



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 17, 2022 – 03:31 PM EDT

PDB ID : 3WTV
Title : Crystal structure of the complex comprised of ETS1(V170G), RUNX1, CBF-BETA, and the tcralpha gene enhancer DNA
Authors : Shiina, M.; Hamada, K.; Ogata, K.
Deposited on : 2014-04-21
Resolution : 2.70 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
Xtriage (Phenix) : 1.13
EDS : 2.29
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) : 2.29

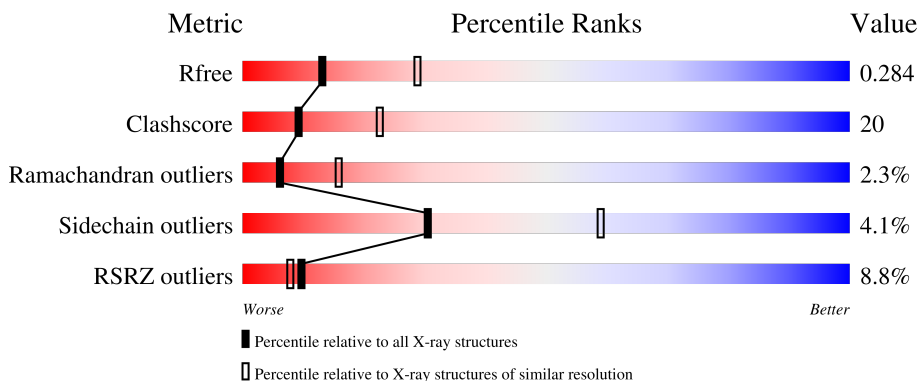
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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.



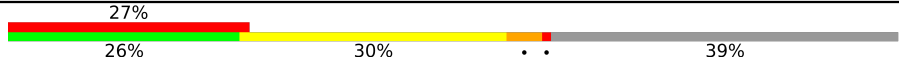
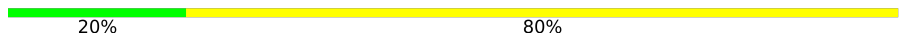
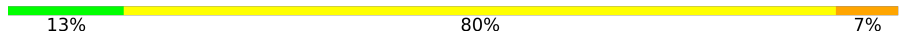


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	A	204	
1	F	204	
2	B	142	
2	G	142	
3	C	166	

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Mol	Chain	Length	Quality of chain
3	H	166	
4	D	15	
4	I	15	
5	E	15	
5	J	15	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7002 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 protein called Runt-related transcription factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	118	911	571	169	167	4	0	0	0
1	F	118	911	571	169	167	4	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	94	LYS	LEU	engineered mutation	UNP Q03347
A	170	GLY	VAL	engineered mutation	UNP Q03347
F	94	LYS	LEU	engineered mutation	UNP Q03347
F	170	GLY	VAL	engineered mutation	UNP Q03347

- Molecule 2 is a protein called Core-binding factor subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	130	1071	671	195	199	6	0	0	0
2	G	129	1062	666	193	197	6	0	0	0

- Molecule 3 is a protein called Protein C-ets-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	118	967	627	165	171	4	0	0	0
3	H	101	853	553	147	149	4	0	0	0

- Molecule 4 is a DNA chain called DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	15	Total	C	N	O	P	0	0	0
			299	144	54	87	14			
4	I	15	Total	C	N	O	P	0	0	0
			299	144	54	87	14			

- Molecule 5 is a DNA chain called DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	15	Total	C	N	O	P	0	0	0
			310	148	59	89	14			
5	J	15	Total	C	N	O	P	0	0	0
			310	148	59	89	14			

