



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 1, 2023 – 11:47 AM EDT

PDB ID : 4IO9
Title : Crystal structure of compound 4d bound to large ribosomal subunit (50S) from *Deinococcus radiodurans*
Authors : Han, S.; Marr, E.S.
Deposited on : 2013-01-07
Resolution : 3.20 Å (reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

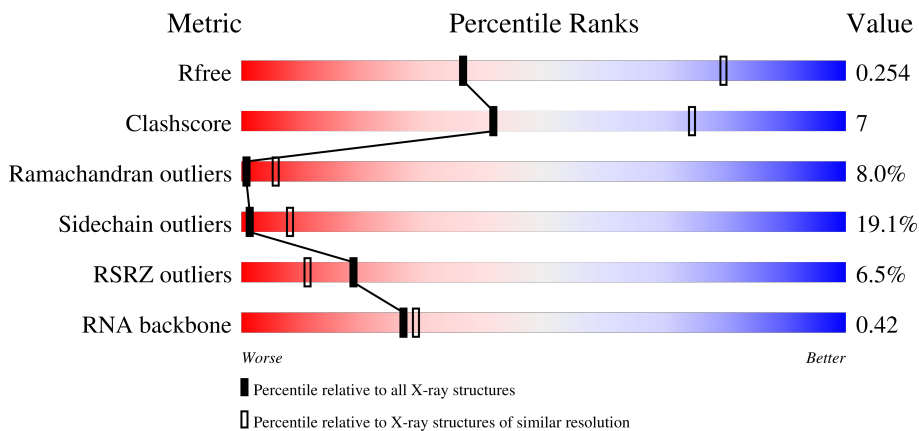
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)
RNA backbone	3102	1010 (3.50-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	X	2880	
2	Y	123	
3	A	274	
4	B	211	

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Mol	Chain	Length	Quality of chain
5	C	205	
6	D	180	
7	E	185	
8	F	144	
9	G	174	
10	H	134	
11	I	156	
12	J	141	
13	K	116	
14	L	114	
15	M	166	
16	N	118	
17	O	100	
18	P	134	
19	Q	95	
20	R	115	
21	S	237	
22	T	91	
23	U	81	
24	V	67	
25	W	55	
26	Z	60	
27	1	55	
28	2	47	
29	3	66	

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Mol	Chain	Length	Quality of chain
30	4	37	<p>97% 76% 22%</p>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
31	MG	X	2902	-	-	-	X
31	MG	X	2903	-	-	-	X
31	MG	X	2912	-	-	-	X
31	MG	Y	201	-	-	-	X
31	MG	Y	203	-	-	-	X

2 Entry composition [i](#)

There are 32 unique types of molecules in this entry. The entry contains 83875 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	X	2686	57651	25718	10642	18606	2685	0	0	0

- Molecule 2 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	Y	122	2598	1161	476	840	121	0	0	0

- Molecule 3 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	A	240	1826	1137	366	321	2	0	0	0

- Molecule 4 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	B	205	1539	965	295	271	8	0	0	0

- Molecule 5 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	C	197	1506	935	287	282	2	0	0	0

- Molecule 6 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	D	177	1400	892	247	254	7	0	0	0

- Molecule 7 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	E	171	1286	812	237	236	1	0	0	0

- Molecule 8 is a protein called 50S ribosomal protein L11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	F	71	503	310	91	99	3	0	0	0

- Molecule 9 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
9	G	142	1114	704	209	198	3	0	0	0

- Molecule 10 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
10	H	134	997	614	198	180	5	0	0	0

- Molecule 11 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
11	I	141	1067	655	216	196	0	0	0

- Molecule 12 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
12	J	136	1090	696	202	185	7	0	0	0

- Molecule 13 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
13	K	113	878	541	178	157	2	0	0	0

- Molecule 14 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
14	L	104	779	476	161	142	0	0	0

- Molecule 15 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
15	M	108	871	543	172	156	0	0	0

- Molecule 16 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
16	N	117	978	608	210	159	1	0	0	0

- Molecule 17 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
17	O	94	741	465	139	137	0	0	0

- Molecule 18 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
18	P	127	1014	639	199	174	2	0	0	0

- Molecule 19 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
19	Q	93	726	458	136	130	2	0	0	0

- Molecule 20 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
20	R	110	825	513	160	151	1	0	0	0

- Molecule 21 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
21	S	175	1345	849	236	254	6	0	0	0

- Molecule 22 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
22	T	84	625	393	122	109	1	0	0	0

- Molecule 23 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
23	U	72	552	341	116	95	0	0	0

- Molecule 24 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
24	V	66	533	327	107	96	3	0	0	0

- Molecule 25 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
25	W	55	424	264	82	76	2	0	0	0

- Molecule 26 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
26	Z	58	457	281	94	77	5	0	0	0

- Molecule 27 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
27	1	53	Total 53 C 53	0	0	53

- Molecule 28 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
28	2	46	Total C 46 46	0	0	46

- Molecule 29 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
29	3	63	Total C 63 63	0	0	63

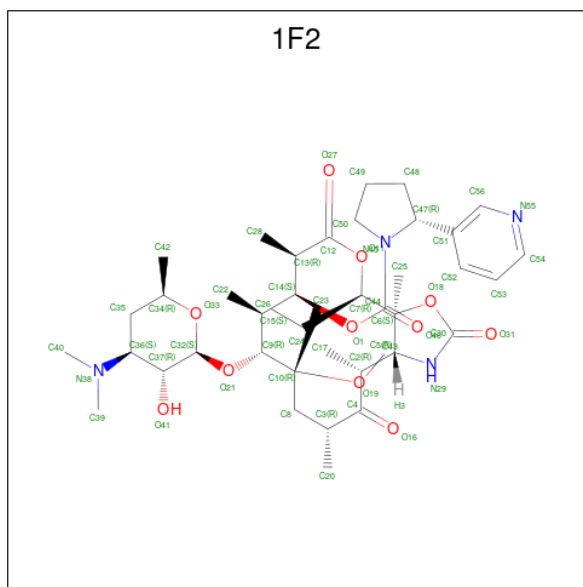
- Molecule 30 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
30	4	37	Total C N O S 297 179 66 47 5	0	0	0

- Molecule 31 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
31	X	28	Total Mg 28 28	0	0
31	Y	5	Total Mg 5 5	0	0
31	J	1	Total Mg 1 1	0	0
31	M	1	Total Mg 1 1	0	0

- Molecule 32 is (3aS,4R,7R,8S,9S,10R,11R,13R,15R,15aR)-4-ethyl-11-methoxy-3a,7,9,11,13,15-hexamethyl-2,6,14-trioxo-10-[[3,4,6-trideoxy-3-(dimethylamino)-beta-D-xylo-hexopyranosyl]oxy]tetradecahydro-2H-oxacyclotetradecino[4,3-d][1,3]oxazol-8-yl (2R)-2-(pyridin-3-yl)pyrrolidine-1-carboxylate (three-letter code: 1F2) (formula: C₄₁H₆₄N₄O₁₁).

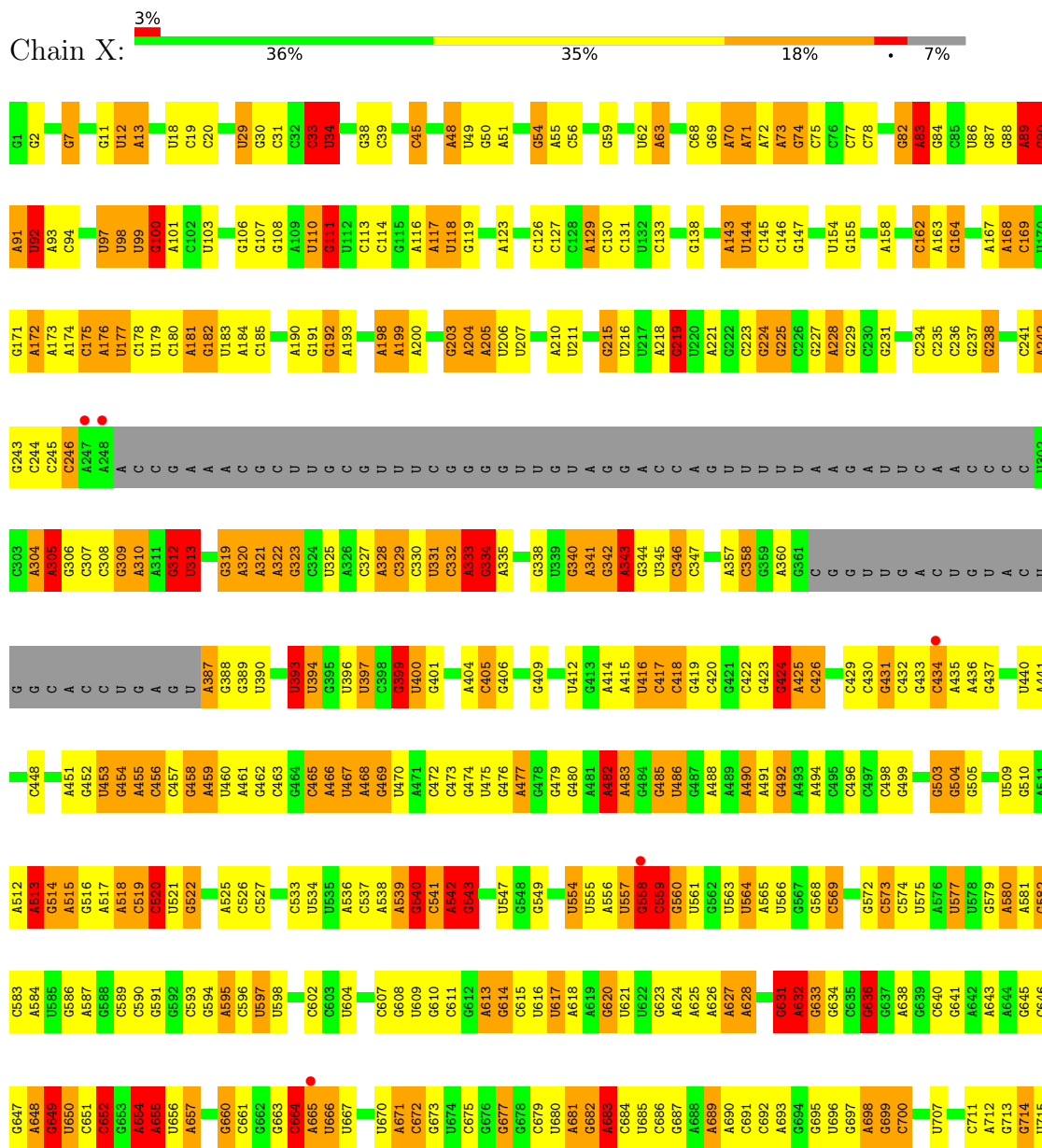


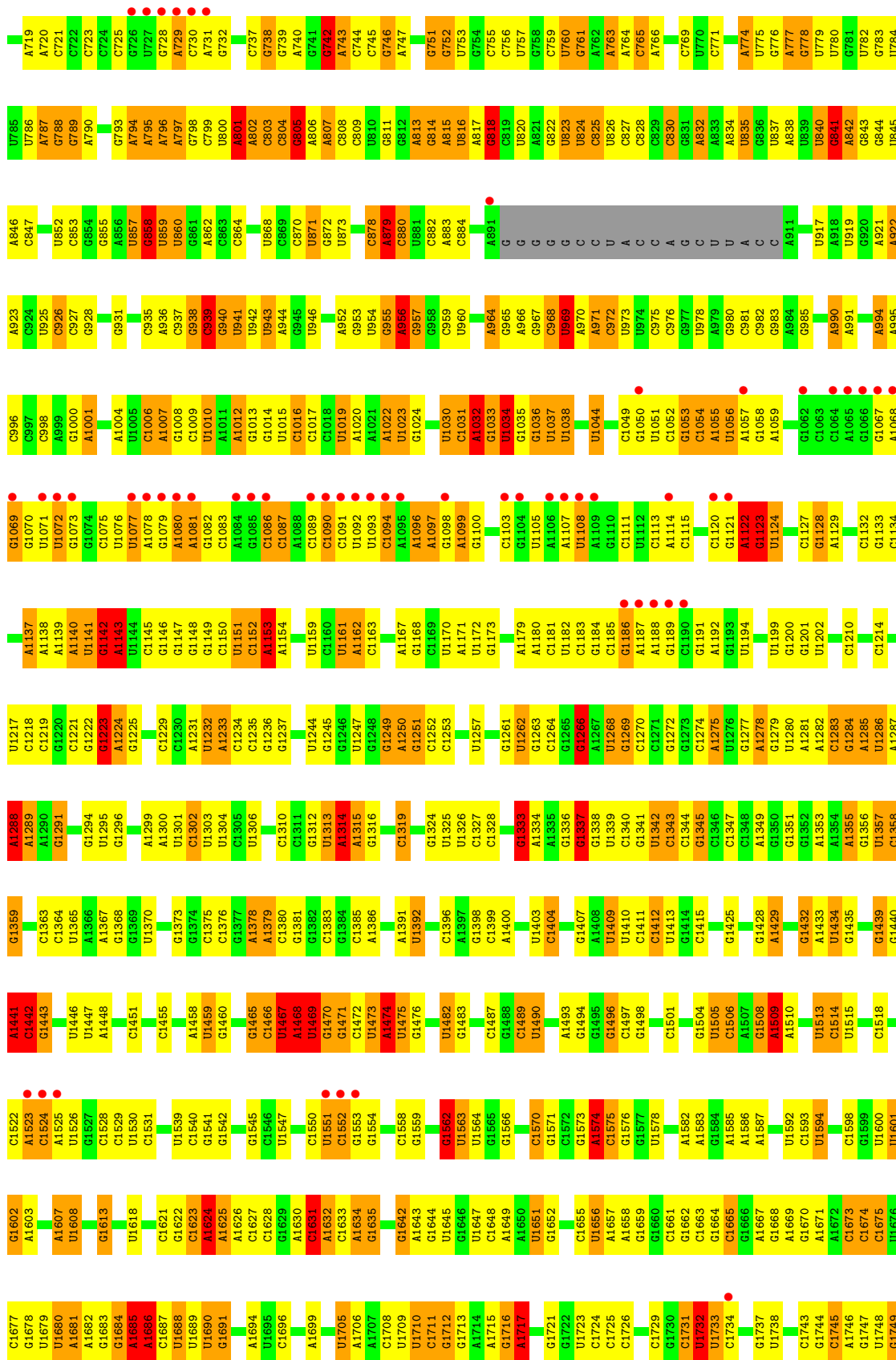
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
32	X	1	56	41	4	11	0	0

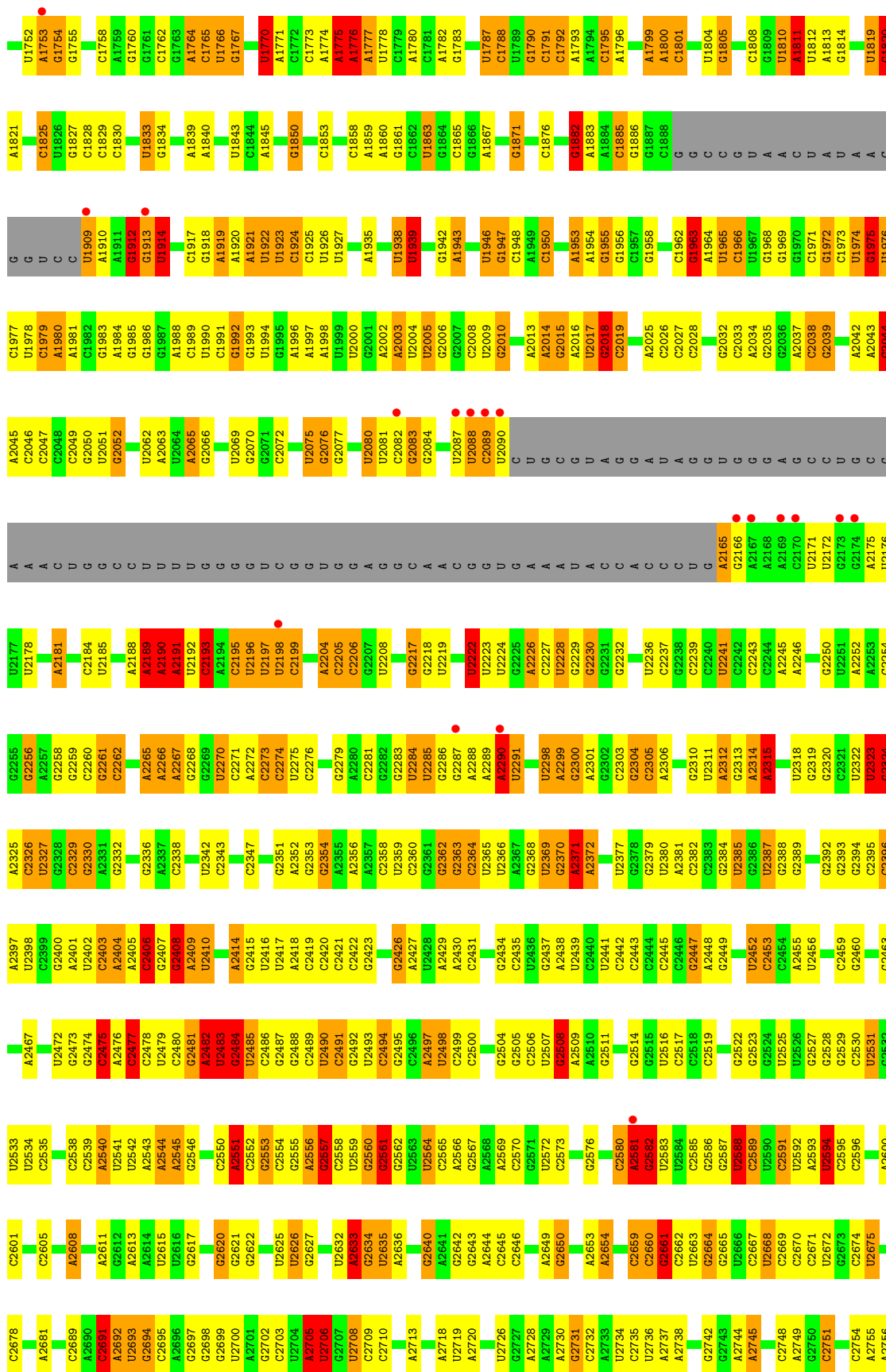
3 Residue-property plots

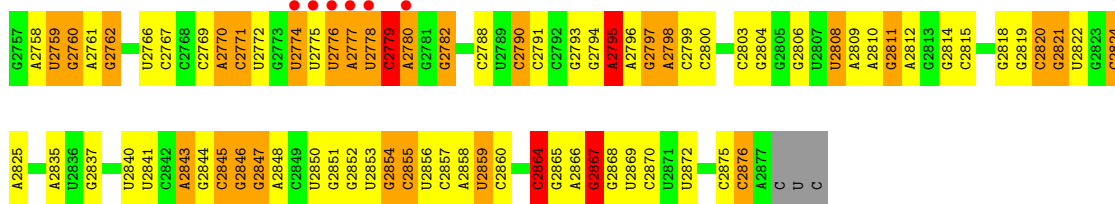
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 23S ribosomal RNA

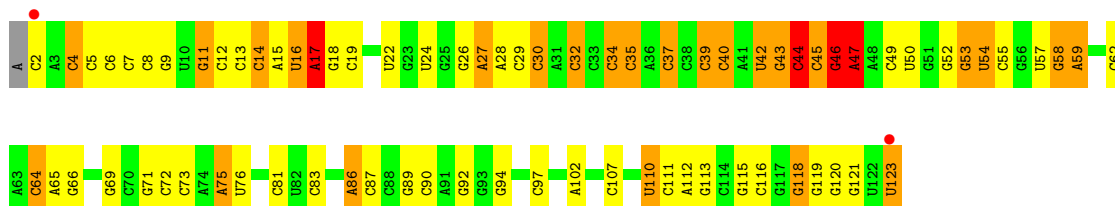




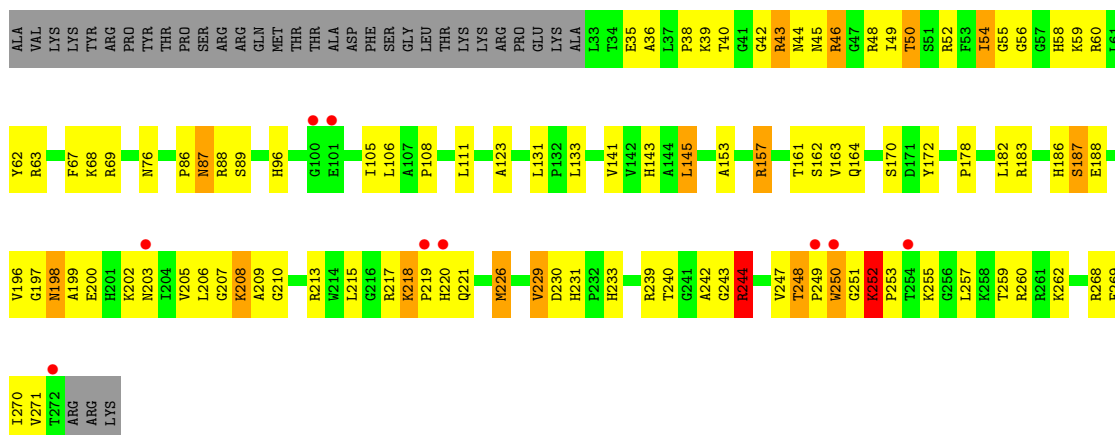




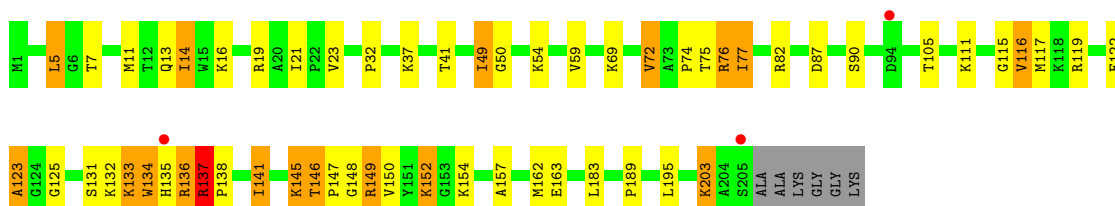
• Molecule 2: 5S ribosomal RNA



• Molecule 3: 50S ribosomal protein L2

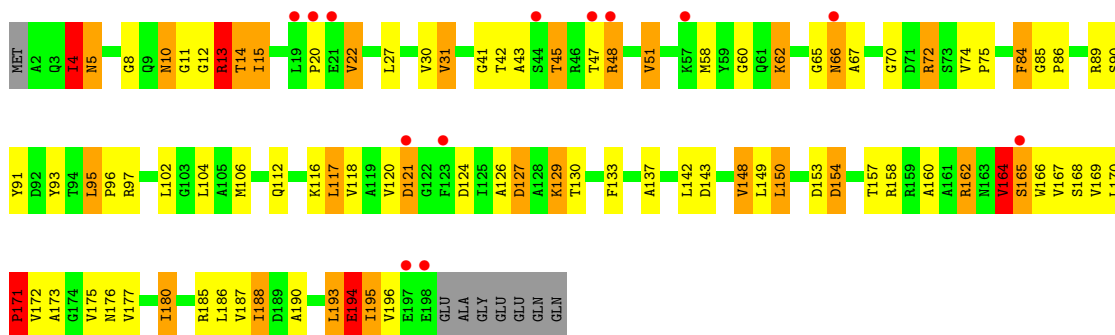


• Molecule 4: 50S ribosomal protein L3

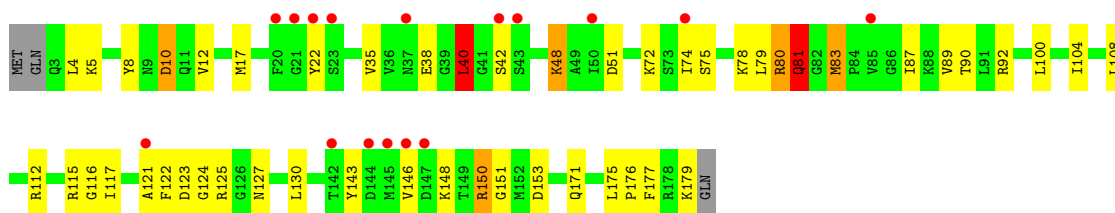


• Molecule 5: 50S ribosomal protein L4

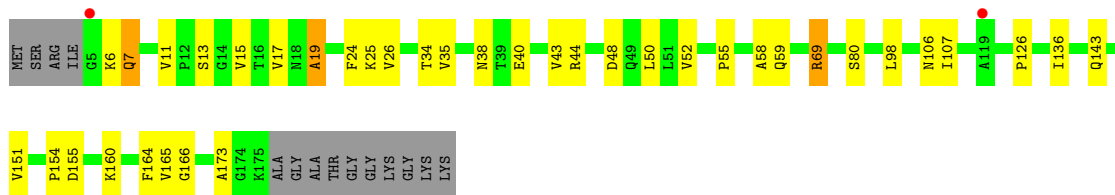




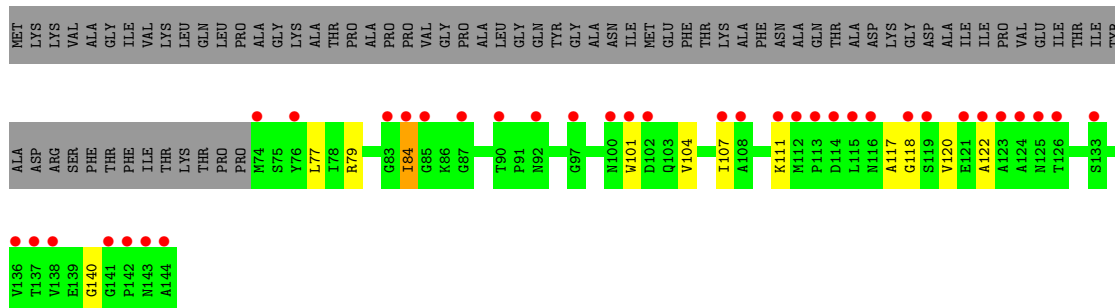
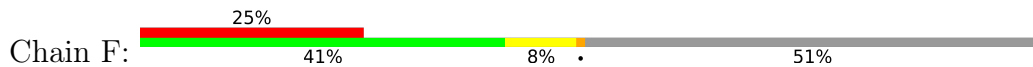
• Molecule 6: 50S ribosomal protein L5



• Molecule 7: 50S ribosomal protein L6

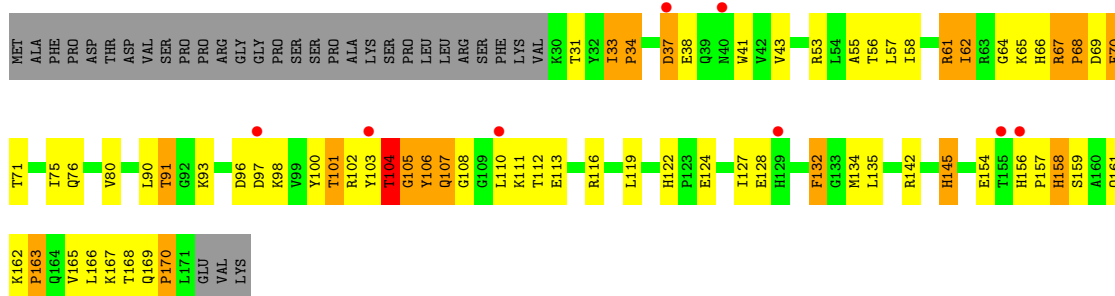


• Molecule 8: 50S ribosomal protein L11



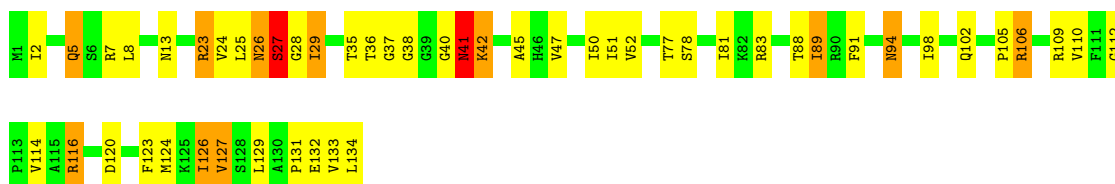
• Molecule 9: 50S ribosomal protein L13





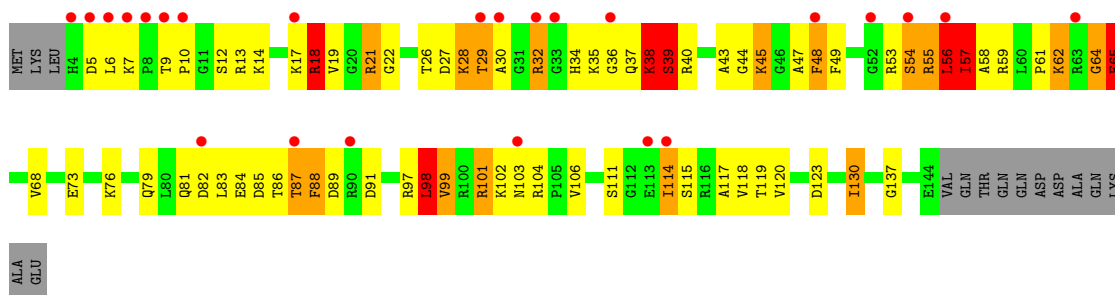
- Molecule 10: 50S ribosomal protein L14

Chain H: 62% 28% 8%



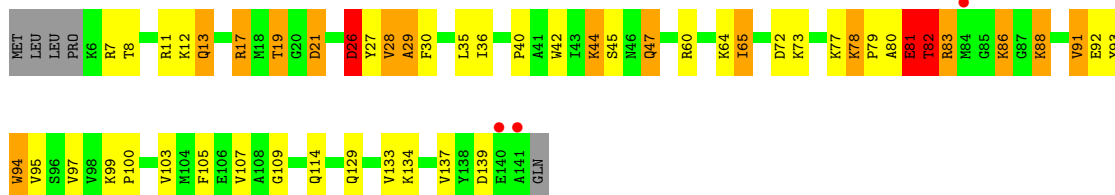
- Molecule 11: 50S ribosomal protein L15

Chain I: 15% 42% 33% 10% 10%



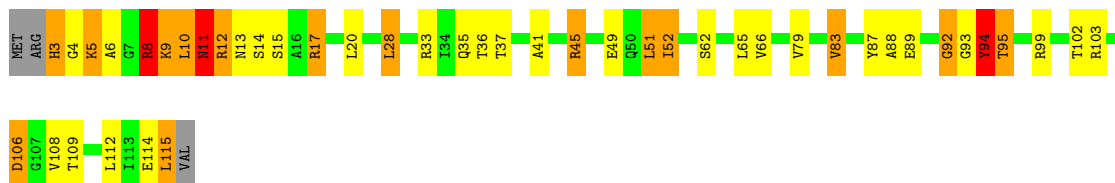
- Molecule 12: 50S ribosomal protein L16

Chain J: 2% 60% 24% 11%

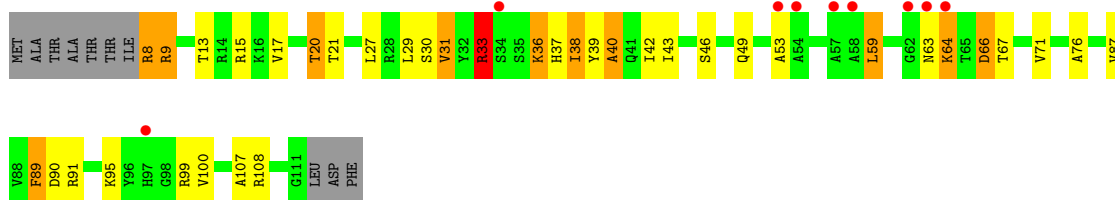


- Molecule 13: 50S ribosomal protein L17

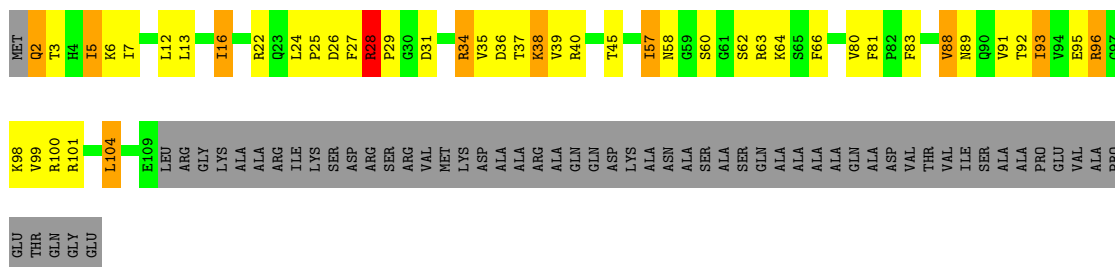
Chain K: 59% 23% 13%



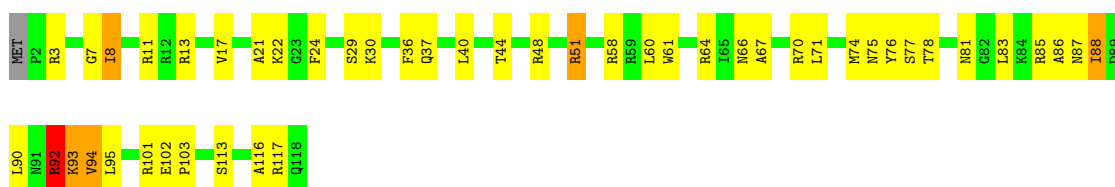
- Molecule 14: 50S ribosomal protein L18



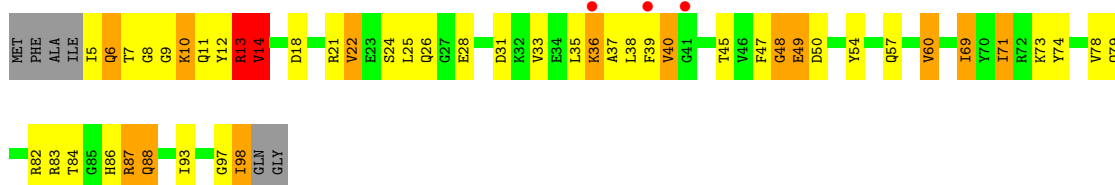
- Molecule 15: 50S ribosomal protein L19



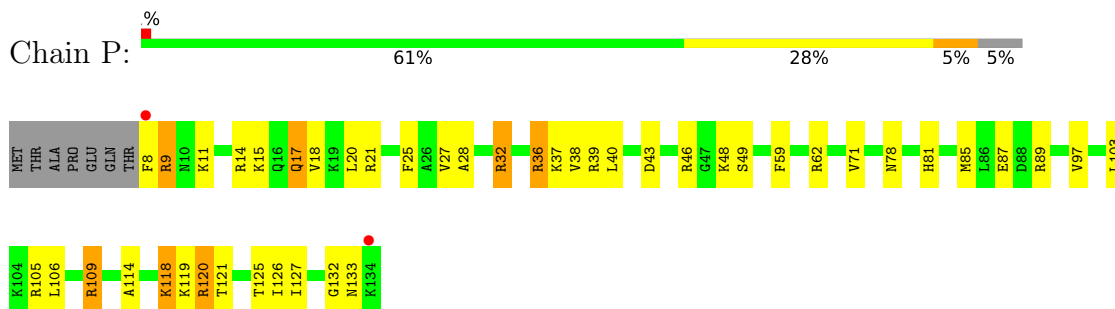
- Molecule 16: 50S ribosomal protein L20



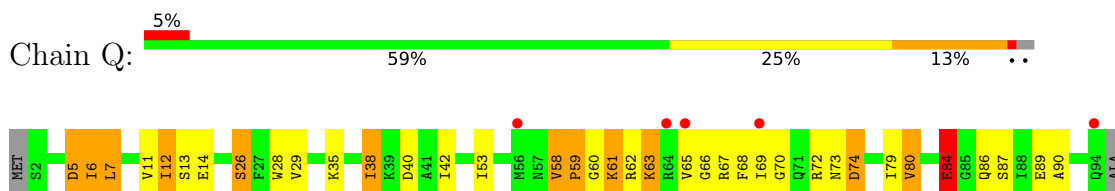
- Molecule 17: 50S ribosomal protein L21



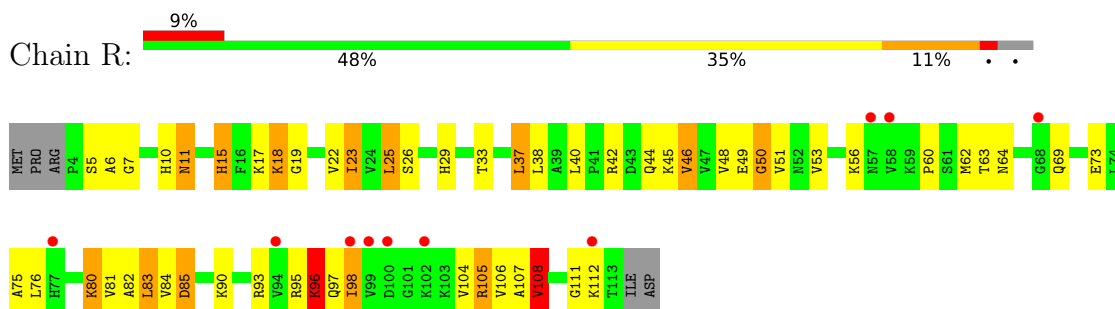
- Molecule 18: 50S ribosomal protein L22



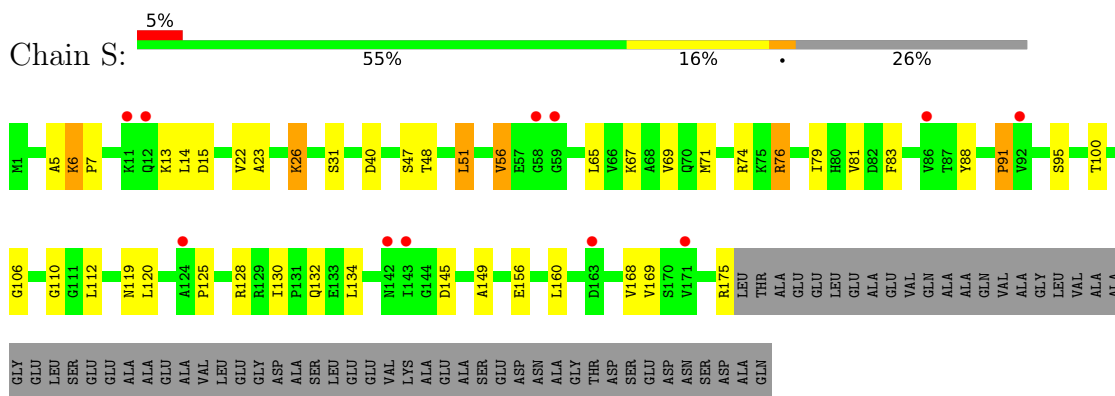
- Molecule 19: 50S ribosomal protein L23



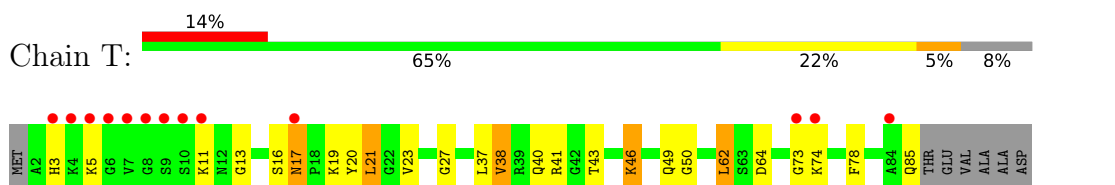
- Molecule 20: 50S ribosomal protein L24



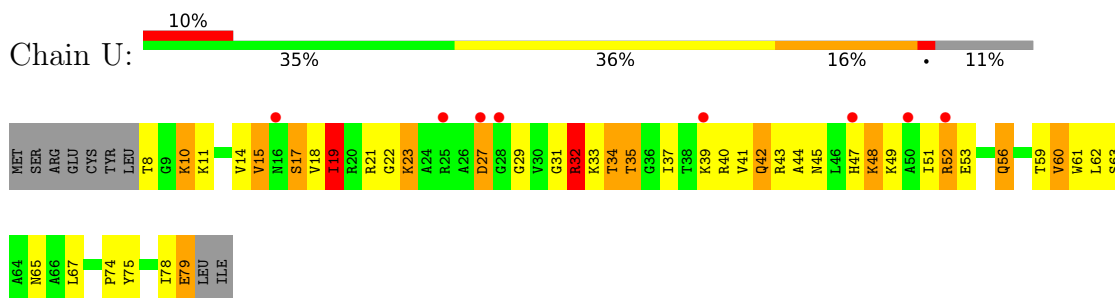
- Molecule 21: 50S ribosomal protein L25



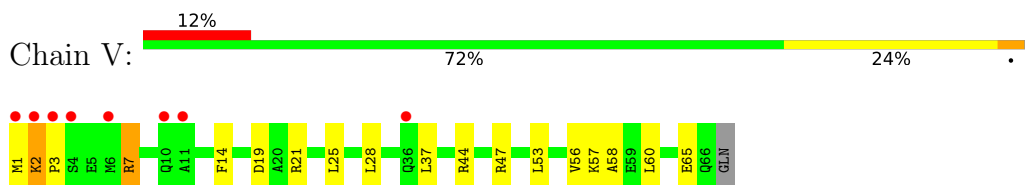
- Molecule 22: 50S ribosomal protein L27



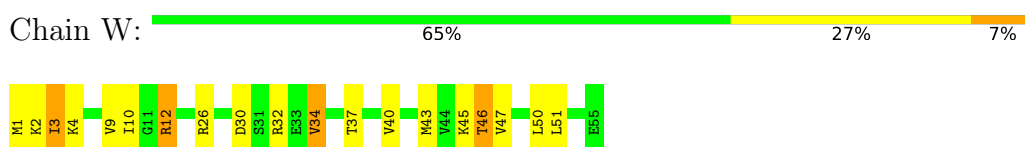
- Molecule 23: 50S ribosomal protein L28



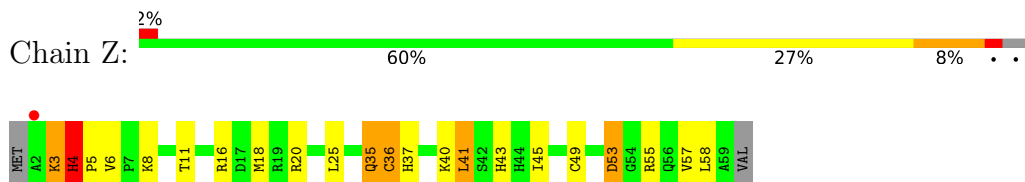
- Molecule 24: 50S ribosomal protein L29



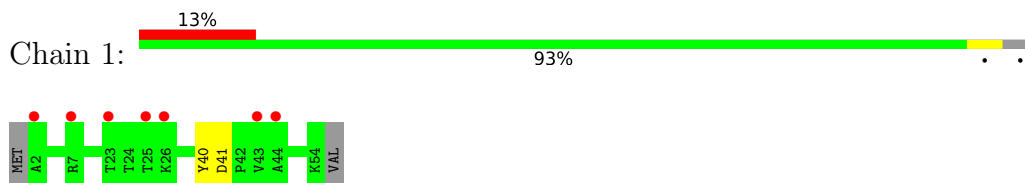
- Molecule 25: 50S ribosomal protein L30



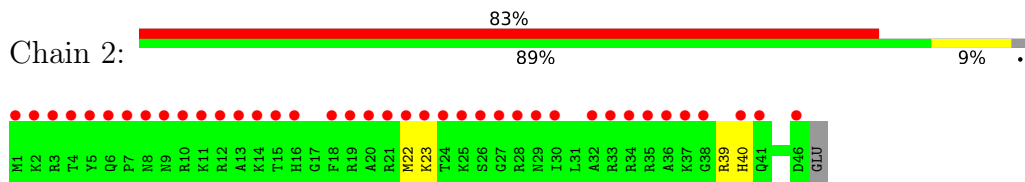
- Molecule 26: 50S ribosomal protein L32



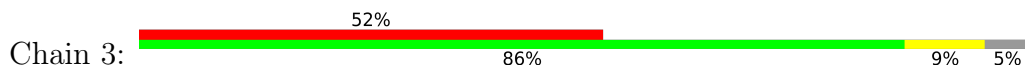
- Molecule 27: 50S ribosomal protein L33

