Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 02:35 AM JST

PDB ID	:	5YZ3
Title	:	Crystal structure of T2R-TTL-28 complex
Authors	:	Yu, Y.; Chen, Q.
Deposited on	:	2017-12-12
Resolution	:	2.54 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:

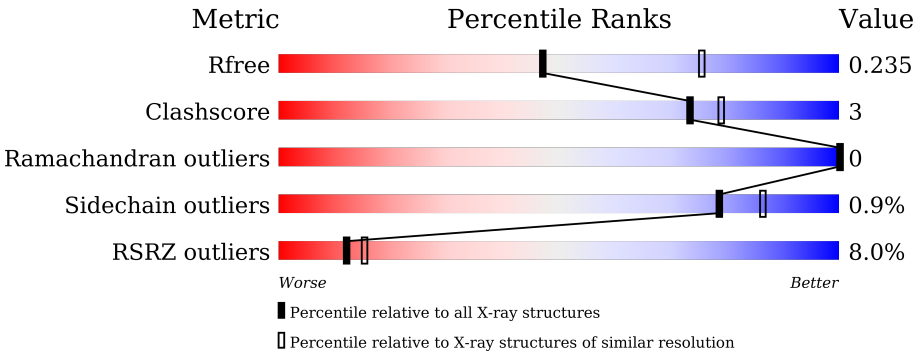
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.54 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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 <=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	A	450	93%
1	C	450	91% 7%
2	B	445	4% 88% 8%
2	D	445	9% 89% 5% 5%
3	E	143	8% 80% 15%
4	F	384	25% 80% 6% 13%

2 Entry composition (i)

There are 15 unique types of molecules in this entry. The entry contains 35117 atoms, of which 16904 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 protein called Tubulin alpha-1B chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	437	Total	C	H	N	O	S	0	1	0
			6743	2166	3321	582	651	23			
1	C	440	Total	C	H	N	O	S	0	1	0
			6783	2178	3340	585	657	23			

- Molecule 2 is a protein called Tubulin beta-2B chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	427	Total	C	H	N	O	S	0	2	0
			6615	2118	3239	578	653	27			
2	D	421	Total	C	H	N	O	S	0	1	0
			6499	2083	3184	563	641	28			

- Molecule 3 is a protein called Stathmin-4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	E	121	Total	C	H	N	O	S	0	0	0
			2014	617	1014	181	197	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	3	MET	-	expression tag	UNP P63043
E	4	ALA	-	expression tag	UNP P63043

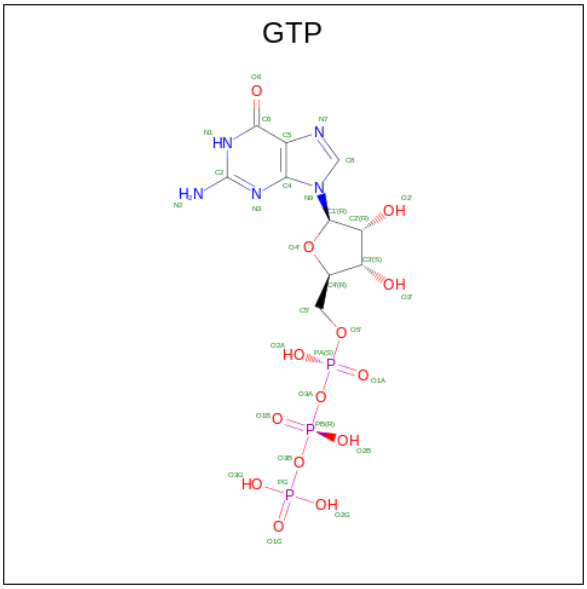
- Molecule 4 is a protein called Tubulin Tyrosine Ligase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	F	334	Total	C	H	N	O	S	0	1	0
			5465	1767	2710	474	500	14			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	379	HIS	-	expression tag	UNP E1BQ43
F	380	HIS	-	expression tag	UNP E1BQ43
F	381	HIS	-	expression tag	UNP E1BQ43
F	382	HIS	-	expression tag	UNP E1BQ43
F	383	HIS	-	expression tag	UNP E1BQ43
F	384	HIS	-	expression tag	UNP E1BQ43

- Molecule 5 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	0	0
			42	10	10	5	14	3		
5	C	1	Total	C	H	N	O	P	0	0
			42	10	10	5	14	3		
5	D	1	Total	C	H	N	O	P	0	0
			42	10	10	5	14	3		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		
6	B	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

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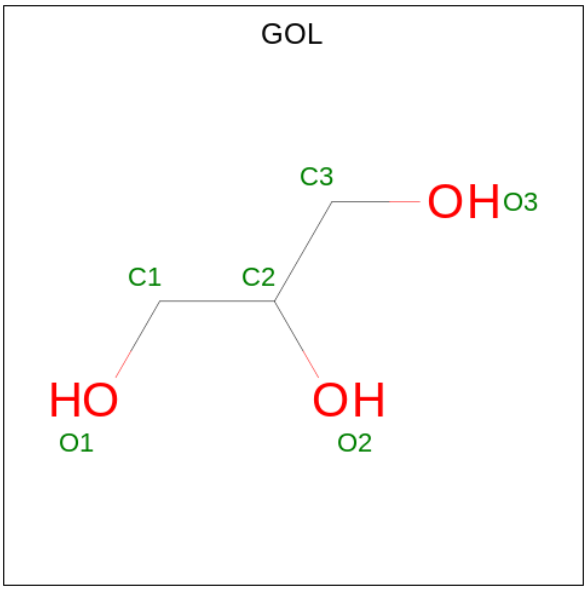
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total Mg 1 1	0	0

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

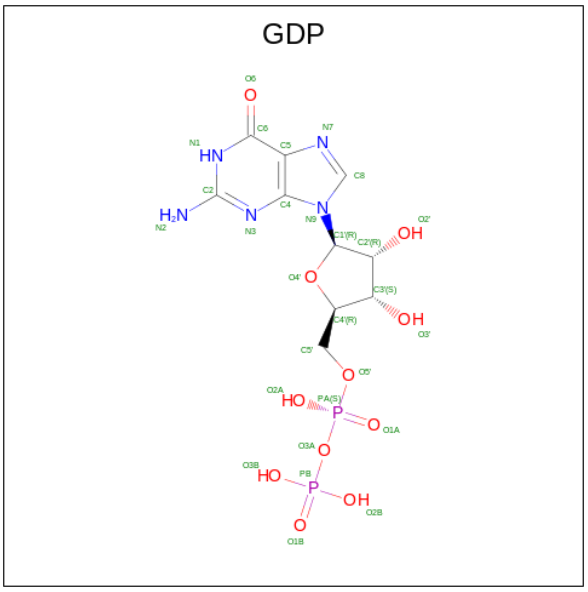
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Ca 1 1	0	0
7	C	1	Total Ca 1 1	0	0

- Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



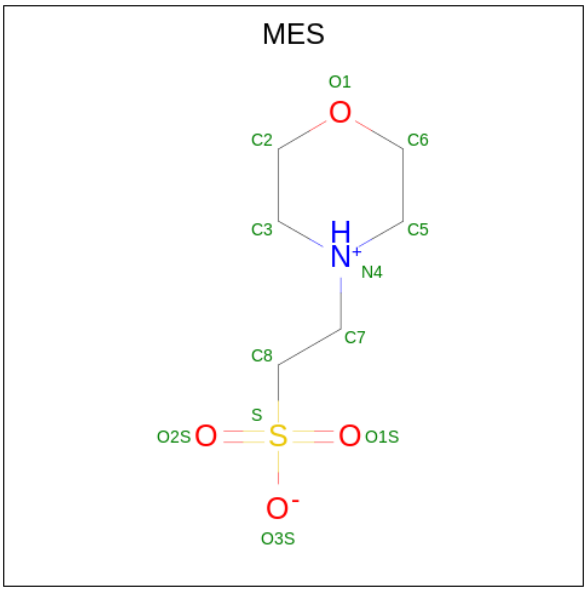
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C H O 14 3 8 3	0	0

- Molecule 9 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: $C_{10}H_{15}N_5O_{11}P_2$).



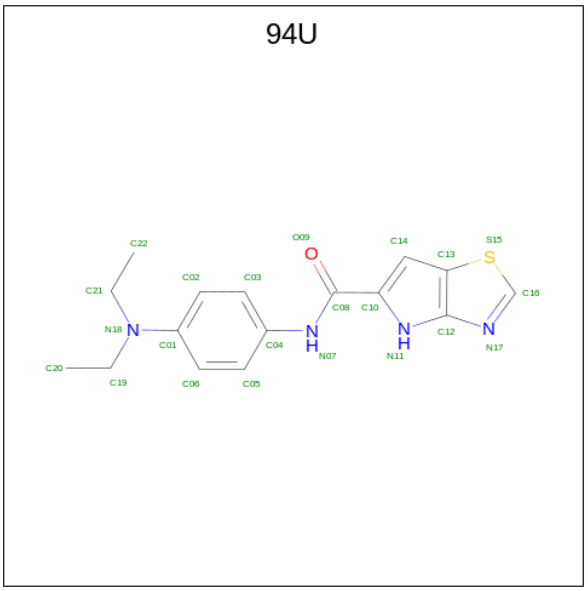
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
9	B	1	Total	C	H	N	O	P	0	0
			38	10	10	5	11	2		

- Molecule 10 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
10	B	1	Total	C	H	N	O	S	0	0
			24	6	12	1	4	1		

- Molecule 11 is N-[4-(diethylamino)phenyl]-4H-pyrrolo[2,3-d][1,3]thiazole-5-carboxamide (three-letter code: 94U) (formula: $C_{16}H_{18}N_4OS$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	B	1	Total C H N O S 40 16 18 4 1 1	0	0
11	D	1	Total C H N O S 40 16 18 4 1 1	0	0

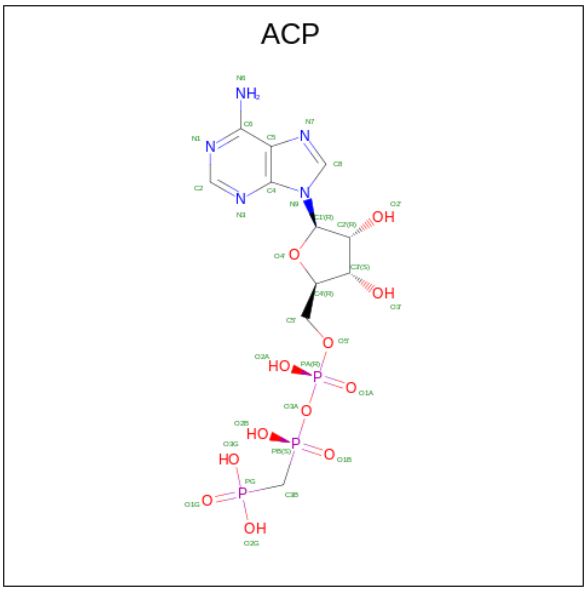
- Molecule 12 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
12	B	1	Total Na 1 1	0	0

- Molecule 13 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
13	D	1	Total Cl 1 1	0	0

- Molecule 14 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (three-letter code: ACP) (formula: $C_{11}H_{18}N_5O_{12}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	F	1	Total	C	N	O	P	0	0
			31	11	5	12	3		

- Molecule 15 is water.