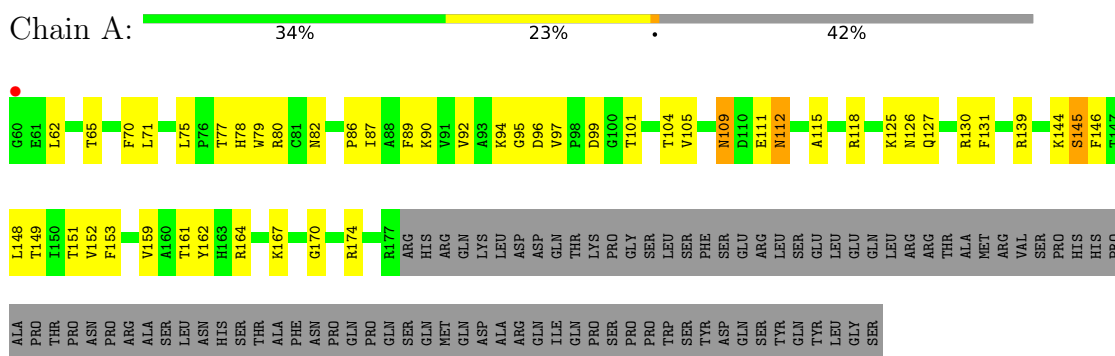
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	3	Total	O	0	0
			3	3		
6	F	4	Total	O	0	0
			4	4		
6	H	1	Total	O	0	0
			1	1		
6	J	1	Total	O	0	0
			1	1		

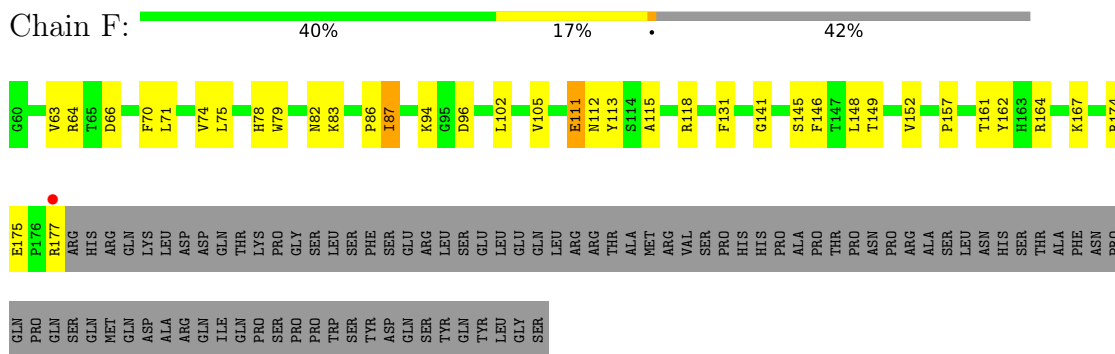
3 Residue-property plots [i](#)

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.

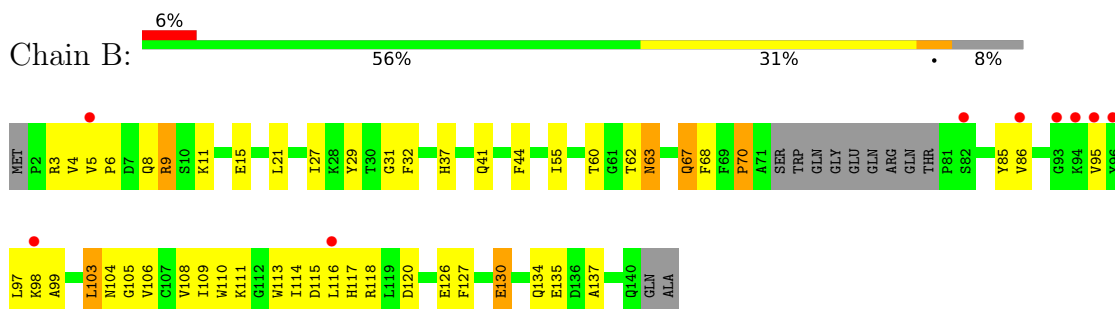
- Molecule 1: Runt-related transcription factor 1



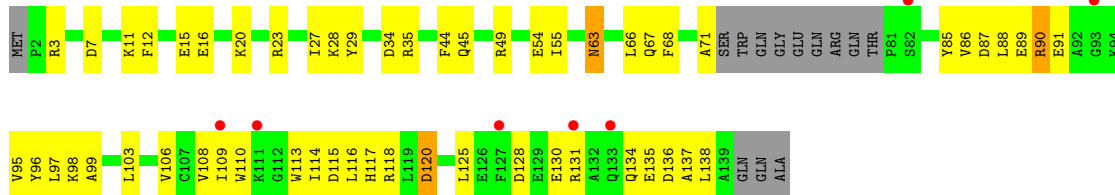
- Molecule 1: Runt-related transcription factor 1