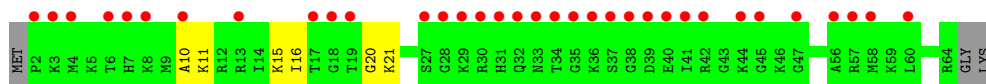


- Molecule 28: 50S ribosomal protein L34

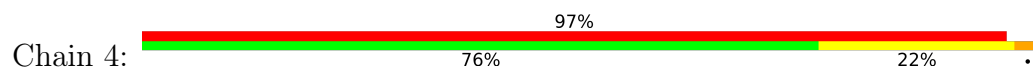


- Molecule 29: 50S ribosomal protein L35





- Molecule 30: 50S ribosomal protein L36



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	169.94Å 409.69Å 694.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.20 30.20 – 3.21	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-3.20) 94.1 (30.20-3.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 3.24Å)	Xtrriage
Refinement program	autobuster	Depositor
R, R_{free}	0.199 , 0.235 0.214 , 0.254	Depositor DCC
R_{free} test set	18481 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	81.2	Xtrriage
Anisotropy	0.747	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 93.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	83875	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.92% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: MG, 1F2

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	X	1.00	46/64561 (0.1%)	1.87	1961/100708 (1.9%)
2	Y	1.05	0/2904	1.78	84/4525 (1.9%)
3	A	0.61	0/1862	0.92	1/2510 (0.0%)
4	B	0.57	0/1567	0.94	1/2105 (0.0%)
5	C	0.62	0/1529	0.98	2/2070 (0.1%)
6	D	0.45	0/1419	0.66	0/1903
7	E	0.45	0/1308	0.67	0/1771
8	F	0.46	0/508	0.64	0/683
9	G	0.58	0/1138	0.94	1/1539 (0.1%)
10	H	0.55	0/1007	0.88	1/1352 (0.1%)
11	I	0.73	1/1081 (0.1%)	1.12	6/1448 (0.4%)
12	J	0.68	1/1113 (0.1%)	0.95	1/1486 (0.1%)
13	K	0.77	2/886 (0.2%)	1.02	3/1188 (0.3%)
14	L	0.53	0/785	0.88	1/1048 (0.1%)
15	M	0.64	0/884	0.98	1/1186 (0.1%)
16	N	0.51	0/994	0.77	0/1323
17	O	0.52	0/750	0.95	1/1000 (0.1%)
18	P	0.56	0/1027	0.85	0/1373
19	Q	0.60	0/737	1.03	5/988 (0.5%)
20	R	0.61	0/835	0.99	0/1121
21	S	0.48	0/1370	0.73	0/1862
22	T	0.55	0/633	0.82	0/838
23	U	0.75	0/556	1.10	1/741 (0.1%)
24	V	0.47	0/537	0.71	0/714
25	W	0.48	0/426	0.81	0/568
26	Z	0.62	0/469	0.97	0/629
30	4	0.44	0/298	0.62	0/390
All	All	0.91	50/91184 (0.1%)	1.69	2070/137069 (1.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected

by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	X	0	6

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	1688	U	C4-O4	9.12	1.30	1.23
1	X	774	A	C5-C4	8.90	1.45	1.38
1	X	1685	A	C3'-O3'	7.74	1.52	1.42
1	X	1468	A	N9-C4	7.70	1.42	1.37
1	X	1333	G	N9-C4	-7.09	1.32	1.38
1	X	1688	U	N3-C4	6.99	1.44	1.38
1	X	1688	U	C2-N3	6.89	1.42	1.37
1	X	2189	A	C3'-O3'	6.77	1.51	1.42
13	K	52	ILE	CG1-CD1	6.46	1.95	1.50
1	X	1980	A	N7-C5	-6.44	1.35	1.39
1	X	1467	U	C1'-N1	6.42	1.58	1.48
1	X	774	A	N7-C5	-6.38	1.35	1.39
1	X	774	A	C6-N1	6.35	1.40	1.35
1	X	346	C	C1'-N1	6.33	1.58	1.48
1	X	2018	G	N9-C8	6.33	1.42	1.37
1	X	559	C	C3'-O3'	6.18	1.50	1.42
1	X	1946	U	C1'-N1	6.16	1.57	1.48
1	X	1288	A	C4'-C3'	-6.11	1.46	1.53
1	X	1975	G	C3'-O3'	6.11	1.50	1.42
11	I	57	ILE	CG1-CD1	6.08	1.92	1.50
1	X	774	A	N1-C2	6.08	1.39	1.34
1	X	1223	G	C2-N3	5.85	1.37	1.32
1	X	699	G	N9-C4	-5.84	1.33	1.38
1	X	796	A	N9-C4	-5.64	1.34	1.37
1	X	759	C	N3-C4	5.62	1.37	1.33
1	X	838	A	C3'-O3'	5.59	1.50	1.42
1	X	661	C	C1'-N1	5.59	1.57	1.48
1	X	646	C	C1'-N1	5.57	1.57	1.48
1	X	343	A	N9-C4	5.56	1.41	1.37
1	X	540	G	C2-N3	5.52	1.37	1.32
1	X	393	U	C1'-N1	5.50	1.57	1.48
1	X	656	U	P-O5'	5.49	1.65	1.59
1	X	559	C	C1'-N1	5.47	1.56	1.48
12	J	19	THR	CA-C	5.42	1.67	1.52
13	K	3	HIS	CA-C	5.41	1.67	1.52
1	X	2735	C	C1'-N1	5.39	1.56	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	796	A	N7-C5	-5.39	1.36	1.39
1	X	540	G	C5-C6	5.37	1.47	1.42
1	X	462	G	C6-O6	5.26	1.28	1.24
1	X	2015	G	N7-C5	5.25	1.42	1.39
1	X	2668	U	C4-C5	5.22	1.48	1.43
1	X	577	U	C4-C5	5.18	1.48	1.43
1	X	927	C	C1'-N1	5.16	1.56	1.48
1	X	2582	G	P-O5'	5.13	1.64	1.59
1	X	774	A	N3-C4	5.13	1.38	1.34
1	X	1688	U	C1'-N1	5.12	1.56	1.48
1	X	2697	G	C5-C4	-5.09	1.34	1.38
1	X	661	C	N1-C2	5.06	1.45	1.40
1	X	757	U	C3'-O3'	-5.03	1.35	1.42
1	X	2800	C	C3'-O3'	-5.00	1.35	1.42