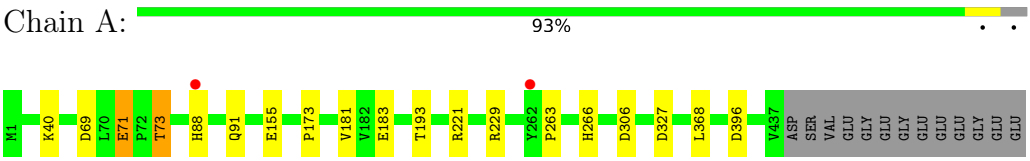
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	151	Total	O	0	0
			151	151		
15	B	127	Total	O	0	0
			127	127		
15	C	243	Total	O	0	0
			243	243		
15	D	75	Total	O	0	0
			75	75		
15	E	27	Total	O	0	0
			27	27		
15	F	54	Total	O	0	0
			54	54		

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.

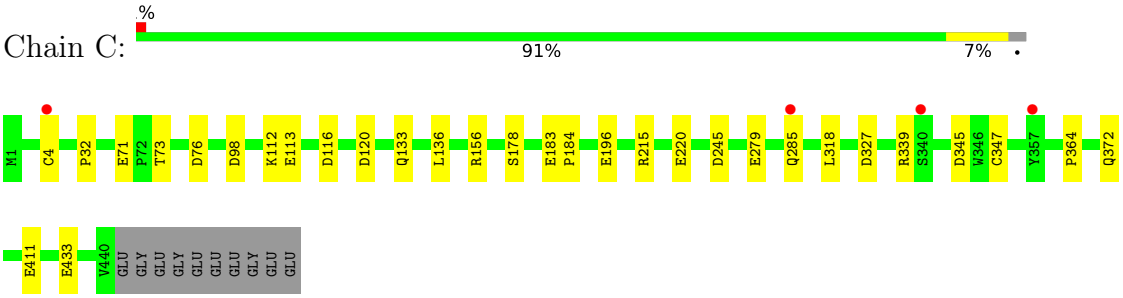
• Molecule 1: Tubulin alpha-1B chain

Chain A:



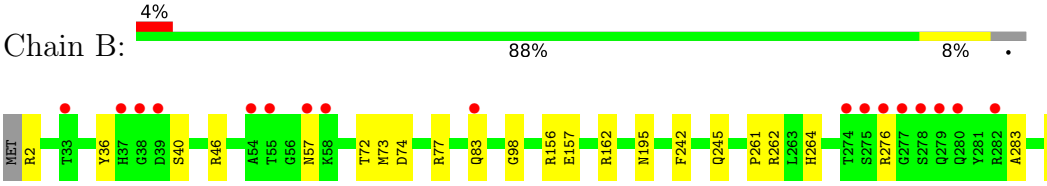
• Molecule 1: Tubulin alpha-1B chain

Chain C:



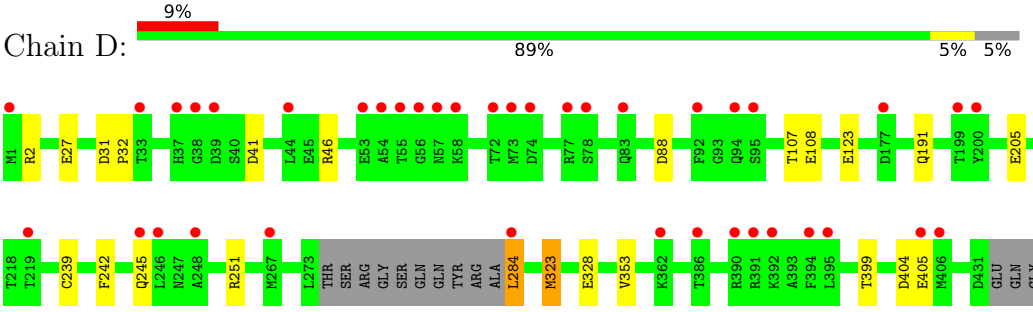
• Molecule 2: Tubulin beta-2B chain

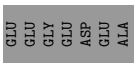
Chain B:



• Molecule 2: Tubulin beta-2B chain

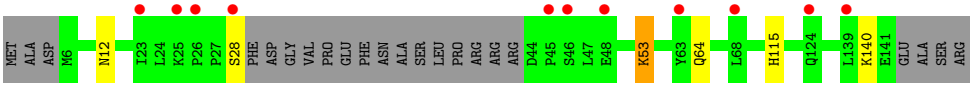
Chain D:





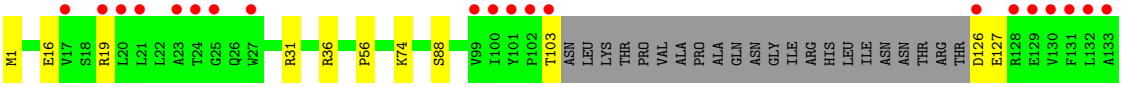
• Molecule 3: Stathmin-4

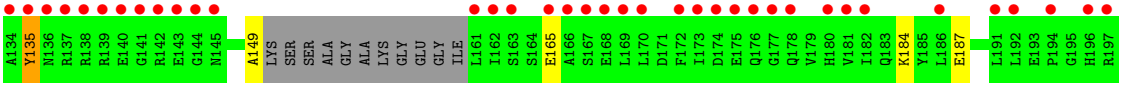
Chain E:

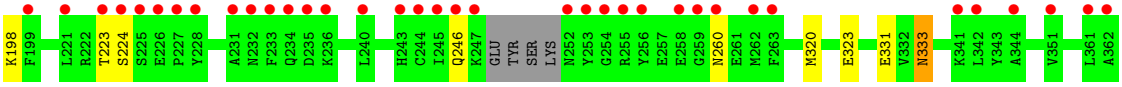


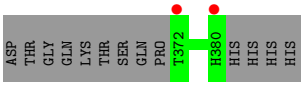
• Molecule 4: Tubulin Tyrosine Ligase

Chain F:









4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	104.89Å 157.50Å 180.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.62 – 2.54 49.62 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.4 (49.62-2.54) 92.1 (49.62-2.54)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.06 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.210 , 0.234 0.212 , 0.235	Depositor DCC
R_{free} test set	1500 reflections (1.54%)	wwPDB-VP
Wilson B-factor (Å^2)	33.3	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 43.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	35117	wwPDB-VP
Average B, all atoms (Å^2)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, GDP, MG, 94U, ACP, CL, MES, GTP, GOL, CA

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	A	0.23	0/3500	0.38	0/4751
1	C	0.24	0/3521	0.38	0/4780
2	B	0.23	0/3451	0.38	0/4674
2	D	0.24	0/3388	0.38	0/4589
3	E	0.23	0/1008	0.34	0/1337
4	F	0.23	0/2817	0.40	1/3805 (0.0%)
All	All	0.24	0/17685	0.38	1/23936 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	F	88	SER	C-N-CA	7.20	139.70	121.70

There are no chirality outliers.

There are no planarity outliers.

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	A	3422	3321	3334	16	0
1	C	3443	3340	3352	27	0
2	B	3376	3239	3247	24	0
2	D	3315	3184	3193	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	1000	1014	1018	6	0
4	F	2755	2710	2721	21	0
5	A	32	10	12	0	0
5	C	32	10	12	0	0
5	D	32	10	12	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	1	0	0	0	0
7	C	1	0	0	0	0
8	A	6	8	8	0	0
9	B	28	10	12	0	0
10	B	12	12	12	0	0
11	B	22	18	0	0	0
11	D	22	18	0	0	0
12	B	1	0	0	0	0
13	D	1	0	0	1	0
14	F	31	0	14	7	0
15	A	151	0	0	11	2
15	B	127	0	0	14	0
15	C	243	0	0	21	2
15	D	75	0	0	12	0
15	E	27	0	0	4	0
15	F	54	0	0	6	0
All	All	18213	16904	16947	114	2

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 3.