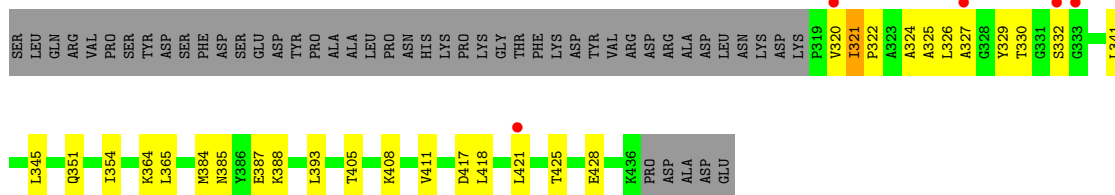
- Molecule 2: Core-binding factor subunit beta



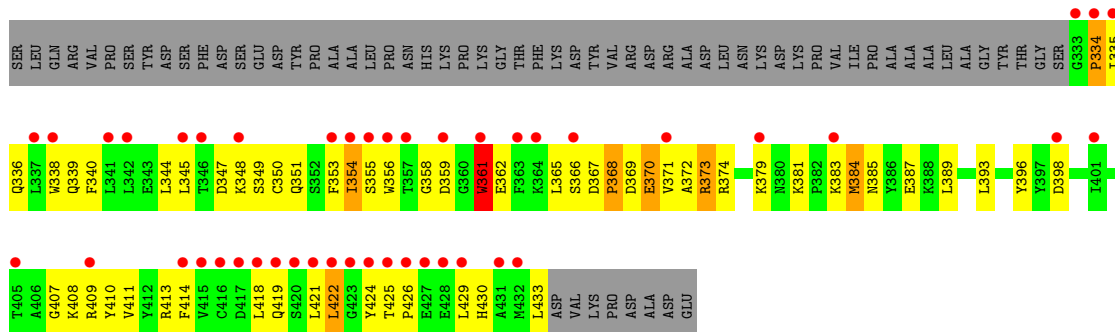
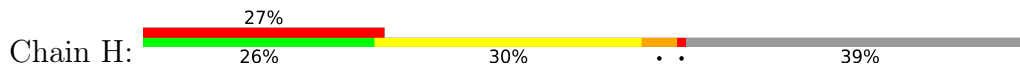
- Molecule 2: Core-binding factor subunit beta



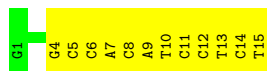
- Molecule 3: Protein C-ets-1



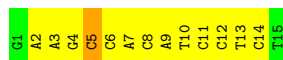
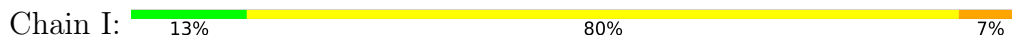
- Molecule 3: Protein C-ets-1




- Molecule 4: DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3')



- Molecule 4: DNA (5'-D(*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3')



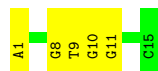
- Molecule 5: DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3')

Chain E:  80% 20%



- Molecule 5: DNA (5'-D(*AP*GP*AP*GP*GP*AP*TP*GP*TP*GP*GP*CP*TP*TP*C)-3')

Chain J:  67% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.22Å 102.25Å 194.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.26 – 2.70 45.26 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.4 (45.26-2.70) 96.5 (45.26-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.62 (at 2.69Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.242 , 0.284 0.242 , 0.284	Depositor DCC
R_{free} test set	4299 reflections (10.08%)	wwPDB-VP
Wilson B-factor (Å ²)	76.5	Xtrriage
Anisotropy	0.365	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 57.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7002	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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](#)