All (2070) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1631	C	C1'-O4'-C4'	-33.72	82.92	109.90
1	X	1288	A	C1'-O4'-C4'	-32.93	83.55	109.90
1	X	1288	A	C5'-C4'-O4'	20.68	133.91	109.10
1	X	1288	A	O4'-C1'-N9	20.52	124.61	108.20
1	X	1019	U	P-O3'-C3'	19.59	143.21	119.70
1	X	774	A	N1-C6-N6	19.05	130.03	118.60
1	X	1631	C	C5'-C4'-O4'	18.47	131.26	109.10
1	X	2808	U	O4'-C1'-N1	17.97	122.57	108.20
1	X	774	A	N7-C8-N9	17.08	122.34	113.80
1	X	2705	A	P-O3'-C3'	17.01	140.11	119.70
1	X	1278	A	O4'-C1'-N9	16.77	121.61	108.20
1	X	1963	G	P-O3'-C3'	16.73	139.78	119.70
1	X	1716	G	P-O3'-C3'	16.57	139.58	119.70
1	X	1631	C	C4'-C3'-C2'	-16.28	86.32	102.60
1	X	1775	A	P-O3'-C3'	16.21	139.15	119.70
1	X	1333	G	N3-C4-N9	-16.10	116.34	126.00
1	X	343	A	O4'-C1'-N9	15.89	120.92	108.20
1	X	994	A	P-O3'-C3'	15.83	138.70	119.70
1	X	788	G	P-O3'-C3'	15.71	138.55	119.70
1	X	1634	A	P-O3'-C3'	15.64	138.47	119.70
1	X	2189	A	P-O3'-C3'	15.53	138.34	119.70
1	X	1475	U	P-O3'-C3'	15.44	138.22	119.70
1	X	1473	U	P-O3'-C3'	15.14	137.87	119.70
1	X	774	A	C5-N7-C8	-15.10	96.35	103.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2497	A	P-O3'-C3'	14.90	137.58	119.70
1	X	2564	U	P-O3'-C3'	14.53	137.13	119.70
1	X	774	A	C6-C5-N7	-14.30	122.29	132.30
1	X	1333	G	O4'-C1'-N9	14.26	119.61	108.20
1	X	594	G	P-O3'-C3'	14.24	136.79	119.70
1	X	399	G	P-O3'-C3'	14.02	136.53	119.70
1	X	469	G	P-O3'-C3'	13.93	136.41	119.70
1	X	1574	A	O4'-C1'-N9	13.91	119.33	108.20
1	X	1355	A	P-O3'-C3'	13.80	136.26	119.70
1	X	558	G	P-O3'-C3'	13.72	136.16	119.70
1	X	73	A	P-O3'-C3'	13.62	136.04	119.70
1	X	1249	G	P-O3'-C3'	13.61	136.03	119.70
1	X	176	A	P-O3'-C3'	13.39	135.77	119.70
1	X	540	G	N1-C6-O6	-13.38	111.87	119.90
1	X	204	A	P-O3'-C3'	13.37	135.74	119.70
1	X	1261	G	P-O3'-C3'	13.07	135.39	119.70
1	X	467	U	P-O3'-C3'	12.91	135.20	119.70
1	X	1938	U	P-O3'-C3'	12.82	135.09	119.70
1	X	2769	C	O4'-C1'-N1	12.79	118.43	108.20
1	X	242	A	O4'-C1'-N9	12.78	118.42	108.20
1	X	2312	A	P-O3'-C3'	12.72	134.96	119.70
1	X	2551	A	P-O3'-C3'	12.67	134.91	119.70
1	X	2018	G	P-O3'-C3'	12.66	134.90	119.70
1	X	2014	A	P-O3'-C3'	12.60	134.82	119.70
2	Y	58	G	P-O3'-C3'	12.56	134.78	119.70
1	X	181	A	P-O3'-C3'	12.54	134.75	119.70
1	X	1036	G	P-O3'-C3'	12.54	134.75	119.70
1	X	100	G	P-O3'-C3'	12.53	134.74	119.70
1	X	99	U	P-O3'-C3'	12.52	134.73	119.70
1	X	71	A	P-O3'-C3'	12.46	134.66	119.70
1	X	1820	G	P-O3'-C3'	12.44	134.63	119.70
2	Y	16	U	P-O3'-C3'	12.29	134.45	119.70
1	X	334	G	P-O3'-C3'	12.27	134.42	119.70
1	X	33	C	O4'-C1'-N1	12.23	117.98	108.20
1	X	2088	U	P-O3'-C3'	12.23	134.37	119.70
1	X	342	G	P-O3'-C3'	12.19	134.33	119.70
1	X	559	C	O4'-C1'-N1	12.09	117.87	108.20
1	X	774	A	C4-C5-N7	12.09	116.74	110.70
1	X	1288	A	C4'-C3'-C2'	-12.04	90.56	102.60
1	X	1688	U	N3-C4-O4	12.02	127.81	119.40
1	X	33	C	P-O3'-C3'	11.96	134.06	119.70
1	X	1790	G	P-O3'-C3'	11.79	133.85	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1142	G	C5'-C4'-O4'	11.74	123.19	109.10
1	X	664	C	P-O3'-C3'	11.72	133.76	119.70
1	X	2298	U	P-O3'-C3'	11.70	133.74	119.70
1	X	537	C	N1-C2-O2	11.68	125.91	118.90
1	X	1333	G	N3-C4-C5	11.66	134.43	128.60
1	X	2190	A	O4'-C1'-N9	11.63	117.50	108.20
1	X	1688	U	N3-C4-C5	-11.60	107.64	114.60
1	X	218	A	P-O3'-C3'	11.56	133.57	119.70
1	X	2736	U	P-O3'-C3'	11.55	133.56	119.70
1	X	454	G	P-O3'-C3'	11.53	133.53	119.70
1	X	2769	C	C1'-O4'-C4'	-11.49	100.70	109.90
1	X	1442	C	P-O3'-C3'	11.48	133.48	119.70
1	X	559	C	C4'-C3'-C2'	-11.37	91.23	102.60
1	X	1467	U	P-O3'-C3'	-11.29	106.15	119.70
1	X	1574	A	C1'-O4'-C4'	-11.15	100.98	109.90
1	X	2668	U	C5-C4-O4	11.13	132.58	125.90
1	X	943	U	O4'-C1'-N1	11.08	117.06	108.20
1	X	2669	C	N1-C2-O2	11.05	125.53	118.90
1	X	518	A	P-O3'-C3'	11.03	132.93	119.70
1	X	774	A	C5-C6-N1	-11.02	112.19	117.70
1	X	969	U	P-O3'-C3'	10.98	132.88	119.70
1	X	1152	C	P-O3'-C3'	10.96	132.86	119.70
1	X	1468	A	O4'-C1'-N9	10.96	116.97	108.20
1	X	774	A	C8-N9-C4	-10.94	101.42	105.80
1	X	1799	A	C1'-O4'-C4'	-10.92	101.17	109.90
1	X	814	G	P-O3'-C3'	10.89	132.76	119.70
1	X	341	A	P-O3'-C3'	10.85	132.72	119.70
1	X	2261	G	P-O3'-C3'	10.85	132.72	119.70
1	X	1467	U	C5-C6-N1	10.85	128.12	122.70
1	X	2691	C	O4'-C1'-N1	10.80	116.84	108.20
1	X	1122	A	P-O3'-C3'	10.79	132.64	119.70
1	X	2204	A	P-O3'-C3'	10.78	132.63	119.70
1	X	683	A	P-O3'-C3'	10.73	132.58	119.70
1	X	48	A	P-O3'-C3'	10.69	132.53	119.70
1	X	1552	C	P-O3'-C3'	10.69	132.53	119.70
1	X	198	A	P-O3'-C3'	10.63	132.46	119.70
1	X	469	G	O4'-C1'-N9	10.60	116.68	108.20
1	X	1429	A	O4'-C1'-N9	10.57	116.66	108.20
1	X	1632	A	P-O3'-C3'	10.57	132.38	119.70
1	X	1850	G	P-O3'-C3'	10.55	132.36	119.70
1	X	1096	A	P-O3'-C3'	10.52	132.32	119.70
1	X	699	G	C5-N7-C8	-10.51	99.05	104.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	803	C	P-O3'-C3'	10.47	132.26	119.70
1	X	2770	A	P-O3'-C3'	10.45	132.24	119.70
1	X	182	G	P-O3'-C3'	10.40	132.18	119.70
1	X	655	A	P-O3'-C3'	10.37	132.14	119.70
1	X	2190	A	P-O5'-C5'	10.35	137.45	120.90
1	X	1053	G	P-O3'-C3'	10.34	132.10	119.70
1	X	1923	U	P-O3'-C3'	10.32	132.09	119.70
1	X	1142	G	O4'-C1'-C2'	-10.30	95.50	105.80
1	X	1186	G	P-O3'-C3'	10.27	132.03	119.70
1	X	638	A	P-O3'-C3'	10.27	132.02	119.70
1	X	1631	C	N1-C1'-C2'	10.22	127.28	114.00
1	X	1574	A	C4'-C3'-C2'	-10.20	92.40	102.60
1	X	2371	A	O4'-C1'-N9	10.16	116.33	108.20
1	X	98	U	P-O3'-C3'	10.16	131.89	119.70
1	X	469	G	C5'-C4'-C3'	-10.15	99.76	116.00
2	Y	54	U	O4'-C1'-N1	10.14	116.31	108.20
1	X	1613	G	C1'-O4'-C4'	-10.14	101.79	109.90
1	X	2018	G	C5'-C4'-O4'	-10.11	96.97	109.10
1	X	34	U	O4'-C1'-N1	10.08	116.27	108.20
1	X	559	C	P-O3'-C3'	10.06	131.77	119.70
1	X	1710	U	P-O3'-C3'	10.01	131.71	119.70
1	X	2498	U	P-O3'-C3'	9.98	131.68	119.70
1	X	418	C	C1'-O4'-C4'	-9.97	101.92	109.90
1	X	1333	G	N3-C2-N2	-9.96	112.93	119.90
1	X	1468	A	C8-N9-C4	-9.96	101.82	105.80
1	X	89	A	P-O3'-C3'	9.94	131.63	119.70
1	X	1631	C	O4'-C4'-C3'	-9.94	94.06	104.00
1	X	1142	G	P-O3'-C3'	9.92	131.60	119.70
1	X	2475	C	O4'-C1'-N1	9.91	116.13	108.20
1	X	1496	G	P-O3'-C3'	9.89	131.57	119.70
1	X	2323	U	O4'-C1'-N1	9.89	116.11	108.20
1	X	1770	U	O4'-C4'-C3'	-9.88	94.12	104.00
1	X	765	C	P-O3'-C3'	9.87	131.54	119.70
1	X	2795	A	P-O3'-C3'	9.85	131.52	119.70
1	X	841	G	O4'-C4'-C3'	-9.84	94.16	104.00
1	X	1223	G	C3'-C2'-C1'	9.78	109.33	101.50
1	X	2016	A	P-O3'-C3'	9.77	131.43	119.70
1	X	1523	A	P-O3'-C3'	9.76	131.41	119.70
1	X	2596	C	O4'-C1'-N1	9.76	116.01	108.20
1	X	242	A	C1'-O4'-C4'	-9.74	102.11	109.90
1	X	1811	A	P-O3'-C3'	9.71	131.35	119.70
1	X	1333	G	C8-N9-C1'	9.70	139.61	127.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	938	G	O4'-C1'-N9	9.68	115.94	108.20
1	X	1266	G	P-O3'-C3'	9.68	131.31	119.70
1	X	1288	A	C3'-C2'-C1'	-9.65	93.78	101.50
1	X	515	A	P-O3'-C3'	9.65	131.28	119.70
1	X	1975	G	C2'-C3'-O3'	9.64	130.71	109.50
1	X	1391	A	P-O3'-C3'	9.60	131.22	119.70
1	X	1055	A	P-O3'-C3'	9.59	131.21	119.70
1	X	1469	U	P-O3'-C3'	9.57	131.19	119.70
1	X	1575	C	P-O3'-C3'	9.57	131.18	119.70
1	X	650	U	O4'-C1'-N1	9.56	115.84	108.20
1	X	2323	U	P-O3'-C3'	9.56	131.17	119.70
1	X	655	A	O4'-C1'-N9	9.55	115.84	108.20
1	X	805	G	O4'-C1'-N9	-9.54	100.56	108.20
1	X	2426	G	P-O3'-C3'	9.54	131.15	119.70
1	X	581	A	P-O3'-C3'	-9.51	108.29	119.70
1	X	1409	U	P-O3'-C3'	9.49	131.09	119.70
1	X	1141	U	P-O3'-C3'	9.48	131.08	119.70
1	X	1753	A	O4'-C1'-N9	9.45	115.76	108.20
1	X	1333	G	N9-C4-C5	9.36	109.14	105.40
2	Y	11	G	C1'-O4'-C4'	-9.36	102.41	109.90
1	X	777	A	P-O3'-C3'	9.35	130.92	119.70
1	X	1732	U	P-O3'-C3'	9.32	130.88	119.70
1	X	1468	A	O4'-C1'-C2'	-9.27	96.53	105.80
1	X	1288	A	O4'-C4'-C3'	-9.25	94.75	104.00
1	X	1412	C	C3'-C2'-C1'	-9.24	94.11	101.50
1	X	2437	G	P-O3'-C3'	9.23	130.77	119.70
1	X	514	G	P-O3'-C3'	9.22	130.76	119.70
1	X	699	G	N3-C4-C5	9.21	133.21	128.60
1	X	1086	C	P-O3'-C3'	9.21	130.75	119.70
1	X	1469	U	N1-C1'-C2'	9.21	125.97	114.00
1	X	540	G	C5-C6-O6	9.20	134.12	128.60
1	X	580	A	P-O3'-C3'	9.16	130.70	119.70
1	X	1345	G	C1'-O4'-C4'	-9.16	102.57	109.90
1	X	1601	U	P-O3'-C3'	9.16	130.69	119.70
1	X	1674	C	O4'-C1'-N1	9.14	115.51	108.20
1	X	68	C	O4'-C1'-N1	9.13	115.50	108.20
1	X	563	U	O4'-C1'-N1	9.12	115.49	108.20
1	X	686	C	O4'-C1'-N1	9.11	115.49	108.20
1	X	2589	C	P-O3'-C3'	9.09	130.61	119.70
1	X	1790	G	O4'-C1'-N9	9.07	115.45	108.20
1	X	175	C	P-O3'-C3'	9.05	130.56	119.70
1	X	2018	G	N3-C4-C5	9.04	133.12	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2481	G	P-O3'-C3'	9.02	130.52	119.70
1	X	632	A	O4'-C1'-N9	9.01	115.41	108.20
1	X	1770	U	C1'-O4'-C4'	-9.00	102.70	109.90
1	X	731	A	P-O3'-C3'	8.98	130.48	119.70
1	X	1975	G	P-O3'-C3'	8.98	130.47	119.70
1	X	957	G	P-O3'-C3'	8.96	130.45	119.70
1	X	2703	C	O4'-C1'-N1	8.94	115.36	108.20
1	X	1233	A	P-O3'-C3'	8.94	130.42	119.70
1	X	346	C	O4'-C1'-N1	8.93	115.35	108.20
1	X	2229	G	P-O3'-C3'	8.92	130.40	119.70
1	X	2491	C	O5'-P-OP2	-8.91	97.68	105.70
1	X	554	U	P-O3'-C3'	8.88	130.36	119.70
1	X	841	G	N9-C1'-C2'	8.88	125.55	114.00
1	X	522	G	O4'-C1'-N9	8.87	115.30	108.20
1	X	579	G	C4-C5-N7	-8.85	107.26	110.80
1	X	2324	G	P-O3'-C3'	8.81	130.28	119.70
1	X	1283	C	P-O3'-C3'	8.80	130.26	119.70
1	X	714	G	O4'-C4'-C3'	-8.78	95.22	104.00
1	X	2530	C	O5'-P-OP2	-8.78	97.80	105.70
1	X	333	A	P-O3'-C3'	8.76	130.21	119.70
1	X	1474	A	P-O3'-C3'	8.74	130.19	119.70
1	X	774	A	C5-C6-N6	-8.73	116.71	123.70
1	X	387	A	P-O3'-C3'	8.73	130.17	119.70
1	X	939	C	P-O3'-C3'	8.69	130.13	119.70
1	X	2493	U	O4'-C1'-N1	8.67	115.14	108.20
1	X	656	U	O4'-C1'-N1	8.66	115.13	108.20
1	X	2044	G	O4'-C1'-C2'	-8.66	97.14	105.80
1	X	2487	G	O4'-C1'-N9	8.64	115.11	108.20
1	X	689	A	C5-N7-C8	-8.63	99.58	103.90
1	X	172	A	P-O3'-C3'	8.62	130.04	119.70
1	X	2633	A	P-O3'-C3'	8.61	130.03	119.70
1	X	2808	U	C1'-O4'-C4'	-8.55	103.06	109.90
1	X	2554	C	N1-C2-O2	8.55	124.03	118.90
1	X	540	G	O4'-C1'-N9	8.54	115.03	108.20
1	X	841	G	C8-N9-C4	-8.52	102.99	106.40
1	X	2671	C	O4'-C1'-N1	8.52	115.02	108.20
1	X	346	C	C6-N1-C2	-8.51	116.89	120.30
1	X	1524	C	P-O3'-C3'	8.49	129.89	119.70
1	X	2018	G	N9-C1'-C2'	8.49	125.04	114.00
1	X	976	C	O4'-C1'-N1	8.49	114.99	108.20
1	X	1359	G	C8-N9-C4	-8.47	103.01	106.40
1	X	574	C	O4'-C1'-N1	8.46	114.97	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1409	U	C1'-O4'-C4'	-8.45	103.14	109.90
1	X	742	G	P-O3'-C3'	8.44	129.83	119.70
2	Y	17	A	O4'-C1'-N9	8.44	114.95	108.20
1	X	2404	A	P-O3'-C3'	8.44	129.83	119.70
1	X	2477	C	O4'-C1'-N1	8.44	114.95	108.20
1	X	699	G	N3-C4-N9	-8.43	120.94	126.00
1	X	636	G	C8-N9-C4	-8.43	103.03	106.40
1	X	2479	U	P-O3'-C3'	8.42	129.81	119.70
1	X	2691	C	C1'-O4'-C4'	-8.39	103.19	109.90
1	X	804	C	O4'-C1'-N1	8.38	114.90	108.20
1	X	2265	A	P-O3'-C3'	8.37	129.75	119.70
1	X	2847	G	C8-N9-C4	-8.35	103.06	106.40
1	X	394	U	O4'-C1'-N1	8.34	114.88	108.20
1	X	2456	U	O4'-C1'-N1	8.34	114.87	108.20
1	X	1006	C	P-O3'-C3'	8.31	129.67	119.70
1	X	2672	U	O4'-C1'-N1	8.30	114.84	108.20
1	X	2812	A	O4'-C1'-N9	8.30	114.84	108.20
1	X	1200	G	O4'-C1'-N9	8.28	114.83	108.20
1	X	1656	U	P-O3'-C3'	8.28	129.63	119.70
1	X	838	A	OP1-P-O3'	8.26	123.37	105.20
1	X	1286	U	O4'-C1'-N1	8.24	114.80	108.20
1	X	2706	U	O4'-C1'-N1	8.24	114.79	108.20
1	X	2854	G	N9-C1'-C2'	8.24	124.72	114.00
1	X	801	A	P-O3'-C3'	8.23	129.57	119.70
1	X	1468	A	P-O3'-C3'	8.21	129.55	119.70
1	X	2744	A	P-O3'-C3'	8.21	129.55	119.70
1	X	2660	C	O4'-C1'-N1	8.20	114.76	108.20
1	X	184	A	O4'-C1'-N9	8.20	114.76	108.20
1	X	2854	G	P-O3'-C3'	8.20	129.53	119.70
1	X	2853	U	O4'-C1'-N1	8.17	114.74	108.20
1	X	490	A	P-O3'-C3'	8.16	129.50	119.70
1	X	2667	C	P-O3'-C3'	8.16	129.50	119.70
1	X	814	G	O4'-C1'-N9	-8.14	101.68	108.20
1	X	1953	A	P-O5'-C5'	-8.14	107.87	120.90
1	X	332	C	O4'-C1'-N1	8.14	114.71	108.20
1	X	953	G	O4'-C1'-N9	8.13	114.71	108.20
1	X	483	A	P-O3'-C3'	-8.13	109.95	119.70
1	X	308	C	O4'-C1'-N1	8.12	114.70	108.20
1	X	1468	A	C3'-C2'-C1'	-8.12	95.01	101.50
1	X	1278	A	C1'-O4'-C4'	-8.11	103.41	109.90
1	X	1333	G	C8-N9-C4	-8.11	103.16	106.40
1	X	2480	C	O4'-C1'-N1	8.10	114.68	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1310	C	O4'-C1'-N1	8.09	114.67	108.20
1	X	1631	C	C5'-C4'-C3'	8.09	128.94	116.00
1	X	1459	U	P-O3'-C3'	8.08	129.39	119.70
1	X	1984	A	P-O5'-C5'	-8.08	107.97	120.90
1	X	1122	A	O4'-C1'-N9	8.08	114.66	108.20
1	X	1250	A	O4'-C1'-N9	-8.07	101.74	108.20
1	X	83	A	P-O3'-C3'	8.07	129.38	119.70
1	X	1199	U	O4'-C1'-N1	8.07	114.65	108.20
1	X	1313	U	O4'-C1'-N1	8.06	114.65	108.20
1	X	2044	G	O4'-C1'-N9	-8.06	101.75	108.20
1	X	656	U	P-O3'-C3'	8.05	129.37	119.70
1	X	2051	U	O4'-C1'-N1	8.05	114.64	108.20
1	X	2032	G	P-O3'-C3'	-8.04	110.05	119.70
1	X	1574	A	C5'-C4'-O4'	8.03	118.74	109.10
1	X	1674	C	C5'-C4'-O4'	-8.03	99.47	109.10
1	X	63	A	C5'-C4'-C3'	-8.01	103.18	116.00
1	X	843	G	P-O3'-C3'	8.00	129.30	119.70
1	X	1278	A	C3'-C2'-C1'	-8.00	95.10	101.50
1	X	990	A	O4'-C4'-C3'	-8.00	96.00	104.00
1	X	2477	C	C5'-C4'-O4'	-8.00	99.50	109.10
1	X	699	G	N7-C8-N9	7.99	117.10	113.10
1	X	1441	A	P-O3'-C3'	7.99	129.29	119.70
1	X	2323	U	N1-C1'-C2'	7.98	124.38	114.00
1	X	514	G	O4'-C1'-N9	-7.98	101.81	108.20
1	X	1938	U	C4'-C3'-C2'	7.97	110.57	102.60
1	X	2481	G	O5'-P-OP1	-7.97	98.53	105.70
1	X	480	G	C5-C6-O6	-7.96	123.83	128.60
1	X	1358	C	O4'-C1'-N1	7.95	114.56	108.20
1	X	1526	U	O4'-C1'-N1	7.95	114.56	108.20
1	X	957	G	N1-C6-O6	-7.94	115.14	119.90
1	X	1250	A	P-O3'-C3'	7.94	129.23	119.70
1	X	2756	A	P-O3'-C3'	7.94	129.22	119.70
1	X	1338	G	P-O3'-C3'	7.93	129.22	119.70
1	X	2237	C	P-O3'-C3'	7.92	129.21	119.70
1	X	343	A	C8-N9-C4	-7.92	102.63	105.80
1	X	1467	U	N1-C2-O2	7.92	128.35	122.80
1	X	2447	G	P-O3'-C3'	7.92	129.21	119.70
1	X	2487	G	C8-N9-C4	-7.91	103.23	106.40
2	Y	26	G	P-O3'-C3'	7.91	129.19	119.70
1	X	1792	C	P-O3'-C3'	7.90	129.18	119.70
1	X	169	C	N1-C2-O2	7.90	123.64	118.90
1	X	1684	G	P-O3'-C3'	7.90	129.18	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	774	A	C4-C5-C6	7.89	120.95	117.00
1	X	2018	G	N3-C4-N9	-7.89	121.27	126.00
1	X	1072	U	P-O3'-C3'	7.88	129.16	119.70
1	X	689	A	N7-C8-N9	7.88	117.74	113.80
1	X	1980	A	C4'-C3'-C2'	-7.87	94.73	102.60
1	X	2013	A	O4'-C1'-N9	-7.87	101.90	108.20
1	X	169	C	O4'-C1'-N1	7.87	114.49	108.20
1	X	2408	G	N3-C4-C5	-7.87	124.67	128.60
1	X	796	A	C5-N7-C8	-7.86	99.97	103.90
1	X	1291	G	O4'-C4'-C3'	-7.86	96.14	104.00
1	X	1607	A	P-O3'-C3'	7.86	129.13	119.70
2	Y	92	G	O4'-C4'-C3'	-7.86	96.14	104.00
1	X	2554	C	O4'-C1'-N1	7.86	114.48	108.20
1	X	975	C	O4'-C1'-N1	7.85	114.48	108.20
1	X	2500	C	O4'-C1'-N1	7.83	114.47	108.20
1	X	2692	A	P-O3'-C3'	7.83	129.10	119.70
19	Q	60	GLY	C-N-CA	7.83	141.28	121.70
1	X	825	C	P-O3'-C3'	-7.83	110.30	119.70
1	X	2481	G	P-O5'-C5'	7.82	133.41	120.90
1	X	2824	C	C2'-C3'-O3'	7.80	126.67	109.50
1	X	1333	G	C4-N9-C1'	-7.79	116.37	126.50
1	X	1631	C	C3'-C2'-C1'	-7.79	95.27	101.50
1	X	1680	U	C2'-C3'-O3'	7.79	126.63	109.50
1	X	2019	C	O5'-P-OP2	-7.78	98.70	105.70
1	X	1509	A	O4'-C1'-N9	7.78	114.42	108.20
1	X	1313	U	C1'-O4'-C4'	-7.77	103.68	109.90
1	X	1665	C	O5'-P-OP2	-7.77	98.70	105.70
1	X	1792	C	N1-C1'-C2'	7.77	124.10	114.00
1	X	2667	C	N1-C2-O2	7.76	123.56	118.90
1	X	2196	U	P-O3'-C3'	7.76	129.01	119.70
2	Y	30	C	O4'-C1'-N1	7.76	114.41	108.20
1	X	1458	A	P-O3'-C3'	7.75	129.01	119.70
1	X	203	G	O4'-C1'-N9	7.75	114.40	108.20
1	X	1559	G	P-O3'-C3'	7.74	128.98	119.70
1	X	2525	U	O4'-C1'-N1	7.72	114.37	108.20
1	X	540	G	C8-N9-C4	-7.71	103.31	106.40
1	X	2859	U	O4'-C1'-N1	7.71	114.37	108.20
1	X	1664	G	O5'-P-OP2	7.70	119.94	110.70
1	X	1142	G	N3-C4-C5	-7.70	124.75	128.60
1	X	1142	G	N3-C4-N9	7.70	130.62	126.00
1	X	738	G	C8-N9-C4	-7.69	103.32	106.40
1	X	1223	G	P-O3'-C3'	7.69	128.92	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1843	U	O4'-C1'-N1	7.69	114.35	108.20
1	X	617	U	N3-C2-O2	-7.69	116.82	122.20
1	X	518	A	N9-C1'-C2'	7.68	123.98	114.00
1	X	864	C	O4'-C1'-N1	7.67	114.34	108.20
1	X	1509	A	C1'-O4'-C4'	-7.67	103.77	109.90
1	X	564	U	O4'-C1'-N1	7.67	114.33	108.20
1	X	1795	C	O4'-C1'-N1	7.66	114.33	108.20
1	X	972	C	C1'-O4'-C4'	-7.66	103.77	109.90
1	X	2867	G	C5-N7-C8	-7.66	100.47	104.30
1	X	1439	G	P-O3'-C3'	7.65	128.88	119.70
1	X	1981	A	N1-C6-N6	7.64	123.19	118.60
1	X	2778	U	P-O3'-C3'	7.64	128.87	119.70
1	X	1770	U	O4'-C1'-N1	7.64	114.31	108.20
1	X	176	A	N9-C1'-C2'	7.62	123.91	114.00
1	X	223	C	O4'-C1'-N1	7.62	114.30	108.20
1	X	516	G	O4'-C1'-N9	7.62	114.30	108.20
2	Y	37	C	O4'-C1'-N1	7.62	114.30	108.20
1	X	636	G	N7-C8-N9	7.62	116.91	113.10
1	X	74	G	O4'-C4'-C3'	-7.61	96.39	104.00
1	X	2018	G	C5-N7-C8	-7.61	100.49	104.30
1	X	809	C	O4'-C1'-N1	7.61	114.29	108.20
1	X	577	U	C4'-C3'-C2'	-7.58	95.02	102.60
1	X	1006	C	P-O5'-C5'	7.58	133.03	120.90
1	X	12	U	N3-C2-O2	-7.57	116.90	122.20
1	X	751	G	O4'-C4'-C3'	-7.57	96.43	104.00
1	X	672	C	O4'-C4'-C3'	-7.57	96.43	104.00
1	X	1791	C	O4'-C1'-N1	7.57	114.25	108.20
1	X	168	A	O4'-C1'-N9	7.56	114.25	108.20
1	X	774	A	N9-C4-C5	-7.55	102.78	105.80
1	X	2439	U	O4'-C1'-N1	7.55	114.24	108.20
1	X	2759	U	P-O3'-C3'	7.54	128.75	119.70
1	X	661	C	N1-C2-O2	7.54	123.43	118.90
1	X	699	G	C4-C5-N7	7.54	113.82	110.80
1	X	542	A	C5-N7-C8	-7.54	100.13	103.90
1	X	1333	G	C2-N3-C4	-7.53	108.14	111.90
1	X	1467	U	N1-C1'-C2'	7.53	123.79	114.00
1	X	1251	G	O4'-C1'-N9	7.52	114.21	108.20
1	X	1184	G	P-O3'-C3'	7.51	128.72	119.70
1	X	661	C	O4'-C1'-N1	7.51	114.21	108.20
1	X	685	U	O4'-C1'-N1	7.51	114.21	108.20
1	X	847	C	O4'-C1'-N1	7.51	114.21	108.20
1	X	1467	U	C4-C5-C6	-7.49	115.20	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1663	C	N1-C2-O2	7.49	123.39	118.90
1	X	1631	C	C2'-C3'-O3'	7.49	125.97	109.50
1	X	2697	G	C5-C6-O6	-7.49	124.11	128.60
1	X	343	A	P-O3'-C3'	7.47	128.67	119.70
1	X	334	G	O4'-C1'-N9	7.47	114.17	108.20
11	I	38	LYS	C-N-CA	7.46	140.36	121.70
1	X	192	G	P-O3'-C3'	7.45	128.64	119.70
1	X	2492	G	O4'-C1'-N9	7.45	114.16	108.20
1	X	661	C	N3-C2-O2	-7.45	116.69	121.90
1	X	1031	C	P-O3'-C3'	7.44	128.63	119.70
1	X	1289	A	P-O5'-C5'	7.44	132.81	120.90
2	Y	11	G	O4'-C1'-N9	7.44	114.15	108.20
1	X	2867	G	C4-C5-N7	7.44	113.78	110.80
1	X	818	G	O4'-C1'-C2'	-7.44	98.36	105.80
1	X	1150	C	P-O3'-C3'	7.44	128.63	119.70
1	X	2274	C	O4'-C1'-N1	7.43	114.15	108.20
1	X	540	G	P-O3'-C3'	7.41	128.59	119.70
1	X	2330	G	P-O3'-C3'	7.41	128.59	119.70
1	X	2668	U	N3-C2-O2	-7.40	117.02	122.20
1	X	725	C	O4'-C1'-N1	7.40	114.12	108.20
1	X	1071	U	P-O3'-C3'	7.40	128.58	119.70
1	X	19	C	O4'-C1'-N1	7.40	114.12	108.20
1	X	579	G	C5-C6-O6	7.40	133.04	128.60
1	X	577	U	N3-C4-C5	-7.39	110.16	114.60
1	X	841	G	N7-C8-N9	7.39	116.80	113.10
1	X	126	C	O4'-C1'-N1	7.39	114.11	108.20
1	X	2347	C	O4'-C1'-N1	7.38	114.10	108.20
1	X	1830	C	P-O3'-C3'	7.37	128.55	119.70
1	X	199	A	P-O3'-C3'	7.36	128.54	119.70
1	X	921	A	P-O3'-C3'	7.36	128.53	119.70
1	X	1038	U	O4'-C1'-N1	7.35	114.08	108.20
1	X	1302	C	O4'-C1'-N1	7.35	114.08	108.20
1	X	2299	A	O4'-C1'-N9	7.35	114.08	108.20
1	X	2190	A	P-O3'-C3'	7.34	128.51	119.70
1	X	2573	C	O4'-C1'-N1	7.34	114.08	108.20
1	X	472	C	O4'-C1'-N1	7.34	114.07	108.20
1	X	788	G	C1'-O4'-C4'	-7.34	104.03	109.90
1	X	1496	G	C3'-C2'-C1'	-7.34	95.63	101.50
2	Y	75	A	P-O3'-C3'	7.34	128.51	119.70
1	X	1137	A	P-O3'-C3'	7.33	128.50	119.70
1	X	2876	C	O4'-C1'-N1	7.33	114.06	108.20
1	X	1991	C	N3-C4-C5	7.32	124.83	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1976	U	O4'-C1'-N1	7.32	114.06	108.20
1	X	2541	U	O4'-C1'-N1	7.32	114.05	108.20
1	X	2841	U	O4'-C1'-N1	7.31	114.05	108.20
1	X	1804	U	O4'-C1'-N1	7.31	114.05	108.20
1	X	541	C	P-O3'-C3'	7.30	128.47	119.70
1	X	2615	U	O4'-C1'-N1	7.30	114.04	108.20
1	X	2814	G	O4'-C1'-N9	7.30	114.04	108.20
1	X	2824	C	P-O3'-C3'	7.30	128.46	119.70
1	X	1627	C	O4'-C1'-N1	7.30	114.04	108.20
1	X	2408	G	C8-N9-C4	-7.30	103.48	106.40
1	X	2697	G	C2-N3-C4	7.29	115.55	111.90
1	X	331	U	O4'-C1'-N1	7.29	114.03	108.20
1	X	559	C	C5'-C4'-O4'	7.28	117.84	109.10
1	X	2222	U	O4'-C1'-N1	7.28	114.03	108.20
1	X	802	A	P-O3'-C3'	7.28	128.43	119.70
1	X	2587	G	C5-C6-O6	-7.27	124.24	128.60
1	X	114	C	O4'-C1'-N1	7.27	114.01	108.20
1	X	959	C	P-O3'-C3'	-7.27	110.98	119.70
1	X	2529	G	O5'-P-OP2	-7.26	99.17	105.70
1	X	2860	C	O4'-C1'-N1	7.26	114.01	108.20
1	X	1153	A	P-O5'-C5'	7.26	132.51	120.90
2	Y	32	C	C6-N1-C2	-7.26	117.40	120.30
1	X	2854	G	N7-C8-N9	7.25	116.72	113.10
1	X	774	A	C6-N1-C2	7.25	122.95	118.60
1	X	978	U	O4'-C1'-N1	7.24	113.99	108.20
1	X	343	A	P-O5'-C5'	7.24	132.48	120.90
1	X	1489	C	P-O3'-C3'	7.23	128.38	119.70
1	X	1731	C	O4'-C1'-N1	7.23	113.98	108.20
1	X	946	U	O4'-C1'-N1	7.22	113.98	108.20
1	X	711	C	O4'-C1'-N1	7.21	113.97	108.20
1	X	558	G	N9-C1'-C2'	7.21	123.37	114.00
1	X	2290	A	P-O3'-C3'	7.21	128.35	119.70
1	X	696	U	O4'-C1'-N1	7.20	113.96	108.20
1	X	610	G	C3'-C2'-C1'	-7.20	95.74	101.50
2	Y	52	G	P-O5'-C5'	7.20	132.42	120.90
1	X	870	C	O4'-C1'-N1	7.20	113.96	108.20
1	X	358	C	O4'-C1'-N1	7.20	113.96	108.20
1	X	2472	U	O4'-C1'-N1	7.19	113.95	108.20
1	X	2015	G	N9-C1'-C2'	7.19	123.35	114.00
1	X	2343	C	O4'-C1'-N1	7.19	113.95	108.20
1	X	2790	C	P-O3'-C3'	-7.19	111.08	119.70
1	X	853	C	C3'-C2'-C1'	-7.18	95.75	101.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1985	G	O4'-C1'-N9	7.18	113.94	108.20
1	X	1505	U	P-O3'-C3'	7.18	128.31	119.70
1	X	691	C	O4'-C1'-N1	7.18	113.94	108.20
1	X	2019	C	O4'-C1'-N1	7.18	113.94	108.20
1	X	2691	C	O3'-P-O5'	-7.17	90.37	104.00
1	X	1986	G	O4'-C4'-C3'	-7.17	96.83	104.00
1	X	1202	U	O4'-C1'-N1	7.17	113.93	108.20
1	X	542	A	C2-N3-C4	-7.17	107.02	110.60
2	Y	81	C	O4'-C1'-N1	7.17	113.93	108.20
1	X	117	A	C1'-O4'-C4'	-7.16	104.17	109.90
1	X	1800	A	P-O3'-C3'	7.16	128.29	119.70
1	X	2587	G	O4'-C1'-N9	7.15	113.92	108.20
1	X	751	G	C2'-C3'-O3'	7.14	125.22	109.50
1	X	617	U	O4'-C1'-N1	7.14	113.91	108.20
1	X	2032	G	C5-C6-O6	-7.14	124.32	128.60
1	X	2400	G	P-O3'-C3'	7.14	128.26	119.70
1	X	738	G	N7-C8-N9	7.13	116.67	113.10
1	X	2735	C	O4'-C1'-N1	7.13	113.91	108.20
1	X	1633	C	O4'-C1'-N1	7.13	113.90	108.20
1	X	1339	U	OP2-P-O3'	7.12	120.88	105.20
1	X	796	A	N7-C8-N9	7.12	117.36	113.80
1	X	955	G	N9-C1'-C2'	7.12	123.25	114.00
1	X	2062	U	O4'-C1'-N1	7.12	113.89	108.20
1	X	2544	A	O4'-C1'-N9	7.11	113.89	108.20
1	X	689	A	O4'-C1'-N9	7.11	113.88	108.20
1	X	2270	U	O4'-C1'-N1	7.11	113.88	108.20
1	X	2314	A	O4'-C1'-N9	7.10	113.88	108.20
1	X	2414	A	P-O3'-C3'	7.10	128.22	119.70
1	X	645	G	P-O3'-C3'	7.10	128.22	119.70
1	X	2567	G	C6-N1-C2	-7.10	120.84	125.10
1	X	2258	G	C4'-C3'-C2'	-7.09	95.51	102.60
1	X	312	G	C1'-O4'-C4'	-7.09	104.23	109.90
1	X	1708	C	O4'-C1'-N1	7.09	113.87	108.20
1	X	2429	A	P-O3'-C3'	-7.09	111.19	119.70
1	X	2564	U	N1-C1'-C2'	7.09	123.22	114.00
1	X	1634	A	C4'-C3'-C2'	7.09	109.69	102.60
1	X	1656	U	O4'-C1'-N1	7.08	113.87	108.20
1	X	1357	U	C1'-O4'-C4'	-7.08	104.24	109.90
1	X	1161	U	O4'-C1'-N1	7.08	113.86	108.20
1	X	1914	U	O4'-C1'-N1	7.08	113.86	108.20
1	X	2705	A	C4'-C3'-C2'	7.08	109.67	102.60
1	X	2372	A	O4'-C1'-N9	7.07	113.86	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1679	U	O4'-C4'-C3'	-7.07	96.93	104.00
1	X	2050	G	P-O3'-C3'	7.07	128.19	119.70
1	X	1621	C	O4'-C1'-N1	7.07	113.85	108.20
1	X	1278	A	N1-C2-N3	7.06	132.83	129.30
1	X	1602	G	P-O3'-C3'	7.06	128.17	119.70
1	X	537	C	P-O3'-C3'	7.05	128.16	119.70
1	X	966	A	P-O3'-C3'	-7.05	111.24	119.70
1	X	483	A	C5'-C4'-O4'	7.05	117.56	109.10
1	X	1412	C	P-O3'-C3'	7.05	128.16	119.70
1	X	1689	U	O4'-C1'-N1	7.04	113.84	108.20
1	X	2478	C	P-O3'-C3'	-7.04	111.25	119.70
1	X	2305	C	P-O3'-C3'	7.04	128.15	119.70
1	X	2018	G	C1'-O4'-C4'	-7.04	104.27	109.90
1	X	1775	A	C4'-C3'-C2'	7.03	109.63	102.60
1	X	59	G	P-O3'-C3'	7.02	128.12	119.70
1	X	683	A	N9-C1'-C2'	7.02	123.13	114.00
1	X	814	G	N9-C1'-C2'	7.02	123.12	114.00
1	X	1467	U	N1-C2-N3	-7.01	110.69	114.90
1	X	393	U	O4'-C1'-N1	7.01	113.81	108.20
1	X	1974	U	O4'-C1'-N1	7.01	113.81	108.20
1	X	2620	G	C4'-C3'-C2'	-7.00	95.60	102.60
1	X	558	G	O4'-C1'-N9	7.00	113.80	108.20
1	X	652	C	P-O5'-C5'	-7.00	109.70	120.90
1	X	1745	C	O4'-C1'-N1	7.00	113.80	108.20
1	X	224	G	C3'-C2'-C1'	7.00	107.10	101.50
1	X	739	G	O4'-C1'-N9	7.00	113.80	108.20
1	X	813	A	P-O3'-C3'	6.99	128.09	119.70
1	X	2311	U	O4'-C1'-N1	6.99	113.79	108.20
1	X	587	A	O4'-C1'-N9	6.99	113.79	108.20
1	X	480	G	C4-C5-N7	6.98	113.59	110.80
1	X	2845	C	O4'-C1'-N1	6.98	113.78	108.20
1	X	575	U	O4'-C1'-N1	6.97	113.78	108.20
1	X	2267	A	P-O3'-C3'	6.97	128.06	119.70
1	X	1963	G	C2'-C3'-O3'	6.96	124.83	113.70
1	X	2478	C	O4'-C1'-N1	6.95	113.76	108.20
1	X	2857	C	O4'-C1'-N1	6.95	113.76	108.20
1	X	2478	C	C6-N1-C2	-6.95	117.52	120.30
1	X	698	A	P-O3'-C3'	6.94	128.03	119.70
1	X	1678	G	O4'-C4'-C3'	-6.93	97.07	104.00
1	X	2291	U	O4'-C1'-N1	6.93	113.74	108.20
1	X	2393	G	O4'-C1'-N9	6.93	113.74	108.20
1	X	2820	C	O4'-C1'-N1	6.92	113.74	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1172	U	O4'-C1'-N1	6.92	113.74	108.20
1	X	2561	G	O4'-C1'-N9	6.92	113.73	108.20
1	X	646	C	O4'-C1'-N1	6.91	113.73	108.20
1	X	422	C	O4'-C1'-N1	6.91	113.72	108.20
1	X	1668	G	P-O3'-C3'	-6.90	111.42	119.70
1	X	2322	U	O4'-C1'-N1	6.90	113.72	108.20
1	X	1489	C	N1-C2-O2	6.89	123.04	118.90
1	X	1980	A	C5'-C4'-O4'	6.89	117.37	109.10
1	X	2366	U	O4'-C1'-N1	6.89	113.71	108.20
1	X	859	U	O4'-C1'-N1	6.88	113.71	108.20
1	X	1151	U	O4'-C1'-N1	6.88	113.70	108.20
1	X	1250	A	C5'-C4'-O4'	6.88	117.36	109.10
2	Y	34	C	O4'-C1'-N1	6.88	113.70	108.20
1	X	1655	C	O4'-C1'-N1	6.87	113.70	108.20
1	X	2208	U	O4'-C1'-N1	6.87	113.70	108.20
1	X	2622	G	C5-C6-O6	-6.86	124.48	128.60
1	X	305	A	O4'-C1'-N9	6.85	113.68	108.20
1	X	526	C	O4'-C1'-N1	6.85	113.68	108.20
1	X	2329	C	O4'-C1'-N1	6.83	113.67	108.20
1	X	1663	C	OP1-P-O3'	6.83	120.23	105.20
1	X	2854	G	O4'-C1'-N9	6.83	113.67	108.20
2	Y	123	U	C2-N1-C1'	6.83	125.89	117.70
1	X	968	C	N1-C2-O2	6.83	123.00	118.90
1	X	1980	A	O4'-C4'-C3'	-6.83	97.17	104.00
1	X	675	C	O4'-C1'-N1	6.82	113.66	108.20
1	X	2605	C	O4'-C1'-N1	6.82	113.66	108.20
1	X	761	G	O5'-P-OP2	-6.81	99.57	105.70
1	X	1279	G	C3'-C2'-C1'	-6.81	96.05	101.50
1	X	1726	C	O4'-C1'-N1	6.80	113.64	108.20
1	X	763	A	P-O3'-C3'	6.80	127.86	119.70
1	X	2080	U	O4'-C1'-N1	6.80	113.64	108.20
1	X	1343	C	O4'-C1'-N1	6.80	113.64	108.20
1	X	2808	U	P-O5'-C5'	6.79	131.77	120.90
1	X	542	A	C3'-C2'-C1'	6.79	106.93	101.50
1	X	1593	C	O4'-C1'-N1	6.79	113.63	108.20
1	X	1770	U	N3-C2-O2	-6.79	117.45	122.20
1	X	2258	G	O4'-C1'-N9	6.79	113.63	108.20
1	X	559	C	N3-C2-O2	-6.78	117.15	121.90
1	X	2303	C	P-O3'-C3'	6.78	127.83	119.70
1	X	2039	G	C8-N9-C4	-6.78	103.69	106.40
1	X	92	U	O4'-C1'-N1	6.77	113.61	108.20
1	X	321	A	P-O3'-C3'	6.77	127.82	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2668	U	N3-C4-O4	-6.76	114.67	119.40
1	X	1882	G	P-O3'-C3'	6.76	127.81	119.70
1	X	689	A	C4-C5-N7	6.75	114.08	110.70
19	Q	61	LYS	N-CA-C	6.75	129.23	111.00
1	X	1218	C	O4'-C1'-N1	6.75	113.60	108.20
1	X	1820	G	C4'-C3'-C2'	6.75	109.35	102.60
1	X	2627	G	C5-C6-O6	-6.75	124.55	128.60
1	X	540	G	N9-C4-C5	6.74	108.09	105.40
1	X	1067	G	P-O3'-C3'	6.74	127.78	119.70
1	X	117	A	P-O3'-C3'	6.73	127.78	119.70
1	X	2032	G	O4'-C4'-C3'	-6.73	97.27	104.00
1	X	2771	C	O4'-C1'-N1	6.73	113.58	108.20
1	X	2482	A	O5'-P-OP1	6.73	118.77	110.70
1	X	322	A	P-O3'-C3'	6.73	127.77	119.70
1	X	1992	G	C5'-C4'-O4'	-6.72	101.04	109.10
1	X	2560	G	C5'-C4'-O4'	6.71	117.16	109.10
1	X	2206	C	O4'-C1'-N1	6.71	113.57	108.20
1	X	466	A	P-O3'-C3'	6.71	127.75	119.70
1	X	1001	A	O4'-C1'-N9	6.71	113.57	108.20
1	X	1631	C	O4'-C1'-N1	6.71	113.57	108.20
1	X	1289	A	O4'-C4'-C3'	-6.71	97.29	104.00
1	X	1530	U	O4'-C1'-N1	6.71	113.57	108.20
1	X	981	C	O4'-C1'-N1	6.71	113.56	108.20
1	X	1679	U	N3-C2-O2	-6.71	117.51	122.20
1	X	2018	G	O5'-P-OP1	-6.71	99.67	105.70
1	X	1340	C	O4'-C1'-N1	6.70	113.56	108.20
1	X	2047	C	O4'-C1'-N1	6.70	113.56	108.20
1	X	2708	U	N3-C2-O2	-6.70	117.51	122.20
1	X	1250	A	O5'-P-OP1	6.69	118.73	110.70
1	X	2476	A	P-O3'-C3'	6.69	127.73	119.70
1	X	2534	U	C1'-O4'-C4'	-6.69	104.55	109.90
1	X	527	C	N1-C2-O2	6.69	122.91	118.90
1	X	387	A	C5'-C4'-O4'	6.68	117.12	109.10
1	X	649	G	O4'-C1'-N9	6.68	113.55	108.20
1	X	313	U	O4'-C1'-N1	6.68	113.54	108.20
1	X	1032	A	C3'-C2'-C1'	-6.68	96.16	101.50
1	X	1439	G	C2'-C3'-O3'	6.68	124.39	113.70
1	X	841	G	O4'-C1'-N9	6.67	113.54	108.20
1	X	2634	G	C1'-O4'-C4'	-6.67	104.56	109.90
1	X	456	C	O4'-C1'-N1	6.67	113.54	108.20
1	X	811	G	O4'-C1'-N9	6.67	113.54	108.20
1	X	796	A	N1-C6-N6	6.67	122.60	118.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1404	C	C1'-O4'-C4'	-6.67	104.56	109.90
1	X	1324	G	O4'-C1'-N9	6.67	113.53	108.20
1	X	1562	G	O4'-C1'-N9	6.67	113.53	108.20
1	X	1771	A	P-O3'-C3'	6.67	127.70	119.70
1	X	1762	C	O4'-C1'-N1	6.66	113.53	108.20
1	X	1716	G	C4'-C3'-C2'	6.66	109.26	102.60
1	X	1319	C	C6-N1-C2	-6.66	117.64	120.30
1	X	13	A	P-O3'-C3'	6.66	127.69	119.70
1	X	31	C	O4'-C1'-N1	6.66	113.52	108.20
1	X	2776	U	P-O3'-C3'	6.66	127.69	119.70
1	X	1469	U	C5'-C4'-O4'	6.65	117.08	109.10
1	X	789	G	P-O3'-C3'	6.65	127.68	119.70
1	X	2441	U	O4'-C1'-N1	6.65	113.52	108.20
1	X	12	U	N1-C2-O2	6.65	127.45	122.80
9	G	106	TYR	N-CA-CB	6.64	122.56	110.60
1	X	2437	G	O4'-C1'-N9	6.64	113.51	108.20
1	X	2794	G	P-O3'-C3'	6.64	127.67	119.70
1	X	1339	U	P-O3'-C3'	6.64	127.67	119.70
1	X	1434	U	C1'-O4'-C4'	-6.64	104.59	109.90
1	X	30	G	C8-N9-C4	-6.64	103.75	106.40
1	X	1470	G	P-O5'-C5'	-6.63	110.29	120.90
1	X	2033	C	P-O3'-C3'	6.63	127.65	119.70
1	X	1712	G	N3-C4-N9	6.62	129.97	126.00
2	Y	72	C	O4'-C1'-N1	6.62	113.50	108.20
1	X	485	G	P-O5'-C5'	6.62	131.49	120.90
1	X	1792	C	N1-C2-O2	6.62	122.87	118.90
1	X	2588	U	O4'-C1'-N1	6.62	113.49	108.20
1	X	2409	A	P-O3'-C3'	6.61	127.64	119.70
1	X	1468	A	N9-C1'-C2'	6.61	122.59	114.00
1	X	1622	G	P-O3'-C3'	6.61	127.63	119.70
1	X	2539	C	O4'-C1'-N1	6.61	113.49	108.20
1	X	1466	C	O4'-C1'-N1	6.61	113.49	108.20
2	Y	9	G	C3'-C2'-C1'	-6.61	96.21	101.50
1	X	700	C	O4'-C1'-N1	6.61	113.48	108.20
1	X	100	G	O4'-C1'-N9	6.60	113.48	108.20
1	X	1249	G	O4'-C1'-N9	6.60	113.48	108.20
1	X	2315	A	P-O5'-C5'	6.60	131.47	120.90
1	X	788	G	O4'-C1'-N9	6.60	113.48	108.20
1	X	955	G	P-O5'-C5'	-6.60	110.34	120.90
1	X	2490	U	O4'-C1'-N1	6.60	113.48	108.20
1	X	1805	G	O4'-C1'-N9	6.59	113.47	108.20
1	X	2804	G	C5-C6-O6	-6.59	124.64	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1469	U	N3-C2-O2	-6.59	117.59	122.20
1	X	1711	C	C1'-O4'-C4'	-6.59	104.63	109.90
1	X	2867	G	N7-C8-N9	6.58	116.39	113.10
1	X	2338	C	O4'-C1'-N1	6.58	113.46	108.20
1	X	2719	U	P-O5'-C5'	6.58	131.42	120.90
1	X	794	A	O4'-C1'-N9	-6.57	102.95	108.20
1	X	2669	C	N3-C2-O2	-6.57	117.30	121.90
1	X	2766	U	O4'-C1'-N1	6.57	113.45	108.20
1	X	2846	G	P-O3'-C3'	6.57	127.58	119.70
1	X	467	U	N3-C2-O2	-6.56	117.61	122.20
1	X	540	G	N9-C1'-C2'	6.56	122.52	114.00
1	X	1446	U	O4'-C1'-N1	6.55	113.44	108.20
1	X	1694	A	O4'-C1'-N9	6.55	113.44	108.20
1	X	308	C	P-O3'-C3'	-6.55	111.84	119.70
1	X	429	C	O4'-C1'-N1	6.55	113.44	108.20
1	X	2081	U	O4'-C1'-N1	6.55	113.44	108.20
1	X	1359	G	N7-C8-N9	6.55	116.37	113.10
1	X	1690	U	O4'-C1'-N1	6.55	113.44	108.20
1	X	399	G	C4'-C3'-C2'	6.54	109.14	102.60
1	X	440	U	P-O3'-C3'	6.54	127.55	119.70
1	X	1652	G	O4'-C1'-N9	-6.54	102.97	108.20
1	X	967	G	P-O5'-C5'	6.54	131.36	120.90
1	X	1647	U	O4'-C1'-N1	6.54	113.43	108.20
1	X	2406	C	O4'-C1'-N1	6.54	113.43	108.20
1	X	2481	G	O3'-P-O5'	-6.54	91.58	104.00
1	X	322	A	O4'-C1'-N9	6.54	113.43	108.20
1	X	2691	C	O4'-C1'-C2'	-6.53	99.27	105.80
1	X	517	A	P-O3'-C3'	6.53	127.53	119.70
23	U	18	VAL	C-N-CA	6.53	138.02	121.70
1	X	1776	A	P-O3'-C3'	6.52	127.53	119.70
1	X	2669	C	O4'-C1'-C2'	-6.52	99.28	105.80
1	X	513	A	P-O3'-C3'	6.52	127.52	119.70
1	X	784	U	O4'-C1'-N1	6.52	113.42	108.20
1	X	2417	U	N3-C2-O2	-6.51	117.64	122.20
1	X	2561	G	C5'-C4'-O4'	-6.51	101.28	109.10
1	X	2437	G	N3-C4-C5	-6.51	125.34	128.60
1	X	241	C	O4'-C4'-C3'	-6.51	97.49	104.00
1	X	242	A	C4'-C3'-C2'	-6.51	96.09	102.60
1	X	1439	G	C3'-C2'-C1'	-6.51	96.29	101.50
1	X	859	U	C5'-C4'-O4'	6.51	116.91	109.10
1	X	1994	U	O4'-C1'-N1	6.51	113.41	108.20
1	X	2384	G	P-O3'-C3'	6.50	127.50	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1749	G	P-O3'-C3'	6.50	127.50	119.70
1	X	320	A	P-O3'-C3'	6.49	127.49	119.70
1	X	424	G	P-O3'-C3'	6.49	127.48	119.70
1	X	631	G	P-O3'-C3'	6.49	127.49	119.70
1	X	1044	U	P-O3'-C3'	6.49	127.49	119.70
1	X	1912	G	P-O3'-C3'	6.49	127.48	119.70
1	X	543	G	C3'-C2'-C1'	-6.49	96.31	101.50
2	Y	47	A	N7-C8-N9	6.49	117.04	113.80
1	X	467	U	N1-C2-O2	6.47	127.33	122.80
1	X	1170	U	O4'-C1'-N1	6.47	113.37	108.20
1	X	2591	C	C2-N3-C4	6.47	123.13	119.90
1	X	2488	G	C5-C6-N1	6.46	114.73	111.50
1	X	312	G	P-O3'-C3'	6.46	127.45	119.70
1	X	2486	C	N1-C2-O2	6.46	122.78	118.90
1	X	1670	G	P-O3'-C3'	6.46	127.45	119.70
1	X	2852	G	O4'-C4'-C3'	-6.46	97.54	104.00
1	X	1925	C	O4'-C1'-N1	6.46	113.36	108.20
1	X	2185	U	O4'-C1'-N1	6.46	113.36	108.20
1	X	580	A	P-O5'-C5'	-6.45	110.58	120.90
1	X	2556	A	P-O3'-C3'	6.45	127.44	119.70
1	X	453	U	O4'-C1'-N1	6.45	113.36	108.20
1	X	2256	G	C8-N9-C4	-6.45	103.82	106.40
1	X	307	C	O4'-C1'-N1	6.44	113.36	108.20
1	X	1200	G	P-O5'-C5'	6.44	131.21	120.90
1	X	2039	G	C5-C6-O6	-6.44	124.73	128.60
1	X	1758	C	O4'-C1'-N1	6.43	113.35	108.20
1	X	1732	U	O4'-C1'-N1	6.43	113.34	108.20
1	X	460	U	P-O3'-C3'	6.43	127.41	119.70
1	X	1685	A	OP1-P-O3'	6.43	119.34	105.20
1	X	469	G	C1'-O4'-C4'	-6.42	104.76	109.90
1	X	2779	C	P-O3'-C3'	6.42	127.41	119.70
1	X	1667	A	O4'-C1'-N9	6.42	113.34	108.20
1	X	608	G	O4'-C4'-C3'	-6.42	97.58	104.00
1	X	1468	A	C5-C6-N1	6.42	120.91	117.70
1	X	2659	C	O4'-C1'-N1	6.42	113.34	108.20
1	X	661	C	C4'-C3'-C2'	-6.42	96.18	102.60
1	X	925	U	P-O3'-C3'	6.41	127.39	119.70
1	X	2044	G	N3-C4-N9	6.41	129.85	126.00
1	X	330	C	O4'-C1'-N1	6.41	113.33	108.20
1	X	1922	U	N3-C2-O2	-6.41	117.71	122.20
1	X	1037	U	C1'-O4'-C4'	-6.41	104.78	109.90
1	X	1031	C	O4'-C1'-N1	6.41	113.33	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2667	C	N3-C2-O2	-6.41	117.42	121.90
1	X	2581	A	P-O3'-C3'	6.40	127.39	119.70
1	X	338	G	O4'-C1'-N9	6.40	113.32	108.20
1	X	458	G	P-O3'-C3'	6.40	127.38	119.70
1	X	1669	A	O4'-C4'-C3'	-6.39	97.61	104.00
1	X	308	C	P-O5'-C5'	6.38	131.11	120.90
1	X	595	A	P-O3'-C3'	6.38	127.36	119.70
1	X	2246	A	N9-C1'-C2'	6.38	122.30	114.00
2	Y	110	U	N3-C2-O2	-6.38	117.73	122.20
19	Q	62	ARG	C-N-CA	6.38	137.66	121.70
1	X	2527	G	O4'-C1'-N9	6.38	113.30	108.20
1	X	2551	A	O3'-P-O5'	-6.38	91.88	104.00
1	X	2691	C	P-O3'-C3'	6.38	127.35	119.70
1	X	1696	C	O4'-C1'-N1	6.37	113.30	108.20
1	X	56	C	O4'-C1'-N1	6.37	113.29	108.20
1	X	1467	U	C4'-C3'-O3'	6.37	125.73	113.00
1	X	2678	C	O4'-C1'-N1	6.37	113.29	108.20
2	Y	6	C	O4'-C1'-N1	6.37	113.29	108.20
1	X	1429	A	C1'-O4'-C4'	-6.36	104.81	109.90
1	X	645	G	O4'-C1'-N9	6.36	113.29	108.20
1	X	1339	U	O4'-C1'-N1	6.36	113.29	108.20
2	Y	53	G	N3-C4-C5	-6.36	125.42	128.60
1	X	2774	U	P-O3'-C3'	6.36	127.33	119.70
1	X	1885	C	N1-C2-O2	6.35	122.71	118.90
1	X	325	U	O4'-C1'-N1	6.35	113.28	108.20
1	X	1679	U	N3-C4-C5	6.35	118.41	114.60
1	X	990	A	C3'-C2'-C1'	-6.35	96.42	101.50
1	X	2523	G	O4'-C1'-N9	6.35	113.28	108.20
1	X	18	U	O4'-C1'-N1	6.35	113.28	108.20
1	X	1979	C	P-O3'-C3'	6.34	127.31	119.70
1	X	618	A	O4'-C1'-N9	6.34	113.27	108.20
2	Y	86	A	C1'-O4'-C4'	-6.34	104.83	109.90
1	X	742	G	C1'-O4'-C4'	-6.34	104.83	109.90
1	X	1882	G	C3'-C2'-C1'	6.34	106.57	101.50
1	X	1664	G	O5'-P-OP1	-6.34	100.00	105.70
1	X	689	A	N1-C6-N6	6.34	122.40	118.60
1	X	593	C	O4'-C1'-N1	6.33	113.26	108.20
1	X	632	A	P-O3'-C3'	6.32	127.29	119.70
1	X	2538	C	O4'-C1'-N1	6.32	113.26	108.20
1	X	2735	C	C6-N1-C2	-6.32	117.77	120.30
10	H	26	ASN	C-N-CA	6.32	137.50	121.70
1	X	1673	C	O4'-C1'-N1	6.31	113.25	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1017	C	O4'-C1'-N1	6.31	113.25	108.20
1	X	1649	A	C2-N3-C4	6.31	113.75	110.60
1	X	1664	G	N9-C1'-C2'	6.30	122.20	114.00
1	X	1716	G	C2'-C3'-O3'	6.30	123.79	113.70
1	X	2217	G	C1'-O4'-C4'	-6.30	104.86	109.90
1	X	566	U	O4'-C1'-N1	6.30	113.24	108.20
1	X	1023	U	O4'-C1'-N1	6.30	113.24	108.20
1	X	1671	A	OP1-P-OP2	6.30	129.05	119.60
1	X	2627	G	N1-C6-O6	6.30	123.68	119.90
1	X	2359	U	P-O3'-C3'	6.30	127.26	119.70
1	X	682	G	C5-C6-N1	6.29	114.65	111.50
1	X	884	C	O4'-C1'-N1	6.29	113.23	108.20
1	X	542	A	N7-C8-N9	6.29	116.94	113.80
1	X	1764	A	N1-C6-N6	6.29	122.37	118.60
2	Y	4	C	O4'-C1'-N1	6.29	113.23	108.20
1	X	2566	A	P-O3'-C3'	6.29	127.25	119.70
1	X	1749	G	C1'-O4'-C4'	-6.29	104.87	109.90
1	X	98	U	O4'-C1'-N1	6.28	113.22	108.20
1	X	467	U	C2-N1-C1'	6.28	125.24	117.70
1	X	2710	C	N1-C2-O2	6.28	122.67	118.90
1	X	1748	U	P-O3'-C3'	6.28	127.23	119.70
1	X	2419	C	O4'-C1'-N1	6.28	113.22	108.20
1	X	2485	U	N1-C2-O2	6.28	127.19	122.80
1	X	2668	U	C5-C6-N1	-6.27	119.56	122.70
1	X	430	C	O4'-C1'-N1	6.27	113.21	108.20
1	X	1142	G	O5'-C5'-C4'	6.27	123.61	111.70
1	X	117	A	O4'-C1'-N9	6.27	113.21	108.20
1	X	926	C	O4'-C1'-N1	6.26	113.21	108.20
1	X	1224	A	P-O3'-C3'	6.26	127.21	119.70
1	X	2593	A	O3'-P-O5'	-6.26	92.11	104.00
1	X	426	C	O4'-C1'-N1	6.26	113.20	108.20
1	X	2760	G	P-O3'-C3'	6.26	127.21	119.70
1	X	2811	G	O4'-C1'-N9	6.25	113.20	108.20
1	X	2582	G	P-O5'-C5'	6.25	130.91	120.90
1	X	2782	G	C6-C5-N7	-6.25	126.65	130.40
1	X	609	U	C3'-C2'-C1'	-6.24	96.50	101.50
1	X	941	U	O4'-C1'-N1	6.24	113.19	108.20
1	X	2875	C	O4'-C1'-N1	6.24	113.19	108.20
1	X	780	U	O4'-C1'-N1	6.24	113.19	108.20
1	X	1558	C	N1-C2-O2	6.24	122.64	118.90
1	X	2645	C	O4'-C1'-N1	6.24	113.19	108.20
1	X	656	U	P-O5'-C5'	6.24	130.88	120.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1468	A	N3-C4-C5	-6.24	122.44	126.80
1	X	1222	G	N3-C4-C5	-6.23	125.48	128.60
1	X	2745	A	P-O3'-C3'	6.23	127.18	119.70
1	X	242	A	P-O3'-C3'	6.23	127.18	119.70
1	X	2506	C	O4'-C1'-N1	6.23	113.19	108.20
1	X	88	G	P-O3'-C3'	6.23	127.17	119.70
1	X	346	C	N1-C1'-C2'	6.23	122.10	114.00
1	X	431	G	O4'-C1'-N9	6.22	113.18	108.20
1	X	917	U	O4'-C1'-N1	6.22	113.18	108.20
1	X	2559	U	C5-C4-O4	-6.22	122.17	125.90
1	X	2359	U	O4'-C1'-N1	6.22	113.17	108.20
1	X	113	C	O4'-C1'-N1	6.21	113.17	108.20
1	X	632	A	C1'-O4'-C4'	-6.21	104.93	109.90
1	X	604	U	O4'-C1'-N1	6.21	113.17	108.20
1	X	1487	C	O4'-C1'-N1	6.21	113.17	108.20
1	X	1551	U	O4'-C1'-N1	6.21	113.17	108.20
1	X	221	A	O4'-C1'-N9	6.21	113.17	108.20
1	X	1746	A	O4'-C1'-N9	6.21	113.17	108.20
1	X	2665	G	C5-C6-O6	-6.21	124.88	128.60
1	X	1652	G	C4-C5-N7	6.21	113.28	110.80
1	X	1675	C	O5'-P-OP1	-6.21	100.11	105.70
1	X	2239	C	O4'-C1'-N1	6.20	113.16	108.20
1	X	526	C	C3'-C2'-C1'	-6.20	96.54	101.50
1	X	2492	G	C3'-C2'-C1'	-6.20	96.54	101.50
2	Y	62	C	O4'-C1'-N1	6.20	113.16	108.20
1	X	577	U	O4'-C1'-N1	6.19	113.15	108.20
1	X	1338	G	N3-C4-C5	-6.19	125.51	128.60
1	X	1016	C	O4'-C1'-N1	6.18	113.15	108.20
1	X	580	A	C5'-C4'-O4'	6.18	116.52	109.10
1	X	2459	C	O4'-C1'-N1	6.18	113.14	108.20
1	X	626	A	P-O3'-C3'	6.18	127.12	119.70
1	X	665	A	O4'-C1'-N9	6.18	113.14	108.20
1	X	1455	C	O4'-C1'-N1	6.17	113.14	108.20
1	X	1279	G	C1'-O4'-C4'	-6.17	104.97	109.90
1	X	2043	A	P-O5'-C5'	-6.17	111.03	120.90
1	X	97	U	O4'-C1'-N1	6.17	113.13	108.20
1	X	175	C	C5-C6-N1	6.17	124.08	121.00
1	X	1328	C	O4'-C1'-N1	6.16	113.13	108.20
1	X	1679	U	N3-C4-O4	-6.16	115.09	119.40
1	X	1865	C	O4'-C1'-N1	6.16	113.13	108.20
1	X	2697	G	P-O3'-C3'	-6.16	112.31	119.70
1	X	1434	U	N1-C1'-C2'	6.16	122.01	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2708	U	O4'-C1'-N1	6.16	113.13	108.20
2	Y	92	G	C3'-C2'-C1'	-6.16	96.57	101.50
1	X	90	G	P-O3'-C3'	6.16	127.09	119.70
1	X	677	G	C4'-C3'-C2'	-6.15	96.45	102.60
1	X	2478	C	C5-C6-N1	6.15	124.08	121.00
1	X	672	C	C3'-C2'-C1'	-6.15	96.58	101.50
1	X	1652	G	C5-C6-O6	-6.15	124.91	128.60
1	X	547	U	O4'-C1'-N1	6.15	113.12	108.20
1	X	774	A	C2-N3-C4	-6.15	107.53	110.60
1	X	1540	C	O4'-C1'-N1	6.15	113.12	108.20
1	X	345	U	O4'-C1'-N1	6.14	113.12	108.20
1	X	1269	G	O4'-C1'-N9	6.14	113.11	108.20
1	X	499	G	P-O3'-C3'	-6.14	112.33	119.70
1	X	1030	U	O4'-C1'-N1	6.14	113.11	108.20
1	X	1171	A	O4'-C1'-N9	6.14	113.11	108.20
1	X	2039	G	N3-C2-N2	-6.14	115.60	119.90
1	X	2821	G	O4'-C1'-N9	6.14	113.11	108.20
1	X	928	G	C5-C6-O6	-6.13	124.92	128.60
1	X	1820	G	P-O5'-C5'	6.13	130.71	120.90
1	X	416	U	O4'-C1'-N1	6.13	113.10	108.20
1	X	1950	C	O4'-C1'-N1	6.13	113.11	108.20
1	X	2246	A	C2-N3-C4	6.13	113.67	110.60
2	Y	24	U	O4'-C1'-N1	6.13	113.11	108.20
1	X	1788	C	O4'-C1'-N1	6.13	113.10	108.20
1	X	1938	U	N1-C1'-C2'	6.12	121.96	114.00
1	X	2840	U	P-O3'-C3'	6.12	127.05	119.70
1	X	483	A	O4'-C1'-N9	6.12	113.09	108.20
1	X	2038	C	OP2-P-O3'	6.11	118.65	105.20
1	X	2371	A	C3'-C2'-C1'	-6.11	96.61	101.50
1	X	1341	G	N3-C4-C5	-6.11	125.55	128.60
1	X	1792	C	N3-C2-O2	-6.11	117.62	121.90
1	X	2540	A	O4'-C1'-N9	6.11	113.09	108.20
1	X	2594	U	C5-C6-N1	6.11	125.75	122.70
1	X	343	A	N7-C8-N9	6.11	116.85	113.80
2	Y	13	C	O4'-C1'-N1	6.11	113.08	108.20
1	X	589	C	O4'-C1'-N1	6.10	113.08	108.20
1	X	1938	U	P-O5'-C5'	6.10	130.66	120.90
1	X	2363	G	C5'-C4'-O4'	6.10	116.42	109.10
1	X	2246	A	P-O3'-C3'	6.09	127.01	119.70
1	X	2431	C	O4'-C1'-N1	6.09	113.08	108.20
1	X	2691	C	C5'-C4'-C3'	-6.09	106.25	116.00
1	X	968	C	C5-C6-N1	6.09	124.05	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1336	G	C4-C5-N7	6.09	113.24	110.80
2	Y	123	U	N1-C1'-C2'	6.09	121.92	114.00
1	X	12	U	C2-N1-C1'	6.09	125.01	117.70
1	X	1412	C	O4'-C4'-C3'	-6.09	97.91	104.00
1	X	2558	C	O4'-C1'-N1	6.09	113.07	108.20
1	X	1993	G	O4'-C1'-N9	6.08	113.07	108.20
1	X	2489	C	O4'-C1'-N1	6.08	113.07	108.20
1	X	2581	A	O4'-C1'-N9	6.08	113.07	108.20
1	X	1086	C	O4'-C1'-N1	6.08	113.06	108.20
1	X	1631	C	C6-N1-C1'	-6.08	113.50	120.80
2	Y	83	C	N1-C2-O2	6.08	122.55	118.90
1	X	557	U	O4'-C1'-N1	6.07	113.06	108.20
2	Y	30	C	P-O5'-C5'	6.07	130.62	120.90
1	X	432	C	O4'-C1'-N1	6.07	113.06	108.20
1	X	480	G	C5-C6-N1	6.07	114.53	111.50
1	X	2230	G	C5-C6-O6	-6.07	124.96	128.60
1	X	2767	C	O4'-C1'-N1	6.07	113.06	108.20
1	X	1468	A	C2-N3-C4	6.07	113.63	110.60
1	X	2661	G	O4'-C1'-N9	6.07	113.05	108.20
1	X	2018	G	C5'-C4'-C3'	-6.07	106.29	116.00
1	X	1717	A	O4'-C1'-N9	6.06	113.05	108.20
1	X	2354	G	O4'-C4'-C3'	-6.06	97.94	104.00
1	X	74	G	C1'-O4'-C4'	-6.06	105.05	109.90
1	X	632	A	C4'-C3'-C2'	-6.06	96.54	102.60
1	X	660	G	N3-C2-N2	-6.06	115.66	119.90
1	X	1442	C	N1-C2-O2	6.05	122.53	118.90
2	Y	35	C	O4'-C1'-N1	6.05	113.04	108.20
1	X	537	C	C6-N1-C1'	-6.05	113.54	120.80
1	X	579	G	N9-C4-C5	6.05	107.82	105.40
1	X	1006	C	N1-C2-O2	6.05	122.53	118.90
2	Y	44	C	O4'-C1'-N1	6.05	113.04	108.20
2	Y	111	C	P-O3'-C3'	6.05	126.96	119.70
1	X	2189	A	O3'-P-O5'	6.05	115.49	104.00
1	X	2245	A	P-O3'-C3'	6.05	126.96	119.70
2	Y	32	C	O4'-C1'-N1	6.05	113.04	108.20
1	X	1981	A	C5-C6-N6	-6.05	118.86	123.70
1	X	455	A	P-O3'-C3'	6.04	126.95	119.70
1	X	63	A	C4'-C3'-C2'	6.04	108.64	102.60
1	X	956	A	C5-C6-N6	-6.04	118.87	123.70
1	X	769	C	O4'-C1'-N1	6.04	113.03	108.20
1	X	1105	U	O4'-C1'-N1	6.03	113.03	108.20
1	X	2484	G	N3-C2-N2	6.03	124.12	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	526	C	O5'-P-OP2	-6.03	100.28	105.70
1	X	236	C	O4'-C1'-N1	6.03	113.02	108.20
1	X	965	G	P-O3'-C3'	-6.03	112.47	119.70
1	X	1191	G	P-O3'-C3'	6.02	126.93	119.70
1	X	2689	C	P-O3'-C3'	6.02	126.93	119.70
1	X	1392	U	P-O3'-C3'	6.02	126.92	119.70
1	X	590	C	O4'-C1'-N1	6.02	113.01	108.20
1	X	1076	U	O4'-C1'-N1	6.02	113.02	108.20
1	X	1154	A	P-O3'-C3'	6.02	126.92	119.70
1	X	1829	C	O4'-C1'-N1	6.02	113.01	108.20
1	X	2645	C	N1-C2-O2	6.02	122.51	118.90
1	X	2854	G	C8-N9-C4	-6.02	103.99	106.40
1	X	652	C	C5'-C4'-C3'	-6.01	106.38	116.00
1	X	1942	G	O4'-C4'-C3'	-6.01	97.99	104.00
1	X	2417	U	O4'-C1'-N1	6.01	113.01	108.20
1	X	2560	G	N3-C4-C5	-6.01	125.59	128.60
1	X	1183	C	O4'-C1'-N1	6.01	113.01	108.20
2	Y	55	C	P-O3'-C3'	6.01	126.91	119.70
1	X	873	U	O4'-C1'-N1	6.01	113.01	108.20
1	X	2395	C	O4'-C1'-N1	6.01	113.01	108.20
1	X	2591	C	N1-C2-O2	6.01	122.50	118.90
1	X	2482	A	P-O5'-C5'	6.00	130.50	120.90
1	X	2695	C	O4'-C1'-N1	6.00	113.00	108.20
1	X	1277	G	N3-C4-C5	-6.00	125.60	128.60
1	X	2314	A	P-O3'-C3'	6.00	126.90	119.70
1	X	1645	U	N3-C2-O2	-6.00	118.00	122.20
1	X	2804	G	C5-C6-N1	6.00	114.50	111.50
1	X	1749	G	O4'-C1'-C2'	-5.99	99.81	105.80
1	X	2193	C	O4'-C1'-N1	5.99	112.99	108.20
1	X	2314	A	P-O5'-C5'	5.99	130.49	120.90
1	X	1268	U	P-O5'-C5'	5.99	130.48	120.90
1	X	2797	G	N3-C4-N9	5.99	129.59	126.00
1	X	2519	C	O4'-C1'-N1	5.98	112.99	108.20
1	X	2591	C	C5-C6-N1	5.98	123.99	121.00
1	X	304	A	P-O5'-C5'	5.98	130.47	120.90
1	X	646	C	C6-N1-C2	-5.98	117.91	120.30
1	X	1515	U	O4'-C1'-N1	5.97	112.98	108.20
1	X	621	U	O4'-C1'-N1	5.97	112.98	108.20
1	X	2258	G	C1'-O4'-C4'	-5.97	105.12	109.90
1	X	2600	A	O4'-C1'-N9	5.97	112.98	108.20
2	Y	14	C	P-O3'-C3'	5.97	126.87	119.70
1	X	611	C	O4'-C1'-N1	5.97	112.98	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	824	U	N1-C1'-C2'	5.97	121.76	114.00