All (114) 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)
4:F:331:GLU:CD	14:F:401:ACP:H3B2	1.73	1.08
4:F:331:GLU:OE2	14:F:401:ACP:H3B2	1.53	1.07
1:A:40:LYS:NZ	15:A:602:HOH:O	1.92	1.02
1:C:339:ARG:O	15:C:601:HOH:O	1.81	0.98
4:F:331:GLU:OE2	14:F:401:ACP:C3B	2.11	0.98
1:A:327:ASP:OD2	15:A:601:HOH:O	1.86	0.93
3:E:28:SER:O	15:E:201:HOH:O	1.87	0.92
13:D:504:CL:CL	15:D:625:HOH:O	2.26	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:41:ASP:OD2	15:D:601:HOH:O	1.92	0.86
4:F:36:ARG:NH2	15:F:502:HOH:O	2.08	0.86
1:A:193:THR:OG1	15:A:603:HOH:O	1.93	0.85
3:E:12:ASN:OD1	15:E:202:HOH:O	1.94	0.84
1:C:120:ASP:OD2	15:C:602:HOH:O	1.96	0.84
1:C:220:GLU:O	15:C:603:HOH:O	1.97	0.82
2:D:27:GLU:OE1	15:D:602:HOH:O	1.98	0.81
1:A:69:ASP:OD2	15:A:604:HOH:O	1.99	0.81
2:D:123:GLU:O	15:D:604:HOH:O	2.01	0.79
2:B:292:GLN:OE1	15:B:601:HOH:O	2.00	0.79
2:D:245:GLN:O	15:D:603:HOH:O	2.00	0.78
1:C:196:GLU:OE1	15:C:604:HOH:O	2.00	0.78
4:F:149:ALA:O	15:F:501:HOH:O	2.02	0.78
1:A:71:GLU:O	15:A:605:HOH:O	2.02	0.77
1:A:181:VAL:O	15:A:606:HOH:O	2.02	0.76
4:F:56:PRO:O	15:F:502:HOH:O	2.04	0.76
4:F:74:LYS:NZ	4:F:331:GLU:OE1	2.20	0.75
1:A:396:ASP:OD2	15:A:607:HOH:O	2.03	0.75
1:C:318:LEU:O	15:C:605:HOH:O	2.05	0.75
4:F:331:GLU:OE2	14:F:401:ACP:PG	2.47	0.72
2:B:291:GLN:O	15:B:602:HOH:O	2.08	0.71
2:B:245:GLN:O	15:B:603:HOH:O	2.09	0.69
1:C:32:PRO:O	15:C:606:HOH:O	2.10	0.68
2:D:245:GLN:OE1	15:D:605:HOH:O	2.11	0.67
2:B:162:ARG:NH2	15:B:614:HOH:O	2.27	0.67
3:E:64:GLN:OE1	15:E:203:HOH:O	2.11	0.67
1:C:279:GLU:OE1	15:C:608:HOH:O	2.13	0.66
1:C:76:ASP:OD2	15:C:607:HOH:O	2.13	0.65
4:F:323:GLU:OE1	15:F:503:HOH:O	2.15	0.63
1:C:285:GLN:NE2	1:C:372:GLN:OE1	2.32	0.61
1:A:221:ARG:NH1	2:B:327:ASP:OD2	2.32	0.61
4:F:224:SER:HA	4:F:246:GLN:HE22	1.64	0.61
2:D:284:LEU:N	15:D:614:HOH:O	2.33	0.60
2:D:251:ARG:NH2	15:D:615:HOH:O	2.34	0.60
4:F:126:ASP:OD1	4:F:127:GLU:N	2.34	0.60
1:C:327:ASP:OD2	15:C:609:HOH:O	2.17	0.58
2:D:245:GLN:NE2	15:D:618:HOH:O	2.37	0.58
1:A:173:PRO:O	15:A:608:HOH:O	2.18	0.57
2:D:88:ASP:OD2	15:D:606:HOH:O	2.17	0.56
2:D:404:ASP:OD1	2:D:405:GLU:N	2.39	0.55
2:B:83:GLN:O	2:B:83:GLN:NE2	2.39	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:345:ASP:OD2	15:C:610:HOH:O	2.18	0.54
1:C:245:ASP:N	15:C:613:HOH:O	2.20	0.54
4:F:246:GLN:OE1	4:F:260:ASN:ND2	2.40	0.54
1:C:279:GLU:OE1	1:C:279:GLU:N	2.42	0.53
1:C:411:GLU:OE1	15:C:611:HOH:O	2.19	0.53
2:B:73:MET:N	15:B:607:HOH:O	2.24	0.52
2:B:334:GLN:NE2	15:B:619:HOH:O	2.31	0.52
4:F:16:GLU:OE2	4:F:19:ARG:NH1	2.42	0.52
2:D:399:THR:O	3:E:140:LYS:NZ	2.43	0.52
1:A:155:GLU:OE1	3:E:53:LYS:NZ	2.43	0.52
2:B:74:ASP:OD1	2:B:77:ARG:NH2	2.43	0.51
2:B:301:ALA:O	2:B:303:CYS:N	2.43	0.51
1:C:133:GLN:NE2	15:C:632:HOH:O	2.35	0.50
2:D:328:GLU:OE1	15:D:607:HOH:O	2.19	0.50
1:C:71:GLU:OE1	1:C:73:THR:N	2.43	0.50
2:D:107:THR:OG1	2:D:108:GLU:N	2.45	0.50
3:E:115:HIS:ND1	15:E:204:HOH:O	2.25	0.50
1:A:263:PRO:O	1:A:266:HIS:ND1	2.40	0.50
2:B:46:ARG:NH2	2:B:242:PHE:O	2.44	0.50
2:D:46:ARG:NH1	2:D:239[B]:CYS:O	2.45	0.49
2:B:283:ALA:N	15:B:622:HOH:O	2.33	0.49
2:D:46:ARG:NH1	2:D:239[A]:CYS:O	2.45	0.49
4:F:331:GLU:OE2	4:F:333:ASN:ND2	2.45	0.49
1:C:364:PRO:HA	15:C:626:HOH:O	2.13	0.49
1:C:98:ASP:OD2	15:C:612:HOH:O	2.19	0.48
1:A:368:LEU:O	15:A:610:HOH:O	2.20	0.48
14:F:401:ACP:O2B	14:F:401:ACP:O3G	2.32	0.48
1:A:71:GLU:OE2	1:A:73:THR:OG1	2.22	0.48
2:B:57:ASN:OD1	15:B:605:HOH:O	2.20	0.48
2:B:72:THR:N	15:B:607:HOH:O	2.47	0.47
1:C:112:LYS:NZ	1:C:113:GLU:OE2	2.44	0.47
1:A:306:ASP:OD1	15:A:611:HOH:O	2.21	0.47
14:F:401:ACP:H3B1	14:F:401:ACP:O1A	2.14	0.47
2:B:157:GLU:OE2	15:B:606:HOH:O	2.21	0.46
4:F:103:THR:N	15:F:511:HOH:O	2.48	0.46
2:D:46:ARG:NH2	2:D:242:PHE:O	2.49	0.46
4:F:135:TYR:OH	4:F:165:GLU:HA	2.16	0.46
2:B:309:ARG:NH1	15:B:631:HOH:O	2.44	0.45
1:C:433:GLU:O	15:C:614:HOH:O	2.21	0.45
2:B:331:LEU:O	2:B:335:ASN:ND2	2.49	0.45
2:B:261:PRO:O	2:B:264:HIS:ND1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:98:GLY:N	15:B:633:HOH:O	2.49	0.44
4:F:223:THR:O	4:F:260:ASN:ND2	2.51	0.44
2:B:379:LYS:NZ	15:B:635:HOH:O	2.51	0.43
1:C:178:SER:OG	1:C:183:GLU:OE1	2.24	0.43
4:F:331:GLU:OE1	15:F:505:HOH:O	2.21	0.42
2:B:262:ARG:NE	2:B:421:GLU:OE2	2.45	0.42
2:B:72:THR:HB	15:B:607:HOH:O	2.20	0.42
1:C:156:ARG:NH1	15:C:647:HOH:O	2.48	0.42
1:A:229:ARG:HD2	15:A:616:HOH:O	2.19	0.41
1:C:116:ASP:O	15:C:616:HOH:O	2.21	0.41
2:D:323:MET:HE2	2:D:353:VAL:HG21	2.02	0.41
1:A:88:HIS:N	1:A:91:GLN:OE1	2.46	0.41
4:F:184:LYS:NZ	4:F:187:GLU:OE2	2.53	0.41
1:C:215:ARG:NH1	15:C:654:HOH:O	2.52	0.41
2:B:36:TYR:OH	2:B:40:SER:O	2.36	0.41
4:F:333:ASN:OD1	14:F:401:ACP:O1G	2.39	0.41
15:C:759:HOH:O	2:D:2:ARG:HG2	2.20	0.41
2:D:191:GLN:HB3	15:D:610:HOH:O	2.20	0.41
2:B:156:ARG:NH1	2:B:195:ASN:OD1	2.50	0.40
2:D:31:ASP:HB2	2:D:32:PRO:HD2	2.03	0.40
1:C:4[B]:CYS:SG	1:C:136:LEU:HG	2.61	0.40
1:C:183:GLU:N	1:C:184:PRO:CD	2.83	0.40
4:F:198:LYS:O	4:F:224:SER:N	2.49	0.40
1:C:215:ARG:NH1	15:C:659:HOH:O	2.55	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
15:A:641:HOH:O	15:C:731:HOH:O[3_545]	1.84	0.36
15:A:691:HOH:O	15:C:772:HOH:O[3_545]	2.19	0.01

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	
1	A	436/450 (97%)	421 (97%)	15 (3%)	0	100	100
1	C	439/450 (98%)	430 (98%)	9 (2%)	0	100	100
2	B	427/445 (96%)	412 (96%)	15 (4%)	0	100	100
2	D	418/445 (94%)	404 (97%)	14 (3%)	0	100	100
3	E	117/143 (82%)	115 (98%)	2 (2%)	0	100	100
4	F	325/384 (85%)	318 (98%)	7 (2%)	0	100	100
All	All	2162/2317 (93%)	2100 (97%)	62 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	
1	A	369/378 (98%)	366 (99%)	3 (1%)	81	88
1	C	372/378 (98%)	371 (100%)	1 (0%)	92	96
2	B	371/383 (97%)	368 (99%)	3 (1%)	81	88
2	D	365/383 (95%)	362 (99%)	3 (1%)	81	88
3	E	109/127 (86%)	108 (99%)	1 (1%)	78	86
4	F	302/342 (88%)	297 (98%)	5 (2%)	60	75
All	All	1888/1991 (95%)	1872 (99%)	16 (1%)	78	88

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	73	THR
1	A	183	GLU
2	B	2	ARG

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Mol	Chain	Res	Type
2	B	276	ARG
2	B	300	MET
1	C	347	CYS
2	D	205	GLU
2	D	284	LEU
2	D	323	MET
3	E	53	LYS
4	F	1	MET
4	F	31	ARG
4	F	135	TYR
4	F	320	MET
4	F	333	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	137	HIS
2	B	190	HIS
2	D	245	GLN
3	E	12	ASN
3	E	71	HIS
4	F	246	GLN
4	F	260	ASN
4	F	333	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 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	$\#\|Z\| > 2$
10	MES	B	503	-	12,12,12	2.20	1 (8%)	14,16,16	2.45	5 (35%)
8	GOL	A	504	-	5,5,5	0.37	0	5,5,5	0.20	0
5	GTP	C	501	6	26,34,34	1.15	2 (7%)	32,54,54	1.40	5 (15%)
5	GTP	D	501	6	26,34,34	1.14	2 (7%)	32,54,54	1.54	6 (18%)
9	GDP	B	501	6	24,30,30	0.96	1 (4%)	30,47,47	1.07	3 (10%)
5	GTP	A	501	6	26,34,34	1.18	2 (7%)	32,54,54	1.38	6 (18%)
11	94U	D	502	-	20,24,24	2.16	9 (45%)	20,33,33	1.76	3 (15%)
14	ACP	F	401	-	27,33,33	2.02	6 (22%)	32,52,52	1.55	6 (18%)
11	94U	B	504	-	20,24,24	1.94	7 (35%)	20,33,33	1.67	3 (15%)

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
10	MES	B	503	-	-	1/6/14/14	0/1/1/1
8	GOL	A	504	-	-	2/4/4/4	-
5	GTP	C	501	6	-	6/18/38/38	0/3/3/3
5	GTP	D	501	6	-	5/18/38/38	0/3/3/3
9	GDP	B	501	6	-	3/12/32/32	0/3/3/3
5	GTP	A	501	6	-	9/18/38/38	0/3/3/3
11	94U	D	502	-	-	4/13/16/16	0/3/3/3
14	ACP	F	401	-	-	5/15/38/38	0/3/3/3
11	94U	B	504	-	-	8/13/16/16	0/3/3/3

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	B	503	MES	C8-S	-7.43	1.66	1.77
14	F	401	ACP	PB-O2B	-4.80	1.45	1.56
14	F	401	ACP	PG-O1G	4.66	1.60	1.50
5	A	501	GTP	C5-C6	-4.25	1.38	1.47
5	C	501	GTP	C5-C6	-4.14	1.39	1.47
11	D	502	94U	C12-N17	4.06	1.42	1.34
5	D	501	GTP	C5-C6	-4.03	1.39	1.47
14	F	401	ACP	PG-O2G	-4.02	1.45	1.54
11	B	504	94U	C12-N17	3.68	1.41	1.34
14	F	401	ACP	C2'-C1'	-3.52	1.48	1.53
11	D	502	94U	C04-N07	3.45	1.48	1.41
14	F	401	ACP	PB-O1B	3.24	1.59	1.51
11	B	504	94U	C04-N07	3.11	1.47	1.41
11	D	502	94U	C12-N11	3.08	1.40	1.34
11	D	502	94U	C03-C02	2.98	1.44	1.38
11	B	504	94U	C03-C02	2.95	1.44	1.38
11	B	504	94U	C12-N11	2.91	1.40	1.34
11	D	502	94U	C08-N07	2.90	1.43	1.35
9	B	501	GDP	C6-N1	-2.90	1.33	1.37
11	B	504	94U	C08-N07	2.77	1.43	1.35
11	D	502	94U	C06-C05	2.74	1.43	1.38
11	B	504	94U	C06-C05	2.60	1.43	1.38
11	D	502	94U	C10-C08	2.36	1.56	1.50
5	D	501	GTP	C2-N3	2.21	1.38	1.33
11	D	502	94U	O09-C08	2.20	1.27	1.23
11	B	504	94U	O09-C08	2.19	1.27	1.23
11	D	502	94U	C13-S15	2.14	1.75	1.74
14	F	401	ACP	C5-N7	-2.04	1.32	1.39
5	A	501	GTP	C2-N3	2.03	1.38	1.33
5	C	501	GTP	C2-N3	2.03	1.38	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	B	503	MES	C5-N4-C3	5.44	121.07	108.83
11	D	502	94U	C10-C08-N07	5.04	126.26	114.04
11	B	504	94U	C10-C08-N07	4.79	125.67	114.04
10	B	503	MES	C7-N4-C5	4.28	122.19	111.23
5	D	501	GTP	PB-O3B-PG	-3.54	120.68	132.83
14	F	401	ACP	N3-C2-N1	-3.35	123.44	128.68
14	F	401	ACP	C3'-C2'-C1'	3.35	106.02	100.98
5	D	501	GTP	PA-O3A-PB	-3.28	121.55	132.83
14	F	401	ACP	PB-O3A-PA	-3.25	122.27	132.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	D	501	GTP	C5-C6-N1	3.22	119.64	113.95
10	B	503	MES	O3S-S-C8	3.22	110.97	105.77
5	A	501	GTP	C5-C6-N1	3.13	119.48	113.95
5	C	501	GTP	PB-O3B-PG	-3.12	122.13	132.83
5	D	501	GTP	C8-N7-C5	3.03	108.77	102.99
11	D	502	94U	O09-C08-N07	-3.03	116.78	123.71
5	A	501	GTP	C8-N7-C5	3.03	108.75	102.99
5	C	501	GTP	C5-C6-N1	3.02	119.28	113.95
5	C	501	GTP	C8-N7-C5	2.99	108.68	102.99
5	A	501	GTP	PB-O3B-PG	-2.98	122.59	132.83
14	F	401	ACP	C4-C5-N7	-2.89	106.39	109.40
11	B	504	94U	O09-C08-N07	-2.87	117.16	123.71
5	D	501	GTP	C2-N1-C6	-2.78	119.97	125.10
5	A	501	GTP	C2-N1-C6	-2.77	120.00	125.10
5	C	501	GTP	C2-N1-C6	-2.67	120.18	125.10
5	C	501	GTP	PA-O3A-PB	-2.63	123.81	132.83
10	B	503	MES	C6-C5-N4	-2.42	106.44	110.10
5	A	501	GTP	PA-O3A-PB	-2.41	124.55	132.83
9	B	501	GDP	PA-O3A-PB	-2.40	124.61	132.83
10	B	503	MES	O1S-S-C8	2.38	109.78	106.92
9	B	501	GDP	C8-N7-C5	2.36	107.48	102.99
14	F	401	ACP	O2'-C2'-C1'	-2.36	102.16	110.85
9	B	501	GDP	C5-C6-N1	2.33	118.07	113.95
11	D	502	94U	C05-C04-C03	-2.31	115.87	119.03
11	B	504	94U	C05-C04-C03	-2.28	115.91	119.03
14	F	401	ACP	O3G-PG-C3B	2.25	111.85	106.40
5	A	501	GTP	O6-C6-C5	-2.23	120.01	124.37
5	D	501	GTP	C3'-C2'-C1'	2.13	104.19	100.98

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	501	GTP	C5'-O5'-PA-O1A
5	A	501	GTP	C5'-O5'-PA-O2A
5	C	501	GTP	C5'-O5'-PA-O1A
5	C	501	GTP	C5'-O5'-PA-O2A
5	D	501	GTP	C5'-O5'-PA-O1A
8	A	504	GOL	O1-C1-C2-C3
9	B	501	GDP	C5'-O5'-PA-O1A
9	B	501	GDP	C5'-O5'-PA-O2A
10	B	503	MES	C8-C7-N4-C5

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Mol	Chain	Res	Type	Atoms
14	F	401	ACP	C5'-O5'-PA-O3A
11	B	504	94U	C22-C21-N18-C01
8	A	504	GOL	O1-C1-C2-O2
11	D	502	94U	C22-C21-N18-C01
11	B	504	94U	C22-C21-N18-C19
11	B	504	94U	C20-C19-N18-C01
11	D	502	94U	C22-C21-N18-C19
11	D	502	94U	C20-C19-N18-C01
11	B	504	94U	C20-C19-N18-C21
5	A	501	GTP	C4'-C5'-O5'-PA
5	D	501	GTP	C4'-C5'-O5'-PA
5	C	501	GTP	PB-O3A-PA-O2A
5	D	501	GTP	C5'-O5'-PA-O2A
14	F	401	ACP	C5'-O5'-PA-O1A
14	F	401	ACP	C5'-O5'-PA-O2A
11	B	504	94U	C02-C01-N18-C21
5	A	501	GTP	PB-O3B-PG-O1G
5	C	501	GTP	C4'-C5'-O5'-PA
5	A	501	GTP	PB-O3A-PA-O2A
5	D	501	GTP	PB-O3A-PA-O2A
11	B	504	94U	C02-C01-N18-C19
11	B	504	94U	C06-C01-N18-C21
11	B	504	94U	C06-C01-N18-C19
11	D	502	94U	C20-C19-N18-C21
14	F	401	ACP	PG-C3B-PB-O1B
14	F	401	ACP	PB-C3B-PG-O1G
5	A	501	GTP	PB-O3B-PG-O2G
5	A	501	GTP	PB-O3B-PG-O3G
5	C	501	GTP	PB-O3B-PG-O2G
5	A	501	GTP	C5'-O5'-PA-O3A
5	C	501	GTP	C5'-O5'-PA-O3A
5	D	501	GTP	C5'-O5'-PA-O3A
9	B	501	GDP	C5'-O5'-PA-O3A
5	A	501	GTP	PB-O3A-PA-O1A

There are no ring outliers.

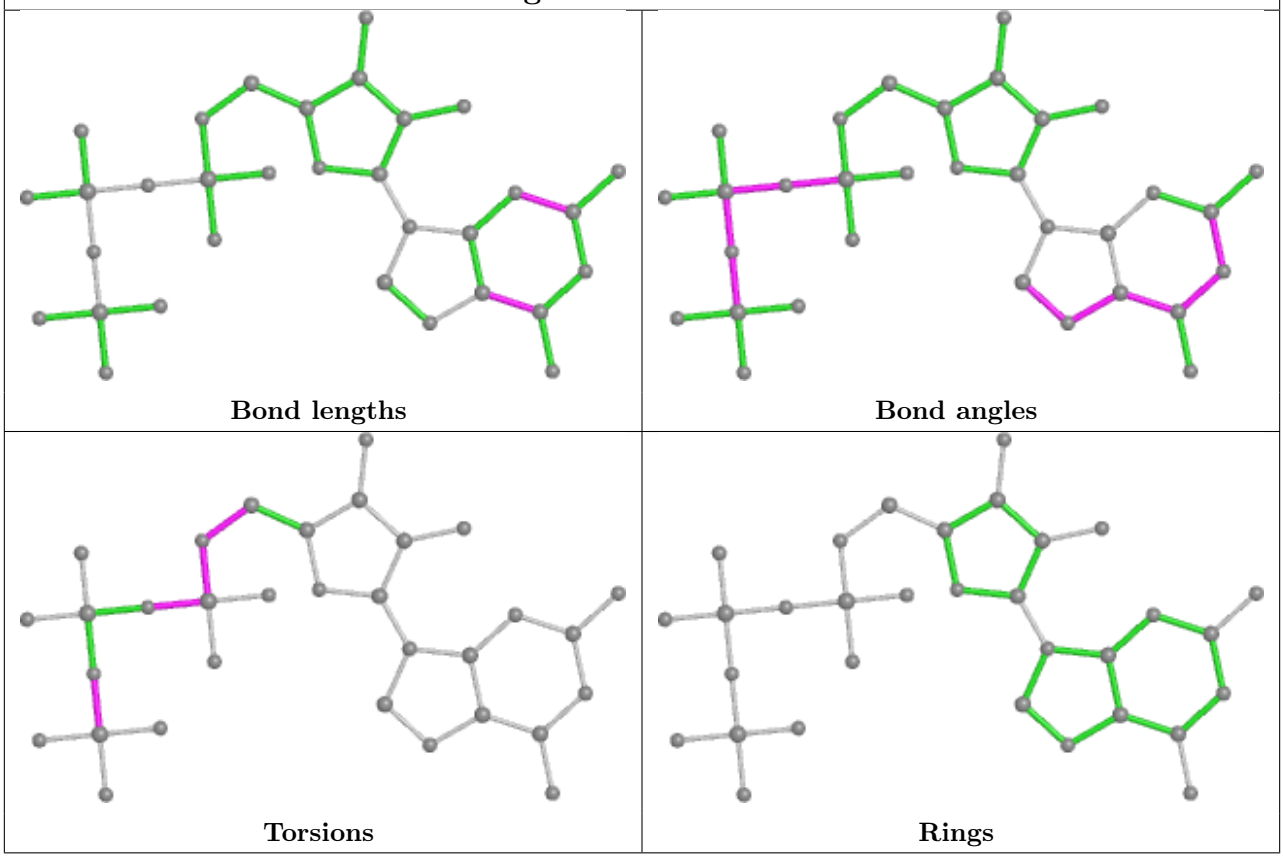
1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	F	401	ACP	7	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 then 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.