1	X	2284	U	O4'-C1'-N1	5.97	112.97	108.20
1	X	2477	C	C6-N1-C2	-5.97	117.91	120.30
1	X	879	A	C5'-C4'-C3'	-5.97	106.45	116.00
1	X	1288	A	P-O3'-C3'	-5.97	112.54	119.70
1	X	826	U	O4'-C1'-N1	5.96	112.97	108.20
1	X	2435	C	N1-C2-O2	5.96	122.48	118.90
1	X	490	A	C5'-C4'-O4'	5.96	116.25	109.10
1	X	1223	G	N7-C8-N9	5.96	116.08	113.10
1	X	1111	C	O4'-C1'-N1	5.96	112.97	108.20
1	X	1746	A	N1-C6-N6	-5.96	115.03	118.60
1	X	2226	A	N1-C6-N6	5.96	122.17	118.60
1	X	30	G	O4'-C1'-N9	5.96	112.97	108.20
1	X	647	G	P-O3'-C3'	5.96	126.85	119.70
1	X	2791	C	O4'-C1'-N1	5.96	112.97	108.20
1	X	1496	G	C2'-C3'-O3'	5.96	123.23	113.70
4	B	132	LYS	C-N-CA	5.96	136.59	121.70
1	X	1396	C	O4'-C1'-N1	5.95	112.96	108.20
1	X	1680	U	O4'-C4'-C3'	-5.95	98.05	104.00
1	X	1787	U	O4'-C1'-N1	5.95	112.96	108.20
1	X	1778	U	O4'-C1'-N1	5.95	112.96	108.20
1	X	162	C	O4'-C1'-N1	5.95	112.96	108.20
1	X	1201	G	N3-C2-N2	-5.95	115.73	119.90
1	X	2392	G	O4'-C1'-N9	5.95	112.96	108.20
1	X	2782	G	O4'-C1'-N9	5.95	112.96	108.20
1	X	448	C	N1-C2-O2	5.95	122.47	118.90
1	X	505	G	N9-C1'-C2'	5.95	121.73	114.00
1	X	1711	C	P-O3'-C3'	5.95	126.83	119.70
1	X	609	U	O4'-C4'-C3'	-5.94	98.06	104.00
1	X	2511	G	P-O5'-C5'	5.94	130.40	120.90
1	X	2855	C	O4'-C1'-N1	5.94	112.95	108.20
1	X	1051	U	O4'-C1'-N1	5.94	112.95	108.20
1	X	437	G	O4'-C1'-N9	5.93	112.95	108.20
1	X	969	U	C4'-C3'-C2'	5.93	108.53	102.60
1	X	2567	G	N3-C4-C5	-5.93	125.63	128.60
1	X	2662	C	N1-C2-O2	5.93	122.46	118.90
1	X	483	A	C4'-C3'-C2'	5.93	108.53	102.60
1	X	397	U	O4'-C1'-N1	5.93	112.94	108.20
1	X	1272	G	O4'-C4'-C3'	-5.93	98.07	104.00
1	X	1221	C	O4'-C1'-N1	5.93	112.94	108.20
1	X	683	A	C2'-C3'-O3'	5.93	123.18	113.70
1	X	456	C	C5'-C4'-O4'	-5.92	102.00	109.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1688	U	N1-C2-O2	-5.92	118.65	122.80
1	X	582	G	OP2-P-O3'	5.92	118.22	105.20
1	X	2844	G	O4'-C1'-N9	5.92	112.94	108.20
1	X	2668	U	N1-C2-N3	5.92	118.45	114.90
1	X	2698	G	C4'-C3'-C2'	-5.92	96.68	102.60
1	X	980	G	O4'-C1'-N9	5.92	112.93	108.20
1	X	1411	C	O4'-C1'-N1	5.92	112.93	108.20
1	X	1706	A	P-O3'-C3'	5.92	126.80	119.70
1	X	1034	U	O4'-C1'-N1	5.91	112.93	108.20
1	X	2072	C	O4'-C1'-N1	5.91	112.93	108.20
1	X	78	C	O4'-C1'-N1	5.91	112.93	108.20
1	X	751	G	P-O3'-C3'	5.91	126.79	119.70
1	X	788	G	N9-C1'-C2'	5.91	121.68	114.00
1	X	1262	U	C5'-C4'-O4'	-5.91	102.01	109.10
1	X	1415	C	O4'-C1'-N1	5.91	112.93	108.20
1	X	2018	G	C4-C5-N7	5.91	113.16	110.80
1	X	959	C	C5-C6-N1	5.91	123.95	121.00
1	X	1349	A	P-O5'-C5'	5.90	130.34	120.90
2	Y	55	C	O4'-C1'-N1	5.90	112.92	108.20
1	X	1624	A	C1'-O4'-C4'	-5.90	105.18	109.90
2	Y	50	U	O4'-C1'-N1	5.90	112.92	108.20
13	K	93	GLY	C-N-CA	-5.90	106.96	121.70
1	X	1664	G	O4'-C1'-N9	-5.89	103.48	108.20
1	X	858	G	C3'-C2'-C1'	5.89	106.21	101.50
1	X	1099	A	P-O3'-C3'	5.89	126.77	119.70
1	X	1214	C	C3'-C2'-C1'	-5.89	96.79	101.50
1	X	1858	C	O4'-C1'-N1	5.89	112.91	108.20
1	X	2487	G	N9-C4-C5	5.89	107.76	105.40
1	X	2377	U	O4'-C1'-N1	5.89	112.91	108.20
17	O	6	GLN	C-N-CA	5.88	136.41	121.70
1	X	2681	A	C4'-C3'-C2'	-5.88	96.72	102.60
1	X	540	G	C3'-C2'-C1'	5.88	106.20	101.50
1	X	1252	C	O4'-C1'-N1	5.88	112.90	108.20
1	X	110	U	O4'-C1'-N1	5.88	112.90	108.20
1	X	1004	A	C2-N3-C4	5.88	113.54	110.60
1	X	1210	C	O4'-C1'-N1	5.88	112.90	108.20
1	X	1263	G	P-O3'-C3'	5.88	126.75	119.70
1	X	1858	C	N1-C2-O2	5.87	122.42	118.90
1	X	2782	G	C5-C6-O6	-5.87	125.08	128.60
1	X	111	G	C5'-C4'-C3'	-5.87	106.61	116.00
2	Y	57	U	O4'-C1'-N1	5.87	112.89	108.20
1	X	731	A	C3'-C2'-C1'	5.86	106.19	101.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2488	G	O4'-C1'-N9	5.86	112.89	108.20
1	X	2671	C	O5'-P-OP2	-5.86	100.42	105.70
2	Y	71	G	C8-N9-C4	-5.86	104.06	106.40
1	X	957	G	N3-C4-C5	-5.86	125.67	128.60
1	X	525	A	O4'-C1'-N9	5.86	112.89	108.20
1	X	469	G	O4'-C1'-C2'	-5.86	99.94	105.80
1	X	1522	C	N1-C2-O2	5.86	122.41	118.90
2	Y	110	U	O4'-C1'-N1	5.86	112.89	108.20
1	X	479	G	C5-C6-O6	-5.86	125.09	128.60
1	X	2082	C	O4'-C1'-N1	5.86	112.89	108.20
13	K	11	ASN	C-N-CA	5.86	136.34	121.70
1	X	2795	A	C3'-C2'-C1'	5.85	106.18	101.50
1	X	960	U	O4'-C1'-N1	5.85	112.88	108.20
1	X	1326	U	N1-C2-O2	5.85	126.89	122.80
1	X	2193	C	O4'-C4'-C3'	-5.85	98.15	104.00
1	X	68	C	N1-C2-O2	5.85	122.41	118.90
1	X	204	A	C2'-C3'-O3'	5.84	123.05	113.70
1	X	1015	U	O4'-C1'-N1	5.84	112.88	108.20
1	X	234	C	O4'-C1'-N1	5.84	112.88	108.20
1	X	2382	C	O4'-C1'-N1	5.84	112.87	108.20
1	X	2668	U	O4'-C1'-N1	5.84	112.87	108.20
1	X	559	C	C1'-O4'-C4'	-5.84	105.23	109.90
1	X	1128	G	P-O3'-C3'	5.84	126.70	119.70
1	X	1663	C	P-O3'-C3'	5.84	126.70	119.70
1	X	2236	U	O4'-C1'-N1	5.84	112.87	108.20
2	Y	53	G	O4'-C1'-N9	5.83	112.87	108.20
1	X	1333	G	N1-C2-N2	5.83	121.45	116.20
1	X	540	G	C4-C5-N7	-5.83	108.47	110.80
1	X	2720	A	P-O3'-C3'	5.83	126.69	119.70
1	X	2742	G	O4'-C4'-C3'	-5.83	98.17	104.00
1	X	1765	C	N1-C2-O2	5.83	122.40	118.90
1	X	509	U	O4'-C1'-N1	5.82	112.86	108.20
1	X	520	C	P-O3'-C3'	5.82	126.69	119.70
1	X	799	C	O4'-C1'-N1	5.82	112.86	108.20
1	X	1010	U	P-O5'-C5'	5.82	130.21	120.90
1	X	93	A	O4'-C1'-N9	5.82	112.86	108.20
1	X	1168	G	P-O5'-C5'	5.82	130.21	120.90
1	X	2808	U	P-O3'-C3'	5.82	126.68	119.70
1	X	1972	G	C5'-C4'-O4'	5.82	116.08	109.10
1	X	1447	U	O4'-C1'-N1	5.82	112.85	108.20
1	X	2051	U	P-O3'-C3'	5.82	126.68	119.70
1	X	2243	C	O4'-C1'-N1	5.82	112.85	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	20	C	O4'-C1'-N1	5.81	112.85	108.20
1	X	1077	U	P-O3'-C3'	5.81	126.67	119.70
1	X	1149	G	P-O3'-C3'	5.81	126.67	119.70
1	X	1336	G	C5-C6-O6	-5.81	125.11	128.60
1	X	1783	G	O4'-C4'-C3'	-5.81	98.19	104.00
1	X	2492	G	N3-C4-C5	-5.81	125.70	128.60
1	X	2852	G	C5-C6-O6	-5.81	125.11	128.60
1	X	729	A	P-O3'-C3'	5.81	126.67	119.70
1	X	1607	A	C2'-C3'-O3'	5.80	122.99	113.70
1	X	1983	G	O4'-C4'-C3'	-5.80	98.19	104.00
1	X	2250	G	P-O5'-C5'	5.80	130.19	120.90
1	X	1353	A	O4'-C1'-N9	5.80	112.84	108.20
1	X	1412	C	C2'-C3'-O3'	5.80	122.98	113.70
1	X	1767	G	C5-C6-O6	-5.80	125.12	128.60
1	X	1811	A	C4'-C3'-C2'	5.80	108.40	102.60
1	X	2015	G	P-O3'-C3'	5.80	126.66	119.70
1	X	1091	C	O4'-C1'-N1	5.80	112.84	108.20
1	X	1935	A	P-O3'-C3'	5.79	126.65	119.70
1	X	2426	G	O4'-C1'-N9	5.79	112.83	108.20
19	Q	60	GLY	N-CA-C	5.79	127.58	113.10
1	X	1314	A	O4'-C1'-C2'	-5.79	100.01	105.80
1	X	2588	U	N3-C2-O2	-5.79	118.15	122.20
1	X	219	G	P-O3'-C3'	5.78	126.64	119.70
1	X	1790	G	C1'-O4'-C4'	-5.78	105.27	109.90
1	X	2198	U	P-O3'-C3'	5.78	126.64	119.70
1	X	1337	G	O4'-C1'-N9	5.78	112.82	108.20
1	X	1232	U	O4'-C1'-N1	5.78	112.82	108.20
1	X	2025	A	O5'-P-OP2	-5.78	100.50	105.70
1	X	2482	A	N1-C2-N3	-5.78	126.41	129.30
1	X	2403	C	N1-C2-O2	5.78	122.36	118.90
2	Y	32	C	C5-C6-N1	5.78	123.89	121.00
1	X	416	U	C1'-O4'-C4'	-5.77	105.28	109.90
1	X	2627	G	C6-C5-N7	-5.77	126.94	130.40
1	X	183	U	O4'-C1'-N1	5.77	112.82	108.20
1	X	707	U	O4'-C1'-N1	5.77	112.82	108.20
1	X	796	A	C8-N9-C4	-5.77	103.49	105.80
1	X	864	C	C6-N1-C2	-5.77	117.99	120.30
1	X	1148	G	O4'-C1'-N9	5.77	112.82	108.20
1	X	2416	U	C3'-C2'-C1'	-5.77	96.89	101.50
1	X	2601	C	C4'-C3'-C2'	-5.77	96.83	102.60
1	X	477	A	P-O3'-C3'	5.77	126.62	119.70
1	X	2710	C	N3-C2-O2	-5.77	117.86	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	860	U	C1'-O4'-C4'	-5.77	105.29	109.90
1	X	216	U	O4'-C1'-N1	5.76	112.81	108.20
1	X	1036	G	C4'-C3'-C2'	5.76	108.36	102.60
1	X	1090	C	O4'-C1'-N1	5.76	112.81	108.20
1	X	1140	A	O5'-P-OP1	-5.76	100.52	105.70
1	X	2731	G	P-O3'-C3'	5.76	126.61	119.70
11	I	28	LYS	C-N-CA	5.76	136.10	121.70
1	X	835	U	N3-C2-O2	-5.76	118.17	122.20
1	X	1313	U	C3'-C2'-C1'	-5.76	96.90	101.50
1	X	1380	C	O4'-C1'-N1	5.76	112.81	108.20
1	X	1850	G	O4'-C1'-N9	5.76	112.80	108.20
1	X	482	A	C8-N9-C4	-5.75	103.50	105.80
1	X	357	A	P-O3'-C3'	5.75	126.61	119.70
1	X	1958	G	C5-C6-O6	-5.75	125.15	128.60
1	X	2749	A	O4'-C1'-N9	5.75	112.80	108.20
1	X	1142	G	N1-C2-N2	-5.75	111.02	116.20
1	X	2481	G	C3'-C2'-C1'	5.75	106.10	101.50
11	I	35	LYS	N-CA-C	-5.75	95.49	111.00
1	X	2327	U	N1-C2-O2	5.74	126.82	122.80
1	X	2195	C	O4'-C1'-N1	5.74	112.79	108.20
1	X	2406	C	P-O5'-C5'	5.74	130.08	120.90
1	X	1770	U	C5-C6-N1	-5.74	119.83	122.70
1	X	1833	U	O4'-C1'-N1	5.74	112.79	108.20
1	X	2256	G	O4'-C1'-N9	5.74	112.79	108.20
1	X	796	A	C6-C5-N7	-5.73	128.29	132.30
1	X	2421	C	O4'-C1'-N1	5.73	112.79	108.20
1	X	1075	C	O4'-C1'-N1	5.73	112.79	108.20
1	X	488	A	O4'-C1'-N9	5.73	112.78	108.20
1	X	943	U	C1'-O4'-C4'	-5.73	105.31	109.90
1	X	771	C	O4'-C1'-N1	5.73	112.78	108.20
1	X	309	G	C8-N9-C4	-5.72	104.11	106.40
1	X	1652	G	C6-C5-N7	-5.72	126.97	130.40
1	X	559	C	N1-C2-O2	5.72	122.33	118.90
1	X	2228	U	N3-C4-C5	-5.72	111.17	114.60
1	X	2038	C	N1-C2-O2	5.72	122.33	118.90
1	X	1625	A	P-O3'-C3'	5.71	126.56	119.70
1	X	2437	G	C3'-C2'-C1'	5.71	106.07	101.50
1	X	957	G	O4'-C1'-N9	5.71	112.77	108.20
1	X	2841	U	P-O3'-C3'	5.71	126.56	119.70
1	X	332	C	P-O3'-C3'	5.71	126.55	119.70
2	Y	45	C	N1-C2-O2	5.71	122.33	118.90
1	X	405	C	N1-C2-O2	5.71	122.33	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2847	G	N9-C1'-C2'	5.71	121.42	114.00
1	X	846	A	O4'-C1'-N9	5.71	112.77	108.20
1	X	1333	G	C5-N7-C8	-5.71	101.45	104.30
1	X	2442	C	P-O3'-C3'	-5.71	112.85	119.70
1	X	2709	C	O4'-C1'-N1	5.71	112.77	108.20
1	X	2846	G	O4'-C1'-N9	5.71	112.77	108.20
1	X	418	C	O4'-C1'-N1	5.71	112.76	108.20
1	X	845	U	O4'-C1'-N1	5.71	112.76	108.20
1	X	2847	G	N9-C4-C5	5.71	107.68	105.40
1	X	615	C	O4'-C1'-N1	5.70	112.76	108.20
1	X	2706	U	P-O3'-C3'	5.70	126.54	119.70
1	X	983	G	P-O3'-C3'	5.70	126.54	119.70
2	Y	29	C	O4'-C1'-N1	5.70	112.76	108.20
1	X	7	G	C5'-C4'-C3'	-5.70	106.89	116.00
1	X	468	A	P-O3'-C3'	5.70	126.53	119.70
1	X	862	A	O4'-C1'-N9	5.70	112.76	108.20
1	X	1466	C	C4'-C3'-C2'	-5.70	96.90	102.60
1	X	1115	C	O4'-C1'-N1	5.69	112.75	108.20
1	X	1459	U	C4'-C3'-C2'	5.69	108.29	102.60
1	X	2276	C	O4'-C1'-N1	5.69	112.75	108.20
1	X	242	A	C5'-C4'-O4'	5.69	115.92	109.10
1	X	1032	A	C5-N7-C8	-5.69	101.06	103.90
1	X	490	A	O4'-C1'-N9	5.69	112.75	108.20
1	X	542	A	C4-C5-N7	5.69	113.54	110.70
1	X	1550	C	O4'-C1'-N1	5.69	112.75	108.20
1	X	2730	A	P-O3'-C3'	5.69	126.52	119.70
1	X	1245	G	C4'-C3'-C2'	-5.68	96.92	102.60
1	X	1278	A	C2-N3-C4	-5.68	107.76	110.60
1	X	1652	G	N9-C4-C5	-5.68	103.13	105.40
1	X	1133	G	O4'-C1'-N9	5.67	112.74	108.20
1	X	1295	U	O4'-C1'-N1	5.67	112.74	108.20
1	X	1412	C	C6-N1-C2	-5.67	118.03	120.30
1	X	1598	C	O4'-C1'-N1	5.67	112.74	108.20
1	X	788	G	C2'-C3'-O3'	5.67	122.78	113.70
1	X	1412	C	O4'-C1'-N1	5.67	112.74	108.20
1	X	1978	U	O4'-C1'-N1	5.67	112.74	108.20
1	X	744	C	O4'-C1'-N1	5.67	112.74	108.20
1	X	837	U	O4'-C1'-N1	5.67	112.74	108.20
1	X	942	U	O4'-C1'-N1	5.67	112.74	108.20
1	X	2037	A	O4'-C1'-N9	5.67	112.74	108.20
1	X	1683	G	P-O3'-C3'	-5.67	112.90	119.70
1	X	1971	C	P-O3'-C3'	-5.67	112.90	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	418	C	C4'-C3'-C2'	-5.67	96.93	102.60
1	X	569	C	N1-C2-O2	5.66	122.30	118.90
1	X	1108	U	O4'-C1'-N1	5.66	112.73	108.20
1	X	2477	C	P-O5'-C5'	5.66	129.96	120.90
1	X	814	G	C5-C6-N1	5.66	114.33	111.50
1	X	1508	G	P-O3'-C3'	5.66	126.49	119.70
1	X	2260	C	O4'-C1'-N1	5.66	112.73	108.20
1	X	796	A	C4-C5-N7	5.66	113.53	110.70
1	X	878	C	N1-C2-O2	5.66	122.29	118.90
1	X	2008	C	C5-C6-N1	5.66	123.83	121.00
1	X	1181	C	O4'-C1'-N1	5.65	112.72	108.20
1	X	805	G	P-O5'-C5'	5.65	129.94	120.90
1	X	2199	C	P-O5'-C5'	5.65	129.94	120.90
1	X	39	C	O4'-C1'-N1	5.65	112.72	108.20
1	X	937	C	O4'-C1'-N1	5.65	112.72	108.20
1	X	2443	C	O4'-C1'-N1	5.65	112.72	108.20
1	X	2449	G	O4'-C1'-N9	5.65	112.72	108.20
2	Y	53	G	C8-N9-C4	-5.65	104.14	106.40
1	X	462	G	C4-C5-C6	5.65	122.19	118.80
1	X	607	C	O4'-C4'-C3'	-5.65	98.35	104.00
1	X	751	G	O4'-C1'-N9	5.65	112.72	108.20
1	X	1513	U	P-O3'-C3'	5.65	126.48	119.70
1	X	1607	A	C3'-C2'-C1'	-5.65	96.98	101.50
1	X	1341	G	P-O3'-C3'	-5.64	112.93	119.70
1	X	2854	G	C5-N7-C8	-5.64	101.48	104.30
1	X	569	C	N3-C2-O2	-5.64	117.95	121.90
1	X	2557	G	C3'-C2'-C1'	5.64	106.02	101.50
1	X	1434	U	O4'-C1'-N1	5.64	112.71	108.20
1	X	817	A	C1'-O4'-C4'	-5.64	105.39	109.90
1	X	1770	U	N1-C2-N3	5.64	118.28	114.90
1	X	2002	A	P-O5'-C5'	5.64	129.92	120.90
1	X	860	U	C5'-C4'-O4'	5.64	115.87	109.10
1	X	2797	G	C6-C5-N7	-5.64	127.02	130.40
1	X	1594	U	O4'-C1'-N1	5.63	112.71	108.20
1	X	868	U	O4'-C1'-N1	5.63	112.70	108.20
1	X	549	G	O4'-C1'-N9	5.63	112.70	108.20
1	X	2483	U	O4'-C1'-N1	5.63	112.70	108.20
1	X	1087	C	O4'-C1'-N1	5.63	112.70	108.20
1	X	1966	C	P-O3'-C3'	-5.63	112.94	119.70
1	X	2219	U	O4'-C1'-N1	5.63	112.70	108.20
1	X	2274	C	N1-C2-O2	5.63	122.28	118.90
1	X	2636	A	P-O3'-C3'	5.63	126.45	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	746	G	N3-C4-C5	-5.63	125.79	128.60
1	X	1142	G	N3-C2-N2	5.63	123.84	119.90
1	X	2726	U	O4'-C1'-N1	5.63	112.70	108.20
2	Y	90	C	N1-C2-O2	5.63	122.28	118.90
1	X	1715	A	P-O3'-C3'	5.62	126.45	119.70
1	X	537	C	N1-C2-N3	-5.62	115.26	119.20
2	Y	118	G	C3'-C2'-C1'	-5.62	97.00	101.50
1	X	2300	G	C2-N3-C4	5.62	114.71	111.90
1	X	103	U	C5-C6-N1	5.62	125.51	122.70
1	X	1245	G	O4'-C1'-N9	5.62	112.69	108.20
1	X	2217	G	P-O3'-C3'	5.61	126.44	119.70
1	X	1683	G	C4-C5-N7	-5.61	108.56	110.80
1	X	327	C	N1-C2-O2	5.61	122.27	118.90
1	X	594	G	C4'-C3'-C2'	-5.61	96.99	102.60
1	X	2362	G	P-O3'-C3'	5.61	126.43	119.70
1	X	956	A	N1-C6-N6	5.61	121.96	118.60
11	I	55	ARG	C-N-CA	5.61	135.72	121.70
1	X	1688	U	O4'-C1'-N1	5.60	112.68	108.20
1	X	2692	A	C3'-C2'-C1'	5.60	105.98	101.50
1	X	2793	G	O4'-C4'-C3'	-5.60	98.40	104.00
1	X	689	A	C8-N9-C4	-5.60	103.56	105.80
1	X	816	U	O4'-C1'-N1	5.60	112.68	108.20
1	X	1753	A	C8-N9-C4	-5.60	103.56	105.80
1	X	465	C	P-O5'-C5'	-5.60	111.94	120.90
1	X	2803	C	P-O5'-C5'	-5.60	111.94	120.90
1	X	2452	U	O4'-C1'-N1	5.60	112.68	108.20
1	X	616	U	O4'-C1'-N1	5.59	112.68	108.20
1	X	940	G	P-O5'-C5'	5.59	129.85	120.90
1	X	2797	G	N3-C4-C5	-5.59	125.80	128.60
1	X	126	C	N1-C1'-C2'	5.59	121.27	114.00
1	X	1288	A	C5'-C4'-C3'	5.59	124.94	116.00
1	X	1490	U	O4'-C1'-N1	5.59	112.67	108.20
1	X	1913	G	O4'-C1'-N9	-5.59	103.73	108.20
1	X	1924	C	N1-C2-O2	5.59	122.25	118.90
1	X	2205	C	O4'-C1'-N1	5.59	112.67	108.20
1	X	1337	G	C8-N9-C4	-5.58	104.17	106.40
1	X	1981	A	C6-C5-N7	-5.58	128.39	132.30
1	X	498	C	N1-C2-O2	5.58	122.25	118.90
1	X	1089	C	P-O3'-C3'	5.58	126.40	119.70
1	X	2252	A	N1-C6-N6	-5.58	115.25	118.60
1	X	2640	G	O4'-C1'-N9	5.58	112.66	108.20
1	X	1712	G	N3-C4-C5	-5.58	125.81	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2410	U	O4'-C1'-N1	5.58	112.66	108.20
2	Y	89	G	O4'-C1'-N9	5.58	112.66	108.20
1	X	462	G	C5-C6-N1	-5.57	108.71	111.50
1	X	1539	U	O4'-C1'-N1	5.57	112.66	108.20
1	X	1977	C	O4'-C1'-N1	5.57	112.66	108.20
1	X	2576	G	C5-C6-O6	-5.57	125.26	128.60
1	X	1150	C	O4'-C1'-N1	5.57	112.65	108.20
1	X	1469	U	O3'-P-O5'	5.57	114.58	104.00
1	X	237	G	O4'-C1'-N9	5.57	112.65	108.20
1	X	465	C	O3'-P-O5'	-5.56	93.43	104.00
1	X	2418	A	P-O3'-C3'	5.56	126.38	119.70
1	X	434	C	C3'-C2'-C1'	5.56	105.95	101.50
1	X	699	G	C8-N9-C4	-5.56	104.17	106.40
1	X	764	A	N9-C1'-C2'	5.56	121.23	114.00
1	X	2479	U	C4'-C3'-C2'	-5.56	97.04	102.60
1	X	737	C	C5-C6-N1	5.56	123.78	121.00
13	K	94	TYR	N-CA-CB	5.56	120.61	110.60
1	X	1237	G	O4'-C1'-N9	5.56	112.65	108.20
1	X	2483	U	C5-C6-N1	5.56	125.48	122.70
1	X	2305	C	O4'-C1'-N1	5.56	112.65	108.20
1	X	2370	G	C1'-O4'-C4'	-5.56	105.45	109.90
1	X	1223	G	C4-C5-N7	5.55	113.02	110.80
1	X	1853	C	O4'-C1'-N1	5.55	112.64	108.20
1	X	2049	C	O4'-C1'-N1	5.55	112.64	108.20
1	X	2050	G	C5-C6-O6	-5.55	125.27	128.60
1	X	1277	G	O4'-C1'-N9	5.55	112.64	108.20
1	X	1398	G	N9-C1'-C2'	5.55	121.21	114.00
1	X	475	U	O4'-C1'-N1	5.54	112.64	108.20
1	X	2564	U	C1'-O4'-C4'	-5.54	105.46	109.90
1	X	2853	U	P-O3'-C3'	5.54	126.35	119.70
1	X	1968	G	O4'-C1'-N9	5.54	112.64	108.20
2	Y	11	G	O4'-C4'-C3'	-5.54	98.46	104.00
1	X	1153	A	P-O3'-C3'	5.54	126.35	119.70
1	X	537	C	C5'-C4'-O4'	5.54	115.75	109.10
1	X	755	C	O4'-C1'-N1	5.54	112.63	108.20
1	X	1056	U	O4'-C1'-N1	5.54	112.63	108.20
1	X	1399	C	O4'-C1'-N1	5.54	112.63	108.20
1	X	1969	G	P-O3'-C3'	-5.54	113.05	119.70
1	X	1514	C	O4'-C1'-N1	5.54	112.63	108.20
1	X	1644	G	O4'-C1'-N9	5.54	112.63	108.20
1	X	1688	U	C4-C5-C6	5.53	123.02	119.70
1	X	2232	G	C5-C6-O6	-5.53	125.28	128.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2312	A	O4'-C1'-N9	5.53	112.63	108.20
1	X	2799	C	O4'-C1'-N1	5.53	112.63	108.20
1	X	2870	C	C6-N1-C2	-5.53	118.09	120.30
1	X	185	C	O4'-C1'-N1	5.53	112.62	108.20
1	X	423	G	C8-N9-C4	-5.52	104.19	106.40
1	X	634	G	P-O5'-C5'	5.52	129.74	120.90
1	X	1608	U	O4'-C1'-N1	5.52	112.62	108.20
1	X	2191	A	O4'-C1'-N9	5.52	112.62	108.20
1	X	1326	U	N3-C2-O2	-5.52	118.33	122.20
1	X	794	A	O4'-C4'-C3'	-5.52	98.48	104.00
1	X	883	A	O4'-C1'-N9	5.52	112.62	108.20
1	X	1249	G	C4'-C3'-C2'	5.52	108.12	102.60
1	X	1662	G	N9-C1'-C2'	5.52	121.17	114.00
1	X	1217	U	O4'-C1'-N1	5.52	112.61	108.20
1	X	820	U	P-O3'-C3'	-5.51	113.08	119.70
1	X	1434	U	P-O3'-C3'	5.51	126.31	119.70
1	X	1863	U	O4'-C1'-N1	5.51	112.61	108.20
1	X	2010	G	O4'-C1'-N9	5.51	112.61	108.20
1	X	522	G	N7-C8-N9	5.51	115.85	113.10
1	X	1992	G	OP1-P-OP2	-5.50	111.34	119.60
1	X	2819	G	N3-C4-C5	-5.50	125.85	128.60
1	X	2835	A	O4'-C1'-N9	5.50	112.60	108.20
1	X	1623	C	N1-C2-O2	5.50	122.20	118.90
1	X	2782	G	C4-C5-N7	5.50	113.00	110.80
1	X	1990	U	O4'-C1'-N1	5.50	112.60	108.20
1	X	1201	G	C8-N9-C4	-5.50	104.20	106.40
1	X	1432	G	C3'-C2'-C1'	5.50	105.90	101.50
1	X	2580	C	P-O3'-C3'	5.50	126.30	119.70
1	X	596	C	P-O5'-C5'	-5.50	112.11	120.90
1	X	1270	C	C6-N1-C2	-5.49	118.10	120.30
1	X	1081	A	P-O3'-C3'	5.49	126.29	119.70
1	X	841	G	C5-N7-C8	-5.49	101.56	104.30
1	X	1468	A	N1-C6-N6	-5.49	115.31	118.60
1	X	1947	G	O4'-C1'-N9	-5.49	103.81	108.20
1	X	633	G	O4'-C1'-N9	5.49	112.59	108.20
1	X	666	U	P-O3'-C3'	5.49	126.28	119.70
1	X	1201	G	P-O3'-C3'	5.48	126.28	119.70
1	X	1792	C	C6-N1-C2	-5.48	118.11	120.30
1	X	308	C	C4'-C3'-C2'	5.48	108.08	102.60
1	X	1341	G	P-O5'-C5'	5.48	129.67	120.90
1	X	1754	G	P-O5'-C5'	5.48	129.67	120.90
1	X	1103	C	O4'-C1'-N1	5.48	112.58	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	45	C	O4'-C1'-N1	5.48	112.58	108.20
1	X	167	A	P-O5'-C5'	5.48	129.66	120.90
1	X	1319	C	C5-C6-N1	5.48	123.74	121.00
2	Y	81	C	C6-N1-C2	-5.48	118.11	120.30
1	X	327	C	O4'-C1'-N1	5.47	112.58	108.20
1	X	818	G	C6-C5-N7	-5.47	127.12	130.40
1	X	1467	U	C4'-C3'-C2'	5.47	108.07	102.60
1	X	2087	U	O4'-C1'-N1	5.47	112.58	108.20
1	X	346	C	C5-C6-N1	5.47	123.74	121.00
1	X	1275	A	P-O3'-C3'	-5.47	113.13	119.70
1	X	1876	C	N1-C2-O2	5.47	122.18	118.90
1	X	2315	A	O4'-C1'-N9	-5.47	103.82	108.20
1	X	2555	G	O4'-C4'-C3'	-5.47	98.53	104.00
1	X	2778	U	N1-C1'-C2'	5.47	121.11	114.00
1	X	503	G	C5-C6-N1	5.47	114.23	111.50
1	X	1725	C	O4'-C1'-N1	5.47	112.58	108.20
1	X	94	C	O4'-C1'-N1	5.47	112.57	108.20
1	X	176	A	C1'-O4'-C4'	-5.47	105.53	109.90
1	X	1341	G	N3-C4-N9	5.46	129.28	126.00
1	X	2224	U	O4'-C1'-N1	5.46	112.57	108.20
1	X	2437	G	C8-N9-C4	-5.46	104.22	106.40
1	X	857	U	O4'-C1'-N1	5.46	112.57	108.20
1	X	1948	C	O4'-C1'-N1	5.46	112.57	108.20
1	X	133	C	O4'-C1'-N1	5.46	112.57	108.20
1	X	1775	A	C2'-C3'-O3'	5.46	122.44	113.70
1	X	765	C	C4'-C3'-C2'	5.46	108.06	102.60
1	X	878	C	O4'-C1'-N1	5.46	112.57	108.20
1	X	1336	G	C5-C6-N1	5.46	114.23	111.50
1	X	2089	C	O4'-C1'-N1	5.46	112.56	108.20
1	X	2342	U	C3'-C2'-C1'	-5.46	97.14	101.50
1	X	2705	A	C4'-C3'-O3'	5.46	123.91	113.00
1	X	1566	G	O4'-C1'-N9	5.46	112.56	108.20
1	X	2032	G	N3-C4-C5	-5.45	125.87	128.60
1	X	2495	G	C5-C6-N1	5.45	114.23	111.50
1	X	309	G	N7-C8-N9	5.45	115.83	113.10
1	X	1142	G	C1'-O4'-C4'	-5.45	105.54	109.90
1	X	1143	A	C3'-C2'-C1'	-5.45	97.14	101.50
1	X	1651	U	O4'-C1'-N1	5.45	112.56	108.20
1	X	1712	G	C4-N9-C1'	5.45	133.59	126.50
1	X	1049	C	O4'-C1'-N1	5.45	112.56	108.20
1	X	1301	U	N3-C2-O2	-5.45	118.39	122.20
2	Y	110	U	N1-C2-O2	5.45	126.61	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	832	A	C5'-C4'-O4'	-5.45	102.56	109.10
1	X	1340	C	O3'-P-O5'	-5.45	93.65	104.00
1	X	2016	A	N1-C2-N3	-5.45	126.58	129.30
1	X	533	C	O4'-C1'-N1	5.44	112.56	108.20
1	X	721	C	O4'-C1'-N1	5.44	112.55	108.20
1	X	2336	G	O5'-P-OP2	-5.44	100.80	105.70
1	X	179	U	O4'-C1'-N1	5.44	112.55	108.20
1	X	2013	A	C1'-O4'-C4'	-5.44	105.55	109.90
1	X	2482	A	C5-C6-N6	-5.44	119.35	123.70
1	X	2646	C	C5-C6-N1	5.44	123.72	121.00
1	X	2808	U	C5'-C4'-O4'	5.44	115.63	109.10
1	X	2407	G	P-O3'-C3'	5.44	126.23	119.70
1	X	1814	G	O4'-C1'-N9	5.44	112.55	108.20
1	X	2675	U	O4'-C1'-N1	5.44	112.55	108.20
1	X	871	U	P-O5'-C5'	5.43	129.60	120.90
1	X	1917	C	O4'-C1'-N1	5.43	112.55	108.20
1	X	1396	C	C6-N1-C2	-5.43	118.13	120.30
1	X	1652	G	N1-C6-O6	5.43	123.16	119.90
1	X	70	A	P-O3'-C3'	5.43	126.22	119.70
1	X	573	C	O4'-C1'-N1	5.43	112.54	108.20
1	X	1015	U	P-O3'-C3'	5.43	126.21	119.70
1	X	1163	C	O4'-C1'-N1	5.43	112.54	108.20
1	X	1278	A	N9-C1'-C2'	5.43	121.06	114.00
1	X	1285	A	P-O3'-C3'	5.43	126.22	119.70
1	X	29	U	O4'-C1'-N1	5.43	112.54	108.20
1	X	2731	G	C3'-C2'-C1'	5.43	105.84	101.50
1	X	840	U	C1'-O4'-C4'	-5.42	105.56	109.90
1	X	2626	U	O4'-C1'-N1	5.42	112.54	108.20
1	X	319	G	N9-C1'-C2'	5.42	121.05	114.00
1	X	2550	C	O4'-C1'-N1	5.42	112.54	108.20
1	X	236	C	N1-C2-O2	5.42	122.15	118.90
2	Y	86	A	O4'-C4'-C3'	-5.42	98.58	104.00
1	X	235	C	N1-C2-O2	5.42	122.15	118.90
1	X	882	C	O4'-C1'-N1	5.42	112.53	108.20
1	X	1364	C	O4'-C1'-N1	5.42	112.53	108.20
1	X	215	G	O4'-C1'-N9	5.42	112.53	108.20
1	X	329	C	C6-N1-C2	-5.42	118.13	120.30
1	X	2076	G	C8-N9-C4	-5.42	104.23	106.40
1	X	1162	A	P-O3'-C3'	5.41	126.20	119.70
1	X	1683	G	N3-C2-N2	-5.41	116.11	119.90
2	Y	49	C	N1-C2-O2	5.41	122.15	118.90
2	Y	7	C	O4'-C1'-N1	5.41	112.53	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	473	C	OP2-P-O3'	5.41	117.10	105.20
1	X	2000	U	O5'-P-OP2	-5.41	100.83	105.70
1	X	2872	U	O4'-C1'-N1	5.41	112.53	108.20
2	Y	5	C	O4'-C1'-N1	5.41	112.53	108.20
1	X	617	U	C2-N1-C1'	5.41	124.19	117.70
1	X	211	U	O4'-C1'-N1	5.41	112.53	108.20
1	X	2229	G	C5'-C4'-O4'	5.41	115.59	109.10
1	X	2323	U	P-O5'-C5'	5.41	129.55	120.90
1	X	2864	C	O4'-C1'-N1	5.41	112.52	108.20
1	X	467	U	N1-C1'-C2'	5.40	121.02	114.00
1	X	1573	G	P-O3'-C3'	5.40	126.18	119.70
1	X	1618	U	P-O3'-C3'	5.40	126.18	119.70
1	X	2285	U	O4'-C1'-N1	5.40	112.52	108.20
1	X	338	G	C8-N9-C4	-5.40	104.24	106.40
1	X	964	A	O4'-C1'-N9	5.40	112.52	108.20
1	X	2273	C	O4'-C1'-N1	5.40	112.52	108.20
1	X	1923	U	C2'-C3'-O3'	5.40	122.34	113.70
1	X	2009	U	P-O3'-C3'	-5.40	113.22	119.70
1	X	178	C	O4'-C1'-N1	5.40	112.52	108.20
1	X	561	U	C3'-C2'-C1'	-5.40	97.18	101.50
1	X	699	G	C8-N9-C1'	5.40	134.02	127.00
1	X	1338	G	N3-C4-N9	5.40	129.24	126.00
1	X	1796	A	P-O3'-C3'	5.40	126.18	119.70
1	X	1939	U	N3-C2-O2	-5.40	118.42	122.20
1	X	2699	G	OP1-P-O3'	5.40	117.08	105.20
1	X	998	C	O4'-C1'-N1	5.40	112.52	108.20
1	X	1325	U	P-O3'-C3'	5.40	126.17	119.70
1	X	175	C	C6-N1-C2	-5.39	118.14	120.30
1	X	537	C	C2-N1-C1'	5.39	124.73	118.80
1	X	2778	U	C3'-C2'-C1'	5.39	105.82	101.50
1	X	90	G	N3-C4-C5	-5.39	125.90	128.60
1	X	346	C	C3'-C2'-C1'	5.39	105.81	101.50
1	X	2586	G	C8-N9-C4	-5.39	104.24	106.40
1	X	753	U	O4'-C1'-N1	5.39	112.51	108.20
1	X	1141	U	O4'-C1'-N1	5.39	112.51	108.20
1	X	1984	A	N1-C6-N6	-5.39	115.36	118.60
1	X	2699	G	C2-N3-C4	5.39	114.59	111.90
1	X	1471	G	O4'-C1'-N9	5.39	112.51	108.20
1	X	145	C	O4'-C1'-N1	5.39	112.51	108.20
1	X	2363	G	O4'-C1'-N9	5.39	112.51	108.20
2	Y	19	C	N1-C2-O2	5.39	122.13	118.90
1	X	1992	G	C2-N3-C4	5.39	114.59	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2018	G	O4'-C1'-N9	5.39	112.51	108.20
1	X	2774	U	O4'-C1'-N1	5.39	112.51	108.20
1	X	2772	U	O4'-C1'-N1	5.38	112.51	108.20
1	X	591	G	O4'-C1'-N9	5.38	112.51	108.20
1	X	820	U	N3-C2-O2	-5.38	118.43	122.20
1	X	1743	C	P-O3'-C3'	-5.38	113.24	119.70
1	X	675	C	C3'-C2'-C1'	-5.38	97.20	101.50
1	X	715	U	O4'-C1'-N1	5.38	112.50	108.20
1	X	1247	U	O4'-C1'-N1	5.38	112.50	108.20
1	X	2533	U	O4'-C1'-N1	5.38	112.50	108.20
1	X	417	C	N1-C2-O2	5.38	122.12	118.90
1	X	1766	U	P-O5'-C5'	-5.38	112.30	120.90
1	X	2176	U	O4'-C1'-N1	5.38	112.50	108.20
1	X	689	A	C1'-O4'-C4'	-5.38	105.60	109.90
1	X	1093	U	O4'-C1'-N1	5.38	112.50	108.20
1	X	2551	A	OP1-P-O3'	5.38	117.03	105.20
1	X	1097	A	P-O3'-C3'	5.37	126.15	119.70
1	X	1274	C	O4'-C1'-N1	5.37	112.50	108.20
2	Y	47	A	C8-N9-C4	-5.37	103.65	105.80
1	X	541	C	O5'-P-OP1	-5.37	100.87	105.70
1	X	1627	C	N1-C2-O2	5.37	122.12	118.90
1	X	660	G	C3'-C2'-C1'	-5.37	97.20	101.50
1	X	613	A	O4'-C1'-N9	5.37	112.50	108.20
1	X	2854	G	P-O5'-C5'	5.37	129.49	120.90
1	X	1752	U	O4'-C1'-N1	5.37	112.49	108.20
1	X	1825	C	O4'-C1'-N1	5.37	112.49	108.20
1	X	2028	C	C5-C6-N1	5.36	123.68	121.00
1	X	2665	G	N3-C4-N9	5.36	129.22	126.00
15	M	28	ARG	N-CA-C	-5.36	96.52	111.00
1	X	2531	U	N3-C2-O2	-5.36	118.45	122.20
1	X	1080	A	C1'-O4'-C4'	-5.36	105.61	109.90
1	X	2275	U	P-O5'-C5'	5.36	129.47	120.90
1	X	238	G	C4'-C3'-C2'	-5.36	97.24	102.60
1	X	697	G	O4'-C1'-N9	5.36	112.49	108.20
1	X	2480	C	N3-C4-C5	5.36	124.04	121.90
1	X	2697	G	C5-C6-N1	5.36	114.18	111.50
1	X	925	U	O4'-C1'-N1	5.36	112.48	108.20
1	X	2840	U	O4'-C1'-N1	5.36	112.48	108.20
2	Y	81	C	C5-C6-N1	5.36	123.68	121.00
1	X	522	G	C5-C6-N1	5.35	114.18	111.50
1	X	652	C	C5-C6-N1	5.35	123.68	121.00
1	X	1262	U	O4'-C1'-N1	5.35	112.48	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2323	U	C5-C6-N1	5.35	125.38	122.70
3	A	242	ALA	N-CA-C	5.35	125.46	111.00
1	X	78	C	C6-N1-C2	-5.35	118.16	120.30
1	X	2788	C	O4'-C1'-N1	5.35	112.48	108.20
1	X	146	C	O4'-C1'-N1	5.35	112.48	108.20
1	X	2008	C	N3-C4-N4	5.35	121.75	118.00
1	X	2815	C	C3'-C2'-C1'	-5.35	97.22	101.50
1	X	1306	U	P-O3'-C3'	-5.35	113.28	119.70
1	X	1270	C	N3-C4-C5	-5.35	119.76	121.90
1	X	1776	A	C4'-C3'-C2'	-5.34	97.25	102.60
1	X	2650	G	P-O3'-C3'	-5.34	113.29	119.70
1	X	2165	A	P-O3'-C3'	5.34	126.11	119.70
1	X	1691	G	C5-C6-O6	-5.34	125.40	128.60
1	X	1828	C	O4'-C1'-N1	5.34	112.47	108.20
1	X	2477	C	O5'-P-OP1	-5.34	100.89	105.70
1	X	2754	C	N1-C2-O2	5.34	122.10	118.90
1	X	2837	G	O4'-C1'-N9	5.34	112.47	108.20
1	X	2088	U	O4'-C1'-N1	5.34	112.47	108.20
1	X	611	C	C6-N1-C2	-5.34	118.17	120.30
1	X	1716	G	C1'-O4'-C4'	5.34	114.17	109.90
1	X	2430	A	O4'-C1'-N9	5.34	112.47	108.20
1	X	129	A	P-O5'-C5'	5.33	129.44	120.90
1	X	1054	C	O4'-C1'-N1	5.33	112.47	108.20
1	X	1315	A	O5'-P-OP2	-5.33	100.90	105.70
1	X	2017	U	P-O3'-C3'	5.33	126.10	119.70
1	X	677	G	P-O3'-C3'	5.33	126.10	119.70
1	X	1092	U	O4'-C1'-N1	5.33	112.47	108.20
1	X	1686	A	OP1-P-OP2	-5.33	111.60	119.60
1	X	2467	A	O4'-C1'-N9	5.33	112.47	108.20
1	X	827	C	O4'-C1'-N1	5.33	112.46	108.20
2	Y	123	U	N1-C2-O2	5.33	126.53	122.80
1	X	579	G	C5-C6-N1	-5.33	108.84	111.50
1	X	190	A	C1'-O4'-C4'	-5.33	105.64	109.90
1	X	1188	A	P-O3'-C3'	5.33	126.09	119.70
1	X	1409	U	O4'-C1'-N1	5.33	112.46	108.20
1	X	440	U	O3'-P-O5'	-5.32	93.89	104.00
1	X	596	C	N3-C4-C5	-5.32	119.77	121.90
1	X	1296	G	P-O3'-C3'	5.32	126.09	119.70
1	X	2559	U	N3-C4-C5	5.32	117.80	114.60
1	X	2856	U	N3-C2-O2	-5.32	118.47	122.20
1	X	696	U	P-O5'-C5'	5.32	129.42	120.90
1	X	54	G	P-O3'-C3'	5.32	126.08	119.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	144	U	O4'-C1'-N1	5.32	112.46	108.20
1	X	687	G	N3-C4-C5	-5.32	125.94	128.60
1	X	1009	C	N1-C2-O2	5.32	122.09	118.90
1	X	1564	U	O4'-C1'-N1	5.32	112.46	108.20
1	X	2075	U	P-O3'-C3'	5.32	126.08	119.70
1	X	2632	U	O4'-C1'-N1	5.32	112.46	108.20
1	X	1871	G	O4'-C1'-N9	5.32	112.46	108.20
1	X	103	U	O4'-C1'-N1	5.32	112.45	108.20
1	X	429	C	C5-C6-N1	5.32	123.66	121.00
1	X	967	G	P-O3'-C3'	5.32	126.08	119.70
1	X	1219	C	O4'-C1'-N1	5.32	112.45	108.20
1	X	1355	A	C4'-C3'-C2'	5.32	107.92	102.60
1	X	1439	G	O4'-C4'-C3'	-5.32	98.69	104.00
1	X	2474	G	C5-C6-N1	5.32	114.16	111.50
1	X	2554	C	N3-C2-O2	-5.32	118.18	121.90
1	X	2844	G	C8-N9-C4	-5.32	104.27	106.40
2	Y	42	U	P-O3'-C3'	5.31	126.08	119.70
1	X	1489	C	C2-N1-C1'	5.31	124.64	118.80
2	Y	64	C	O4'-C1'-N1	5.31	112.45	108.20
1	X	246	C	O4'-C1'-N1	5.31	112.44	108.20
1	X	682	G	C3'-C2'-C1'	5.31	105.75	101.50
1	X	2027	C	C3'-C2'-C1'	-5.31	97.25	101.50
1	X	586	G	O4'-C4'-C3'	-5.30	98.69	104.00
1	X	756	C	O4'-C1'-N1	5.30	112.44	108.20
1	X	2507	U	O4'-C1'-N1	5.30	112.44	108.20
1	X	536	A	P-O3'-C3'	5.30	126.06	119.70
19	Q	73	ASN	C-N-CA	5.30	134.95	121.70
1	X	747	A	N1-C6-N6	5.30	121.78	118.60
1	X	1506	C	O4'-C1'-N1	5.30	112.44	108.20
1	X	2281	C	O4'-C1'-N1	5.30	112.44	108.20
1	X	580	A	N9-C1'-C2'	5.30	120.89	114.00
1	X	650	U	P-O5'-C5'	5.30	129.38	120.90
1	X	1275	A	C2-N3-C4	5.30	113.25	110.60
1	X	1400	A	P-O3'-C3'	-5.30	113.34	119.70
1	X	2299	A	C1'-O4'-C4'	-5.30	105.66	109.90
1	X	1069	G	C3'-C2'-C1'	5.29	105.73	101.50
1	X	2195	C	C6-N1-C2	-5.29	118.18	120.30
1	X	2662	C	O4'-C1'-N1	5.29	112.43	108.20
12	J	82	THR	C-N-CA	5.29	134.93	121.70
1	X	1336	G	C6-C5-N7	-5.29	127.23	130.40
1	X	1349	A	C2-N3-C4	5.29	113.25	110.60
1	X	75	C	O4'-C1'-N1	5.29	112.43	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2625	U	C5-C4-O4	-5.29	122.73	125.90
2	Y	2	C	O4'-C1'-N1	5.29	112.43	108.20
2	Y	87	C	O4'-C1'-N1	5.29	112.43	108.20
1	X	2528	G	O4'-C1'-N9	-5.29	103.97	108.20
1	X	2594	U	C4-C5-C6	-5.29	116.53	119.70
1	X	70	A	P-O5'-C5'	-5.28	112.45	120.90
1	X	1147	G	O4'-C1'-N9	5.28	112.43	108.20
1	X	1489	C	N3-C2-O2	-5.28	118.20	121.90
1	X	2434	G	N3-C4-C5	-5.28	125.96	128.60
1	X	2484	G	C8-N9-C4	-5.28	104.29	106.40
1	X	943	U	N3-C2-O2	-5.28	118.50	122.20
1	X	1712	G	O4'-C1'-N9	5.28	112.42	108.20
1	X	2044	G	N3-C4-C5	-5.28	125.96	128.60
1	X	1467	U	C5-C4-O4	-5.28	122.73	125.90
1	X	2522	G	O4'-C1'-N9	5.28	112.42	108.20
1	X	2702	G	P-O3'-C3'	5.28	126.03	119.70
1	X	2642	G	P-O5'-C5'	5.27	129.34	120.90
1	X	574	C	OP2-P-O3'	5.27	116.80	105.20
1	X	660	G	C8-N9-C4	-5.27	104.29	106.40
1	X	2776	U	O4'-C1'-N1	5.27	112.42	108.20
1	X	2304	G	P-O3'-C3'	5.27	126.03	119.70
1	X	1055	A	O4'-C1'-N9	5.27	112.42	108.20
1	X	527	C	O4'-C1'-N1	5.27	112.41	108.20
1	X	2228	U	N3-C4-O4	5.27	123.09	119.40
1	X	2556	A	C5'-C4'-O4'	5.27	115.42	109.10
1	X	549	G	C3'-C2'-C1'	-5.27	97.29	101.50
1	X	434	C	O4'-C1'-N1	5.26	112.41	108.20
1	X	687	G	N3-C4-N9	5.26	129.16	126.00
1	X	11	G	C8-N9-C4	-5.26	104.30	106.40
1	X	1022	A	P-O3'-C3'	-5.26	113.39	119.70
1	X	1162	A	C4'-C3'-C2'	-5.26	97.34	102.60
1	X	859	U	N3-C2-O2	-5.26	118.52	122.20
1	X	1306	U	O4'-C1'-N1	5.26	112.41	108.20
1	X	2858	A	C5-C6-N6	-5.26	119.49	123.70
1	X	514	G	N9-C1'-C2'	5.26	120.83	114.00
11	I	32	ARG	N-CA-C	-5.26	96.81	111.00
1	X	1032	A	N7-C8-N9	5.25	116.43	113.80
1	X	1375	C	O4'-C1'-N1	5.25	112.40	108.20
1	X	1524	C	N1-C2-O2	5.25	122.05	118.90
1	X	1767	G	C5-C6-N1	5.25	114.13	111.50
1	X	759	C	N3-C4-C5	-5.25	119.80	121.90
1	X	828	C	N1-C2-O2	5.25	122.05	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	522	G	C5-N7-C8	-5.25	101.68	104.30
1	X	579	G	O4'-C1'-N9	5.25	112.40	108.20
1	X	2332	G	O4'-C1'-N9	5.25	112.40	108.20
2	Y	12	C	O4'-C4'-C3'	-5.25	98.75	104.00
1	X	607	C	C3'-C2'-C1'	-5.25	97.30	101.50
1	X	2528	G	OP1-P-O3'	5.25	116.74	105.20
1	X	2275	U	P-O3'-C3'	5.24	125.99	119.70
1	X	2649	A	P-O3'-C3'	-5.24	113.41	119.70
1	X	1642	G	O4'-C1'-N9	-5.24	104.01	108.20
1	X	2522	G	C8-N9-C4	-5.24	104.30	106.40
1	X	597	U	O4'-C4'-C3'	-5.24	98.76	104.00
11	I	48	PHE	C-N-CA	5.24	134.80	121.70
1	X	1277	G	C5-C6-N1	5.24	114.12	111.50
1	X	1729	C	C6-N1-C2	-5.24	118.20	120.30
1	X	2570	C	O4'-C1'-N1	5.24	112.39	108.20
1	X	1303	U	N1-C2-O2	5.23	126.46	122.80
1	X	2491	C	C3'-C2'-C1'	-5.23	97.31	101.50
1	X	344	G	N9-C1'-C2'	5.23	120.80	114.00
1	X	2241	U	N3-C4-O4	5.23	123.06	119.40
1	X	787	A	P-O3'-C3'	5.23	125.98	119.70
1	X	1319	C	O4'-C1'-N1	5.23	112.38	108.20
1	X	542	A	P-O3'-C3'	5.23	125.97	119.70
1	X	1635	G	C8-N9-C4	-5.23	104.31	106.40
1	X	1733	U	P-O3'-C3'	5.23	125.97	119.70
1	X	2489	C	P-O3'-C3'	-5.23	113.43	119.70
2	Y	46	G	C3'-C2'-C1'	5.23	105.68	101.50
2	Y	44	C	P-O3'-C3'	5.23	125.97	119.70
1	X	620	G	P-O3'-C3'	5.22	125.97	119.70
1	X	982	C	O4'-C1'-N1	5.22	112.38	108.20
1	X	2184	C	O4'-C1'-N1	5.22	112.38	108.20
1	X	1562	G	C4'-C3'-C2'	5.22	107.82	102.60
1	X	1918	G	P-O3'-C3'	5.22	125.97	119.70
1	X	698	A	C1'-O4'-C4'	-5.22	105.72	109.90
1	X	795	A	N1-C6-N6	5.22	121.73	118.60
1	X	1407	G	C4-N9-C1'	5.22	133.28	126.50
1	X	1986	G	N3-C4-C5	-5.22	125.99	128.60
1	X	1148	G	C5-C6-O6	-5.21	125.47	128.60
1	X	542	A	N1-C6-N6	5.21	121.73	118.60
1	X	652	C	C4'-C3'-C2'	5.21	107.81	102.60
1	X	1249	G	N1-C6-O6	-5.21	116.77	119.90
1	X	1363	C	O4'-C1'-N1	5.21	112.37	108.20
1	X	1721	G	O4'-C4'-C3'	-5.21	98.79	104.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	859	U	P-O3'-C3'	5.21	125.95	119.70
1	X	2812	A	C8-N9-C4	-5.21	103.72	105.80
1	X	1501	C	O4'-C1'-N1	5.21	112.37	108.20
2	Y	22	U	O4'-C1'-N1	5.21	112.37	108.20
1	X	200	A	P-O3'-C3'	5.21	125.95	119.70
1	X	661	C	C6-N1-C2	-5.21	118.22	120.30
1	X	746	G	N3-C4-N9	5.21	129.12	126.00
1	X	1529	C	O4'-C1'-N1	5.21	112.37	108.20
1	X	2646	C	O4'-C1'-N1	5.21	112.36	108.20
2	Y	14	C	C3'-C2'-C1'	5.21	105.67	101.50
1	X	1123	G	P-O3'-C3'	5.20	125.94	119.70
1	X	2669	C	C2-N1-C1'	5.20	124.52	118.80
1	X	2748	C	P-O3'-C3'	-5.20	113.46	119.70
2	Y	50	U	C3'-C2'-C1'	-5.20	97.34	101.50
1	X	841	G	C1'-O4'-C4'	-5.20	105.74	109.90
1	X	503	G	O4'-C4'-C3'	-5.20	98.80	104.00
1	X	786	U	O4'-C1'-N1	5.20	112.36	108.20
1	X	1470	G	C8-N9-C4	-5.20	104.32	106.40
1	X	1679	U	C2-N3-C4	-5.20	123.88	127.00
1	X	1774	A	P-O3'-C3'	5.20	125.94	119.70
1	X	2065	A	O4'-C1'-N9	5.20	112.36	108.20
1	X	1351	G	O4'-C1'-N9	5.20	112.36	108.20
1	X	2380	U	O4'-C1'-N1	5.20	112.36	108.20
1	X	89	A	N9-C1'-C2'	5.20	120.75	114.00
1	X	943	U	N1-C1'-C2'	5.20	120.75	114.00
1	X	2692	A	O3'-P-O5'	-5.20	94.13	104.00
1	X	623	G	O4'-C1'-N9	5.19	112.36	108.20
1	X	518	A	P-O5'-C5'	5.19	129.21	120.90
1	X	991	A	N9-C1'-C2'	5.19	120.75	114.00
1	X	2541	U	N3-C2-O2	-5.19	118.57	122.20
1	X	420	C	O4'-C1'-N1	5.19	112.35	108.20
1	X	594	G	O4'-C1'-N9	5.18	112.35	108.20
1	X	1277	G	N1-C6-O6	-5.18	116.79	119.90
1	X	1467	U	C2-N3-C4	5.18	130.11	127.00
1	X	71	A	C4'-C3'-C2'	5.18	107.78	102.60
1	X	2533	U	N1-C2-O2	5.18	126.43	122.80
1	X	1819	U	N3-C2-O2	-5.18	118.58	122.20
1	X	2032	G	C2'-C3'-O3'	5.18	121.99	113.70
1	X	521	U	C2-N1-C1'	5.17	123.91	117.70
1	X	1006	C	C1'-O4'-C4'	-5.17	105.76	109.90
1	X	2421	C	P-O5'-C5'	5.17	129.18	120.90
1	X	1219	C	N1-C2-O2	5.17	122.00	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1770	U	C4-C5-C6	5.17	122.80	119.70
1	X	422	C	C6-N1-C2	-5.17	118.23	120.30
1	X	1685	A	P-O3'-C3'	5.17	125.90	119.70
1	X	2569	A	P-O5'-C5'	-5.17	112.63	120.90
1	X	683	A	C4'-C3'-C2'	5.17	107.77	102.60
1	X	1132	C	C5-C6-N1	5.16	123.58	121.00
2	Y	73	C	O4'-C1'-N1	5.16	112.33	108.20
1	X	2226	A	C5-C6-N6	-5.16	119.57	123.70
2	Y	119	G	C3'-C2'-C1'	-5.16	97.37	101.50
1	X	1799	A	O4'-C1'-N9	5.16	112.33	108.20
1	X	1222	G	P-O3'-C3'	5.16	125.89	119.70
1	X	2018	G	C3'-C2'-C1'	-5.16	97.37	101.50
2	Y	53	G	N3-C4-N9	5.16	129.09	126.00
1	X	460	U	O5'-P-OP1	-5.16	101.06	105.70
1	X	1482	U	C1'-O4'-C4'	-5.16	105.77	109.90
1	X	2731	G	O4'-C1'-N9	5.16	112.33	108.20
1	X	822	G	C4'-C3'-C2'	-5.16	97.44	102.60
1	X	1312	G	N7-C8-N9	5.16	115.68	113.10
2	Y	107	C	P-O3'-C3'	5.16	125.89	119.70
1	X	2751	C	O4'-C1'-N1	5.15	112.32	108.20
2	Y	92	G	C5-C6-N1	5.15	114.08	111.50
1	X	1152	C	O4'-C1'-N1	5.15	112.32	108.20
1	X	1223	G	N3-C2-N2	5.15	123.51	119.90
1	X	1244	U	C5-C6-N1	5.15	125.28	122.70
1	X	1628	C	C6-N1-C2	-5.15	118.24	120.30
1	X	2525	U	N3-C2-O2	-5.15	118.60	122.20
1	X	1294	G	O4'-C1'-N9	5.15	112.32	108.20
1	X	1327	C	C5-C6-N1	5.15	123.57	121.00
1	X	1677	C	C4'-C3'-C2'	-5.15	97.45	102.60
1	X	2782	G	N1-C6-O6	5.14	122.99	119.90
1	X	2500	C	N1-C2-O2	5.14	121.98	118.90
1	X	2700	U	OP1-P-OP2	-5.14	111.89	119.60
1	X	1674	C	OP1-P-O3'	5.14	116.51	105.20
1	X	927	C	N1-C2-O2	5.14	121.98	118.90
1	X	983	G	C8-N9-C4	-5.14	104.34	106.40
1	X	1764	A	C5-C6-N6	-5.14	119.59	123.70
1	X	2699	G	N3-C4-C5	-5.14	126.03	128.60
1	X	1383	C	N1-C2-O2	5.14	121.98	118.90
1	X	2460	G	O4'-C1'-N9	5.14	112.31	108.20
1	X	779	U	O4'-C1'-N1	5.14	112.31	108.20
1	X	1574	A	P-O5'-C5'	5.14	129.12	120.90
1	X	2858	A	O4'-C1'-N9	5.14	112.31	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	761	G	C1'-O4'-C4'	-5.13	105.79	109.90
1	X	1988	A	N9-C4-C5	-5.13	103.75	105.80
1	X	2523	G	C6-C5-N7	-5.13	127.32	130.40
1	X	1709	U	P-O3'-C3'	5.13	125.86	119.70
1	X	1705	U	O4'-C1'-N1	5.13	112.31	108.20
1	X	1805	G	N3-C4-C5	-5.13	126.03	128.60
1	X	1547	U	O4'-C1'-N1	5.13	112.30	108.20
1	X	2077	G	O4'-C1'-N9	5.13	112.30	108.20
1	X	1801	C	P-O3'-C3'	5.13	125.85	119.70
1	X	2430	A	N1-C6-N6	-5.13	115.52	118.60
1	X	957	G	C4-C5-N7	-5.12	108.75	110.80
1	X	191	G	O4'-C1'-N9	5.12	112.30	108.20
1	X	2165	A	C3'-C2'-C1'	5.12	105.60	101.50
1	X	30	G	N3-C4-C5	-5.12	126.04	128.60
1	X	509	U	P-O3'-C3'	5.12	125.85	119.70
1	X	540	G	N3-C4-C5	-5.12	126.04	128.60
1	X	1223	G	N3-C4-N9	5.12	129.07	126.00
1	X	459	A	P-O3'-C3'	5.12	125.84	119.70
1	X	1631	C	O5'-C5'-C4'	-5.12	101.97	111.70
1	X	2847	G	N7-C8-N9	5.12	115.66	113.10
1	X	776	G	N9-C1'-C2'	5.12	120.65	114.00
1	X	2535	C	O4'-C1'-N1	5.12	112.29	108.20
1	X	2621	G	OP2-P-O3'	5.12	116.45	105.20
1	X	825	C	P-O5'-C5'	5.11	129.08	120.90
1	X	1245	G	P-O3'-C3'	5.11	125.84	119.70
1	X	1306	U	C3'-C2'-C1'	-5.11	97.41	101.50
1	X	1094	C	O4'-C1'-N1	5.11	112.29	108.20
1	X	1711	C	C5'-C4'-O4'	5.11	115.23	109.10
1	X	2408	G	C2-N3-C4	5.11	114.45	111.90
1	X	2635	U	O4'-C1'-N1	5.11	112.29	108.20
1	X	729	A	C1'-O4'-C4'	-5.11	105.81	109.90
1	X	1312	G	P-O3'-C3'	5.11	125.83	119.70
1	X	538	A	P-O3'-C3'	5.11	125.83	119.70
1	X	559	C	C6-N1-C2	-5.11	118.26	120.30
1	X	2851	G	C5-C6-O6	-5.11	125.54	128.60
1	X	328	A	C5'-C4'-O4'	5.10	115.22	109.10
2	Y	54	U	N3-C2-O2	-5.10	118.63	122.20
2	Y	120	G	O4'-C1'-N9	5.10	112.28	108.20
1	X	89	A	C5'-C4'-O4'	5.10	115.22	109.10
1	X	1149	G	C4'-C3'-C2'	-5.10	97.50	102.60
2	Y	27	A	P-O3'-C3'	5.10	125.82	119.70
1	X	1012	A	O4'-C1'-N9	5.10	112.28	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	1946	U	N3-C2-O2	-5.10	118.63	122.20
1	X	2806	G	O4'-C1'-N9	5.10	112.28	108.20
1	X	2850	U	P-O5'-C5'	5.10	129.06	120.90
1	X	2867	G	C6-C5-N7	-5.10	127.34	130.40
1	X	1682	A	C8-N9-C4	-5.10	103.76	105.80
1	X	2508	G	N9-C1'-C2'	5.10	120.63	114.00
1	X	743	A	C3'-C2'-C1'	-5.10	97.42	101.50
1	X	858	G	N7-C8-N9	5.10	115.65	113.10
1	X	1083	C	O4'-C1'-N1	5.10	112.28	108.20
1	X	2385	U	O4'-C1'-N1	5.09	112.28	108.20
1	X	1690	U	N1-C1'-C2'	-5.09	106.40	112.00
1	X	2033	C	O4'-C1'-N1	5.09	112.28	108.20
1	X	312	G	O4'-C1'-N9	5.09	112.27	108.20
1	X	473	C	O4'-C1'-N1	5.09	112.27	108.20
1	X	830	C	N3-C2-O2	-5.09	118.34	121.90
1	X	430	C	C5-C6-N1	5.09	123.54	121.00
1	X	1980	A	N1-C6-N6	5.09	121.65	118.60
1	X	2279	G	C8-N9-C4	-5.09	104.37	106.40
1	X	2369	U	O4'-C1'-N1	5.09	112.27	108.20
1	X	2654	A	P-O3'-C3'	-5.09	113.59	119.70
5	C	194	GLU	C-N-CA	5.09	134.42	121.70
1	X	971	A	C2-N3-C4	5.08	113.14	110.60
1	X	1805	G	C2-N3-C4	5.08	114.44	111.90
1	X	2033	C	N1-C1'-C2'	5.08	120.61	114.00
1	X	2463	G	O4'-C1'-N9	5.08	112.27	108.20
1	X	745	C	N1-C2-O2	5.08	121.95	118.90
1	X	820	U	O4'-C1'-N1	5.08	112.27	108.20
1	X	2178	U	O4'-C1'-N1	5.08	112.27	108.20
2	Y	123	U	N3-C2-O2	-5.08	118.64	122.20
1	X	1681	A	C2-N3-C4	-5.08	108.06	110.60
1	X	2396	C	P-O3'-C3'	-5.08	113.61	119.70
1	X	2720	A	O4'-C4'-C3'	-5.08	98.92	104.00
1	X	2843	A	P-O5'-C5'	-5.08	112.77	120.90
2	Y	113	G	O4'-C1'-N9	5.08	112.26	108.20
1	X	496	C	P-O3'-C3'	-5.08	113.61	119.70
1	X	1504	G	P-O3'-C3'	5.08	125.79	119.70
1	X	2003	A	C2-N3-C4	5.08	113.14	110.60
1	X	1518	C	O4'-C1'-N1	5.08	112.26	108.20
1	X	1578	U	O4'-C1'-N1	5.08	112.26	108.20
1	X	2818	G	C4'-C3'-C2'	-5.08	97.52	102.60
1	X	855	G	O4'-C1'-N9	5.07	112.26	108.20
1	X	2181	A	O4'-C1'-N9	5.07	112.26	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	180	C	O4'-C1'-N1	5.07	112.26	108.20
1	X	177	U	O4'-C1'-N1	5.07	112.26	108.20
1	X	1253	C	O4'-C1'-N1	5.07	112.26	108.20
1	X	2777	A	P-O3'-C3'	5.07	125.78	119.70
1	X	1796	A	O4'-C1'-N9	5.07	112.25	108.20
1	X	164	G	C8-N9-C4	-5.07	104.37	106.40
1	X	2734	U	O4'-C1'-N1	5.07	112.25	108.20
1	X	2365	U	O4'-C1'-N1	5.07	112.25	108.20
1	X	2482	A	C5'-C4'-O4'	5.07	115.18	109.10
1	X	2560	G	C8-N9-C4	-5.07	104.37	106.40
1	X	2858	A	N1-C6-N6	5.07	121.64	118.60
1	X	991	A	N1-C2-N3	-5.06	126.77	129.30
5	C	171	PRO	C-N-CA	5.06	134.36	121.70
1	X	244	C	N1-C2-O2	5.06	121.94	118.90
1	X	430	C	C6-N1-C2	-5.06	118.28	120.30
1	X	1753	A	N7-C8-N9	5.06	116.33	113.80
1	X	756	C	C3'-C2'-C1'	-5.06	97.45	101.50
1	X	1120	C	P-O3'-C3'	5.06	125.77	119.70
1	X	1753	A	C2-N3-C4	5.06	113.13	110.60
1	X	1006	C	O4'-C1'-N1	5.06	112.25	108.20
1	X	598	U	C3'-C2'-C1'	-5.05	97.46	101.50
1	X	740	A	N9-C1'-C2'	5.05	120.57	114.00
1	X	1688	U	C5-C6-N1	5.05	125.23	122.70
1	X	172	A	P-O5'-C5'	5.05	128.99	120.90
1	X	611	C	C3'-C2'-C1'	-5.05	97.46	101.50
1	X	738	G	N3-C4-C5	-5.05	126.07	128.60
1	X	1124	U	O4'-C1'-N1	5.05	112.24	108.20
1	X	174	A	O5'-P-OP2	-5.05	101.15	105.70
1	X	1159	U	O4'-C1'-N1	5.05	112.24	108.20
1	X	2002	A	C5-C6-N6	-5.05	119.66	123.70
1	X	2229	G	C2-N3-C4	5.05	114.43	111.90
1	X	2869	U	O4'-C1'-N1	5.05	112.24	108.20
1	X	131	C	O4'-C1'-N1	5.05	112.24	108.20
1	X	223	C	C6-N1-C2	-5.05	118.28	120.30
1	X	2762	G	C6-C5-N7	-5.05	127.37	130.40
1	X	542	A	N1-C2-N3	5.05	131.82	129.30
1	X	972	C	N1-C1'-C2'	5.05	120.56	114.00
1	X	1712	G	C6-C5-N7	-5.05	127.37	130.40
1	X	2002	A	N1-C6-N6	5.05	121.63	118.60
1	X	2608	A	C1'-O4'-C4'	-5.05	105.86	109.90
1	X	203	G	C4'-C3'-C2'	-5.05	97.55	102.60
1	X	654	A	O4'-C1'-N9	5.05	112.24	108.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	778	G	C4'-C3'-C2'	-5.05	97.55	102.60
1	X	1764	A	C4-C5-N7	5.05	113.22	110.70
1	X	334	G	N3-C4-C5	-5.04	126.08	128.60
1	X	1467	U	O4'-C1'-C2'	-5.04	100.75	105.80
1	X	469	G	N3-C4-C5	-5.04	126.08	128.60
1	X	1613	G	C5'-C4'-O4'	5.04	115.15	109.10
1	X	2371	A	N7-C8-N9	5.04	116.32	113.80
1	X	2670	C	N3-C2-O2	-5.04	118.37	121.90
1	X	2422	C	O4'-C1'-N1	5.04	112.23	108.20
1	X	539	A	P-O3'-C3'	5.04	125.75	119.70
1	X	699	G	C2-N3-C4	-5.04	109.38	111.90
1	X	1570	C	N1-C2-O2	5.04	121.92	118.90
1	X	1592	U	O4'-C1'-N1	5.04	112.23	108.20
1	X	2387	U	O4'-C1'-N1	5.04	112.23	108.20
1	X	1229	C	O4'-C1'-N1	5.04	112.23	108.20
1	X	774	A	C5'-C4'-O4'	5.04	115.14	109.10
1	X	1407	G	N9-C1'-C2'	5.04	120.55	114.00
1	X	1541	G	C4'-C3'-C2'	-5.04	97.56	102.60
2	Y	97	C	N1-C2-O2	5.04	121.92	118.90
1	X	1280	U	P-O3'-C3'	5.03	125.74	119.70
1	X	2261	G	C5-C6-N1	5.03	114.02	111.50
1	X	1716	G	C8-N9-C4	-5.03	104.39	106.40
1	X	344	G	C8-N9-C4	-5.03	104.39	106.40
1	X	1304	U	O4'-C1'-N1	5.03	112.22	108.20
1	X	469	G	C2'-C3'-O3'	5.03	121.75	113.70
1	X	1376	C	N1-C2-O2	5.03	121.92	118.90
1	X	2299	A	P-O3'-C3'	5.03	125.73	119.70
1	X	2347	C	C3'-C2'-C1'	-5.03	97.48	101.50
1	X	2044	G	C6-C5-N7	-5.03	127.38	130.40
1	X	2069	U	O4'-C1'-N1	5.03	112.22	108.20
1	X	2592	U	O4'-C1'-N1	5.03	112.22	108.20
1	X	2090	U	O4'-C1'-N1	5.03	112.22	108.20
1	X	2230	G	C3'-C2'-C1'	-5.03	97.48	101.50
1	X	2453	C	N1-C2-O2	5.03	121.92	118.90
1	X	2553	G	C5-N7-C8	-5.02	101.79	104.30
1	X	2800	C	C5'-C4'-C3'	-5.02	107.97	116.00
1	X	404	A	O4'-C1'-N9	5.02	112.22	108.20
1	X	2587	G	N1-C6-O6	5.02	122.91	119.90
2	Y	39	C	N1-C2-O2	5.02	121.91	118.90
1	X	1244	U	O4'-C1'-N1	5.02	112.21	108.20
1	X	1347	C	OP2-P-O3'	5.02	116.24	105.20
1	X	1828	C	N1-C2-O2	5.02	121.91	118.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	2032	G	N3-C4-N9	5.02	129.01	126.00
1	X	2222	U	N3-C2-O2	-5.02	118.69	122.20
1	X	2567	G	O4'-C1'-N9	5.02	112.21	108.20
1	X	2572	U	N3-C4-O4	5.02	122.91	119.40
1	X	823	U	OP1-P-O3'	5.01	116.23	105.20
1	X	1201	G	N1-C6-O6	5.01	122.91	119.90
2	Y	14	C	N1-C2-O2	5.01	121.91	118.90
1	X	7	G	C4'-C3'-C2'	5.01	107.61	102.60
1	X	2327	U	N3-C2-O2	-5.01	118.69	122.20
1	X	610	G	O3'-P-O5'	-5.01	94.48	104.00
1	X	752	G	O3'-P-O5'	-5.01	94.48	104.00
1	X	1665	C	P-O5'-C5'	5.01	128.91	120.90
1	X	1956	G	C5-C6-O6	-5.01	125.59	128.60
1	X	130	C	O4'-C1'-N1	5.01	112.21	108.20
1	X	1333	G	C6-C5-N7	5.01	133.41	130.40
1	X	1729	C	O4'-C1'-N1	5.01	112.21	108.20
1	X	2239	C	N1-C2-O2	5.01	121.91	118.90
1	X	462	G	C4-C5-N7	-5.01	108.80	110.80
1	X	1683	G	N9-C4-C5	5.01	107.40	105.40
1	X	2423	G	O5'-P-OP2	-5.01	101.19	105.70
1	X	2561	G	C4-C5-N7	5.01	112.80	110.80
1	X	614	G	O4'-C1'-N9	5.00	112.20	108.20
1	X	2621	G	N3-C4-C5	-5.00	126.10	128.60
14	L	20	THR	C-N-CA	5.00	134.21	121.70
1	X	225	G	O4'-C1'-N9	5.00	112.20	108.20
1	X	2039	G	N7-C8-N9	5.00	115.60	113.10
1	X	2075	U	O4'-C1'-N1	5.00	112.20	108.20
1	X	2188	A	O4'-C1'-N9	5.00	112.20	108.20
1	X	2193	C	N1-C2-O2	5.00	121.90	118.90
1	X	454	G	C2'-C3'-O3'	5.00	121.70	113.70