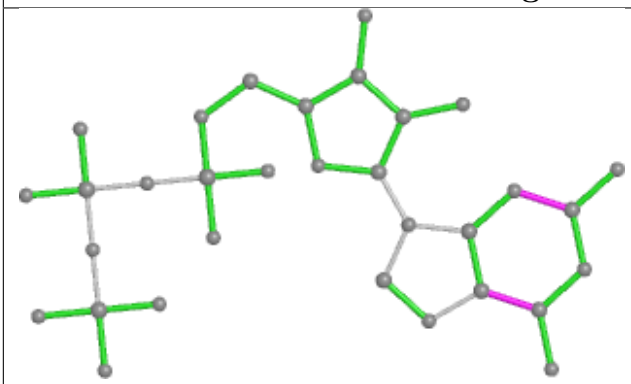
Ligand GTP C 501



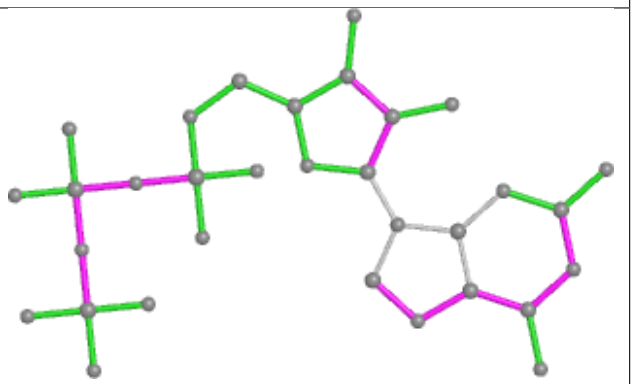
Bond lengths **Bond angles**

Torsions **Rings**

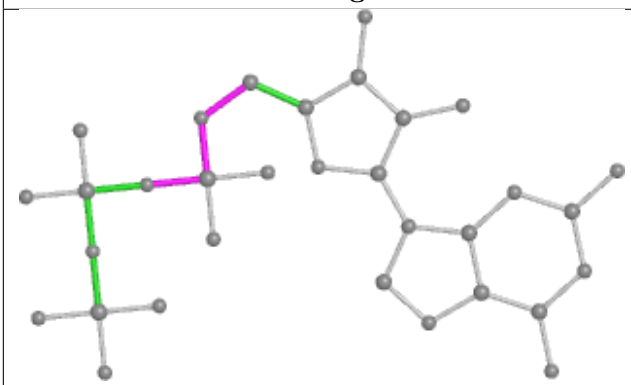
Ligand GTP D 501



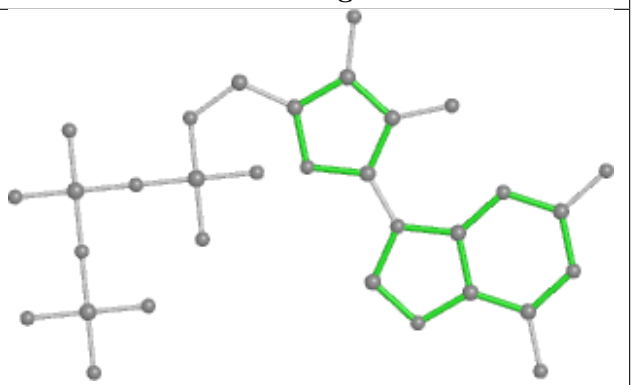
Bond lengths



Bond angles

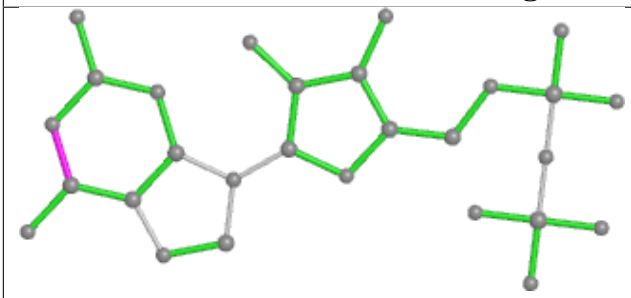


Torsions

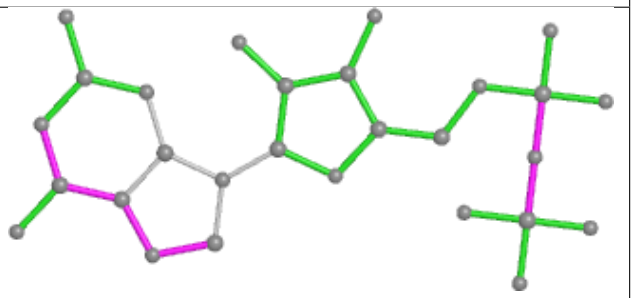


Rings

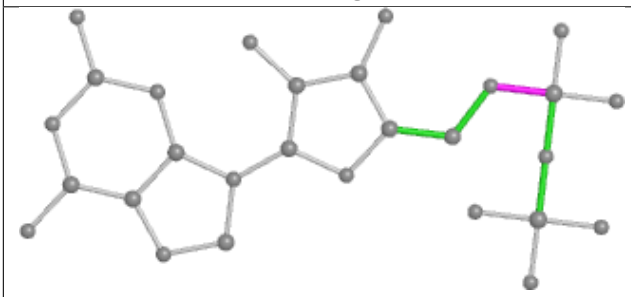
Ligand GDP B 501



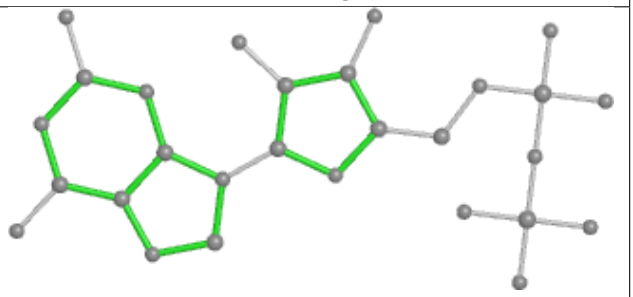
Bond lengths



Bond angles

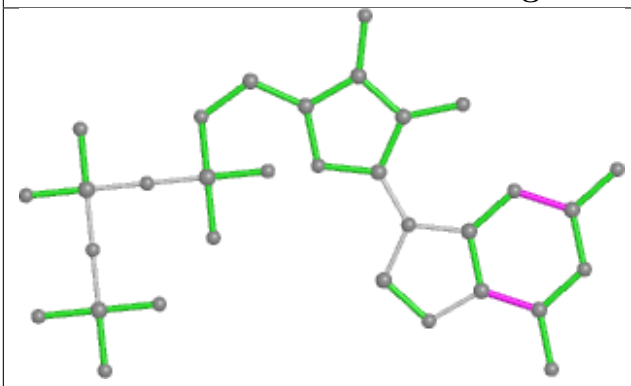


Torsions

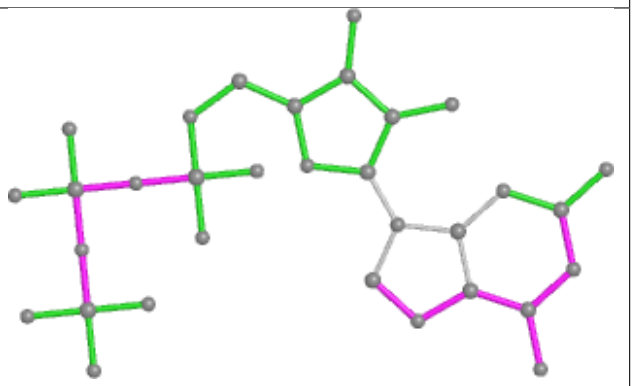


Rings

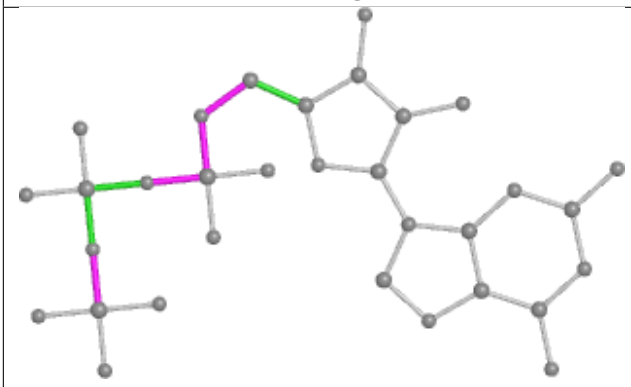
Ligand GTP A 501



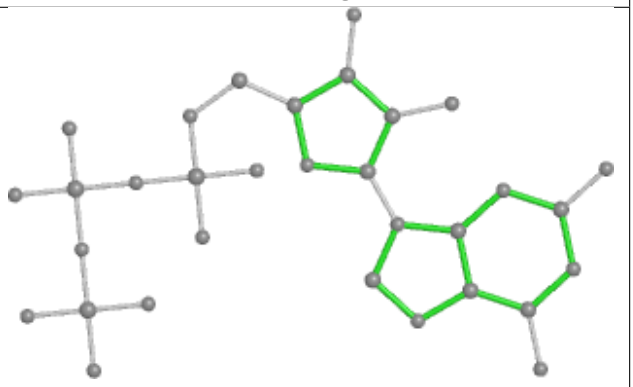
Bond lengths



Bond angles

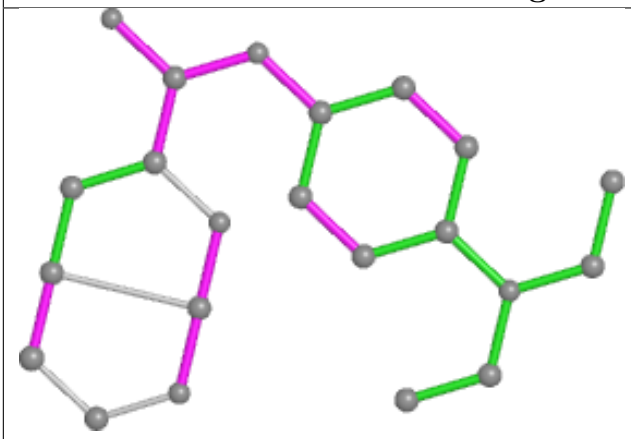


Torsions

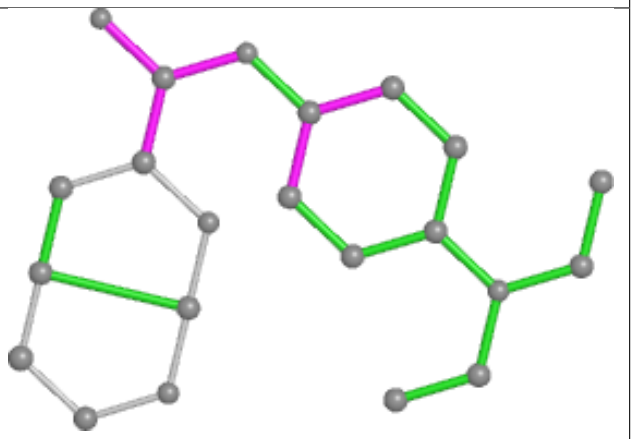


Rings

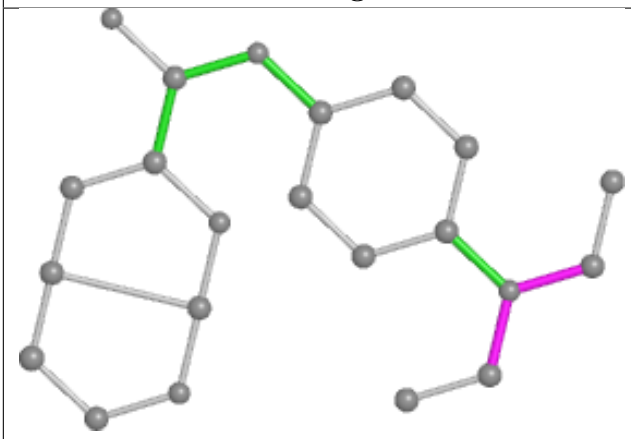
Ligand 94U D 502