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	X	1143	A	Sidechain
1	X	1337	G	Sidechain
1	X	474	G	Sidechain
1	X	671	A	Sidechain
1	X	683	A	Sidechain
1	X	805	G	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	X	57651	0	29049	404	0
2	Y	2598	0	1328	17	0
3	A	1826	0	1885	61	0
4	B	1539	0	1600	61	0
5	C	1506	0	1525	49	0
6	D	1400	0	1481	22	0
7	E	1286	0	1336	10	0
8	F	503	0	520	5	0
9	G	1114	0	1144	68	0
10	H	997	0	1046	31	0
11	I	1067	0	1103	37	0
12	J	1090	0	1125	36	0
13	K	878	0	930	36	0
14	L	779	0	820	25	0
15	M	871	0	894	25	0
16	N	978	0	1020	33	0
17	O	741	0	756	34	0
18	P	1014	0	1096	20	0
19	Q	726	0	753	15	0
20	R	825	0	881	27	0
21	S	1345	0	1372	18	0
22	T	625	0	655	11	0
23	U	552	0	604	26	0
24	V	533	0	558	5	0
25	W	424	0	470	8	0
26	Z	457	0	462	12	0
27	1	53	0	0	1	0
28	2	46	0	0	2	0
29	3	63	0	0	3	0
30	4	297	0	330	5	0
31	J	1	0	0	0	0
31	M	1	0	0	0	0
31	X	28	0	0	0	0
31	Y	5	0	0	0	0
32	X	56	0	64	3	0
All	All	83875	0	54807	954	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (954) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:M:57:ILE:CD1	15:M:57:ILE:CG1	1.79	1.58
11:I:57:ILE:CD1	11:I:57:ILE:CG1	1.92	1.45
13:K:52:ILE:CD1	13:K:52:ILE:CG1	1.95	1.45
9:G:100:TYR:HB2	9:G:116:ARG:NH1	1.66	1.08
11:I:62:LYS:HE2	11:I:64:GLY:HA2	1.34	1.03
1:X:787:A:H2	1:X:800:U:HO2'	1.06	1.03
1:X:1333:G:N2	1:X:1344:C:H41	1.55	1.03
1:X:1882:G:N2	1:X:1885:C:H41	1.55	1.03
23:U:31:GLY:HA2	23:U:32:ARG:HH11	1.18	1.02
4:B:152:LYS:HB2	9:G:106:TYR:HB3	1.42	1.01
1:X:617:U:H5	1:X:632:A:C2	1.82	0.97
3:A:43:ARG:N	3:A:43:ARG:HD2	1.82	0.95
11:I:62:LYS:CE	11:I:64:GLY:HA2	1.97	0.95
1:X:1919:A:H2	1:X:1926:U:H3	0.97	0.94
26:Z:4:HIS:HB3	26:Z:5:PRO:HD3	1.49	0.94
17:O:87:ARG:HG2	17:O:87:ARG:HH11	1.32	0.94
1:X:1030:U:H3	1:X:1153:A:H62	1.12	0.91
3:A:247:VAL:HG23	3:A:248:THR:HG23	1.52	0.91
1:X:1033:G:H22	1:X:1153:A:H2	1.20	0.90
1:X:2371:A:H2	1:X:2403:C:H42	1.19	0.90
1:X:1466:C:H2'	1:X:1467:U:O4'	1.73	0.89
9:G:61:ARG:HH11	9:G:66:HIS:H	1.15	0.89
9:G:100:TYR:HB2	9:G:116:ARG:HH12	1.35	0.89
4:B:116:VAL:HG22	4:B:136:ARG:HE	1.36	0.88
14:L:38:ILE:HG13	14:L:39:TYR:H	1.37	0.87
17:O:5:ILE:HD12	17:O:6:GLN:H	1.41	0.86
1:X:1542:G:H22	1:X:1562:G:H1	0.92	0.86
3:A:43:ARG:HD2	3:A:43:ARG:H	1.39	0.85
4:B:32:PRO:HB3	4:B:72:VAL:HG11	1.55	0.85
13:K:17:ARG:NH1	13:K:20:LEU:HD23	1.91	0.85
32:X:2929:1F2:H9	32:X:2929:1F2:H54	1.58	0.85
3:A:231:HIS:HD2	3:A:233:HIS:H	1.24	0.85
1:X:1030:U:H3	1:X:1153:A:N6	1.75	0.84
3:A:210:GLY:HA2	3:A:213:ARG:HB2	1.59	0.84
1:X:1919:A:H2	1:X:1926:U:N3	1.74	0.84
1:X:559:C:H2'	1:X:560:G:C1'	2.08	0.83
19:Q:61:LYS:H	19:Q:72:ARG:HA	1.42	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:215:G:H21	1:X:632:A:H8	1.25	0.83
1:X:1468:A:H5'	1:X:1472:C:N4	1.94	0.83
1:X:1007:A:H1'	17:O:6:GLN:HG2	1.61	0.83
1:X:1333:G:H22	1:X:1344:C:H41	1.26	0.82
1:X:1494:G:HO2'	1:X:1574:A:H2	1.26	0.82
9:G:33:ILE:HB	9:G:34:PRO:HD3	1.61	0.82
9:G:108:GLY:H	9:G:110:LEU:HG	1.45	0.82
1:X:1333:G:H22	1:X:1344:C:N4	1.78	0.82
13:K:17:ARG:HH11	13:K:20:LEU:HD23	1.41	0.81
4:B:116:VAL:HG22	4:B:136:ARG:NE	1.95	0.81
1:X:1333:G:N2	1:X:1344:C:N4	2.28	0.81
13:K:3:HIS:HB3	13:K:5:LYS:HD2	1.62	0.79
1:X:774:A:O5'	1:X:774:A:H8	1.65	0.78
1:X:559:C:H2'	1:X:560:G:H1'	1.66	0.78
9:G:93:LYS:HG2	9:G:96:ASP:HB3	1.65	0.78
1:X:2266:A:H62	1:X:2323:U:H3	1.29	0.77
3:A:244:ARG:HB3	3:A:252:LYS:HZ1	1.49	0.77
3:A:231:HIS:CD2	3:A:233:HIS:H	2.03	0.77
9:G:106:TYR:O	9:G:110:LEU:HG	1.85	0.77
1:X:1142:G:H21	9:G:101:THR:CG2	1.99	0.76
10:H:13:ASN:HD21	10:H:109:ARG:HG2	1.50	0.76
23:U:48:LYS:HG2	23:U:49:LYS:H	1.50	0.76
5:C:148:VAL:HG13	5:C:185:ARG:HB2	1.66	0.76
9:G:61:ARG:NH1	9:G:66:HIS:H	1.84	0.76
1:X:2797:G:OP2	13:K:3:HIS:NE2	2.18	0.76
13:K:11:ASN:OD1	13:K:17:ARG:NH2	2.18	0.76
4:B:147:PRO:C	4:B:149:ARG:H	1.88	0.75
12:J:78:LYS:HA	12:J:88:LYS:HZ3	1.49	0.75
20:R:107:ALA:HB1	20:R:111:GLY:HA2	1.68	0.75
23:U:31:GLY:HA2	23:U:32:ARG:NH1	1.99	0.74
1:X:2042:A:H5''	5:C:65:GLY:HA2	1.70	0.74
1:X:83:A:H5''	20:R:17:LYS:HG2	1.70	0.74
9:G:132:PHE:CZ	9:G:145:HIS:HB2	2.23	0.74
4:B:54:LYS:HB2	4:B:75:THR:O	1.88	0.74
1:X:1882:G:H22	1:X:1885:C:H41	1.31	0.73
1:X:631:G:H4'	1:X:632:A:H5'	1.70	0.73
4:B:152:LYS:CB	9:G:106:TYR:HB3	2.18	0.73
16:N:61:TRP:HH2	16:N:94:VAL:H	1.37	0.73
25:W:12:ARG:CG	25:W:12:ARG:HH11	2.01	0.73
16:N:66:ASN:HB3	16:N:76:TYR:HB2	1.70	0.73
1:X:1770:U:H5	1:X:1775:A:N7	1.86	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1373:G:H22	1:X:2192:U:H3	1.37	0.72
12:J:40:PRO:HB3	12:J:99:LYS:HD2	1.71	0.72
9:G:100:TYR:HB2	9:G:116:ARG:HH11	1.52	0.72
1:X:1266:G:N7	11:I:32:ARG:NH1	2.38	0.71
17:O:87:ARG:HG2	17:O:87:ARG:NH1	2.01	0.71
1:X:1142:G:H21	9:G:101:THR:HG22	1.54	0.71
5:C:43:ALA:HB1	5:C:86:PRO:HB2	1.72	0.71
1:X:2691:C:OP1	1:X:2694:G:H4'	1.90	0.71
12:J:27:TYR:HB2	12:J:137:VAL:HG21	1.72	0.71
1:X:1542:G:N2	1:X:1562:G:H1	1.78	0.70
14:L:8:ARG:HG3	14:L:9:ARG:N	2.06	0.70
1:X:512:A:H4'	18:P:15:LYS:HB3	1.73	0.70
1:X:823:U:OP1	11:I:32:ARG:NH1	2.24	0.70
6:D:4:LEU:HG	6:D:5:LYS:H	1.57	0.70
1:X:1448:A:H61	1:X:1574:A:H61	1.37	0.70
29:3:15:LYS:CA	29:3:16:ILE:CA	2.70	0.70
1:X:2241:U:H5	22:T:17:ASN:OD1	1.75	0.70
4:B:7:THR:HG21	15:M:5:ILE:HD11	1.74	0.70
1:X:542:A:C2	1:X:2004:U:H2'	2.27	0.69
1:X:227:G:H2'	1:X:228:A:C8	2.27	0.69
4:B:152:LYS:HB2	9:G:106:TYR:CB	2.22	0.69
9:G:104:THR:OG1	9:G:110:LEU:HB3	1.93	0.69
1:X:2561:G:H5'	1:X:2561:G:H8	1.55	0.69
3:A:172:TYR:HA	3:A:186:HIS:HA	1.73	0.69
7:E:107:ILE:HD11	7:E:151:VAL:HG12	1.74	0.69
1:X:2368:G:H5''	1:X:2369:U:H5'	1.72	0.69
25:W:12:ARG:HH11	25:W:12:ARG:HG2	1.57	0.69
12:J:27:TYR:CB	12:J:137:VAL:HG21	2.22	0.69
14:L:8:ARG:HG3	14:L:9:ARG:H	1.57	0.69
1:X:558:G:H4'	1:X:559:C:H5'	1.76	0.68
1:X:1673:C:H5''	4:B:136:ARG:HD3	1.76	0.68
17:O:88:GLN:HE21	17:O:88:GLN:HA	1.58	0.68
1:X:1745:C:P	15:M:101:ARG:HH22	2.16	0.68
15:M:34:ARG:NH2	15:M:88:VAL:HG13	2.08	0.68
17:O:12:TYR:HB2	17:O:40:VAL:H	1.58	0.68
22:T:23:VAL:HA	22:T:38:VAL:HG23	1.74	0.68
26:Z:4:HIS:HB3	26:Z:5:PRO:CD	2.24	0.68
11:I:62:LYS:NZ	11:I:64:GLY:HA2	2.09	0.68
23:U:32:ARG:HG2	23:U:33:LYS:H	1.58	0.68
1:X:1675:C:OP1	4:B:134:TRP:CE2	2.46	0.68
1:X:2653:A:O2'	10:H:41:ASN:ND2	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:504:G:H21	18:P:78:ASN:HD21	1.42	0.67
1:X:2561:G:H5'	1:X:2561:G:C8	2.29	0.67
20:R:23:ILE:HG22	20:R:33:THR:HB	1.75	0.67
5:C:45:THR:HG21	5:C:85:GLY:HA3	1.75	0.67
10:H:13:ASN:ND2	10:H:109:ARG:HG2	2.09	0.67
11:I:32:ARG:HD2	17:O:79:GLN:NE2	2.09	0.67
1:X:652:C:H42	1:X:657:A:H61	1.42	0.67
1:X:2387:U:H2'	1:X:2388:G:H8	1.58	0.67
9:G:67:ARG:HG2	9:G:70:PHE:HA	1.77	0.67
4:B:16:LYS:HB2	4:B:21:ILE:HD11	1.77	0.66
1:X:1257:U:H5''	11:I:17:LYS:HG3	1.77	0.66
1:X:1882:G:H21	1:X:1885:C:H41	1.43	0.66
1:X:323:G:OP1	1:X:343:A:H5''	1.94	0.66
1:X:1673:C:H5''	4:B:136:ARG:CD	2.26	0.66
1:X:1673:C:C5'	4:B:136:ARG:HD2	2.25	0.66
3:A:218:LYS:HE3	3:A:221:GLN:HB2	1.76	0.66
5:C:137:ALA:HB1	5:C:142:LEU:HB2	1.77	0.66
5:C:112:GLN:HA	5:C:116:LYS:HD3	1.78	0.66
17:O:57:GLN:H	17:O:97:GLY:HA3	1.61	0.66
1:X:1673:C:C5'	4:B:136:ARG:CD	2.74	0.66
1:X:1744:G:OP1	15:M:100:ARG:HD2	1.96	0.66
1:X:1770:U:C5	1:X:1775:A:N7	2.64	0.65
1:X:617:U:C5	1:X:632:A:C2	2.74	0.65
1:X:1673:C:H5'	4:B:136:ARG:HD2	1.79	0.65
1:X:2319:G:H2'	1:X:2320:G:H8	1.61	0.65
1:X:2594:U:H2'	1:X:2595:C:H6	1.61	0.65
20:R:90:LYS:HB2	20:R:108:VAL:HG21	1.78	0.65
1:X:640:C:H4'	1:X:660:G:H21	1.62	0.65
1:X:320:A:N3	1:X:340:G:O2'	2.29	0.64
14:L:38:ILE:HG13	14:L:39:TYR:N	2.11	0.64
1:X:1753:A:O5'	1:X:1753:A:H8	1.81	0.64
4:B:134:TRP:HD1	4:B:134:TRP:H	1.45	0.64
9:G:33:ILE:HB	9:G:34:PRO:CD	2.27	0.64
20:R:10:HIS:HD2	20:R:44:GLN:NE2	1.96	0.64
3:A:200:GLU:HG3	3:A:202:LYS:HB2	1.80	0.64
3:A:86:PRO:O	3:A:87:ASN:HB2	1.97	0.64
6:D:80:ARG:HD3	6:D:83:MET:HB2	1.80	0.63
23:U:52:ARG:HG3	23:U:79:GLU:HA	1.80	0.63
1:X:1468:A:H5'	1:X:1472:C:H42	1.64	0.63
1:X:482:A:H2'	1:X:483:A:O4'	1.98	0.63
1:X:564:U:H2'	1:X:565:A:C8	2.34	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:K:3:HIS:CG	13:K:5:LYS:HZ2	2.15	0.63
1:X:1113:C:H2'	1:X:1114:A:H8	1.64	0.63
17:O:87:ARG:HH11	17:O:87:ARG:CG	2.10	0.62
5:C:148:VAL:O	5:C:167:VAL:HA	2.00	0.62
13:K:11:ASN:ND2	13:K:12:ARG:HE	1.97	0.62
23:U:22:GLY:HA3	23:U:39:LYS:HD2	1.82	0.62
9:G:107:GLN:HA	9:G:110:LEU:HB2	1.81	0.62
1:X:2477:C:H6	1:X:2477:C:H5'	1.65	0.62
1:X:2042:A:H5''	5:C:65:GLY:CA	2.30	0.62
2:Y:30:C:OP1	14:L:37:HIS:HB3	2.00	0.62
11:I:18:ARG:HB3	11:I:21:ARG:HB2	1.81	0.62
1:X:504:G:H4'	18:P:27:VAL:HG13	1.81	0.62
1:X:971:A:H61	12:J:83:ARG:HH22	1.47	0.62
10:H:98:ILE:HG22	10:H:106:ARG:HG3	1.81	0.62
13:K:79:VAL:HA	13:K:83:VAL:HG13	1.81	0.62
19:Q:53:ILE:HD13	19:Q:80:VAL:HG13	1.80	0.62
1:X:1737:G:H2'	1:X:1738:U:C6	2.36	0.61
1:X:2659:C:H5'	4:B:189:PRO:HA	1.82	0.61
1:X:2790:C:O2'	26:Z:43:HIS:HD2	1.83	0.61
11:I:17:LYS:O	11:I:18:ARG:HB2	2.00	0.61
1:X:2779:C:H2'	1:X:2780:A:C8	2.36	0.61
1:X:2222:U:H2'	1:X:2223:U:C6	2.35	0.61
11:I:62:LYS:HG2	11:I:64:GLY:H	1.66	0.61
1:X:559:C:H2'	1:X:560:G:O4'	2.01	0.60
4:B:152:LYS:H	9:G:106:TYR:HB3	1.65	0.60
17:O:73:LYS:HB2	17:O:82:ARG:HB2	1.82	0.60
16:N:93:LYS:HD3	17:O:5:ILE:HB	1.84	0.60
23:U:22:GLY:HA3	23:U:39:LYS:CD	2.31	0.60
3:A:226:MET:HG2	3:A:230:ASP:HB2	1.84	0.60
2:Y:45:C:H2'	6:D:92:ARG:NH1	2.17	0.60
32:X:2929:1F2:H9	32:X:2929:1F2:C43	2.29	0.60
1:X:857:U:H3'	1:X:858:G:C8	2.37	0.60
3:A:58:HIS:O	3:A:58:HIS:ND1	2.35	0.60
4:B:75:THR:O	4:B:76:ARG:HB2	2.00	0.60
14:L:38:ILE:HG21	14:L:71:VAL:HG21	1.84	0.60
1:X:954:U:OP2	11:I:38:LYS:HG2	2.01	0.60
1:X:636:G:H5''	1:X:636:G:H8	1.67	0.59
1:X:689:A:H8	1:X:2052:G:H21	1.50	0.59
10:H:83:ARG:HH21	10:H:89:ILE:HD11	1.67	0.59
1:X:2845:C:H5''	13:K:65:LEU:HD11	1.84	0.59
1:X:2044:G:OP1	5:C:62:LYS:NZ	2.36	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:45:ASN:CG	3:A:46:ARG:H	2.05	0.59
1:X:1811:A:H5''	3:A:161:THR:HG21	1.83	0.59
12:J:19:THR:HG23	12:J:99:LYS:HD3	1.84	0.59
1:X:2795:A:H4'	13:K:5:LYS:HG2	1.83	0.59
4:B:122:PHE:O	4:B:123:ALA:HB2	2.02	0.59
6:D:75:SER:HB2	6:D:79:LEU:HD13	1.82	0.59
1:X:793:G:H21	1:X:796:A:H62	1.50	0.59
1:X:172:A:H61	1:X:175:C:H3'	1.68	0.59
1:X:1173:G:H21	17:O:88:GLN:HE22	1.50	0.59
14:L:33:ARG:CZ	14:L:38:ILE:HB	2.33	0.59
1:X:1466:C:C2'	1:X:1467:U:O4'	2.47	0.59
1:X:2310:G:H4'	22:T:43:THR:H	1.68	0.59
5:C:164:VAL:C	5:C:166:TRP:H	2.05	0.59
13:K:3:HIS:CG	13:K:5:LYS:NZ	2.70	0.59
16:N:37:GLN:HA	16:N:40:LEU:HD12	1.85	0.59
1:X:1448:A:H61	1:X:1574:A:N6	2.01	0.59
14:L:38:ILE:HG13	14:L:40:ALA:H	1.67	0.59
1:X:692:C:H2'	1:X:693:A:H8	1.68	0.58
23:U:27:ASP:H	23:U:32:ARG:HH21	1.51	0.58
1:X:746:G:N7	1:X:774:A:C6	2.71	0.58
1:X:1142:G:OP1	9:G:107:GLN:O	2.21	0.58
1:X:333:A:H2'	5:C:162:ARG:HH12	1.68	0.58
9:G:132:PHE:HB2	9:G:145:HIS:CE1	2.38	0.58
16:N:93:LYS:CE	17:O:5:ILE:HD13	2.34	0.58
16:N:93:LYS:HE3	17:O:5:ILE:HD13	1.85	0.58
19:Q:29:VAL:HG11	19:Q:38:ILE:HD11	1.85	0.58
20:R:105:ARG:HH12	20:R:112:LYS:HA	1.69	0.58
10:H:132:GLU:HG2	10:H:134:LEU:HG	1.84	0.58
3:A:133:LEU:HB2	3:A:187:SER:HA	1.86	0.58
1:X:1675:C:OP1	4:B:134:TRP:NE1	2.36	0.58
11:I:38:LYS:HD2	11:I:40:ARG:O	2.04	0.58
1:X:2594:U:H2'	1:X:2595:C:C6	2.39	0.57
12:J:36:ILE:HD11	12:J:103:VAL:HG22	1.86	0.57
23:U:17:SER:HB2	23:U:44:ALA:HA	1.86	0.57
1:X:405:C:H2'	1:X:406:G:H8	1.69	0.57
1:X:2387:U:H2'	1:X:2388:G:C8	2.39	0.57
13:K:87:TYR:HE1	13:K:94:TYR:HD2	1.51	0.57
16:N:66:ASN:HB3	16:N:76:TYR:CB	2.35	0.57
1:X:1342:U:H5''	1:X:1343:C:H5	1.68	0.57
12:J:44:LYS:HD2	12:J:47:GLN:HE22	1.68	0.57
1:X:692:C:H2'	1:X:693:A:C8	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:939:C:H5'	1:X:939:C:H6	1.70	0.57
1:X:1467:U:H2'	1:X:1468:A:OP1	2.05	0.57
1:X:1773:C:H1'	1:X:2588:U:H5''	1.87	0.57
9:G:67:ARG:CG	9:G:70:PHE:HA	2.35	0.57
11:I:97:ARG:O	11:I:98:LEU:HB2	2.05	0.56
18:P:28:ALA:HB2	18:P:71:VAL:HG22	1.87	0.56
1:X:1142:G:C1'	9:G:103:TYR:HD2	2.18	0.56
1:X:794:A:H5'	3:A:218:LYS:HZ2	1.71	0.56
18:P:14:ARG:HA	18:P:17:GLN:HG2	1.87	0.56
21:S:95:SER:HB3	21:S:119:ASN:HB3	1.87	0.56
1:X:333:A:H2'	5:C:162:ARG:NH1	2.20	0.56
20:R:25:LEU:HD12	20:R:81:VAL:N	2.20	0.56
23:U:31:GLY:C	23:U:32:ARG:HE	2.08	0.56
6:D:150:ARG:HE	6:D:151:GLY:H	1.54	0.56
8:F:117:ALA:HB1	8:F:122:ALA:HB1	1.87	0.56
13:K:13:ASN:ND2	13:K:15:SER:OG	2.38	0.56
26:Z:35:GLN:O	26:Z:37:HIS:N	2.39	0.56
1:X:1805:G:H1'	3:A:50:THR:CG2	2.36	0.56
1:X:1673:C:H5'	4:B:136:ARG:CD	2.36	0.56
9:G:103:TYR:CG	9:G:111:LYS:HA	2.40	0.56
1:X:485:G:C6	1:X:520:C:N4	2.74	0.56
21:S:6:LYS:H	21:S:7:PRO:HD3	1.71	0.56
1:X:38:G:H21	5:C:42:THR:HG21	1.71	0.56
1:X:794:A:H2	1:X:1767:G:N3	2.04	0.56
9:G:162:LYS:H	9:G:163:PRO:HD3	1.71	0.56
1:X:794:A:H5'	3:A:218:LYS:NZ	2.21	0.55
1:X:2490:U:H2'	1:X:2491:C:O4'	2.06	0.55
4:B:147:PRO:HB2	4:B:149:ARG:HD2	1.87	0.55
9:G:105:GLY:O	9:G:106:TYR:C	2.42	0.55
4:B:149:ARG:NH2	9:G:106:TYR:HD1	2.04	0.55
1:X:2319:G:H2'	1:X:2320:G:C8	2.42	0.55
1:X:1673:C:C5'	4:B:136:ARG:HD3	2.36	0.55
1:X:2620:G:H5''	9:G:104:THR:HB	1.89	0.55
11:I:10:PRO:HA	11:I:14:LYS:HB2	1.89	0.55
23:U:62:LEU:HD23	23:U:67:LEU:HD12	1.89	0.55
3:A:86:PRO:O	3:A:87:ASN:CB	2.54	0.55
1:X:1122:A:O2'	1:X:1123:G:H4'	2.07	0.55
9:G:68:PRO:HD2	9:G:76:GLN:HB3	1.88	0.55
25:W:1:MET:HB3	25:W:34:VAL:HG12	1.89	0.55
30:4:19:ARG:HD2	30:4:24:LEU:HD22	1.89	0.55
1:X:38:G:H1	1:X:453:U:H3	1.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:2516:U:H2'	1:X:2517:C:C6	2.42	0.55
3:A:43:ARG:HB3	3:A:54:ILE:HG13	1.89	0.55
16:N:66:ASN:HB3	16:N:76:TYR:H	1.71	0.55
1:X:1008:G:H5''	16:N:92:ARG:HB3	1.89	0.54
1:X:1268:U:C2	5:C:66:ASN:HA	2.42	0.54
1:X:1278:A:H61	1:X:1996:A:H5''	1.72	0.54
2:Y:53:G:H21	2:Y:54:U:H5''	1.72	0.54
9:G:157:PRO:C	9:G:159:SER:H	2.10	0.54
16:N:101:ARG:O	16:N:103:PRO:HD3	2.08	0.54
3:A:243:GLY:C	3:A:244:ARG:HD3	2.28	0.54
17:O:10:LYS:HG3	17:O:11:GLN:HG2	1.90	0.54
1:X:742:G:N1	3:A:208:LYS:HD3	2.23	0.54
1:X:654:A:H2'	1:X:655:A:H5'	1.89	0.54
1:X:1962:C:H2'	1:X:1963:G:H5'	1.90	0.54
5:C:158:ARG:HA	5:C:169:VAL:HG21	1.90	0.54
1:X:494:A:C8	20:R:56:LYS:HD2	2.43	0.54
8:F:79:ARG:HG2	8:F:84:ILE:HB	1.89	0.54
1:X:219:G:N2	1:X:231:G:H2'	2.22	0.54
1:X:1337:G:OP2	18:P:105:ARG:NH1	2.41	0.54
9:G:106:TYR:O	9:G:110:LEU:CG	2.56	0.54
24:V:25:LEU:HD21	24:V:47:ARG:HG2	1.90	0.54
1:X:224:G:H4'	1:X:399:G:C5	2.42	0.53
1:X:797:A:C5	3:A:229:VAL:HG21	2.43	0.53
1:X:969:U:C4	12:J:17:ARG:HB2	2.43	0.53
4:B:116:VAL:HG13	4:B:136:ARG:HH21	1.73	0.53
13:K:49:GLU:O	13:K:52:ILE:HG12	2.08	0.53
23:U:53:GLU:HB2	23:U:56:GLN:O	2.09	0.53
5:C:133:PHE:HB2	5:C:160:ALA:HB1	1.89	0.53
11:I:58:ALA:O	11:I:59:ARG:HB3	2.08	0.53
1:X:2761:A:H5''	1:X:2762:G:H5'	1.89	0.53
1:X:617:U:H5	1:X:632:A:N1	2.04	0.53
1:X:1033:G:N2	1:X:1153:A:H2	1.98	0.53
1:X:1810:U:H2'	3:A:157:ARG:HD3	1.90	0.53
4:B:149:ARG:CZ	9:G:106:TYR:HD1	2.21	0.53
5:C:41:GLY:HA3	5:C:89:ARG:O	2.09	0.53
12:J:28:VAL:HG12	12:J:29:ALA:H	1.72	0.53
16:N:21:ALA:HB1	16:N:29:SER:HA	1.91	0.53
1:X:2241:U:C5	22:T:17:ASN:OD1	2.59	0.53
6:D:40:LEU:HD21	6:D:87:ILE:HD12	1.91	0.53
12:J:28:VAL:HG23	12:J:137:VAL:HB	1.91	0.53
1:X:617:U:H5	1:X:632:A:H2	1.50	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:O:69:ILE:HG22	17:O:86:HIS:HB3	1.90	0.53
18:P:38:VAL:HG12	18:P:97:VAL:HG21	1.90	0.53
21:S:91:PRO:HG2	21:S:125:PRO:HD2	1.90	0.53
1:X:1373:G:N2	1:X:2192:U:H3	2.04	0.53
4:B:152:LYS:H	9:G:106:TYR:CB	2.22	0.53
14:L:76:ALA:HB2	14:L:107:ALA:HA	1.91	0.53
2:Y:32:C:H1'	2:Y:59:A:H61	1.73	0.53
12:J:109:GLY:HA3	21:S:112:LEU:HD21	1.90	0.52
16:N:66:ASN:HB2	16:N:70:ARG:HH11	1.73	0.52
1:X:1033:G:N2	1:X:1153:A:C2	2.75	0.52
3:A:42:GLY:C	3:A:43:ARG:HH11	2.12	0.52
4:B:183:LEU:HD11	15:M:16:ILE:HG21	1.90	0.52
1:X:760:U:C6	26:Z:3:LYS:HG3	2.44	0.52
1:X:2545:A:H61	10:H:40:GLY:HA3	1.74	0.52
1:X:1468:A:H5'	1:X:1472:C:H41	1.73	0.52
1:X:1765:C:N3	3:A:208:LYS:HE2	2.24	0.52
2:Y:16:U:H3'	2:Y:17:A:H5''	1.90	0.52
10:H:41:ASN:ND2	10:H:41:ASN:H	2.08	0.52
17:O:38:LEU:HD23	17:O:47:PHE:HB3	1.91	0.52
19:Q:5:ASP:O	19:Q:7:LEU:N	2.43	0.52
10:H:27:SER:HA	10:H:50:ILE:HD12	1.91	0.52
22:T:50:GLY:O	22:T:62:LEU:HB2	2.10	0.52
1:X:879:A:H2'	1:X:879:A:N3	2.25	0.52
1:X:2397:A:H2'	1:X:2398:U:O4'	2.10	0.52
9:G:62:ILE:HG22	9:G:135:LEU:HD21	1.90	0.52
12:J:12:LYS:O	12:J:13:GLN:HB2	2.08	0.52
12:J:100:PRO:HB2	21:S:74:ARG:HG2	1.91	0.52
15:M:99:VAL:HG11	15:M:104:LEU:HD22	1.90	0.52
1:X:415:A:H61	1:X:436:A:H61	1.58	0.52
1:X:1845:A:N1	1:X:2070:G:H1'	2.25	0.52
5:C:30:VAL:HG11	5:C:177:VAL:HG21	1.92	0.52
10:H:2:ILE:HB	10:H:45:ALA:HB3	1.92	0.52
18:P:25:PHE:HD1	18:P:127:ILE:HD11	1.74	0.52
1:X:1939:U:H1'	1:X:2531:U:OP1	2.10	0.52
1:X:1943:A:H8	1:X:1943:A:H5''	1.74	0.52
10:H:24:VAL:HA	10:H:51:ILE:HG22	1.92	0.52
14:L:29:LEU:HB3	14:L:89:PHE:HA	1.92	0.52
1:X:1686:A:H5''	1:X:1687:C:OP2	2.09	0.52
9:G:98:LYS:HB3	9:G:116:ARG:HB2	1.92	0.52
19:Q:66:GLY:O	19:Q:68:PHE:N	2.30	0.52
1:X:1630:A:N1	18:P:114:ALA:HB2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1643:A:H61	1:X:1656:U:H3	1.59	0.51
1:X:1805:G:H1'	3:A:50:THR:HG21	1.90	0.51
1:X:2368:G:H5''	1:X:2369:U:C5'	2.39	0.51
16:N:24:PHE:O	16:N:29:SER:HB3	2.11	0.51
1:X:118:U:H4'	1:X:119:G:H5''	1.93	0.51
1:X:305:A:H5'	1:X:305:A:H8	1.75	0.51
1:X:1342:U:H5''	1:X:1343:C:C5	2.46	0.51
1:X:1586:A:H2'	1:X:1587:A:C8	2.45	0.51
1:X:2551:A:H62	4:B:145:LYS:HG3	1.74	0.51
4:B:147:PRO:C	4:B:149:ARG:N	2.58	0.51
13:K:10:LEU:HA	13:K:17:ARG:HG2	1.93	0.51
14:L:33:ARG:NH1	14:L:38:ILE:HB	2.25	0.51
3:A:182:LEU:HB2	3:A:268:ARG:O	2.10	0.51
6:D:8:TYR:O	6:D:12:VAL:HB	2.11	0.51
6:D:123:ASP:HB3	6:D:127:ASN:HB2	1.92	0.51
9:G:57:LEU:HD22	9:G:170:PRO:HA	1.92	0.51
14:L:36:LYS:HB3	14:L:64:LYS:HB2	1.93	0.51
16:N:81:ASN:HD22	16:N:117:ARG:HH21	1.58	0.51
21:S:51:LEU:HB3	21:S:65:LEU:HD12	1.92	0.51
1:X:486:U:H4'	1:X:519:C:H2'	1.92	0.51
20:R:51:VAL:HG21	20:R:76:LEU:HD21	1.92	0.51
1:X:504:G:N2	18:P:78:ASN:HD21	2.08	0.51
11:I:21:ARG:HH11	11:I:22:GLY:HA3	1.76	0.51
1:X:1737:G:H2'	1:X:1738:U:H6	1.75	0.51
9:G:70:PHE:HB2	16:N:64:ARG:HG2	1.93	0.51
13:K:45:ARG:HD2	13:K:95:THR:HG22	1.92	0.51
14:L:8:ARG:CG	14:L:9:ARG:H	2.22	0.51
1:X:2475:C:OP1	12:J:83:ARG:HB3	2.11	0.50
5:C:8:GLY:HA3	5:C:120:VAL:HB	1.93	0.50
9:G:103:TYR:HB3	9:G:107:GLN:HE21	1.76	0.50
24:V:7:ARG:HB2	24:V:60:LEU:HD11	1.92	0.50
3:A:43:ARG:N	3:A:43:ARG:CD	2.64	0.50
6:D:78:LYS:HG2	6:D:80:ARG:HH11	1.76	0.50
1:X:2083:G:H1	1:X:2172:U:H3	1.58	0.50
9:G:162:LYS:N	9:G:163:PRO:CD	2.74	0.50
23:U:48:LYS:HG2	23:U:49:LYS:N	2.21	0.50
23:U:49:LYS:HA	23:U:62:LEU:H	1.75	0.50
3:A:36:ALA:HB1	3:A:62:TYR:O	2.12	0.50
5:C:148:VAL:HB	5:C:167:VAL:HG12	1.94	0.50
10:H:25:LEU:HD11	10:H:52:VAL:HG23	1.93	0.50
19:Q:66:GLY:C	19:Q:68:PHE:H	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:954:U:P	11:I:38:LYS:HG2	2.52	0.50
12:J:92:GLU:HG3	12:J:93:TYR:CD2	2.46	0.50
18:P:17:GLN:HG3	18:P:18:VAL:HG23	1.94	0.50
6:D:143:TYR:HA	6:D:146:VAL:HG22	1.94	0.50
9:G:124:GLU:O	9:G:128:GLU:HB2	2.11	0.50
13:K:17:ARG:NH1	13:K:20:LEU:CD2	2.71	0.50
20:R:29:HIS:CD2	20:R:51:VAL:HG22	2.47	0.50
1:X:490:A:N3	1:X:492:G:H5''	2.27	0.50
1:X:670:U:H2'	1:X:671:A:C8	2.46	0.50
1:X:1142:G:N2	9:G:101:THR:HG22	2.22	0.50
1:X:654:A:H2'	1:X:655:A:C5'	2.42	0.50
16:N:44:THR:O	16:N:48:ARG:HG3	2.12	0.50
21:S:69:VAL:HG22	21:S:81:VAL:HG13	1.93	0.50
1:X:922:A:N1	1:X:2256:G:H1'	2.27	0.50
1:X:2034:A:O4'	4:B:141:ILE:HD12	2.12	0.50
4:B:50:GLY:HA3	4:B:75:THR:HG21	1.94	0.50
4:B:146:THR:OG1	4:B:147:PRO:HD3	2.12	0.50
7:E:154:PRO:HA	7:E:160:LYS:O	2.11	0.50
11:I:28:LYS:NZ	11:I:36:GLY:HA3	2.27	0.50
12:J:65:ILE:HG23	12:J:107:VAL:HG12	1.94	0.50
1:X:2272:A:H5''	14:L:15:ARG:HH21	1.77	0.49
1:X:2508:G:C8	1:X:2508:G:H5'	2.47	0.49
2:Y:45:C:H2'	6:D:92:ARG:HH11	1.77	0.49
4:B:147:PRO:O	4:B:149:ARG:N	2.45	0.49
5:C:22:VAL:HG13	5:C:106:MET:HG2	1.94	0.49
21:S:6:LYS:HB2	21:S:31:SER:HB3	1.94	0.49
23:U:14:VAL:O	23:U:15:VAL:HG22	2.12	0.49
1:X:742:G:N7	3:A:209:ALA:O	2.45	0.49
1:X:1134:C:H1'	30:4:18:ARG:HH12	1.77	0.49
1:X:1467:U:H6	1:X:1467:U:H3'	1.77	0.49
1:X:1493:A:H2'	1:X:1494:G:O4'	2.12	0.49
4:B:5:LEU:HG	4:B:195:LEU:HD11	1.94	0.49
9:G:132:PHE:HZ	9:G:142:ARG:HA	1.77	0.49
14:L:31:VAL:HG21	14:L:100:VAL:HG23	1.93	0.49
1:X:2352:A:H2'	1:X:2353:G:C8	2.46	0.49
7:E:6:LYS:HB3	7:E:69:ARG:HD3	1.92	0.49
7:E:11:VAL:HB	7:E:15:VAL:HG21	1.93	0.49
17:O:12:TYR:CB	17:O:40:VAL:H	2.23	0.49
17:O:71:ILE:HD11	17:O:86:HIS:HB2	1.94	0.49
23:U:51:ILE:HG23	23:U:59:THR:HA	1.94	0.49
1:X:2266:A:H2	1:X:2325:A:H62	1.58	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:I:117:ALA:HA	11:I:137:GLY:O	2.13	0.49
1:X:649:G:H1	1:X:660:G:H1	1.60	0.49
10:H:78:SER:HA	10:H:91:PHE:O	2.11	0.49
23:U:48:LYS:CG	23:U:49:LYS:H	2.24	0.49
23:U:49:LYS:HB2	23:U:61:TRP:CE3	2.48	0.49
1:X:426:C:HO2'	1:X:1863:U:HO2'	1.59	0.49
1:X:2867:G:O5'	1:X:2867:G:H8	1.95	0.49
12:J:77:LYS:O	12:J:79:PRO:HD3	2.12	0.49
13:K:11:ASN:HD22	13:K:12:ARG:HE	1.60	0.49
18:P:36:ARG:HA	18:P:39:ARG:HD2	1.93	0.49
1:X:29:U:H4'	16:N:11:ARG:HH22	1.78	0.49
1:X:2017:U:H2'	1:X:2018:G:H5''	1.95	0.49
10:H:110:VAL:HG23	10:H:129:LEU:HB2	1.94	0.49
16:N:88:ILE:HG12	17:O:49:GLU:HB2	1.94	0.49
18:P:32:ARG:HA	18:P:121:THR:HG22	1.94	0.49
23:U:23:LYS:HD2	23:U:35:THR:HG21	1.95	0.49
1:X:389:G:H2'	1:X:390:U:C6	2.47	0.49
12:J:27:TYR:HB3	12:J:137:VAL:HG21	1.93	0.49
1:X:2080:U:H3	1:X:2175:A:H61	1.60	0.49
1:X:2661:G:O6	1:X:2708:U:H1'	2.13	0.49
10:H:28:GLY:HA3	10:H:35:THR:OG1	2.13	0.49
1:X:91:A:H2'	1:X:92:U:C6	2.48	0.48
1:X:649:G:N2	1:X:660:G:N2	2.60	0.48
1:X:818:G:H1'	1:X:844:G:O2'	2.13	0.48
11:I:54:SER:C	11:I:56:LEU:H	2.17	0.48
14:L:15:ARG:HH11	14:L:15:ARG:HA	1.78	0.48
1:X:2505:G:H1'	30:4:1:MET:HB2	1.93	0.48
4:B:13:GLN:O	4:B:14:ILE:HD12	2.13	0.48
4:B:195:LEU:H	15:M:2:GLN:HG2	1.78	0.48
13:K:10:LEU:HD23	13:K:17:ARG:HB2	1.95	0.48
16:N:83:LEU:HD12	16:N:113:SER:HB2	1.93	0.48
18:P:103:LEU:HB2	18:P:119:LYS:HB2	1.96	0.48
21:S:6:LYS:N	21:S:7:PRO:HD3	2.27	0.48
1:X:12:U:O2	1:X:12:U:H2'	2.13	0.48
6:D:72:LYS:HE2	6:D:81:GLN:HE21	1.79	0.48
7:E:38:ASN:HB3	7:E:40:GLU:HG2	1.95	0.48
1:X:2352:A:H2'	1:X:2353:G:H8	1.78	0.48
1:X:2484:G:HO2'	1:X:2485:U:H6	1.59	0.48
14:L:27:LEU:HB2	14:L:87:VAL:HG22	1.95	0.48
1:X:89:A:H4'	1:X:90:G:H5''	1.96	0.48
5:C:164:VAL:O	5:C:166:TRP:N	2.40	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:1:40:TYR:CA	27:1:41:ASP:CA	2.91	0.48
1:X:1562:G:H5''	1:X:1563:U:H5'	1.94	0.48
1:X:1705:U:O2	1:X:1717:A:H8	1.97	0.48
9:G:157:PRO:O	9:G:159:SER:N	2.46	0.48
11:I:32:ARG:HD2	17:O:79:GLN:HE22	1.77	0.48
1:X:1278:A:H2	1:X:1997:A:H62	1.61	0.48
1:X:2270:U:H2'	1:X:2271:C:C6	2.47	0.48
5:C:5:ASN:HB2	5:C:10:ASN:HA	1.95	0.48
10:H:116:ARG:HH11	15:M:38:LYS:HD3	1.78	0.48
17:O:10:LYS:HD2	17:O:37:ALA:HB3	1.95	0.48
1:X:617:U:C5	1:X:632:A:N1	2.81	0.48
1:X:627:A:H2'	1:X:628:A:C8	2.49	0.48
1:X:1685:A:H5''	10:H:5:GLN:HG2	1.96	0.48
17:O:33:VAL:HG12	17:O:57:GLN:HE21	1.79	0.48
1:X:1997:A:H2'	1:X:1998:A:C8	2.49	0.48
4:B:133:LYS:HD2	4:B:137:ARG:HB3	1.96	0.48
23:U:51:ILE:HA	23:U:59:THR:O	2.13	0.48
1:X:1833:U:H2'	1:X:1834:G:C8	2.49	0.48
1:X:2289:A:H3'	1:X:2290:A:H8	1.78	0.48
3:A:43:ARG:H	3:A:43:ARG:CD	2.18	0.48
30:4:19:ARG:HB2	30:4:24:LEU:HD13	1.96	0.48
1:X:1777:A:H1'	1:X:1921:A:N6	2.28	0.47
4:B:117:MET:H	4:B:136:ARG:HG3	1.79	0.47
9:G:55:ALA:C	9:G:134:MET:HE1	2.35	0.47
20:R:105:ARG:HH21	20:R:107:ALA:HB2	1.77	0.47
1:X:712:A:H2'	1:X:713:G:O4'	2.15	0.47
1:X:1820:G:OP2	3:A:239:ARG:NH1	2.47	0.47
8:F:101:TRP:HZ3	8:F:140:GLY:HA3	1.78	0.47
1:X:1776:A:OP1	1:X:1965:U:H5'	2.15	0.47
3:A:45:ASN:CG	3:A:46:ARG:N	2.67	0.47
11:I:88:PHE:HE2	11:I:119:THR:HB	1.79	0.47
18:P:40:LEU:HB3	26:Z:25:LEU:HD13	1.96	0.47
1:X:738:G:O5'	1:X:738:G:H8	1.98	0.47
1:X:1787:U:H2'	1:X:1788:C:C6	2.49	0.47
12:J:81:GLU:HG2	12:J:82:THR:H	1.79	0.47
1:X:1992:G:H1'	13:K:106:ASP:O	2.13	0.47
26:Z:16:ARG:HD3	26:Z:20:ARG:CZ	2.44	0.47
1:X:334:G:C8	5:C:164:VAL:HG13	2.49	0.47
1:X:1859:A:H2'	1:X:1860:A:C8	2.50	0.47
4:B:59:VAL:HG21	4:B:74:PRO:HB2	1.96	0.47
4:B:137:ARG:HG3	4:B:138:PRO:HD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:118:VAL:HG22	5:C:188:ILE:HD12	1.97	0.47
5:C:149:LEU:HD23	5:C:180:ILE:HG22	1.96	0.47
18:P:59:PHE:HD2	26:Z:41:LEU:HD22	1.79	0.47
23:U:27:ASP:N	23:U:32:ARG:HH21	2.12	0.47
1:X:33:C:O2'	1:X:34:U:H5''	2.14	0.47
1:X:143:A:H2'	1:X:144:U:C6	2.49	0.47
1:X:1137:A:H4'	1:X:1138:A:O5'	2.15	0.47
1:X:2189:A:C2	1:X:2190:A:C4	3.02	0.47
9:G:106:TYR:O	9:G:110:LEU:CD1	2.62	0.47
24:V:56:VAL:C	24:V:58:ALA:H	2.18	0.47
1:X:86:U:H5''	1:X:87:G:H5'	1.95	0.47
1:X:774:A:O5'	1:X:774:A:C8	2.56	0.47
1:X:879:A:H5'	1:X:880:C:OP2	2.14	0.47
11:I:62:LYS:NZ	11:I:64:GLY:CA	2.76	0.47
16:N:61:TRP:CZ3	16:N:94:VAL:N	2.82	0.47
1:X:203:G:H21	1:X:205:A:H62	1.62	0.47
1:X:1586:A:H5'	3:A:38:PRO:HG3	1.95	0.47
1:X:2654:A:H5'	10:H:42:LYS:H	1.80	0.47
22:T:3:HIS:CD2	22:T:5:LYS:HB2	2.50	0.47
23:U:19:ILE:HA	23:U:42:GLN:HA	1.95	0.47
1:X:935:C:H2'	1:X:936:A:C8	2.50	0.47
1:X:1173:G:H4'	17:O:22:VAL:HG23	1.97	0.47
4:B:122:PHE:O	4:B:123:ALA:CB	2.63	0.47
10:H:77:THR:HA	10:H:94:ASN:HB3	1.96	0.47
10:H:83:ARG:NH2	10:H:89:ILE:HD11	2.28	0.47
20:R:105:ARG:HH22	20:R:111:GLY:C	2.17	0.47
1:X:577:U:OP1	11:I:40:ARG:NH2	2.48	0.46
3:A:106:LEU:H	3:A:106:LEU:HD12	1.79	0.46
5:C:186:LEU:HG	5:C:188:ILE:HG12	1.95	0.46
6:D:35:VAL:HG22	6:D:90:THR:HG23	1.97	0.46
1:X:2289:A:H3'	1:X:2290:A:C8	2.49	0.46
9:G:103:TYR:CZ	9:G:111:LYS:HB2	2.51	0.46
10:H:112:GLY:O	10:H:131:PRO:HD2	2.15	0.46
12:J:26:ASP:O	12:J:27:TYR:HD1	1.97	0.46
1:X:650:U:H2'	1:X:651:C:C6	2.50	0.46
1:X:923:A:C6	12:J:12:LYS:HG2	2.50	0.46
1:X:1469:U:OP2	1:X:1471:G:OP2	2.33	0.46
1:X:2581:A:H2'	1:X:2582:G:C4'	2.46	0.46
6:D:72:LYS:HG2	6:D:81:GLN:HG2	1.98	0.46
18:P:118:LYS:HD2	18:P:120:ARG:HH21	1.80	0.46
20:R:23:ILE:HG23	20:R:84:VAL:HG21	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
20:R:40:LEU:HB2	20:R:45:LYS:HB2	1.97	0.46
1:X:542:A:C2	1:X:2004:U:C2'	2.97	0.46
1:X:1030:U:N3	1:X:1153:A:N6	2.54	0.46
11:I:76:LYS:HB3	11:I:79:GLN:HG2	1.98	0.46
12:J:77:LYS:HG3	12:J:78:LYS:H	1.80	0.46
25:W:46:THR:HG22	25:W:47:VAL:HG13	1.97	0.46
1:X:1142:G:O5'	9:G:107:GLN:HB3	2.15	0.46
1:X:2705:A:H1'	1:X:2706:U:H2'	1.97	0.46
8:F:77:LEU:HD13	8:F:107:ILE:HG23	1.98	0.46
9:G:67:ARG:HE	9:G:70:PHE:HA	1.80	0.46
9:G:69:ASP:H	9:G:76:GLN:HE21	1.62	0.46
12:J:27:TYR:HB2	12:J:137:VAL:HG11	1.98	0.46
13:K:87:TYR:CE1	13:K:94:TYR:HD2	2.30	0.46
7:E:164:PHE:O	7:E:166:GLY:N	2.48	0.46
3:A:270:ILE:HG13	3:A:271:VAL:H	1.81	0.46
10:H:28:GLY:O	10:H:29:ILE:HB	2.16	0.46
1:X:334:G:H5'	5:C:162:ARG:NH2	2.31	0.46
1:X:673:G:H5'	5:C:93:TYR:CD1	2.51	0.46
1:X:719:A:H2'	1:X:720:A:O4'	2.16	0.46
16:N:74:MET:HB3	16:N:75:ASN:H	1.59	0.46
1:X:641:G:N2	1:X:643:A:H3'	2.31	0.46
9:G:156:HIS:N	9:G:157:PRO:CD	2.79	0.46
13:K:3:HIS:HB3	13:K:5:LYS:CD	2.42	0.46
1:X:510:G:N1	1:X:513:A:OP2	2.45	0.45
1:X:1469:U:P	1:X:1471:G:OP2	2.74	0.45
1:X:2326:C:H2'	1:X:2327:U:H6	1.81	0.45
8:F:104:VAL:HA	8:F:107:ILE:HD12	1.98	0.45
13:K:3:HIS:ND1	13:K:5:LYS:NZ	2.47	0.45
13:K:36:THR:HG23	13:K:41:ALA:HB2	1.98	0.45
20:R:105:ARG:HH22	20:R:112:LYS:HA	1.80	0.45
1:X:568:G:H2'	1:X:569:C:O4'	2.15	0.45
1:X:652:C:H5''	1:X:652:C:H6	1.81	0.45
1:X:834:A:H5'	1:X:835:U:H6	1.81	0.45
1:X:1744:G:H5''	15:M:100:ARG:HD3	1.98	0.45
3:A:182:LEU:HD12	3:A:269:PHE:HB2	1.97	0.45
12:J:42:TRP:CG	12:J:95:VAL:HG11	2.52	0.45
20:R:45:LYS:HA	20:R:76:LEU:O	2.16	0.45
28:2:39:ARG:CA	28:2:40:HIS:CA	2.95	0.45
5:C:4:ILE:HG23	5:C:13:ARG:HH22	1.80	0.45
1:X:597:U:O4	1:X:683:A:H1'	2.16	0.45
1:X:2372:A:H5''	11:I:61:PRO:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Y:46:G:H4'	6:D:92:ARG:HH12	1.82	0.45
5:C:74:VAL:HB	5:C:75:PRO:HD2	1.97	0.45
5:C:170:LEU:HA	5:C:171:PRO:HD3	1.84	0.45
24:V:21:ARG:HH11	24:V:53:LEU:HD11	1.81	0.45
1:X:2790:C:O2'	26:Z:43:HIS:CD2	2.65	0.45
10:H:123:PHE:HB3	10:H:126:ILE:HG13	1.98	0.45
20:R:22:VAL:HG11	20:R:80:LYS:HZ3	1.80	0.45
1:X:312:G:HO2'	1:X:313:U:H6	1.63	0.45
1:X:1974:U:H2'	1:X:1975:G:H5''	1.99	0.45
1:X:2273:C:H2'	1:X:2274:C:H6	1.82	0.45
5:C:127:ASP:HB2	5:C:129:LYS:HG3	1.99	0.45
12:J:26:ASP:O	12:J:27:TYR:CD1	2.70	0.45
1:X:347:C:H4'	20:R:15:HIS:CD2	2.52	0.45
1:X:1012:A:H2'	1:X:1013:G:O4'	2.16	0.45
1:X:2226:A:H2'	1:X:2227:C:H6	1.81	0.45
1:X:2821:G:H2'	1:X:2822:U:C6	2.51	0.45
4:B:59:VAL:HG21	4:B:74:PRO:CB	2.46	0.45
25:W:3:ILE:HD12	25:W:51:LEU:HD13	1.99	0.45
26:Z:16:ARG:O	26:Z:20:ARG:HD2	2.17	0.45
1:X:1882:G:H22	1:X:1885:C:N4	2.08	0.45
15:M:93:ILE:HG21	15:M:96:ARG:HG2	1.99	0.45
25:W:47:VAL:HB	25:W:50:LEU:HD12	1.99	0.45
1:X:1922:U:OP1	1:X:2583:U:O2'	2.33	0.45
1:X:2189:A:H2	1:X:2190:A:C5	2.35	0.45
1:X:2691:C:O2'	1:X:2693:U:H5'	2.17	0.45
3:A:252:LYS:HD2	3:A:253:PRO:HD3	1.98	0.45
4:B:23:VAL:HG21	4:B:183:LEU:HD13	1.98	0.45
9:G:66:HIS:HA	16:N:67:ALA:HB1	1.98	0.45
12:J:27:TYR:CZ	21:S:76:ARG:HB3	2.52	0.45
4:B:146:THR:CB	4:B:147:PRO:HD3	2.47	0.45
14:L:38:ILE:CG1	14:L:39:TYR:H	2.18	0.45
1:X:310:A:N1	1:X:333:A:O2'	2.42	0.44
1:X:649:G:H22	1:X:660:G:N2	2.15	0.44
1:X:763:A:OP1	1:X:1631:C:N4	2.43	0.44
1:X:1167:A:C8	16:N:51:ARG:HG3	2.52	0.44
1:X:1474:A:H2'	1:X:1474:A:N3	2.32	0.44
1:X:1688:U:HO2'	1:X:1690:U:H5	1.63	0.44
13:K:33:ARG:HD3	13:K:112:LEU:HD22	2.00	0.44
15:M:5:ILE:H	15:M:5:ILE:HG13	1.62	0.44
1:X:1142:G:O4'	9:G:107:GLN:HG2	2.17	0.44
2:Y:40:C:H42	2:Y:46:G:H1	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:C:45:THR:HG22	5:C:47:THR:OG1	2.17	0.44
7:E:25:LYS:HG3	7:E:34:THR:HG22	1.98	0.44
12:J:77:LYS:HG3	12:J:78:LYS:N	2.32	0.44
17:O:36:LYS:HD2	17:O:54:TYR:HB2	1.99	0.44
20:R:96:LYS:CG	20:R:97:GLN:H	2.31	0.44
1:X:679:C:H2'	1:X:680:U:C6	2.52	0.44
1:X:1833:U:H2'	1:X:1834:G:H8	1.80	0.44
4:B:115:GLY:HA2	4:B:157:ALA:CB	2.47	0.44
6:D:116:GLY:HA2	6:D:176:PRO:HB2	1.99	0.44
7:E:43:VAL:HB	7:E:52:VAL:HG13	1.99	0.44
9:G:58:ILE:HG12	9:G:80:VAL:HG11	1.99	0.44
16:N:75:ASN:ND2	16:N:77:SER:HB3	2.33	0.44
25:W:12:ARG:CG	25:W:12:ARG:NH1	2.72	0.44
1:X:577:U:H5''	1:X:956:A:N6	2.32	0.44
1:X:857:U:H3'	1:X:858:G:H8	1.79	0.44
1:X:2751:C:H5'	4:B:203:LYS:HD3	1.98	0.44
19:Q:11:VAL:HB	19:Q:26:SER:HB2	1.98	0.44
1:X:680:U:H2'	1:X:681:A:H5''	1.99	0.44
1:X:1032:A:C8	1:X:1032:A:H3'	2.52	0.44
3:A:67:PHE:HD2	3:A:153:ALA:HB3	1.83	0.44
3:A:145:LEU:HD13	3:A:163:VAL:HG11	2.00	0.44
1:X:573:C:HO2'	1:X:1266:G:H1	1.65	0.44
1:X:1509:A:H8	1:X:1510:A:C8	2.36	0.44
1:X:1827:G:H1'	1:X:1914:U:C2	2.53	0.44
2:Y:47:A:H8	6:D:92:ARG:CZ	2.30	0.44
3:A:123:ALA:HB3	3:A:131:LEU:HB3	2.00	0.44
9:G:158:HIS:HA	9:G:161:GLN:HB2	2.00	0.44
1:X:1007:A:H4'	16:N:93:LYS:HB3	2.00	0.44
1:X:1465:G:H2'	1:X:1466:C:C6	2.53	0.44
1:X:2445:C:H5''	30:4:6:SER:HB3	2.00	0.44
5:C:164:VAL:HB	5:C:165:SER:H	1.42	0.44
22:T:38:VAL:HG13	22:T:40:GLN:HG2	1.98	0.44
1:X:168:A:H2'	1:X:169:C:C6	2.53	0.44
1:X:400:U:H5	23:U:21:ARG:HH12	1.64	0.44
1:X:1050:G:H1	1:X:1127:C:H42	1.65	0.44
1:X:1385:C:H2'	1:X:1386:A:O4'	2.17	0.44
1:X:457:C:H5''	16:N:3:ARG:HB3	1.99	0.44
1:X:663:G:C5	1:X:664:C:H1'	2.52	0.44
1:X:1583:A:H3'	3:A:86:PRO:HG3	2.00	0.44
2:Y:116:C:H4'	14:L:49:GLN:HG2	2.00	0.44
3:A:67:PHE:CE2	3:A:106:LEU:HD11	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:67:PHE:HB3	3:A:153:ALA:H	1.83	0.44
4:B:11:MET:HA	4:B:23:VAL:O	2.17	0.44
4:B:32:PRO:CB	4:B:72:VAL:HG11	2.38	0.44
10:H:105:PRO:HG3	10:H:126:ILE:HD13	1.99	0.44
20:R:22:VAL:HG13	20:R:82:ALA:HA	2.00	0.44
14:L:42:ILE:HD12	14:L:87:VAL:HG21	2.00	0.43
1:X:834:A:H5'	1:X:835:U:C6	2.53	0.43
1:X:1811:A:H3'	3:A:178:PRO:HB2	1.99	0.43
2:Y:118:G:H21	14:L:39:TYR:HH	1.66	0.43
12:J:73:LYS:H	12:J:94:TRP:HD1	1.66	0.43
1:X:465:C:O2'	1:X:483:A:N6	2.52	0.43
5:C:47:THR:HG23	5:C:84:PHE:H	1.83	0.43
5:C:154:ASP:HB2	5:C:157:THR:OG1	2.17	0.43
12:J:86:LYS:H	12:J:86:LYS:HG2	1.68	0.43
1:X:800:U:H5''	1:X:801:A:H5'	2.01	0.43
1:X:1113:C:H2'	1:X:1114:A:C8	2.48	0.43
1:X:1630:A:C2	18:P:114:ALA:HB2	2.54	0.43
1:X:1745:C:OP1	15:M:101:ARG:NH2	2.51	0.43
1:X:1955:G:OP2	3:A:239:ARG:NH1	2.49	0.43
1:X:1973:C:H2'	1:X:1974:U:O4'	2.18	0.43
10:H:28:GLY:HA3	10:H:35:THR:H	1.84	0.43
11:I:81:GLN:HG2	11:I:114:ILE:HG22	1.99	0.43
13:K:11:ASN:OD1	13:K:17:ARG:CZ	2.66	0.43
20:R:56:LYS:HD3	20:R:69:GLN:HG2	1.99	0.43
1:X:534:U:H4'	1:X:564:U:H4'	2.00	0.43
1:X:1731:C:H2'	1:X:1732:U:H3'	1.98	0.43
5:C:58:MET:HB2	5:C:70:GLY:O	2.18	0.43
9:G:103:TYR:CB	9:G:107:GLN:HE21	2.32	0.43
11:I:38:LYS:HB3	11:I:39:SER:H	1.65	0.43
22:T:20:TYR:HB3	22:T:21:LEU:H	1.72	0.43
23:U:33:LYS:O	23:U:34:THR:HB	2.18	0.43
1:X:82:G:N2	1:X:100:G:O2'	2.51	0.43
1:X:224:G:H4'	1:X:399:G:C6	2.53	0.43
1:X:841:G:H2'	1:X:842:A:C8	2.53	0.43
1:X:1674:C:H2'	1:X:1675:C:C6	2.54	0.43
29:3:20:GLY:CA	29:3:21:LYS:CA	2.97	0.43
1:X:163:A:H2'	1:X:164:G:C8	2.54	0.43
1:X:2197:U:H2'	1:X:2198:U:C6	2.54	0.43
1:X:1223:G:H5'	1:X:1225:G:O4'	2.18	0.43
1:X:1314:A:H2	1:X:1642:G:N3	2.16	0.43
1:X:1674:C:H2'	1:X:1675:C:H6	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:2543:A:C2	1:X:2626:U:H4'	2.54	0.43
9:G:67:ARG:HB3	9:G:70:PHE:HA	2.01	0.43
20:R:37:LEU:HD21	20:R:49:GLU:HG3	2.01	0.43
1:X:476:G:H2'	1:X:477:A:C8	2.54	0.43
1:X:2189:A:H2	1:X:2190:A:C4	2.37	0.43
1:X:2356:A:H1'	14:L:89:PHE:CZ	2.54	0.43
3:A:250:TRP:HB3	3:A:251:GLY:H	1.61	0.43
13:K:8:ARG:O	13:K:9:LYS:HB3	2.19	0.43
14:L:89:PHE:CD1	14:L:89:PHE:N	2.71	0.43
28:2:22:MET:CA	28:2:23:LYS:CA	2.97	0.43
1:X:1966:C:H4'	1:X:2585:C:H4'	2.01	0.43
1:X:2633:A:N1	1:X:2644:A:H5''	2.34	0.43
2:Y:43:G:H5'	2:Y:44:C:H5'	2.00	0.43
7:E:17:VAL:HG22	7:E:26:VAL:HG22	2.00	0.43
9:G:70:PHE:CB	16:N:64:ARG:HG2	2.49	0.43
19:Q:58:VAL:HA	19:Q:59:PRO:HD3	1.81	0.43
19:Q:68:PHE:O	19:Q:70:GLY:N	2.51	0.43
1:X:77:C:H42	1:X:106:G:H1	1.66	0.42
1:X:807:A:H2'	1:X:808:C:C6	2.53	0.42
1:X:1699:A:H61	1:X:1723:U:H3	1.66	0.42
4:B:115:GLY:HA2	4:B:157:ALA:HB2	2.00	0.42
10:H:124:MET:O	10:H:127:VAL:HG12	2.18	0.42
1:X:331:U:O2'	5:C:162:ARG:NH1	2.52	0.42
1:X:830:C:O2'	1:X:852:U:H5''	2.19	0.42
1:X:1367:A:H2'	1:X:1368:G:O4'	2.19	0.42
1:X:1623:C:H4'	1:X:1624:A:O5'	2.19	0.42
4:B:125:GLY:H	4:B:135:HIS:HA	1.84	0.42
10:H:88:THR:HB	15:M:80:VAL:HB	2.00	0.42
17:O:36:LYS:HD3	17:O:39:PHE:HB3	2.01	0.42
1:X:540:G:C5	1:X:2005:U:H5''	2.54	0.42
1:X:2324:G:N3	1:X:2360:C:H2'	2.35	0.42
1:X:2394:G:H4'	11:I:65:PHE:HB3	2.02	0.42
1:X:2581:A:H2'	1:X:2582:G:H4'	2.00	0.42
2:Y:65:A:H2'	2:Y:66:G:H8	1.84	0.42
17:O:13:ARG:HB3	17:O:14:VAL:H	1.71	0.42
19:Q:84:GLU:HA	19:Q:86:GLN:HE21	1.84	0.42
1:X:1989:C:O2'	1:X:2798:A:N3	2.49	0.42
17:O:98:ILE:H	17:O:98:ILE:HG13	1.75	0.42
1:X:334:G:H8	5:C:164:VAL:HG13	1.83	0.42
1:X:1142:G:H5'	9:G:111:LYS:HB3	2.02	0.42
1:X:1658:A:H2'	1:X:1659:G:O4'	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Y:34:C:H2'	2:Y:35:C:C6	2.54	0.42
2:Y:94:G:H5'	21:S:74:ARG:HH12	1.84	0.42
4:B:16:LYS:HB2	4:B:21:ILE:CD1	2.46	0.42
10:H:27:SER:HB3	10:H:50:ILE:HG13	2.01	0.42
11:I:130:ILE:H	11:I:130:ILE:HG13	1.65	0.42
12:J:27:TYR:CE2	21:S:76:ARG:HB3	2.55	0.42
12:J:27:TYR:HE2	21:S:76:ARG:HD3	1.84	0.42
1:X:82:G:N1	1:X:100:G:H2'	2.34	0.42
1:X:636:G:H8	1:X:636:G:C5'	2.33	0.42
1:X:1030:U:C4	1:X:1031:C:H5	2.37	0.42
1:X:2191:A:OP1	1:X:2193:C:N4	2.52	0.42
3:A:186:HIS:HB2	3:A:188:GLU:HG3	2.01	0.42
3:A:247:VAL:HG23	3:A:248:THR:N	2.34	0.42
5:C:27:LEU:O	5:C:31:VAL:HG22	2.19	0.42
5:C:150:LEU:HB3	5:C:169:VAL:HG23	2.00	0.42
5:C:194:GLU:O	5:C:195:ILE:HG12	2.20	0.42
9:G:69:ASP:H	9:G:76:GLN:NE2	2.17	0.42
9:G:106:TYR:CE2	9:G:108:GLY:CA	3.03	0.42
10:H:26:ASN:HB3	10:H:38:GLY:H	1.84	0.42
11:I:43:ALA:C	11:I:45:LYS:H	2.23	0.42
13:K:28:LEU:HD13	13:K:115:LEU:HD23	2.00	0.42
19:Q:28:TRP:HZ3	19:Q:58:VAL:HG21	1.84	0.42
20:R:46:VAL:HG21	20:R:80:LYS:HE3	2.02	0.42
21:S:23:ALA:HA	21:S:83:PHE:O	2.18	0.42
1:X:110:U:H3'	1:X:111:G:H5''	2.01	0.42
1:X:451:A:H2'	1:X:452:G:C8	2.55	0.42
1:X:1909:U:P	1:X:1912:G:H1	2.42	0.42
1:X:2540:A:O2'	10:H:23:ARG:HG3	2.19	0.42
1:X:2663:U:C4	1:X:2664:G:N7	2.88	0.42
3:A:43:ARG:N	3:A:43:ARG:HH11	2.17	0.42
3:A:108:PRO:HB3	3:A:143:HIS:HE1	1.84	0.42
5:C:95:LEU:HD23	5:C:96:PRO:HD2	2.00	0.42
6:D:17:MET:HG2	6:D:22:TYR:HB2	2.01	0.42
7:E:19:ALA:HB1	7:E:24:PHE:HD2	1.84	0.42
9:G:65:LYS:HD2	9:G:66:HIS:CD2	2.54	0.42
14:L:63:ASN:HB3	14:L:66:ASP:HB2	2.01	0.42
15:M:39:VAL:HA	15:M:45:THR:HA	2.02	0.42
20:R:105:ARG:HH22	20:R:112:LYS:CA	2.33	0.42
26:Z:36:CYS:HB3	26:Z:49:CYS:HB3	1.91	0.42
29:3:10:ALA:CA	29:3:11:LYS:CA	2.98	0.42
1:X:503:G:H2'	1:X:504:G:O4'	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:648:A:H4'	1:X:649:G:O4'	2.20	0.42
1:X:1441:A:H4'	1:X:1442:C:O5'	2.19	0.42
1:X:1766:U:O4	1:X:1780:A:H2	2.02	0.42
1:X:2617:G:O2'	1:X:2755:A:N1	2.45	0.42
19:Q:59:PRO:HA	19:Q:74:ASP:OD1	2.19	0.42
1:X:2542:U:O2	1:X:2544:A:H8	2.03	0.42
3:A:231:HIS:HD2	3:A:233:HIS:N	2.04	0.42
19:Q:35:LYS:HA	19:Q:38:ILE:HG22	2.01	0.42
20:R:80:LYS:O	20:R:80:LYS:NZ	2.45	0.42
1:X:543:G:H5'	16:N:24:PHE:CE1	2.55	0.42
1:X:760:U:O2	1:X:1997:A:H1'	2.20	0.42
1:X:2674:C:H2'	1:X:2675:U:H6	1.85	0.42
9:G:43:VAL:HB	9:G:167:LYS:HG2	2.02	0.42
1:X:1151:U:H3'	9:G:91:THR:HG21	2.01	0.41
1:X:1378:A:H3'	1:X:1379:A:H8	1.85	0.41
1:X:1939:U:O2	1:X:1939:U:H2'	2.18	0.41
1:X:2315:A:H2	1:X:2364:C:O2	2.03	0.41
2:Y:8:C:H4'	2:Y:30:C:H5'	2.01	0.41
3:A:208:LYS:HE3	3:A:208:LYS:HA	2.01	0.41
4:B:183:LEU:HD11	15:M:16:ILE:CG2	2.50	0.41
12:J:36:ILE:HD12	12:J:133:VAL:HG21	2.01	0.41
15:M:34:ARG:HB2	15:M:91:VAL:HG23	2.02	0.41
21:S:168:VAL:HG12	21:S:169:VAL:HG23	2.01	0.41
1:X:171:G:H2'	1:X:172:A:O4'	2.20	0.41
1:X:1300:A:H5'	13:K:103:ARG:HD2	2.01	0.41
1:X:2065:A:H2'	1:X:2066:G:O4'	2.20	0.41
1:X:2326:C:H2'	1:X:2327:U:C6	2.55	0.41
1:X:2326:C:C2	1:X:2327:U:C5	3.08	0.41
15:M:66:PHE:CE2	15:M:81:PHE:HB2	2.55	0.41
1:X:577:U:C5'	1:X:956:A:N6	2.84	0.41
4:B:134:TRP:CD1	4:B:134:TRP:N	2.83	0.41
6:D:38:GLU:HG3	6:D:40:LEU:HD23	2.01	0.41
9:G:162:LYS:H	9:G:163:PRO:CD	2.33	0.41
22:T:46:LYS:HB2	22:T:78:PHE:CE2	2.56	0.41
23:U:49:LYS:HB2	23:U:61:TRP:CD2	2.56	0.41
1:X:1231:A:H2'	1:X:1232:U:C6	2.55	0.41
1:X:2820:C:H5''	15:M:60:SER:HB3	2.02	0.41
5:C:62:LYS:HE2	5:C:62:LYS:HB3	1.72	0.41
14:L:30:SER:O	14:L:40:ALA:HA	2.21	0.41
17:O:60:VAL:HA	17:O:93:ILE:HG22	2.03	0.41
4:B:5:LEU:HD22	4:B:49:ILE:HG22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:D:48:LYS:H	6:D:48:LYS:HG3	1.67	0.41
11:I:87:THR:HB	11:I:97:ARG:HD3	2.02	0.41
24:V:2:LYS:H	24:V:3:PRO:CD	2.34	0.41
1:X:558:G:H4'	1:X:559:C:C5'	2.48	0.41
1:X:636:G:C5'	1:X:636:G:C8	3.04	0.41
1:X:1442:C:O2'	1:X:1443:G:H5'	2.19	0.41
1:X:2406:C:C5'	1:X:2408:G:H5'	2.50	0.41
3:A:141:VAL:HG13	3:A:162:SER:HB3	2.02	0.41
5:C:12:GLY:O	5:C:14:THR:N	2.49	0.41
13:K:51:LEU:HD12	13:K:66:VAL:HG22	2.02	0.41
15:M:39:VAL:HG12	15:M:45:THR:HG23	2.01	0.41
1:X:346:C:O2	1:X:347:C:C5	2.73	0.41
1:X:564:U:H2'	1:X:565:A:H8	1.83	0.41
1:X:1283:C:H5''	1:X:1284:G:O5'	2.21	0.41
1:X:1813:A:O5'	1:X:1813:A:H8	2.04	0.41
4:B:149:ARG:CZ	9:G:106:TYR:CD1	3.03	0.41
6:D:104:ILE:HA	6:D:108:LEU:HD12	2.03	0.41
15:M:24:LEU:HD12	15:M:83:PHE:CG	2.56	0.41
15:M:58:ASN:HD22	15:M:58:ASN:H	1.67	0.41
16:N:93:LYS:HE2	17:O:5:ILE:HD13	2.03	0.41
18:P:43:ASP:OD2	18:P:46:ARG:NH2	2.54	0.41
21:S:47:SER:OG	21:S:48:THR:N	2.51	0.41
1:X:393:U:H2'	1:X:394:U:H6	1.85	0.41
1:X:1919:A:C2	1:X:1926:U:N3	2.67	0.41
4:B:183:LEU:HD21	15:M:16:ILE:HD13	2.02	0.41
5:C:173:ALA:HB1	5:C:193:LEU:HD13	2.02	0.41
23:U:10:LYS:HD2	23:U:60:VAL:HG11	2.02	0.41
1:X:162:C:H2'	1:X:163:A:H8	1.85	0.41
1:X:172:A:N6	1:X:175:C:H3'	2.34	0.41
1:X:405:C:H2'	1:X:406:G:C8	2.53	0.41
1:X:1032:A:H2'	1:X:1034:U:H5''	2.03	0.41
1:X:1179:A:H2'	1:X:1180:A:C8	2.56	0.41
1:X:1687:C:O5'	1:X:1687:C:H6	2.03	0.41
1:X:2206:C:H1'	3:A:262:LYS:HE2	2.03	0.41
5:C:117:LEU:HB3	5:C:187:VAL:HA	2.02	0.41
10:H:114:VAL:HG22	10:H:133:VAL:HG22	2.02	0.41
11:I:102:LYS:O	11:I:104:ARG:N	2.37	0.41
13:K:106:ASP:OD1	13:K:108:VAL:HB	2.21	0.41
17:O:40:VAL:HG13	17:O:45:THR:HG22	2.01	0.41
21:S:149:ALA:HB1	21:S:160:LEU:HD13	2.02	0.41
1:X:1467:U:C6	1:X:1467:U:C5'	3.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:1795:C:H5'	3:A:257:LEU:HD13	2.02	0.41
1:X:2189:A:C2	1:X:2190:A:C5	3.09	0.41
11:I:30:ALA:HB3	11:I:34:HIS:CE1	2.56	0.41
11:I:54:SER:O	11:I:56:LEU:N	2.50	0.41
13:K:3:HIS:CE1	13:K:5:LYS:HZ2	2.31	0.41
16:N:17:VAL:HG11	16:N:36:PHE:HB2	2.02	0.41
17:O:48:GLY:C	17:O:50:ASP:H	2.24	0.41
19:Q:72:ARG:H	19:Q:72:ARG:HG2	1.65	0.41
22:T:21:LEU:HD11	22:T:41:ARG:CZ	2.51	0.41
1:X:82:G:H1	1:X:100:G:H2'	1.86	0.40
1:X:573:C:H5''	17:O:74:TYR:OH	2.21	0.40
1:X:689:A:H2	1:X:815:A:H61	1.68	0.40
1:X:2262:C:C2	1:X:2368:G:C2	3.10	0.40
3:A:205:VAL:O	3:A:207:GLY:N	2.54	0.40
11:I:76:LYS:HG2	11:I:111:SER:HB2	2.03	0.40
12:J:35:LEU:HD23	12:J:105:PHE:HD2	1.86	0.40
12:J:133:VAL:HG12	21:S:76:ARG:CZ	2.51	0.40
13:K:49:GLU:OE1	13:K:95:THR:HB	2.21	0.40
16:N:93:LYS:H	16:N:93:LYS:HG3	1.69	0.40
20:R:93:ARG:HE	20:R:108:VAL:HG12	1.86	0.40
22:T:3:HIS:HD2	22:T:5:LYS:HB2	1.85	0.40
1:X:1235:C:H2'	1:X:1236:G:C8	2.57	0.40
1:X:1467:U:C2'	1:X:1468:A:OP1	2.68	0.40
1:X:2855:C:H1'	13:K:92:GLY:O	2.22	0.40
1:X:2864:C:H2'	1:X:2865:G:C8	2.57	0.40
2:Y:64:C:H2'	2:Y:65:A:H8	1.85	0.40
4:B:77:ILE:HD13	15:M:3:THR:HG22	2.02	0.40
9:G:119:LEU:HD23	9:G:122:HIS:HD2	1.85	0.40
12:J:26:ASP:HB3	12:J:27:TYR:H	1.60	0.40
16:N:86:ALA:HB2	16:N:116:ALA:HB2	2.03	0.40
1:X:424:G:H4'	1:X:425:A:O5'	2.21	0.40
1:X:636:G:H5''	1:X:636:G:C8	2.53	0.40
1:X:1142:G:OP2	1:X:2494:C:H5''	2.21	0.40
1:X:2551:A:N7	4:B:145:LYS:HB2	2.37	0.40
15:M:27:PHE:CE2	15:M:91:VAL:HG11	2.56	0.40
1:X:309:G:H5''	1:X:310:A:OP1	2.21	0.40
1:X:2556:A:H5''	1:X:2557:G:H5'	2.04	0.40
3:A:217:ARG:O	3:A:218:LYS:C	2.60	0.40
6:D:177:PHE:HB2	6:D:179:LYS:HE3	2.02	0.40
9:G:62:ILE:CG2	9:G:135:LEU:HD21	2.50	0.40
16:N:93:LYS:O	16:N:94:VAL:HB	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
17:O:10:LYS:HA	17:O:36:LYS:HA	2.03	0.40
1:X:89:A:C4'	1:X:90:G:H5''	2.52	0.40
1:X:393:U:H2'	1:X:394:U:C6	2.57	0.40
1:X:761:G:OP2	18:P:109:ARG:HG3	2.21	0.40
1:X:1151:U:H4'	1:X:1153:A:H5''	2.04	0.40
1:X:1287:A:H2'	1:X:1288:A:H5''	2.03	0.40
1:X:2482:A:H4'	1:X:2483:U:OP1	2.21	0.40
32:X:2929:1F2:H46	32:X:2929:1F2:O41	2.21	0.40
3:A:60:ARG:HD3	3:A:86:PRO:HB2	2.04	0.40
5:C:48:ARG:HB2	5:C:51:VAL:HG22	2.02	0.40
13:K:17:ARG:HD3	13:K:17:ARG:HA	2.00	0.40
19:Q:6:ILE:HG22	19:Q:7:LEU:HD23	2.01	0.40
20:R:48:VAL:HG12	20:R:50:GLY:H	1.86	0.40
25:W:40:VAL:HG22	25:W:43:MET:HE3	2.04	0.40
26:Z:45:ILE:HG21	26:Z:57:VAL:HG23	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	A	238/274 (87%)	178 (75%)	39 (16%)	21 (9%)	1	4
4	B	203/211 (96%)	171 (84%)	24 (12%)	8 (4%)	3	22
5	C	195/205 (95%)	129 (66%)	45 (23%)	21 (11%)	0	2
6	D	175/180 (97%)	141 (81%)	27 (15%)	7 (4%)	3	21
7	E	169/185 (91%)	139 (82%)	20 (12%)	10 (6%)	1	12
8	F	69/144 (48%)	57 (83%)	10 (14%)	2 (3%)	4	28
9	G	140/174 (80%)	105 (75%)	21 (15%)	14 (10%)	0	3
10	H	132/134 (98%)	115 (87%)	11 (8%)	6 (4%)	2	18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
11	I	139/156 (89%)	85 (61%)	28 (20%)	26 (19%)	0	0
12	J	134/141 (95%)	101 (75%)	19 (14%)	14 (10%)	0	3
13	K	111/116 (96%)	92 (83%)	11 (10%)	8 (7%)	1	7
14	L	102/114 (90%)	75 (74%)	19 (19%)	8 (8%)	1	6
15	M	106/166 (64%)	89 (84%)	13 (12%)	4 (4%)	3	22
16	N	115/118 (98%)	92 (80%)	17 (15%)	6 (5%)	2	15
17	O	92/100 (92%)	67 (73%)	13 (14%)	12 (13%)	0	1
18	P	125/134 (93%)	108 (86%)	12 (10%)	5 (4%)	3	21
19	Q	91/95 (96%)	64 (70%)	14 (15%)	13 (14%)	0	1
20	R	108/115 (94%)	66 (61%)	24 (22%)	18 (17%)	0	0
21	S	173/237 (73%)	135 (78%)	28 (16%)	10 (6%)	1	13
22	T	82/91 (90%)	64 (78%)	11 (13%)	7 (8%)	1	4
23	U	70/81 (86%)	43 (61%)	16 (23%)	11 (16%)	0	1
24	V	64/67 (96%)	58 (91%)	4 (6%)	2 (3%)	4	26
25	W	53/55 (96%)	50 (94%)	3 (6%)	0	100	100
26	Z	56/60 (93%)	48 (86%)	5 (9%)	3 (5%)	2	14
30	4	35/37 (95%)	31 (89%)	3 (9%)	1 (3%)	4	28
All	All	2977/3390 (88%)	2303 (77%)	437 (15%)	237 (8%)	1	6

All (237) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	56	GLY
3	A	89	SER
3	A	198	ASN
3	A	199	ALA
3	A	249	PRO
3	A	250	TRP
4	B	116	VAL
4	B	146	THR
5	C	13	ARG
5	C	20	PRO
5	C	60	GLY
5	C	67	ALA
5	C	121	ASP
5	C	129	LYS

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Mol	Chain	Res	Type
5	C	164	VAL
5	C	165	SER
5	C	172	VAL
5	C	195	ILE
6	D	81	GLN
7	E	55	PRO
7	E	126	PRO
7	E	165	VAL
9	G	34	PRO
9	G	67	ARG
9	G	91	THR
9	G	104	THR
9	G	107	GLN
9	G	158	HIS
10	H	27	SER
10	H	29	ILE
11	I	18	ARG
11	I	37	GLN
11	I	39	SER
11	I	55	ARG
11	I	62	LYS
11	I	98	LEU
11	I	99	VAL
11	I	103	ASN
12	J	13	GLN
12	J	21	ASP
12	J	26	ASP
13	K	6	ALA
13	K	9	LYS
14	L	21	THR
15	M	26	ASP
15	M	29	PRO
16	N	7	GLY
16	N	87	ASN
17	O	7	THR
17	O	9	GLY
17	O	10	LYS
17	O	14	VAL
17	O	31	ASP
17	O	48	GLY
19	Q	6	ILE
19	Q	12	ILE