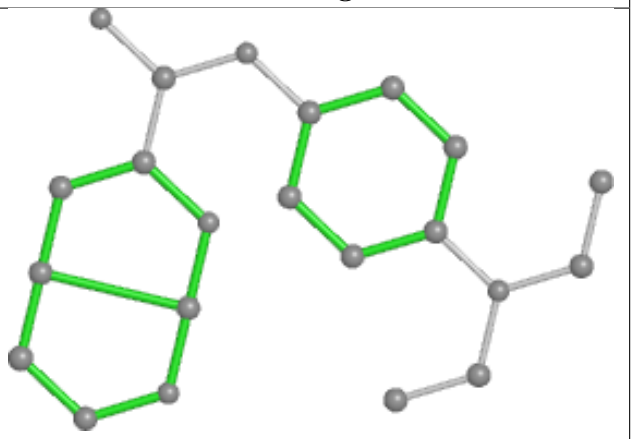
Bond lengths



Bond angles

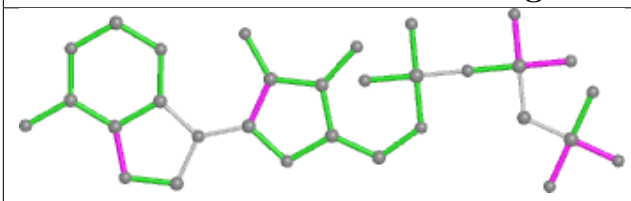


Torsions

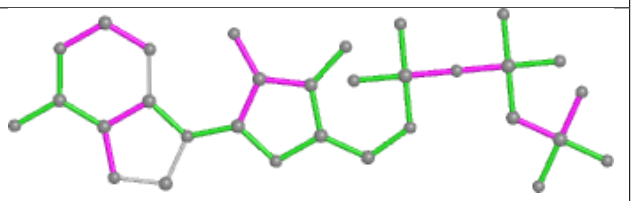


Rings

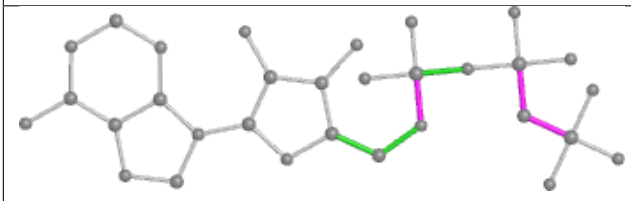
Ligand ACP F 401



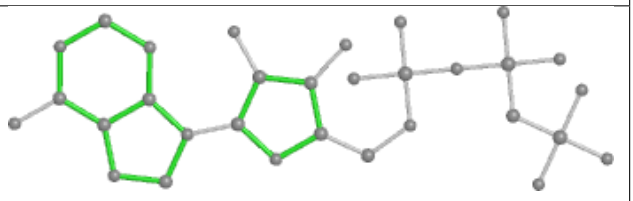
Bond lengths



Bond angles

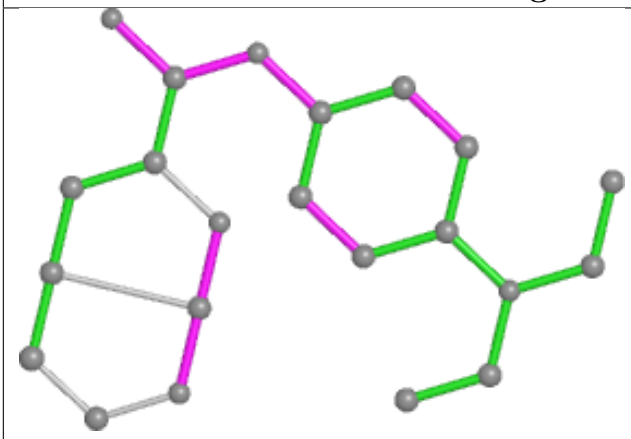


Torsions

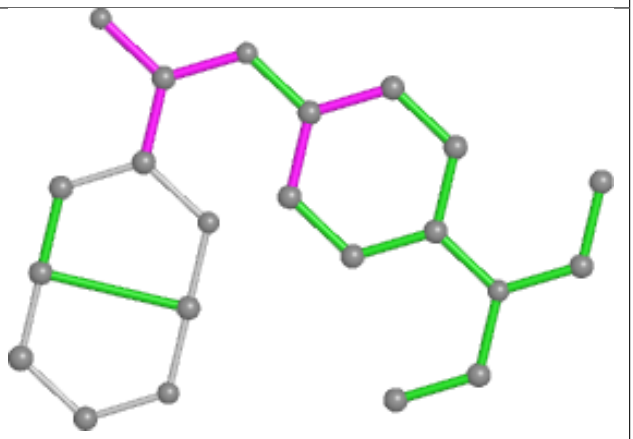


Rings

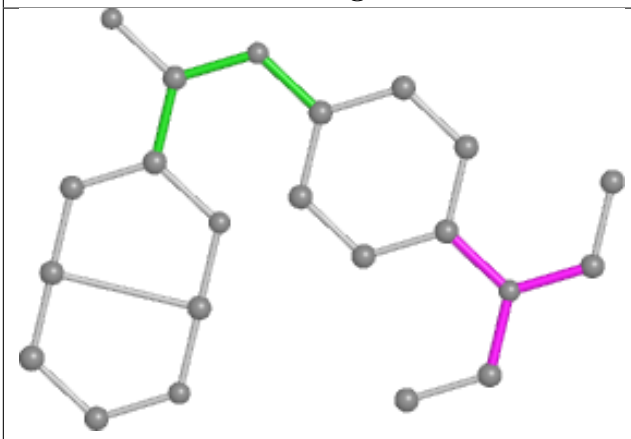
Ligand 94U B 504



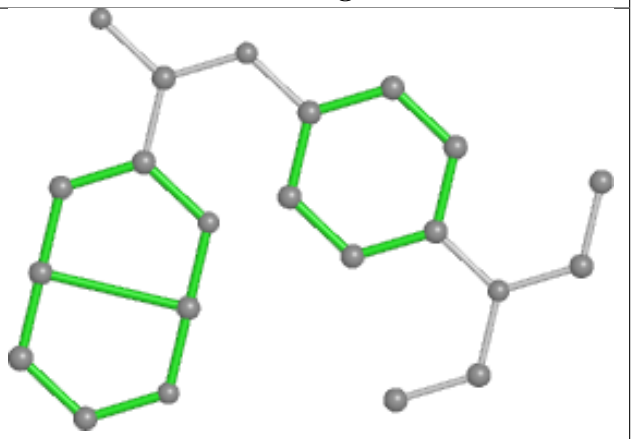
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	A	437/450 (97%)	0.16	2 (0%) 91 94	23, 37, 64, 82	0
1	C	440/450 (97%)	-0.08	4 (0%) 84 88	17, 29, 51, 72	0
2	B	427/445 (95%)	0.28	20 (4%) 31 38	19, 37, 75, 115	0
2	D	421/445 (94%)	0.71	42 (9%) 7 9	27, 54, 85, 111	0
3	E	121/143 (84%)	0.60	11 (9%) 9 11	26, 51, 85, 103	0
4	F	334/384 (86%)	1.38	95 (28%) 0 0	32, 64, 126, 150	0
All	All	2180/2317 (94%)	0.45	174 (7%) 12 15	17, 42, 89, 150	0