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Mol	Chain	Res	Type
19	Q	67	ARG
19	Q	69	ILE
19	Q	90	ALA
20	R	5	SER
20	R	11	ASN
21	S	26	LYS
22	T	19	LYS
23	U	15	VAL
23	U	19	ILE
23	U	29	GLY
23	U	60	VAL
24	V	2	LYS
26	Z	4	HIS
26	Z	36	CYS
26	Z	53	ASP
3	A	87	ASN
3	A	206	LEU
3	A	248	THR
4	B	76	ARG
4	B	123	ALA
4	B	148	GLY
5	C	22	VAL
5	C	84	PHE
6	D	124	GLY
7	E	13	SER
9	G	64	GLY
9	G	97	ASP
9	G	170	PRO
11	I	47	ALA
11	I	49	PHE
11	I	54	SER
11	I	56	LEU
11	I	64	GLY
11	I	86	THR
12	J	11	ARG
12	J	80	ALA
12	J	83	ARG
13	K	4	GLY
13	K	92	GLY
14	L	33	ARG
14	L	40	ALA
14	L	46	SER

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Mol	Chain	Res	Type
17	O	8	GLY
17	O	60	VAL
18	P	9	ARG
19	Q	74	ASP
19	Q	84	GLU
19	Q	87	SER
19	Q	89	GLU
20	R	7	GLY
20	R	60	PRO
20	R	83	LEU
20	R	85	ASP
20	R	98	ILE
20	R	108	VAL
21	S	5	ALA
21	S	88	TYR
21	S	91	PRO
21	S	156	GLU
22	T	21	LEU
22	T	74	LYS
23	U	27	ASP
23	U	32	ARG
23	U	41	VAL
23	U	56	GLN
24	V	57	LYS
3	A	170	SER
3	A	220	HIS
3	A	244	ARG
4	B	90	SER
5	C	15	ILE
5	C	190	ALA
6	D	122	PHE
7	E	19	ALA
7	E	58	ALA
7	E	173	ALA
8	F	120	VAL
9	G	33	ILE
9	G	37	ASP
10	H	5	GLN
10	H	37	GLY
11	I	29	THR
11	I	65	PHE
11	I	82	ASP

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Mol	Chain	Res	Type
11	I	88	PHE
11	I	115	SER
12	J	60	ARG
12	J	81	GLU
13	K	11	ASN
14	L	20	THR
16	N	92	ARG
16	N	95	LEU
17	O	13	ARG
17	O	36	LYS
17	O	49	GLU
18	P	132	GLY
19	Q	5	ASP
19	Q	59	PRO
19	Q	63	LYS
20	R	6	ALA
20	R	18	LYS
20	R	26	SER
20	R	50	GLY
20	R	62	MET
20	R	63	THR
20	R	96	LYS
21	S	76	ARG
21	S	106	GLY
3	A	59	LYS
4	B	137	ARG
5	C	72	ARG
5	C	196	VAL
6	D	42	SER
7	E	7	GLN
9	G	163	PRO
11	I	91	ASP
13	K	8	ARG
13	K	14	SER
13	K	88	ALA
14	L	59	LEU
15	M	25	PRO
16	N	94	VAL
19	Q	13	SER
21	S	6	LYS
22	T	13	GLY
22	T	73	GLY

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Mol	Chain	Res	Type
23	U	34	THR
23	U	48	LYS
3	A	35	GLU
3	A	54	ILE
3	A	55	GLY
3	A	187	SER
3	A	219	PRO
5	C	4	ILE
5	C	11	GLY
5	C	66	ASN
5	C	126	ALA
6	D	10	ASP
6	D	40	LEU
7	E	80	SER
9	G	68	PRO
10	H	41	ASN
10	H	42	LYS
11	I	19	VAL
11	I	101	ARG
12	J	29	ALA
12	J	30	PHE
12	J	78	LYS
14	L	53	ALA
16	N	8	ILE
18	P	20	LEU
18	P	81	HIS
18	P	85	MET
21	S	56	VAL
22	T	11	LYS
22	T	27	GLY
30	4	13	ASN
3	A	252	LYS
3	A	255	LYS
5	C	171	PRO
6	D	121	ALA
9	G	105	GLY
11	I	9	THR
11	I	84	GLU
12	J	28	VAL
12	J	82	THR
20	R	75	ALA
21	S	110	GLY

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Mol	Chain	Res	Type
23	U	74	PRO
4	B	72	VAL
8	F	118	GLY
11	I	44	GLY
11	I	68	VAL
12	J	91	VAL
20	R	19	GLY
14	L	38	ILE
15	M	28	ARG
17	O	22	VAL
20	R	64	ASN
3	A	197	GLY
7	E	136	ILE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
3	A	185/215 (86%)	151 (82%)	34 (18%)	1 8
4	B	155/157 (99%)	128 (83%)	27 (17%)	2 10
5	C	157/163 (96%)	120 (76%)	37 (24%)	1 3
6	D	153/156 (98%)	133 (87%)	20 (13%)	4 19
7	E	136/144 (94%)	125 (92%)	11 (8%)	11 42
8	F	51/107 (48%)	49 (96%)	2 (4%)	32 67
9	G	118/146 (81%)	93 (79%)	25 (21%)	1 5
10	H	103/103 (100%)	87 (84%)	16 (16%)	2 12
11	I	108/121 (89%)	76 (70%)	32 (30%)	0 1
12	J	110/115 (96%)	88 (80%)	22 (20%)	1 6
13	K	90/93 (97%)	68 (76%)	22 (24%)	0 2
14	L	74/82 (90%)	56 (76%)	18 (24%)	0 2
15	M	94/134 (70%)	66 (70%)	28 (30%)	0 1
16	N	96/97 (99%)	81 (84%)	15 (16%)	2 12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	O	75/79 (95%)	57 (76%)	18 (24%)	0	3
18	P	109/115 (95%)	89 (82%)	20 (18%)	1	8
19	Q	75/76 (99%)	62 (83%)	13 (17%)	2	10
20	R	91/96 (95%)	70 (77%)	21 (23%)	1	3
21	S	149/192 (78%)	130 (87%)	19 (13%)	4	20
22	T	62/67 (92%)	53 (86%)	9 (14%)	3	15
23	U	57/66 (86%)	37 (65%)	20 (35%)	0	0
24	V	54/55 (98%)	46 (85%)	8 (15%)	3	14
25	W	48/48 (100%)	35 (73%)	13 (27%)	0	1
26	Z	51/53 (96%)	39 (76%)	12 (24%)	1	3
30	4	35/35 (100%)	31 (89%)	4 (11%)	5	24
All	All	2436/2715 (90%)	1970 (81%)	466 (19%)	1	8

All (466) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	39	LYS
3	A	40	THR
3	A	43	ARG
3	A	44	ASN
3	A	46	ARG
3	A	48	ARG
3	A	49	ILE
3	A	50	THR
3	A	52	ARG
3	A	63	ARG
3	A	68	LYS
3	A	69	ARG
3	A	76	ASN
3	A	88	ARG
3	A	96	HIS
3	A	105	ILE
3	A	111	LEU
3	A	145	LEU
3	A	157	ARG
3	A	164	GLN
3	A	183	ARG
3	A	196	VAL

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Mol	Chain	Res	Type
3	A	198	ASN
3	A	203	ASN
3	A	208	LYS
3	A	215	LEU
3	A	218	LYS
3	A	226	MET
3	A	229	VAL
3	A	240	THR
3	A	244	ARG
3	A	252	LYS
3	A	259	THR
3	A	260	ARG
4	B	5	LEU
4	B	14	ILE
4	B	19	ARG
4	B	37	LYS
4	B	41	THR
4	B	49	ILE
4	B	69	LYS
4	B	77	ILE
4	B	82	ARG
4	B	87	ASP
4	B	105	THR
4	B	111	LYS
4	B	119	ARG
4	B	131	SER
4	B	133	LYS
4	B	134	TRP
4	B	136	ARG
4	B	137	ARG
4	B	141	ILE
4	B	145	LYS
4	B	149	ARG
4	B	150	VAL
4	B	152	LYS
4	B	154	LYS
4	B	162	MET
4	B	163	GLU
4	B	203	LYS
5	C	4	ILE
5	C	5	ASN
5	C	10	ASN

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Mol	Chain	Res	Type
5	C	13	ARG
5	C	14	THR
5	C	15	ILE
5	C	31	VAL
5	C	45	THR
5	C	48	ARG
5	C	51	VAL
5	C	62	LYS
5	C	72	ARG
5	C	90	SER
5	C	91	TYR
5	C	95	LEU
5	C	97	ARG
5	C	102	LEU
5	C	104	LEU
5	C	117	LEU
5	C	121	ASP
5	C	124	ASP
5	C	127	ASP
5	C	130	THR
5	C	143	ASP
5	C	148	VAL
5	C	150	LEU
5	C	153	ASP
5	C	154	ASP
5	C	162	ARG
5	C	164	VAL
5	C	168	SER
5	C	175	VAL
5	C	176	ASN
5	C	180	ILE
5	C	188	ILE
5	C	193	LEU
5	C	194	GLU
6	D	10	ASP
6	D	40	LEU
6	D	48	LYS
6	D	51	ASP
6	D	74	ILE
6	D	80	ARG
6	D	81	GLN
6	D	83	MET

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Mol	Chain	Res	Type
6	D	89	VAL
6	D	100	LEU
6	D	112	ARG
6	D	115	ARG
6	D	117	ILE
6	D	125	ARG
6	D	130	LEU
6	D	148	LYS
6	D	150	ARG
6	D	153	ASP
6	D	171	GLN
6	D	175	LEU
7	E	7	GLN
7	E	35	VAL
7	E	44	ARG
7	E	48	ASP
7	E	50	LEU
7	E	59	GLN
7	E	69	ARG
7	E	98	LEU
7	E	106	ASN
7	E	143	GLN
7	E	155	ASP
8	F	84	ILE
8	F	111	LYS
9	G	31	THR
9	G	37	ASP
9	G	38	GLU
9	G	41	TRP
9	G	53	ARG
9	G	56	THR
9	G	61	ARG
9	G	62	ILE
9	G	70	PHE
9	G	71	THR
9	G	75	ILE
9	G	90	LEU
9	G	101	THR
9	G	102	ARG
9	G	104	THR
9	G	112	THR
9	G	113	GLU

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Mol	Chain	Res	Type
9	G	127	ILE
9	G	132	PHE
9	G	145	HIS
9	G	154	GLU
9	G	165	VAL
9	G	166	LEU
9	G	168	THR
9	G	169	GLN
10	H	7	ARG
10	H	8	LEU
10	H	23	ARG
10	H	27	SER
10	H	36	THR
10	H	41	ASN
10	H	47	VAL
10	H	81	ILE
10	H	89	ILE
10	H	94	ASN
10	H	102	GLN
10	H	106	ARG
10	H	116	ARG
10	H	120	ASP
10	H	126	ILE
10	H	127	VAL
11	I	5	ASP
11	I	6	LEU
11	I	7	LYS
11	I	12	SER
11	I	13	ARG
11	I	18	ARG
11	I	21	ARG
11	I	26	THR
11	I	27	ASP
11	I	29	THR
11	I	38	LYS
11	I	39	SER
11	I	45	LYS
11	I	48	PHE
11	I	53	ARG
11	I	56	LEU
11	I	57	ILE
11	I	65	PHE

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Mol	Chain	Res	Type
11	I	73	GLU
11	I	83	LEU
11	I	85	ASP
11	I	87	THR
11	I	89	ASP
11	I	98	LEU
11	I	99	VAL
11	I	101	ARG
11	I	106	VAL
11	I	114	ILE
11	I	118	VAL
11	I	120	VAL
11	I	123	ASP
11	I	130	ILE
12	J	7	ARG
12	J	8	THR
12	J	17	ARG
12	J	21	ASP
12	J	26	ASP
12	J	44	LYS
12	J	45	SER
12	J	47	GLN
12	J	64	LYS
12	J	65	ILE
12	J	72	ASP
12	J	81	GLU
12	J	82	THR
12	J	86	LYS
12	J	88	LYS
12	J	91	VAL
12	J	94	TRP
12	J	97	VAL
12	J	114	GLN
12	J	129	GLN
12	J	134	LYS
12	J	139	ASP
13	K	5	LYS
13	K	8	ARG
13	K	10	LEU
13	K	11	ASN
13	K	12	ARG
13	K	17	ARG

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Mol	Chain	Res	Type
13	K	28	LEU
13	K	35	GLN
13	K	37	THR
13	K	45	ARG
13	K	51	LEU
13	K	62	SER
13	K	83	VAL
13	K	89	GLU
13	K	94	TYR
13	K	95	THR
13	K	99	ARG
13	K	102	THR
13	K	106	ASP
13	K	109	THR
13	K	114	GLU
13	K	115	LEU
14	L	8	ARG
14	L	9	ARG
14	L	13	THR
14	L	17	VAL
14	L	31	VAL
14	L	33	ARG
14	L	36	LYS
14	L	43	ILE
14	L	59	LEU
14	L	64	LYS
14	L	66	ASP
14	L	67	THR
14	L	89	PHE
14	L	90	ASP
14	L	91	ARG
14	L	95	LYS
14	L	99	ARG
14	L	108	ARG
15	M	2	GLN
15	M	5	ILE
15	M	6	LYS
15	M	7	ILE
15	M	12	LEU
15	M	13	LEU
15	M	16	ILE
15	M	22	ARG

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Mol	Chain	Res	Type
15	M	28	ARG
15	M	31	ASP
15	M	34	ARG
15	M	35	VAL
15	M	36	ASP
15	M	37	THR
15	M	38	LYS
15	M	40	ARG
15	M	57	ILE
15	M	62	SER
15	M	63	ARG
15	M	64	LYS
15	M	88	VAL
15	M	89	ASN
15	M	92	THR
15	M	93	ILE
15	M	95	GLU
15	M	96	ARG
15	M	98	LYS
15	M	104	LEU
16	N	8	ILE
16	N	13	ARG
16	N	22	LYS
16	N	30	LYS
16	N	51	ARG
16	N	58	ARG
16	N	60	LEU
16	N	71	LEU
16	N	78	THR
16	N	85	ARG
16	N	88	ILE
16	N	90	LEU
16	N	92	ARG
16	N	93	LYS
16	N	102	GLU
17	O	13	ARG
17	O	14	VAL
17	O	18	ASP
17	O	21	ARG
17	O	24	SER
17	O	25	LEU
17	O	26	GLN

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Mol	Chain	Res	Type
17	O	28	GLU
17	O	35	LEU
17	O	40	VAL
17	O	69	ILE
17	O	71	ILE
17	O	78	VAL
17	O	83	ARG
17	O	84	THR
17	O	87	ARG
17	O	88	GLN
17	O	98	ILE
18	P	8	PHE
18	P	9	ARG
18	P	11	LYS
18	P	17	GLN
18	P	21	ARG
18	P	32	ARG
18	P	36	ARG
18	P	37	LYS
18	P	48	LYS
18	P	49	SER
18	P	62	ARG
18	P	87	GLU
18	P	89	ARG
18	P	106	LEU
18	P	109	ARG
18	P	118	LYS
18	P	120	ARG
18	P	125	THR
18	P	126	ILE
18	P	133	ASN
19	Q	7	LEU
19	Q	12	ILE
19	Q	14	GLU
19	Q	26	SER
19	Q	38	ILE
19	Q	40	ASP
19	Q	42	ILE
19	Q	58	VAL
19	Q	63	LYS
19	Q	65	VAL
19	Q	79	ILE

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Mol	Chain	Res	Type
19	Q	80	VAL
19	Q	84	GLU
20	R	11	ASN
20	R	15	HIS
20	R	18	LYS
20	R	23	ILE
20	R	25	LEU
20	R	37	LEU
20	R	38	LEU
20	R	42	ARG
20	R	46	VAL
20	R	53	VAL
20	R	73	GLU
20	R	80	LYS
20	R	83	LEU
20	R	85	ASP
20	R	95	ARG
20	R	96	LYS
20	R	98	ILE
20	R	104	VAL
20	R	105	ARG
20	R	106	VAL
20	R	108	VAL
21	S	13	LYS
21	S	14	LEU
21	S	15	ASP
21	S	22	VAL
21	S	26	LYS
21	S	40	ASP
21	S	51	LEU
21	S	56	VAL
21	S	67	LYS
21	S	71	MET
21	S	79	ILE
21	S	100	THR
21	S	120	LEU
21	S	128	ARG
21	S	130	ILE
21	S	132	GLN
21	S	134	LEU
21	S	145	ASP
21	S	175	ARG

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Mol	Chain	Res	Type
22	T	16	SER
22	T	17	ASN
22	T	37	LEU
22	T	38	VAL
22	T	46	LYS
22	T	49	GLN
22	T	62	LEU
22	T	64	ASP
22	T	85	GLN
23	U	8	THR
23	U	10	LYS
23	U	11	LYS
23	U	17	SER
23	U	19	ILE
23	U	23	LYS
23	U	32	ARG
23	U	35	THR
23	U	37	ILE
23	U	40	ARG
23	U	42	GLN
23	U	43	ARG
23	U	45	ASN
23	U	47	HIS
23	U	52	ARG
23	U	63	SER
23	U	65	ASN
23	U	75	TYR
23	U	78	ILE
23	U	79	GLU
24	V	1	MET
24	V	7	ARG
24	V	14	PHE
24	V	19	ASP
24	V	28	LEU
24	V	37	LEU
24	V	44	ARG
24	V	65	GLU
25	W	2	LYS
25	W	3	ILE
25	W	4	LYS
25	W	9	VAL
25	W	10	ILE

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Mol	Chain	Res	Type
25	W	12	ARG
25	W	26	ARG
25	W	30	ASP
25	W	32	ARG
25	W	34	VAL
25	W	37	THR
25	W	45	LYS
25	W	46	THR
26	Z	3	LYS
26	Z	4	HIS
26	Z	6	VAL
26	Z	8	LYS
26	Z	11	THR
26	Z	18	MET
26	Z	35	GLN
26	Z	40	LYS
26	Z	41	LEU
26	Z	53	ASP
26	Z	55	ARG
26	Z	58	LEU
30	4	2	LYS
30	4	9	LYS
30	4	13	ASN
30	4	30	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
3	A	231	HIS
4	B	129	HIS
4	B	180	ASN
5	C	61	GLN
6	D	9	ASN
6	D	81	GLN
6	D	129	ASN
7	E	20	GLN
7	E	106	ASN
7	E	143	GLN
9	G	76	GLN
9	G	122	HIS
10	H	41	ASN
11	I	66	ASN

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Mol	Chain	Res	Type
12	J	47	GLN
12	J	58	HIS
13	K	13	ASN
14	L	41	GLN
14	L	49	GLN
15	M	58	ASN
16	N	31	GLN
16	N	72	HIS
16	N	91	ASN
17	O	11	GLN
17	O	79	GLN
17	O	88	GLN
18	P	78	ASN
18	P	81	HIS
18	P	115	ASN
19	Q	8	GLN
19	Q	73	ASN
19	Q	86	GLN
20	R	10	HIS
20	R	29	HIS
20	R	44	GLN
22	T	3	HIS
22	T	17	ASN
22	T	49	GLN
22	T	57	HIS
25	W	49	HIS
25	W	54	GLN
26	Z	29	ASN
26	Z	43	HIS
26	Z	44	HIS
30	4	36	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	X	2683/2880 (93%)	672 (25%)	252 (9%)
2	Y	121/123 (98%)	26 (21%)	7 (5%)
All	All	2804/3003 (93%)	698 (24%)	259 (9%)

All (698) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	X	2	G
1	X	7	G
1	X	34	U
1	X	45	C
1	X	48	A
1	X	49	U
1	X	50	G
1	X	51	A
1	X	54	G
1	X	55	A
1	X	63	A
1	X	69	G
1	X	70	A
1	X	71	A
1	X	72	A
1	X	74	G
1	X	82	G
1	X	83	A
1	X	84	G
1	X	89	A
1	X	90	G
1	X	91	A
1	X	92	U
1	X	97	U
1	X	98	U
1	X	99	U
1	X	100	G
1	X	101	A
1	X	107	G
1	X	108	G
1	X	111	G
1	X	116	A
1	X	117	A
1	X	118	U
1	X	123	A
1	X	127	C
1	X	129	A
1	X	138	G
1	X	143	A
1	X	147	G
1	X	154	U
1	X	155	G
1	X	158	A

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Mol	Chain	Res	Type
1	X	173	A
1	X	176	A
1	X	177	U
1	X	181	A
1	X	182	G
1	X	192	G
1	X	193	A
1	X	198	A
1	X	199	A
1	X	204	A
1	X	205	A
1	X	206	U
1	X	207	U
1	X	210	A
1	X	219	G
1	X	225	G
1	X	228	A
1	X	229	G
1	X	238	G
1	X	242	A
1	X	243	G
1	X	245	C
1	X	246	C
1	X	304	A
1	X	305	A
1	X	306	G
1	X	310	A
1	X	312	G
1	X	313	U
1	X	319	G
1	X	321	A
1	X	322	A
1	X	323	G
1	X	329	C
1	X	332	C
1	X	333	A
1	X	334	G
1	X	335	A
1	X	340	G
1	X	341	A
1	X	342	G
1	X	343	A

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Mol	Chain	Res	Type
1	X	358	C
1	X	360	A
1	X	388	G
1	X	393	U
1	X	396	U
1	X	397	U
1	X	399	G
1	X	400	U
1	X	401	G
1	X	409	G
1	X	412	U
1	X	414	A
1	X	416	U
1	X	418	C
1	X	419	G
1	X	424	G
1	X	425	A
1	X	431	G
1	X	433	G
1	X	435	A
1	X	441	A
1	X	454	G
1	X	455	A
1	X	456	C
1	X	458	G
1	X	459	A
1	X	461	A
1	X	463	C
1	X	467	U
1	X	468	A
1	X	469	G
1	X	470	U
1	X	482	A
1	X	486	U
1	X	491	A
1	X	492	G
1	X	504	G
1	X	514	G
1	X	515	A
1	X	518	A
1	X	519	C
1	X	520	C

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Mol	Chain	Res	Type
1	X	539	A
1	X	540	G
1	X	541	C
1	X	542	A
1	X	543	G
1	X	554	U
1	X	555	U
1	X	556	A
1	X	557	U
1	X	558	G
1	X	559	C
1	X	560	G
1	X	572	G
1	X	580	A
1	X	582	G
1	X	583	C
1	X	584	A
1	X	595	A
1	X	602	C
1	X	613	A
1	X	614	G
1	X	620	G
1	X	624	A
1	X	625	A
1	X	627	A
1	X	628	A
1	X	631	G
1	X	632	A
1	X	633	G
1	X	636	G
1	X	648	A
1	X	649	G
1	X	652	C
1	X	654	A
1	X	655	A
1	X	657	A
1	X	664	C
1	X	665	A
1	X	666	U
1	X	667	U
1	X	681	A
1	X	682	G

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Mol	Chain	Res	Type
1	X	683	A
1	X	684	C
1	X	690	A
1	X	695	G
1	X	698	A
1	X	699	G
1	X	700	C
1	X	723	C
1	X	728	G
1	X	729	A
1	X	730	C
1	X	732	G
1	X	742	G
1	X	743	A
1	X	751	G
1	X	752	G
1	X	760	U
1	X	766	A
1	X	778	G
1	X	782	U
1	X	783	G
1	X	788	G
1	X	789	G
1	X	790	A
1	X	795	A
1	X	797	A
1	X	798	G
1	X	801	A
1	X	802	A
1	X	803	C
1	X	804	C
1	X	805	G
1	X	806	A
1	X	807	A
1	X	814	G
1	X	815	A
1	X	816	U
1	X	818	G
1	X	824	U
1	X	825	C
1	X	832	A
1	X	840	U

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Mol	Chain	Res	Type
1	X	841	G
1	X	859	U
1	X	860	U
1	X	871	U
1	X	872	G
1	X	878	C
1	X	879	A
1	X	880	C
1	X	919	U
1	X	922	A
1	X	926	C
1	X	931	G
1	X	938	G
1	X	939	C
1	X	940	G
1	X	941	U
1	X	943	U
1	X	944	A
1	X	952	A
1	X	955	G
1	X	956	A
1	X	957	G
1	X	964	A
1	X	968	C
1	X	969	U
1	X	970	A
1	X	972	C
1	X	973	U
1	X	985	G
1	X	994	A
1	X	995	A
1	X	996	C
1	X	1000	G
1	X	1001	A
1	X	1006	C
1	X	1007	A
1	X	1010	U
1	X	1014	G
1	X	1016	C
1	X	1019	U
1	X	1020	A
1	X	1022	A

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Mol	Chain	Res	Type
1	X	1023	U
1	X	1024	G
1	X	1033	G
1	X	1034	U
1	X	1035	G
1	X	1036	G
1	X	1037	U
1	X	1038	U
1	X	1044	U
1	X	1052	C
1	X	1053	G
1	X	1054	C
1	X	1055	A
1	X	1056	U
1	X	1057	A
1	X	1058	G
1	X	1059	A
1	X	1068	A
1	X	1069	G
1	X	1070	G
1	X	1072	U
1	X	1073	G
1	X	1077	U
1	X	1078	A
1	X	1079	G
1	X	1081	A
1	X	1082	G
1	X	1086	C
1	X	1087	C
1	X	1090	C
1	X	1094	C
1	X	1097	A
1	X	1098	G
1	X	1099	A
1	X	1100	G
1	X	1107	A
1	X	1108	U
1	X	1121	G
1	X	1122	A
1	X	1123	G
1	X	1124	U
1	X	1128	G

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Mol	Chain	Res	Type
1	X	1129	A
1	X	1139	A
1	X	1140	A
1	X	1141	U
1	X	1142	G
1	X	1143	A
1	X	1145	C
1	X	1146	G
1	X	1152	C
1	X	1153	A
1	X	1161	U
1	X	1162	A
1	X	1182	U
1	X	1185	C
1	X	1187	A
1	X	1189	G
1	X	1192	A
1	X	1194	U
1	X	1224	A
1	X	1233	A
1	X	1234	C
1	X	1250	A
1	X	1251	G
1	X	1262	U
1	X	1264	C
1	X	1266	G
1	X	1269	G
1	X	1275	A
1	X	1281	A
1	X	1282	A
1	X	1284	G
1	X	1285	A
1	X	1286	U
1	X	1288	A
1	X	1289	A
1	X	1302	C
1	X	1313	U
1	X	1314	A
1	X	1315	A
1	X	1316	G
1	X	1319	C
1	X	1334	A

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Mol	Chain	Res	Type
1	X	1342	U
1	X	1345	G
1	X	1356	G
1	X	1357	U
1	X	1358	C
1	X	1359	G
1	X	1365	U
1	X	1370	U
1	X	1378	A
1	X	1379	A
1	X	1381	G
1	X	1392	U
1	X	1404	C
1	X	1409	U
1	X	1410	U
1	X	1412	C
1	X	1413	U
1	X	1425	G
1	X	1428	G
1	X	1429	A
1	X	1432	G
1	X	1433	A
1	X	1434	U
1	X	1435	G
1	X	1440	G
1	X	1441	A
1	X	1442	C
1	X	1443	G
1	X	1451	C
1	X	1459	U
1	X	1460	G
1	X	1465	G
1	X	1467	U
1	X	1468	A
1	X	1469	U
1	X	1470	G
1	X	1474	A
1	X	1475	U
1	X	1476	G
1	X	1483	G
1	X	1489	C
1	X	1490	U

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Mol	Chain	Res	Type
1	X	1497	C
1	X	1498	G
1	X	1505	U
1	X	1506	C
1	X	1508	G
1	X	1509	A
1	X	1513	U
1	X	1514	C
1	X	1523	A
1	X	1524	C
1	X	1525	A
1	X	1528	C
1	X	1531	C
1	X	1545	G
1	X	1551	U
1	X	1552	C
1	X	1553	G
1	X	1554	G
1	X	1562	G
1	X	1563	U
1	X	1570	C
1	X	1571	G
1	X	1574	A
1	X	1575	C
1	X	1576	G
1	X	1582	A
1	X	1585	A
1	X	1594	U
1	X	1600	U
1	X	1601	U
1	X	1602	G
1	X	1603	A
1	X	1607	A
1	X	1608	U
1	X	1613	G
1	X	1624	A
1	X	1625	A
1	X	1626	A
1	X	1631	C
1	X	1632	A
1	X	1634	A
1	X	1635	G

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Mol	Chain	Res	Type
1	X	1648	C
1	X	1651	U
1	X	1657	A
1	X	1661	C
1	X	1665	C
1	X	1681	A
1	X	1685	A
1	X	1686	A
1	X	1691	G
1	X	1710	U
1	X	1711	C
1	X	1712	G
1	X	1713	G
1	X	1716	G
1	X	1717	A
1	X	1724	C
1	X	1732	U
1	X	1733	U
1	X	1734	C
1	X	1747	G
1	X	1749	G
1	X	1754	G
1	X	1755	G
1	X	1760	G
1	X	1764	A
1	X	1775	A
1	X	1776	A
1	X	1777	A
1	X	1782	A
1	X	1790	G
1	X	1791	C
1	X	1792	C
1	X	1793	A
1	X	1799	A
1	X	1800	A
1	X	1801	C
1	X	1808	C
1	X	1810	U
1	X	1811	A
1	X	1812	U
1	X	1819	U
1	X	1821	A

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Mol	Chain	Res	Type
1	X	1825	C
1	X	1840	A
1	X	1850	G
1	X	1861	G
1	X	1867	A
1	X	1883	A
1	X	1886	G
1	X	1910	A
1	X	1912	G
1	X	1913	G
1	X	1914	U
1	X	1919	A
1	X	1920	A
1	X	1921	A
1	X	1923	U
1	X	1924	C
1	X	1927	U
1	X	1938	U
1	X	1939	U
1	X	1943	A
1	X	1946	U
1	X	1947	G
1	X	1950	C
1	X	1953	A
1	X	1954	A
1	X	1955	G
1	X	1964	A
1	X	1965	U
1	X	1972	G
1	X	1976	U
1	X	1979	C
1	X	1980	A
1	X	2003	A
1	X	2005	U
1	X	2006	G
1	X	2010	G
1	X	2014	A
1	X	2015	G
1	X	2018	G
1	X	2019	C
1	X	2026	C
1	X	2035	G

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Mol	Chain	Res	Type
1	X	2038	C
1	X	2039	G
1	X	2044	G
1	X	2045	A
1	X	2046	C
1	X	2052	G
1	X	2063	A
1	X	2076	G
1	X	2084	G
1	X	2089	C
1	X	2166	G
1	X	2171	U
1	X	2181	A
1	X	2189	A
1	X	2190	A
1	X	2191	A
1	X	2195	C
1	X	2196	U
1	X	2197	U
1	X	2199	C
1	X	2204	A
1	X	2205	C
1	X	2217	G
1	X	2218	G
1	X	2222	U
1	X	2230	G
1	X	2259	G
1	X	2262	C
1	X	2265	A
1	X	2266	A
1	X	2267	A
1	X	2268	G
1	X	2283	G
1	X	2284	U
1	X	2285	U
1	X	2286	G
1	X	2287	G
1	X	2288	A
1	X	2290	A
1	X	2291	U
1	X	2298	U
1	X	2299	A

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Mol	Chain	Res	Type
1	X	2300	G
1	X	2301	A
1	X	2304	G
1	X	2305	C
1	X	2306	A
1	X	2313	G
1	X	2314	A
1	X	2315	A
1	X	2318	U
1	X	2323	U
1	X	2324	G
1	X	2326	C
1	X	2329	C
1	X	2330	G
1	X	2351	G
1	X	2358	C
1	X	2362	G
1	X	2363	G
1	X	2364	C
1	X	2371	A
1	X	2379	G
1	X	2381	A
1	X	2385	U
1	X	2389	G
1	X	2396	C
1	X	2401	A
1	X	2402	U
1	X	2404	A
1	X	2405	A
1	X	2406	C
1	X	2408	G
1	X	2409	A
1	X	2410	U
1	X	2414	A
1	X	2415	G
1	X	2420	C
1	X	2426	G
1	X	2427	A
1	X	2438	A
1	X	2447	G
1	X	2448	A
1	X	2452	U

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Mol	Chain	Res	Type
1	X	2453	C
1	X	2455	A
1	X	2473	G
1	X	2475	C
1	X	2477	C
1	X	2481	G
1	X	2482	A
1	X	2483	U
1	X	2484	G
1	X	2494	C
1	X	2497	A
1	X	2498	U
1	X	2499	C
1	X	2504	G
1	X	2508	G
1	X	2509	A
1	X	2514	G
1	X	2545	A
1	X	2546	G
1	X	2552	C
1	X	2553	G
1	X	2557	G
1	X	2561	G
1	X	2562	G
1	X	2565	C
1	X	2581	A
1	X	2582	G
1	X	2588	U
1	X	2589	C
1	X	2591	C
1	X	2594	U
1	X	2608	A
1	X	2611	A
1	X	2613	A
1	X	2633	A
1	X	2635	U
1	X	2640	G
1	X	2643	G
1	X	2650	G
1	X	2661	G
1	X	2664	G
1	X	2668	U

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Mol	Chain	Res	Type
1	X	2691	C
1	X	2692	A
1	X	2693	U
1	X	2694	G
1	X	2705	A
1	X	2706	U
1	X	2713	A
1	X	2718	A
1	X	2728	A
1	X	2731	G
1	X	2732	C
1	X	2737	A
1	X	2745	A
1	X	2758	A
1	X	2759	U
1	X	2760	G
1	X	2770	A
1	X	2771	C
1	X	2774	U
1	X	2775	U
1	X	2776	U
1	X	2777	A
1	X	2778	U
1	X	2779	C
1	X	2780	A
1	X	2782	G
1	X	2795	A
1	X	2796	A
1	X	2798	A
1	X	2808	U
1	X	2809	A
1	X	2810	A
1	X	2811	G
1	X	2825	A
1	X	2843	A
1	X	2847	G
1	X	2854	G
1	X	2859	U
1	X	2864	C
1	X	2866	A
1	X	2868	G
2	Y	4	C

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Mol	Chain	Res	Type
2	Y	11	G
2	Y	14	C
2	Y	15	A
2	Y	17	A
2	Y	18	G
2	Y	27	A
2	Y	28	A
2	Y	37	C
2	Y	39	C
2	Y	40	C
2	Y	42	U
2	Y	43	G
2	Y	44	C
2	Y	46	G
2	Y	47	A
2	Y	59	A
2	Y	69	G
2	Y	75	A
2	Y	76	U
2	Y	102	A
2	Y	110	U
2	Y	112	A
2	Y	115	G
2	Y	121	G
2	Y	123	U

All (259) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	X	13	A
1	X	33	C
1	X	34	U
1	X	48	A
1	X	49	U
1	X	50	G
1	X	62	U
1	X	70	A
1	X	71	A
1	X	73	A
1	X	74	G
1	X	82	G
1	X	83	A

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Mol	Chain	Res	Type
1	X	89	A
1	X	98	U
1	X	99	U
1	X	100	G
1	X	117	A
1	X	154	U
1	X	173	A
1	X	176	A
1	X	181	A
1	X	182	G
1	X	198	A
1	X	199	A
1	X	204	A
1	X	242	A
1	X	312	G
1	X	321	A
1	X	322	A
1	X	328	A
1	X	332	C
1	X	334	G
1	X	341	A
1	X	342	G
1	X	343	A
1	X	387	A
1	X	396	U
1	X	399	G
1	X	416	U
1	X	417	C
1	X	418	C
1	X	419	G
1	X	434	C
1	X	454	G
1	X	458	G
1	X	466	A
1	X	467	U
1	X	469	G
1	X	504	G
1	X	513	A
1	X	522	G
1	X	539	A
1	X	540	G
1	X	542	A

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Mol	Chain	Res	Type
1	X	557	U
1	X	558	G
1	X	559	C
1	X	580	A
1	X	631	G
1	X	648	A
1	X	652	C
1	X	664	C
1	X	672	C
1	X	677	G
1	X	681	A
1	X	682	G
1	X	683	A
1	X	698	A
1	X	714	G
1	X	751	G
1	X	765	C
1	X	775	U
1	X	777	A
1	X	788	G
1	X	803	C
1	X	806	A
1	X	813	A
1	X	814	G
1	X	818	G
1	X	824	U
1	X	840	U
1	X	842	A
1	X	858	G
1	X	859	U
1	X	860	U
1	X	871	U
1	X	872	G
1	X	878	C
1	X	938	G
1	X	939	C
1	X	955	G
1	X	969	U
1	X	972	C
1	X	990	A
1	X	994	A
1	X	995	A

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Mol	Chain	Res	Type
1	X	1000	G
1	X	1006	C
1	X	1007	A
1	X	1019	U
1	X	1032	A
1	X	1036	G
1	X	1038	U
1	X	1053	G
1	X	1055	A
1	X	1056	U
1	X	1057	A
1	X	1069	G
1	X	1072	U
1	X	1080	A
1	X	1081	A
1	X	1086	C
1	X	1096	A
1	X	1099	A
1	X	1122	A
1	X	1141	U
1	X	1142	G
1	X	1152	C
1	X	1185	C
1	X	1186	G
1	X	1223	G
1	X	1233	A
1	X	1249	G
1	X	1266	G
1	X	1288	A
1	X	1289	A
1	X	1291	G
1	X	1299	A
1	X	1314	A
1	X	1315	A
1	X	1333	G
1	X	1345	G
1	X	1355	A
1	X	1357	U
1	X	1403	U
1	X	1404	C
1	X	1409	U
1	X	1412	C

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Mol	Chain	Res	Type
1	X	1429	A
1	X	1433	A
1	X	1434	U
1	X	1439	G
1	X	1441	A
1	X	1442	C
1	X	1459	U
1	X	1473	U
1	X	1475	U
1	X	1482	U
1	X	1496	G
1	X	1508	G
1	X	1513	U
1	X	1552	C
1	X	1562	G
1	X	1570	C
1	X	1574	A
1	X	1575	C
1	X	1601	U
1	X	1602	G
1	X	1607	A
1	X	1613	G
1	X	1624	A
1	X	1631	C
1	X	1632	A
1	X	1634	A
1	X	1680	U
1	X	1684	G
1	X	1685	A
1	X	1710	U
1	X	1711	C
1	X	1716	G
1	X	1732	U
1	X	1749	G
1	X	1770	U
1	X	1775	A
1	X	1777	A
1	X	1790	G
1	X	1799	A
1	X	1800	A
1	X	1810	U
1	X	1811	A

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Mol	Chain	Res	Type
1	X	1820	G
1	X	1839	A
1	X	1871	G
1	X	1882	G
1	X	1883	A
1	X	1909	U
1	X	1920	A
1	X	1923	U
1	X	1938	U
1	X	1953	A
1	X	1963	G
1	X	1975	G
1	X	1980	A
1	X	2005	U
1	X	2014	A
1	X	2018	G
1	X	2045	A
1	X	2075	U
1	X	2083	G
1	X	2088	U
1	X	2165	A
1	X	2189	A
1	X	2193	C
1	X	2204	A
1	X	2217	G
1	X	2228	U
1	X	2254	C
1	X	2261	G
1	X	2265	A
1	X	2290	A
1	X	2298	U
1	X	2299	A
1	X	2305	C
1	X	2312	A
1	X	2314	A
1	X	2323	U
1	X	2354	G
1	X	2363	G
1	X	2370	G
1	X	2381	A
1	X	2401	A
1	X	2447	G

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Mol	Chain	Res	Type
1	X	2477	C
1	X	2482	A
1	X	2497	A
1	X	2498	U
1	X	2508	G
1	X	2551	A
1	X	2560	G
1	X	2561	G
1	X	2564	U
1	X	2565	C
1	X	2580	C
1	X	2634	G
1	X	2660	C
1	X	2691	C
1	X	2705	A
1	X	2731	G
1	X	2738	A
1	X	2758	A
1	X	2770	A
1	X	2777	A
1	X	2778	U
1	X	2808	U
1	X	2810	A
1	X	2824	C
1	X	2846	G
1	X	2848	A
1	X	2854	G
1	X	2867	G
1	X	2876	C
2	Y	4	C
2	Y	11	G
2	Y	14	C
2	Y	46	G
2	Y	58	G
2	Y	59	A
2	Y	86	A

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 35 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
32	1F2	X	2929	-	58,60,60	1.98	13 (22%)	81,90,90	2.09	26 (32%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
32	1F2	X	2929	-	-	1/74/115/115	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
32	X	2929	1F2	C30-N29	9.08	1.46	1.33
32	X	2929	1F2	C8-C10	4.43	1.59	1.52
32	X	2929	1F2	C5-N29	3.51	1.52	1.45
32	X	2929	1F2	O11-C12	-3.49	1.26	1.34
32	X	2929	1F2	C17-C2	3.22	1.60	1.53
32	X	2929	1F2	C10-C9	-3.18	1.46	1.55
32	X	2929	1F2	O11-C7	3.02	1.51	1.46
32	X	2929	1F2	O18-C6	-2.89	1.43	1.47
32	X	2929	1F2	C13-C12	-2.73	1.45	1.51
32	X	2929	1F2	C25-C6	2.39	1.57	1.52
32	X	2929	1F2	C15-C14	-2.34	1.48	1.54
32	X	2929	1F2	C15-C9	-2.30	1.49	1.55
32	X	2929	1F2	O18-C30	-2.10	1.33	1.36

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
32	X	2929	1F2	C24-C7-C6	-5.42	107.76	115.23
32	X	2929	1F2	C25-C6-C5	-4.99	109.41	116.42
32	X	2929	1F2	C48-C47-N45	4.90	107.56	101.94
32	X	2929	1F2	O16-C4-C2	-4.29	114.51	120.60
32	X	2929	1F2	O31-C30-N29	-4.26	124.25	129.22
32	X	2929	1F2	O46-C44-N45	-4.00	119.63	124.26
32	X	2929	1F2	O11-C12-C13	3.65	119.57	111.56
32	X	2929	1F2	C3-C4-C2	3.65	125.44	119.10
32	X	2929	1F2	C22-C15-C14	3.53	117.73	111.40
32	X	2929	1F2	C22-C15-C9	3.32	119.72	112.55
32	X	2929	1F2	C8-C10-C9	-2.92	106.72	110.25
32	X	2929	1F2	C50-N45-C47	-2.83	108.54	111.83
32	X	2929	1F2	C10-C9-C15	2.83	117.95	113.61
32	X	2929	1F2	O19-C10-C9	2.79	113.59	107.85
32	X	2929	1F2	C32-C37-C36	2.73	113.66	109.19
32	X	2929	1F2	C54-N55-C56	2.67	121.46	116.85
32	X	2929	1F2	O18-C6-C5	-2.65	100.53	103.21
32	X	2929	1F2	C32-O21-C9	-2.55	111.82	116.25
32	X	2929	1F2	O1-C14-C13	-2.51	103.12	107.55
32	X	2929	1F2	O18-C30-O31	2.49	124.53	121.66
32	X	2929	1F2	O27-C12-C13	-2.43	117.57	124.08
32	X	2929	1F2	C7-O11-C12	2.36	122.38	118.18
32	X	2929	1F2	O19-C10-C8	2.28	110.72	105.71
32	X	2929	1F2	C28-C13-C14	-2.12	108.12	112.92
32	X	2929	1F2	O11-C7-C6	2.08	110.18	105.48
32	X	2929	1F2	O41-C37-C36	2.03	113.38	109.77

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
32	X	2929	1F2	C35-C36-N38-C40

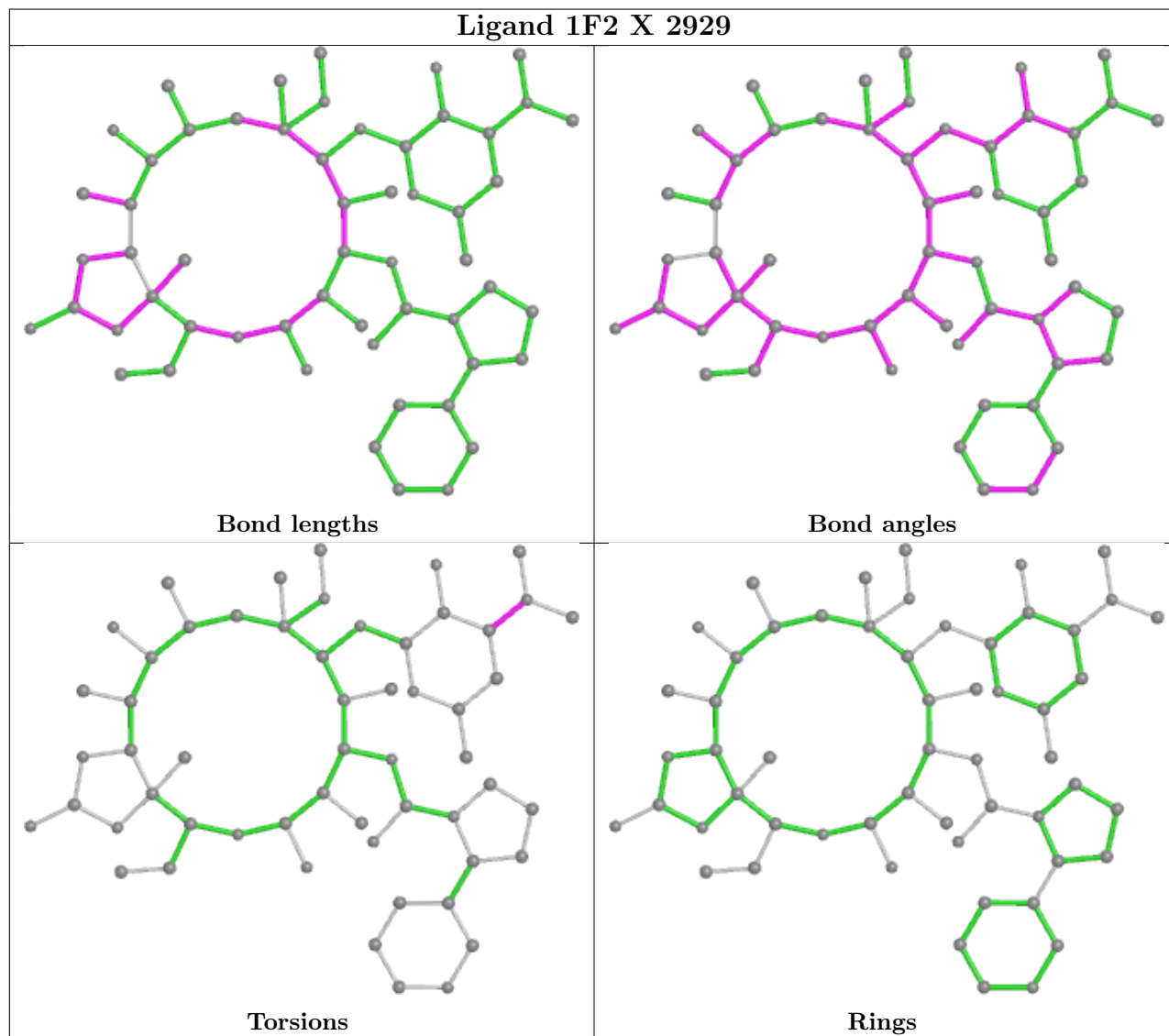
There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
32	X	2929	1F2	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	X	2686/2880 (93%)	-0.18	85 (3%) 47 31	43, 92, 197, 276	0
2	Y	122/123 (99%)	-0.05	2 (1%) 72 59	83, 136, 170, 192	0
3	A	240/274 (87%)	0.06	9 (3%) 40 26	68, 116, 146, 173	0
4	B	205/211 (97%)	-0.29	3 (1%) 73 61	45, 73, 106, 154	0
5	C	197/205 (96%)	0.17	13 (6%) 18 11	57, 114, 155, 187	0
6	D	177/180 (98%)	0.53	16 (9%) 9 5	146, 183, 216, 227	0
7	E	171/185 (92%)	-0.20	2 (1%) 79 67	92, 143, 192, 206	0
8	F	71/144 (49%)	2.27	36 (50%) 0 0	211, 236, 252, 257	0
9	G	142/174 (81%)	0.08	8 (5%) 24 13	73, 97, 145, 161	0
10	H	134/134 (100%)	-0.42	0 100 100	50, 70, 96, 120	0
11	I	141/156 (90%)	0.81	24 (17%) 1 1	67, 129, 174, 204	0
12	J	136/141 (96%)	0.11	3 (2%) 62 48	74, 103, 149, 184	0
13	K	113/116 (97%)	-0.42	0 100 100	35, 60, 79, 91	0
14	L	104/114 (91%)	0.39	9 (8%) 10 5	98, 134, 156, 168	0
15	M	108/166 (65%)	-0.41	0 100 100	50, 73, 111, 144	0
16	N	117/118 (99%)	-0.29	0 100 100	60, 90, 127, 160	0
17	O	94/100 (94%)	-0.22	3 (3%) 47 31	67, 115, 156, 173	0
18	P	127/134 (94%)	-0.39	2 (1%) 72 59	50, 67, 108, 158	0
19	Q	93/95 (97%)	-0.04	5 (5%) 25 14	73, 106, 162, 195	0
20	R	110/115 (95%)	0.22	10 (9%) 9 5	88, 117, 170, 178	0
21	S	175/237 (73%)	0.14	11 (6%) 20 11	121, 155, 175, 190	0
22	T	84/91 (92%)	0.58	13 (15%) 2 1	80, 108, 186, 199	0
23	U	72/81 (88%)	0.47	8 (11%) 5 3	92, 128, 153, 162	0
24	V	66/67 (98%)	0.55	8 (12%) 4 2	100, 132, 211, 216	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
25	W	55/55 (100%)	-0.39	0 100 100	81, 98, 126, 152	0
26	Z	58/60 (96%)	-0.14	1 (1%) 70 57	49, 71, 105, 113	0
27	1	53/55 (96%)	0.84	7 (13%) 3 2	8, 32, 61, 96	0
28	2	46/47 (97%)	3.75	39 (84%) 0 0	3, 16, 37, 59	0
29	3	63/66 (95%)	2.54	34 (53%) 0 0	3, 25, 40, 60	0
30	4	37/37 (100%)	6.83	36 (97%) 0 0	227, 254, 265, 269	0
All	All	5997/6561 (91%)	0.07	387 (6%) 18 11	3, 100, 196, 276	0

All (387) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
30	4	24	LEU	18.2
30	4	17	VAL	15.6
27	1	7	ARG	15.5
30	4	25	VAL	14.0
8	F	125	ASN	12.0
28	2	27	GLY	11.0
24	V	1	MET	10.7
30	4	11	CYS	10.4
22	T	9	SER	10.4
30	4	34	GLN	10.1
30	4	35	ARG	9.7
30	4	28	SER	9.6
30	4	20	HIS	9.4
30	4	6	SER	9.2
2	Y	123	U	9.0
1	X	731	A	8.6
29	3	38	GLY	8.6
30	4	19	ARG	8.6
30	4	29	ASN	8.4
26	Z	2	ALA	8.4
22	T	8	GLY	8.3
30	4	10	MET	8.3
30	4	21	GLY	8.2
30	4	12	ASP	7.9
29	3	40	GLU	7.9
8	F	114	ASP	7.9
11	I	9	THR	7.8
29	3	33	ASN	7.8
30	4	7	VAL	7.8

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Mol	Chain	Res	Type	RSRZ
28	2	32	ALA	7.7
30	4	16	VAL	7.5
11	I	8	PRO	7.4
29	3	37	SER	7.4
23	U	27	ASP	7.2
24	V	3	PRO	7.2
11	I	6	LEU	7.1
11	I	5	ASP	7.1
24	V	2	LYS	7.0
29	3	31	HIS	6.9
3	A	203	ASN	6.9
30	4	22	ARG	6.8
30	4	5	SER	6.5
30	4	23	VAL	6.5
28	2	26	SER	6.4
28	2	9	ASN	6.3
6	D	43	SER	6.2
30	4	26	ILE	6.2
9	G	97	ASP	6.0
1	X	1095	A	5.9
28	2	7	PRO	5.9
28	2	8	ASN	5.8
30	4	14	CYS	5.8
8	F	126	THR	5.8
8	F	85	GLY	5.7
1	X	248	A	5.7
28	2	36	ALA	5.7
30	4	36	GLN	5.7
23	U	28	GLY	5.7
11	I	29	THR	5.6
11	I	52	GLY	5.4
29	3	45	GLY	5.3
28	2	24	THR	5.3
8	F	121	GLU	5.3
22	T	10	SER	5.3
1	X	1085	G	5.3
29	3	30	ARG	5.3
5	C	165	SER	5.2
5	C	19	LEU	5.2
6	D	42	SER	5.1
5	C	44	SER	5.1
27	1	2	ALA	5.1

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Mol	Chain	Res	Type	RSRZ
28	2	34	ARG	5.1
1	X	1077	U	5.1
1	X	1086	C	5.0
30	4	13	ASN	5.0
12	J	84	MET	5.0
1	X	1106	A	5.0
2	Y	2	C	4.9
30	4	4	ARG	4.9
11	I	63	ARG	4.8
1	X	1068	A	4.8
1	X	1090	C	4.8
14	L	58	ALA	4.8
11	I	4	HIS	4.7
6	D	145	MET	4.7
30	4	27	CYS	4.7
28	2	22	MET	4.7
30	4	1	MET	4.6
28	2	35	ARG	4.6
5	C	47	THR	4.6
11	I	114	ILE	4.6
28	2	29	ASN	4.6
28	2	4	THR	4.6
11	I	33	GLY	4.5
8	F	115	LEU	4.5
1	X	1080	A	4.5
28	2	11	LYS	4.5
29	3	7	HIS	4.5
29	3	60	LEU	4.4
8	F	144	ALA	4.4
14	L	64	LYS	4.4
5	C	197	GLU	4.4
8	F	84	ILE	4.4
4	B	205	SER	4.4
30	4	30	VAL	4.3
1	X	1523	A	4.3
11	I	30	ALA	4.3
28	2	30	ILE	4.3
28	2	33	ARG	4.3
6	D	147	ASP	4.3
8	F	102	ASP	4.3
1	X	1107	A	4.3
29	3	42	ARG	4.3

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Mol	Chain	Res	Type	RSRZ
1	X	665	A	4.2
28	2	16	HIS	4.2
6	D	20	PHE	4.2
28	2	37	LYS	4.2
28	2	6	GLN	4.2
1	X	1089	C	4.1
28	2	2	LYS	4.1
11	I	7	LYS	4.1
27	1	43	VAL	4.1
11	I	10	PRO	4.1
22	T	6	GLY	4.1
1	X	1072	U	4.1
1	X	728	G	4.0
1	X	1081	A	4.0
22	T	11	LYS	4.0
21	S	58	GLY	4.0
1	X	727	U	4.0
1	X	2776	U	4.0
19	Q	64	ARG	4.0
20	R	57	ASN	4.0
28	2	25	LYS	3.9
29	3	2	PRO	3.9
28	2	41	GLN	3.9
3	A	250	TRP	3.9
8	F	137	THR	3.9
28	2	20	ALA	3.9
6	D	85	VAL	3.9
27	1	44	ALA	3.9
24	V	4	SER	3.9
29	3	10	ALA	3.9
20	R	99	VAL	3.8
1	X	730	C	3.8
14	L	97	HIS	3.8
28	2	14	LYS	3.8
1	X	1103	C	3.8
12	J	140	GLU	3.8
6	D	146	VAL	3.8
29	3	39	ASP	3.8
27	1	25	THR	3.8
23	U	47	HIS	3.8
1	X	1552	C	3.8
8	F	142	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	X	2166	G	3.7
30	4	15	LYS	3.7
20	R	94	VAL	3.7
30	4	37	GLY	3.7
29	3	8	LYS	3.7
28	2	1	MET	3.7
5	C	198	GLU	3.7
8	F	113	PRO	3.7
29	3	34	THR	3.7
1	X	1524	C	3.6
1	X	1093	U	3.6
22	T	7	VAL	3.6
1	X	558	G	3.6
1	X	1186	G	3.6
8	F	122	ALA	3.6
22	T	73	GLY	3.6
29	3	32	GLN	3.5
8	F	116	ASN	3.5
9	G	129	HIS	3.5
6	D	22	TYR	3.5
8	F	136	VAL	3.5
29	3	28	GLY	3.5
1	X	1065	A	3.5
1	X	1066	G	3.5
1	X	1091	C	3.5
9	G	156	HIS	3.5
20	R	102	LYS	3.5
27	1	26	LYS	3.5
11	I	82	ASP	3.4
1	X	2774	U	3.4
20	R	100	ASP	3.4
21	S	59	GLY	3.4
1	X	2777	A	3.4
1	X	1187	A	3.4
28	2	15	THR	3.4
21	S	143	ILE	3.4
29	3	35	GLY	3.4
8	F	97	GLY	3.3
3	A	272	THR	3.3
1	X	1551	U	3.3
1	X	2088	U	3.3
29	3	36	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
3	A	249	PRO	3.2
6	D	74	ILE	3.2
4	B	135	HIS	3.2
6	D	142	THR	3.2
30	4	32	HIS	3.2
1	X	1084	A	3.1
30	4	18	ARG	3.1
1	X	891	A	3.1
1	X	2169	A	3.1
1	X	1104	G	3.1
14	L	54	ALA	3.1
30	4	8	LYS	3.1
1	X	1098	G	3.1
17	O	39	PHE	3.1
1	X	1120	C	3.1
20	R	98	ILE	3.1
6	D	37	ASN	3.0
1	X	2780	A	3.0
1	X	2170	C	3.0
29	3	27	SER	3.0
1	X	1094	C	3.0
11	I	36	GLY	3.0
1	X	1078	A	3.0
1	X	1188	A	3.0
1	X	1108	U	3.0
1	X	1189	G	2.9
14	L	63	ASN	2.9
8	F	101	TRP	2.9
28	2	18	PHE	2.9
1	X	1734	C	2.9
8	F	100	ASN	2.9
8	F	133	SER	2.9
4	B	94	ASP	2.9
24	V	36	GLN	2.9
1	X	1109	A	2.9
21	S	86	VAL	2.9
8	F	123	ALA	2.9
1	X	2089	C	2.9
11	I	48	PHE	2.9
28	2	10	ARG	2.8
29	3	47	GLY	2.8
1	X	2082	C	2.8

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Mol	Chain	Res	Type	RSRZ
1	X	1067	G	2.8
6	D	23	SER	2.8
28	2	46	ASP	2.8
8	F	108	ALA	2.8
1	X	2775	U	2.8
6	D	144	ASP	2.8
1	X	2290	A	2.8
11	I	54	SER	2.8
20	R	77	HIS	2.7
21	S	12	GLN	2.7
1	X	2778	U	2.7
24	V	11	ALA	2.7
5	C	123	PHE	2.7
22	T	74	LYS	2.7
1	X	1913	G	2.7
30	4	3	VAL	2.7
1	X	1079	G	2.7
14	L	57	ALA	2.7
29	3	6	THR	2.7
7	E	5	GLY	2.7
18	P	8	PHE	2.7
1	X	2287	G	2.7
20	R	58	VAL	2.6
29	3	41	ILE	2.6
8	F	76	TYR	2.6
3	A	220	HIS	2.6
14	L	62	GLY	2.6
8	F	107	ILE	2.6
14	L	53	ALA	2.6
28	2	12	ARG	2.6
1	X	2167	A	2.6
19	Q	94	GLN	2.6
21	S	163	ASP	2.6
19	Q	65	VAL	2.6
1	X	2087	U	2.6
24	V	10	GLN	2.6
8	F	111	LYS	2.6
1	X	1071	U	2.5
1	X	1553	G	2.5
1	X	2090	U	2.5
30	4	33	LYS	2.5
17	O	41	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
8	F	119	SER	2.5
30	4	31	LYS	2.5
5	C	121	ASP	2.5
1	X	1114	A	2.5
20	R	112	LYS	2.5
22	T	17	ASN	2.5
30	4	9	LYS	2.5
23	U	50	ALA	2.5
3	A	254	THR	2.5
29	3	13	ARG	2.5
21	S	171	VAL	2.5
28	2	5	TYR	2.5
23	U	16	ASN	2.5
29	3	29	LYS	2.5
8	F	87	GLY	2.4
5	C	48	ARG	2.4
3	A	219	PRO	2.4
19	Q	69	ILE	2.4
21	S	11	LYS	2.4
1	X	1057	A	2.4
11	I	32	ARG	2.4
1	X	1121	G	2.4
1	X	2174	G	2.4
29	3	3	LYS	2.4
21	S	142	ASN	2.4
8	F	83	GLY	2.4
21	S	92	VAL	2.4
12	J	141	ALA	2.4
22	T	3	HIS	2.4
22	T	84	ALA	2.4
28	2	13	ALA	2.4
14	L	34	SER	2.4
3	A	101	GLU	2.4
28	2	28	ARG	2.3
28	2	38	GLY	2.3
29	3	58	MET	2.3
11	I	113	GLU	2.3
28	2	19	ARG	2.3
23	U	25	ARG	2.3
9	G	40	ASN	2.3
22	T	4	LYS	2.3
9	G	155	THR	2.3

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Mol	Chain	Res	Type	RSRZ
8	F	118	GLY	2.3
28	2	40	HIS	2.3
21	S	124	ALA	2.3
3	A	100	GLY	2.3
23	U	39	LYS	2.3
29	3	56	ALA	2.3
29	3	19	THR	2.3
11	I	56	LEU	2.3
28	2	21	ARG	2.3
8	F	92	ASN	2.2
9	G	37	ASP	2.2
1	X	1062	G	2.2
1	X	729	A	2.2
29	3	4	MET	2.2
8	F	90	THR	2.2
20	R	68	GLY	2.2
8	F	141	GLY	2.2
9	G	103	TYR	2.2
24	V	6	MET	2.2
5	C	21	GLU	2.2
1	X	1092	U	2.2
1	X	2581	A	2.2
1	X	726	G	2.2
1	X	1050	G	2.2
29	3	17	THR	2.2
5	C	20	PRO	2.2
11	I	87	THR	2.2
11	I	103	ASN	2.2
6	D	121	ALA	2.2
6	D	21	GLY	2.2
1	X	1073	G	2.1
1	X	434	C	2.1
17	O	36	LYS	2.1
1	X	2198	U	2.1
29	3	18	GLY	2.1
8	F	143	ASN	2.1
28	2	23	LYS	2.1
7	E	119	ALA	2.1
11	I	17	LYS	2.1
28	2	3	ARG	2.1
8	F	112	MET	2.1
5	C	57	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
29	3	44	LYS	2.1
8	F	74	MET	2.1
22	T	5	LYS	2.1
1	X	2173	G	2.1
18	P	134	LYS	2.1
6	D	50	ILE	2.1
1	X	1753	A	2.1
27	1	23	THR	2.1
19	Q	56	MET	2.1
1	X	1069	G	2.0
1	X	1909	U	2.0
1	X	247	A	2.0
8	F	138	VAL	2.0
9	G	110	LEU	2.0
5	C	66	ASN	2.0
29	3	57	ARG	2.0
8	F	124	ALA	2.0
1	X	1525	A	2.0
23	U	52	ARG	2.0
1	X	1064	C	2.0
1	X	1190	C	2.0
11	I	90	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

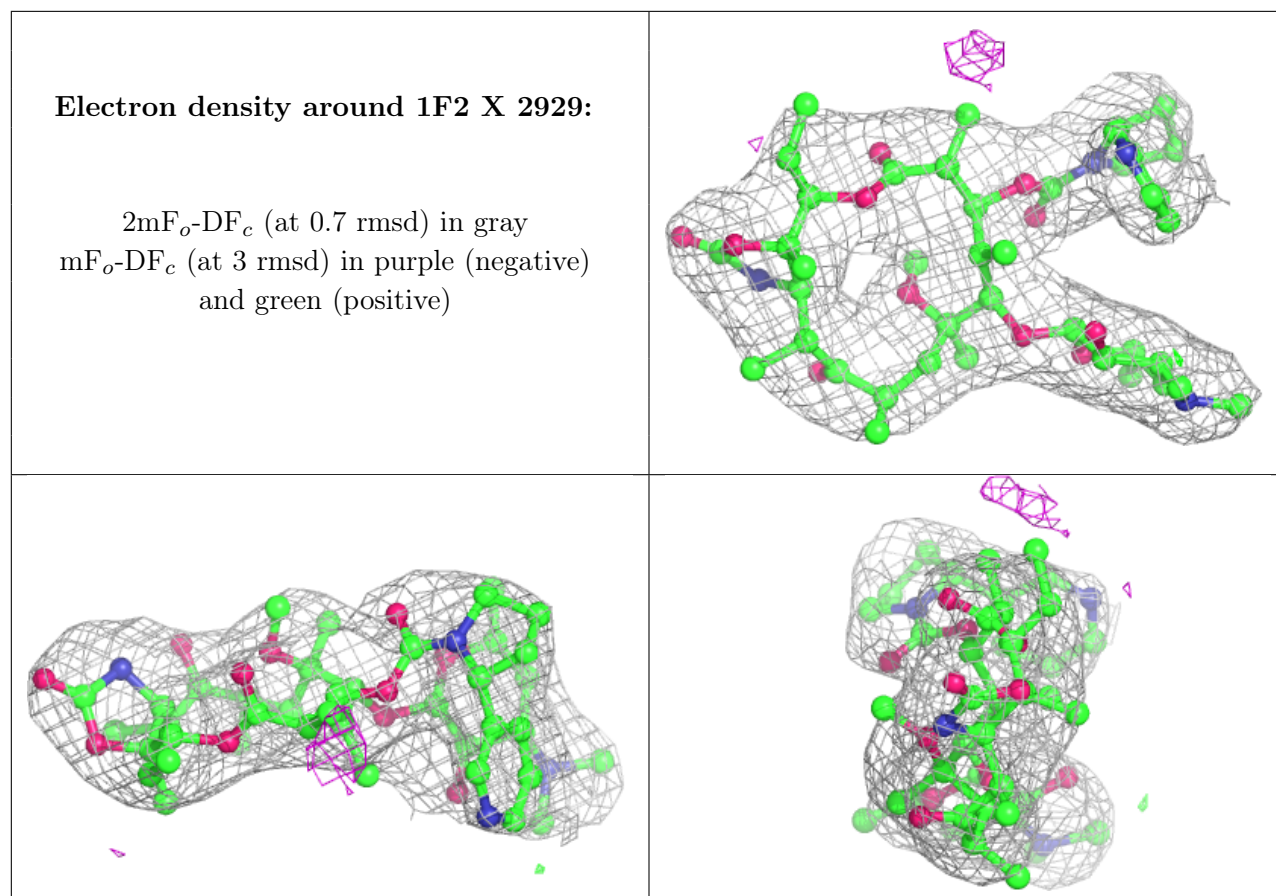
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
31	MG	Y	203	1/1	0.45	0.72	87,87,87,87	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
31	MG	X	2902	1/1	0.60	0.63	94,94,94,94	0
31	MG	X	2903	1/1	0.64	0.67	89,89,89,89	0
31	MG	Y	201	1/1	0.66	0.82	89,89,89,89	0
31	MG	X	2905	1/1	0.78	0.27	65,65,65,65	0
31	MG	X	2912	1/1	0.78	0.50	71,71,71,71	0
31	MG	J	201	1/1	0.79	0.30	100,100,100,100	0
31	MG	X	2928	1/1	0.84	0.53	62,62,62,62	0
31	MG	Y	205	1/1	0.85	0.30	79,79,79,79	0
31	MG	X	2909	1/1	0.90	0.51	96,96,96,96	0
31	MG	X	2924	1/1	0.91	1.20	69,69,69,69	0
31	MG	X	2907	1/1	0.91	0.65	51,51,51,51	0
31	MG	X	2904	1/1	0.92	0.29	107,107,107,107	0
31	MG	X	2917	1/1	0.92	0.54	55,55,55,55	0
31	MG	X	2925	1/1	0.94	0.53	122,122,122,122	0
31	MG	X	2906	1/1	0.94	0.41	58,58,58,58	0
31	MG	X	2918	1/1	0.95	0.67	42,42,42,42	0
31	MG	X	2919	1/1	0.95	0.47	30,30,30,30	0
31	MG	X	2920	1/1	0.95	0.20	115,115,115,115	0
31	MG	X	2915	1/1	0.95	0.68	57,57,57,57	0
31	MG	X	2911	1/1	0.95	0.31	68,68,68,68	0
31	MG	X	2913	1/1	0.96	0.63	61,61,61,61	0
31	MG	Y	204	1/1	0.96	0.31	82,82,82,82	0
31	MG	X	2922	1/1	0.96	0.65	44,44,44,44	0
31	MG	Y	202	1/1	0.96	0.51	58,58,58,58	0
31	MG	X	2914	1/1	0.97	0.40	27,27,27,27	0
31	MG	X	2927	1/1	0.97	0.50	62,62,62,62	0
31	MG	X	2921	1/1	0.97	0.42	81,81,81,81	0
32	1F2	X	2929	56/56	0.97	0.19	42,68,77,83	0
31	MG	X	2916	1/1	0.98	0.62	37,37,37,37	0
31	MG	X	2910	1/1	0.98	0.44	42,42,42,42	0
31	MG	X	2908	1/1	0.98	0.92	37,37,37,37	0
31	MG	X	2923	1/1	0.98	0.38	34,34,34,34	0
31	MG	M	201	1/1	0.98	0.55	23,23,23,23	0
31	MG	X	2901	1/1	0.98	0.36	50,50,50,50	0
31	MG	X	2926	1/1	0.99	0.97	45,45,45,45	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.