All (174) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	F	136	ASN	11.6
4	F	142	ARG	10.6
2	B	279	GLN	9.3
2	B	280	GLN	7.7
4	F	101	TYR	6.2
4	F	133	ALA	6.1
4	F	169	LEU	6.1
4	F	103	THR	5.7
4	F	173	ILE	5.7
4	F	259	GLY	5.6
4	F	253	TYR	5.6
4	F	244	CYS	5.5
4	F	170	LEU	5.5
4	F	172	PHE	5.4
4	F	176	GLN	5.3
4	F	130	VAL	5.3
4	F	372	THR	5.2
4	F	143	GLU	5.1
4	F	254	GLY	5.0

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Mol	Chain	Res	Type	RSRZ
4	F	178	GLN	4.9
4	F	256	TYR	4.8
4	F	141	GLY	4.7
4	F	137	ARG	4.7
4	F	131	PHE	4.7
4	F	132	LEU	4.6
4	F	139	ARG	4.6
4	F	233	PHE	4.5
4	F	100	ILE	4.5
2	B	276	ARG	4.5
4	F	129	GLU	4.5
4	F	194	PRO	4.5
2	D	1	MET	4.5
4	F	231	ALA	4.3
4	F	252	ASN	4.3
2	D	57	ASN	4.2
4	F	167	SER	4.1
4	F	177	GLY	4.1
4	F	232	ASN	4.1
4	F	223	THR	4.1
4	F	361	LEU	4.1
4	F	225	SER	4.1
4	F	255	ARG	4.0
2	D	55	THR	4.0
4	F	247	LYS	3.9
4	F	17	VAL	3.9
4	F	140	GLU	3.9
2	D	92	PHE	3.9
4	F	20	LEU	3.8
2	D	394	PHE	3.8
4	F	102	PRO	3.8
2	D	73	MET	3.8
4	F	236	LYS	3.8
2	D	77	ARG	3.7
2	B	57	ASN	3.7
2	D	39	ASP	3.6
4	F	240	LEU	3.6
2	B	275	SER	3.6
4	F	224	SER	3.6
4	F	246	GLN	3.5
2	D	83	GLN	3.5
4	F	182	ILE	3.5

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Mol	Chain	Res	Type	RSRZ
2	B	282	ARG	3.4
3	E	46	SER	3.4
4	F	234	GLN	3.4
2	D	391	ARG	3.4
4	F	24	THR	3.4
2	D	245	GLN	3.3
2	B	277	GLY	3.3
4	F	99	VAL	3.3
4	F	191	LEU	3.3
2	D	219	THR	3.3
3	E	48	GLU	3.3
4	F	144	GLY	3.3
1	A	262	TYR	3.3
2	D	284	LEU	3.2
4	F	380	HIS	3.2
4	F	138	ARG	3.2
4	F	27	TRP	3.2
4	F	126	ASP	3.2
4	F	362	ALA	3.1
4	F	162	ILE	3.1
4	F	161	LEU	3.1
4	F	25	GLY	3.1
4	F	258	GLU	3.1
2	D	390	ARG	3.1
2	B	55	THR	3.0
4	F	21	LEU	3.0
4	F	135	TYR	3.0
3	E	25	LYS	3.0
4	F	197	ARG	3.0
2	D	72	THR	3.0
2	D	54	ALA	2.9
4	F	235	ASP	2.9
4	F	245	ILE	2.9
2	B	83	GLN	2.8
2	B	274	THR	2.8
2	D	94	GLN	2.8
2	D	37	HIS	2.8
3	E	139	LEU	2.7
4	F	186	LEU	2.7
4	F	168	GLU	2.7
4	F	243	HIS	2.7
4	F	134	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	217	LEU	2.7
4	F	23	ALA	2.7
4	F	192	LEU	2.7
4	F	342	LEU	2.7
2	B	39	ASP	2.7
2	D	246	LEU	2.6
2	D	216	LYS	2.6
4	F	145	ASN	2.6
2	D	248	ALA	2.6
2	D	405	GLU	2.6
2	D	386	THR	2.5
4	F	166	ALA	2.5
4	F	174	ASP	2.5
2	D	212	PHE	2.5
3	E	28	SER	2.5
1	C	4[A]	CYS	2.5
2	D	95	SER	2.5
4	F	263	PHE	2.4
2	D	74	ASP	2.4
2	D	406	MET	2.4
3	E	26	PRO	2.4
3	E	68	LEU	2.4
2	B	54	ALA	2.4
2	D	33	THR	2.4
2	B	427	ASP	2.4
1	C	357	TYR	2.4
4	F	199	PHE	2.4
2	D	200	TYR	2.4
4	F	128	ARG	2.4
4	F	181	VAL	2.3
3	E	63	TYR	2.3
4	F	163	SER	2.3
2	D	38	GLY	2.3
1	A	88	HIS	2.3
3	E	45	PRO	2.3
2	B	33	THR	2.3
2	D	392	LYS	2.3
1	C	340	SER	2.3
4	F	260	ASN	2.3
4	F	165	GLU	2.3
2	D	58	LYS	2.3
4	F	19	ARG	2.3

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Mol	Chain	Res	Type	RSRZ
2	B	38	GLY	2.2
4	F	221	LEU	2.2
2	D	56	GLY	2.2
2	B	278	SER	2.2
4	F	175	GLU	2.2
4	F	226	GLU	2.2
2	B	58	LYS	2.2
4	F	351	VAL	2.2
2	B	37	HIS	2.2
3	E	23	ILE	2.2
4	F	228	TYR	2.2
4	F	262	MET	2.2
4	F	227	PRO	2.2
2	B	338	SER	2.2
2	B	428	ALA	2.2
2	D	53	GLU	2.2
3	E	124	GLN	2.2
2	D	78	SER	2.2
4	F	196	HIS	2.1
2	D	199	THR	2.1
4	F	180	HIS	2.1
4	F	344	ALA	2.1
4	F	341	LYS	2.1
2	D	362	LYS	2.1
2	D	395	LEU	2.1
2	D	267	MET	2.0
2	D	177	ASP	2.0
2	D	44	LEU	2.0
1	C	285	GLN	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)

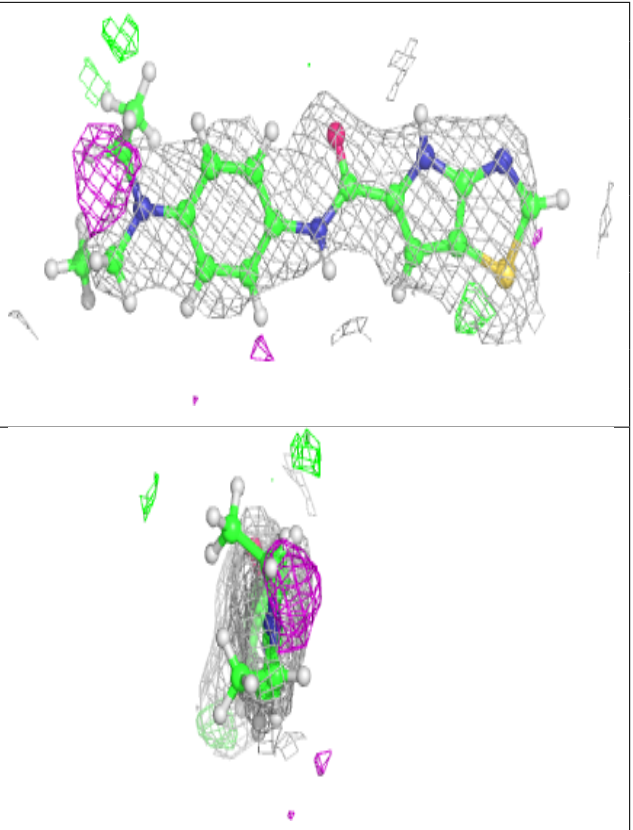
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' 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
12	NA	B	505	1/1	0.78	0.18	55,55,55,55	0
13	CL	D	504	1/1	0.79	0.26	43,43,43,43	0
6	MG	D	503	1/1	0.88	0.06	57,57,57,57	0
11	94U	D	502	22/22	0.89	0.28	39,50,63,73	0
14	ACP	F	401	31/31	0.91	0.20	48,59,85,98	0
11	94U	B	504	22/22	0.93	0.22	28,36,54,54	0
8	GOL	A	504	6/6	0.94	0.17	42,55,58,67	0
6	MG	A	502	1/1	0.94	0.22	21,21,21,21	0
6	MG	B	502	1/1	0.95	0.18	27,27,27,27	0
5	GTP	D	501	32/32	0.95	0.14	40,50,61,65	0
10	MES	B	503	12/12	0.96	0.16	28,35,47,48	0
7	CA	A	503	1/1	0.96	0.06	50,50,50,50	0
9	GDP	B	501	28/28	0.97	0.23	21,26,36,40	0
7	CA	C	503	1/1	0.98	0.07	37,37,37,37	0
5	GTP	C	501	32/32	0.98	0.17	14,22,28,33	0
5	GTP	A	501	32/32	0.98	0.22	18,27,34,40	0
6	MG	C	502	1/1	0.99	0.21	22,22,22,22	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.

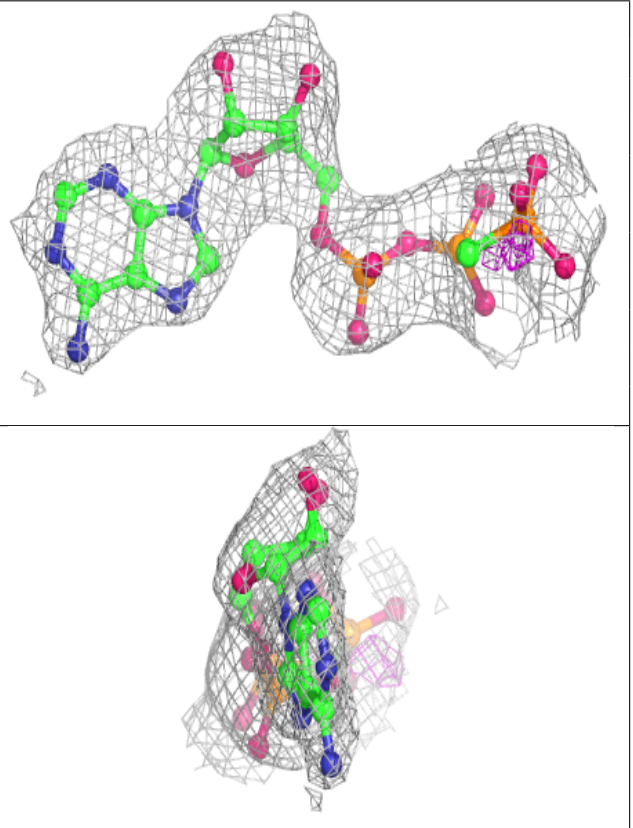
Electron density around 94U D 502:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



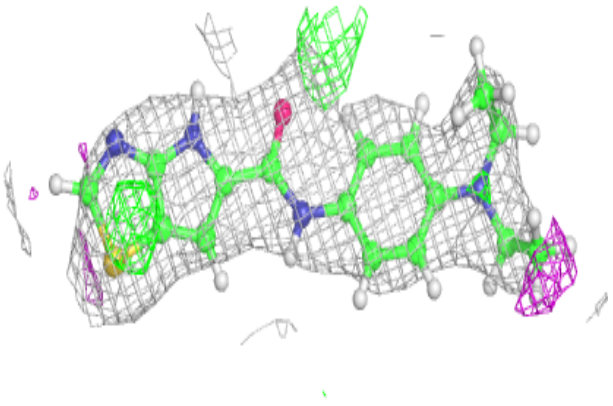
Electron density around ACP F 401:

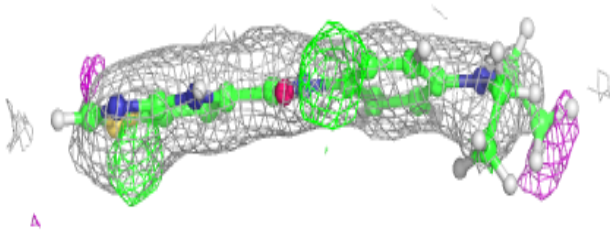
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

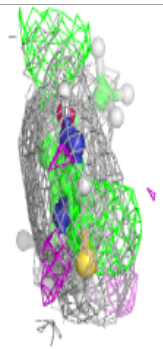


Electron density around 94U B 504:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

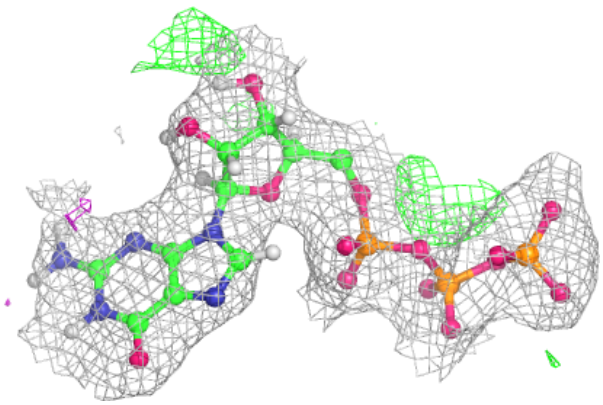


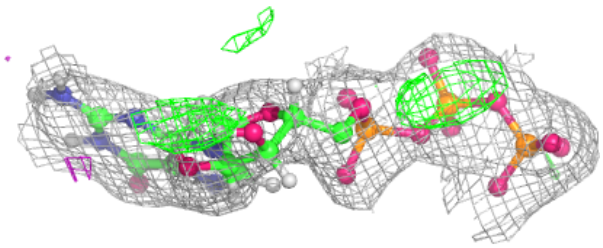


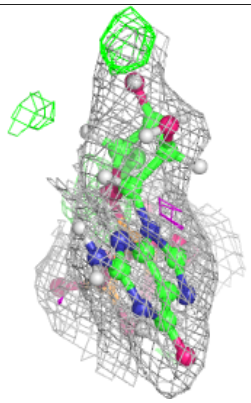


Electron density around GTP D 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

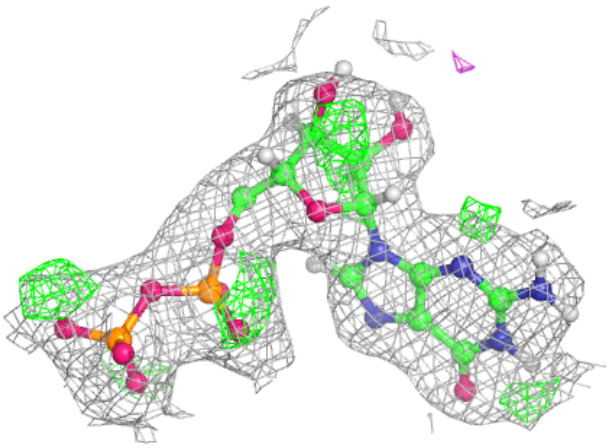


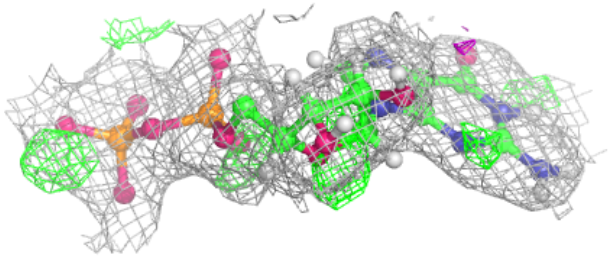


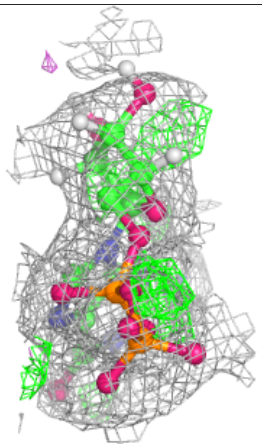


Electron density around GDP B 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

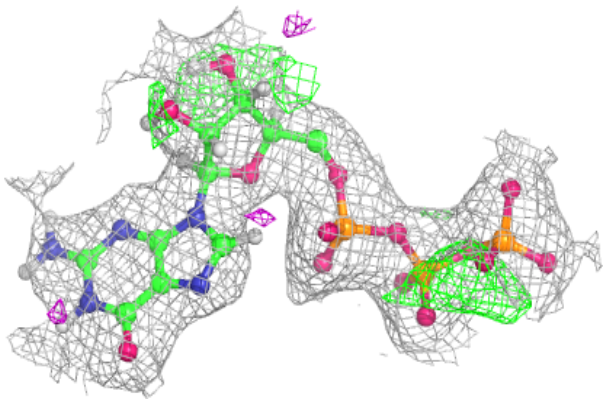


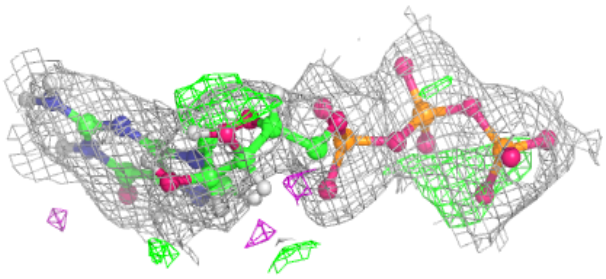


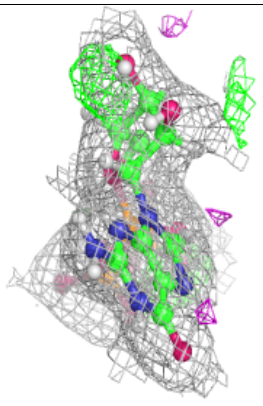


Electron density around GTP C 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

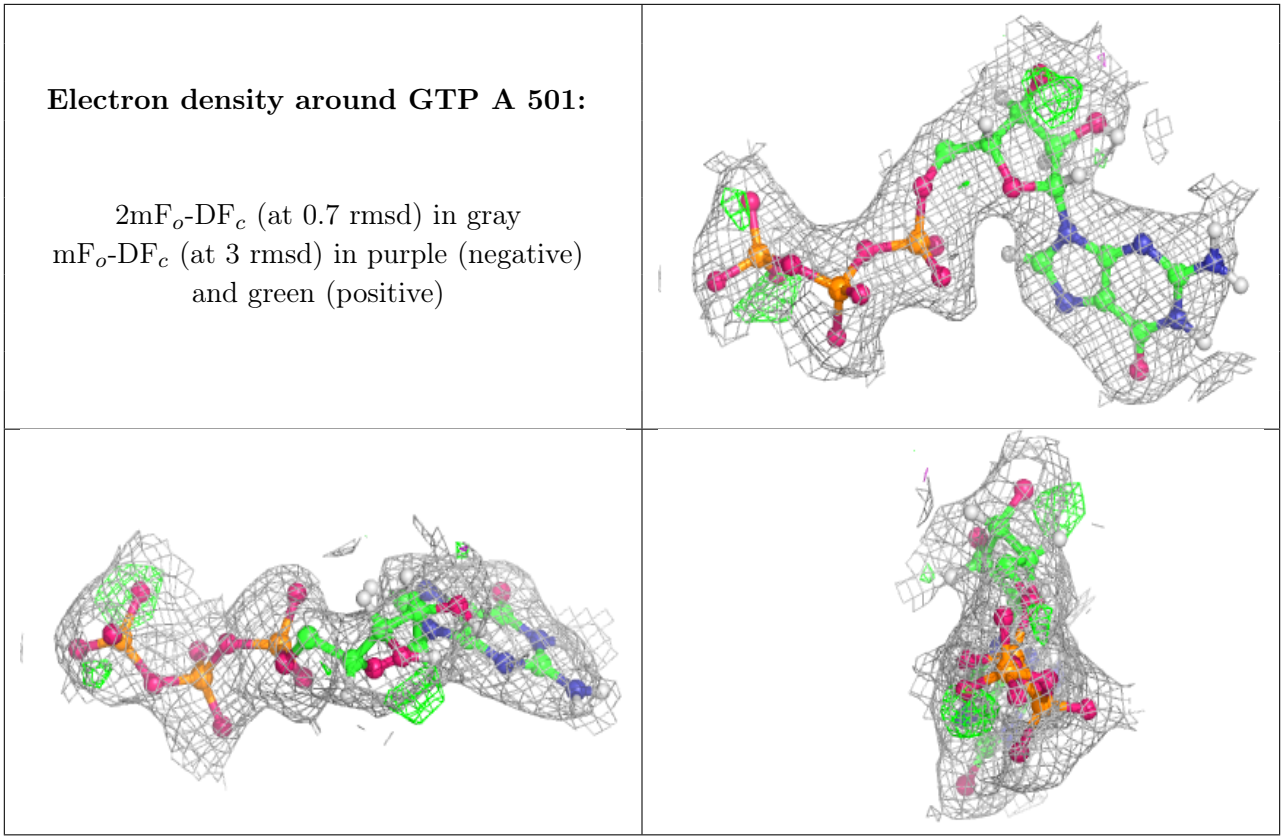






Electron density around GTP A 501:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.