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	0.42	0/930	0.58	0/1263
1	F	0.48	0/930	0.65	0/1263
2	B	0.34	0/1093	0.49	0/1466
2	G	0.37	0/1084	0.50	0/1454
3	C	0.44	0/994	0.56	0/1342
3	H	0.32	0/877	0.45	0/1181
4	D	0.54	0/334	0.87	0/512
4	I	0.57	0/334	0.83	0/512
5	E	0.50	0/348	0.81	0/537
5	J	0.55	0/348	0.85	0/537
All	All	0.43	0/7272	0.62	0/10067

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
4	D	0	2
4	I	0	2
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	4	DG	Sidechain
4	D	5	DC	Sidechain
4	I	4	DG	Sidechain
4	I	5	DC	Sidechain

5.2 Too-close contacts [i](#)

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	911	0	916	41	0
1	F	911	0	916	26	0
2	B	1071	0	1034	45	0
2	G	1062	0	1026	47	0
3	C	967	0	966	23	0
3	H	853	0	847	57	0
4	D	299	0	170	13	0
4	I	299	0	170	20	0
5	E	310	0	171	3	0
5	J	310	0	171	6	0
6	C	3	0	0	0	0
6	F	4	0	0	0	0
6	H	1	0	0	0	0
6	J	1	0	0	0	0
All	All	7002	0	6387	266	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 20.

All (266) 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:I:2:DA:H2''	4:I:3:DA:H5'	1.31	1.10
4:I:9:DA:H2''	4:I:10:DT:H5'	1.36	1.03
4:I:11:DC:H2''	4:I:12:DC:H5'	1.39	1.02
3:H:424:TYR:HB2	3:H:429:LEU:HG	1.53	0.91
1:A:82:ASN:HD21	1:A:118:ARG:HH11	1.19	0.90
4:I:2:DA:H2''	4:I:3:DA:C5'	2.04	0.87
1:A:97:VAL:H	1:A:127:GLN:HE22	1.24	0.84
1:A:87:ILE:O	1:A:87:ILE:HD12	1.80	0.82
1:A:144:LYS:HE3	3:H:407:GLY:HA3	1.62	0.81
2:B:9:ARG:HB3	2:B:9:ARG:HH11	1.48	0.78
2:B:60:THR:HG23	2:B:62:THR:H	1.51	0.74
1:F:71:LEU:HD11	1:F:94:LYS:HE3	1.69	0.73
3:H:367:ASP:HB3	3:H:370:GLU:HB2	1.68	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:322:PRO:HB2	3:C:325:ALA:HB3	1.70	0.73
4:D:10:DT:H2''	4:D:11:DC:H5'	1.70	0.73
4:I:10:DT:H2''	4:I:11:DC:O5'	1.89	0.72
3:H:372:ALA:HB1	3:H:385:ASN:HA	1.73	0.71
3:H:365:LEU:HD23	3:H:410:TYR:HD2	1.56	0.70
1:A:139:ARG:HG3	1:A:139:ARG:HH11	1.56	0.69
3:H:335:ILE:HD12	3:H:339:GLN:NE2	2.09	0.68
1:F:63:VAL:HB	1:F:74:VAL:HG22	1.75	0.67
2:G:63:ASN:N	2:G:63:ASN:HD22	1.91	0.67
2:B:115:ASP:OD1	2:B:118:ARG:HD3	1.95	0.67
1:F:148:LEU:HB2	1:F:162:TYR:HB3	1.75	0.67
2:B:67:GLN:HE21	2:B:67:GLN:HA	1.61	0.66
2:G:55:ILE:HD13	2:G:66:LEU:HD12	1.77	0.66
1:F:63:VAL:HG22	1:F:64:ARG:H	1.60	0.66
3:H:345:LEU:HB3	3:H:356:TRP:HE1	1.61	0.66
4:I:13:DT:H6	4:I:13:DT:H5'	1.60	0.66
1:A:104:THR:HG22	1:A:151:THR:HB	1.78	0.65
3:H:409:ARG:O	3:H:411:VAL:HG23	1.96	0.65
1:A:109:ASN:ND2	1:A:112:ASN:H	1.94	0.65
1:A:82:ASN:ND2	1:A:118:ARG:HH11	1.92	0.64
2:G:90:ARG:O	2:G:90:ARG:HD3	1.98	0.64
2:G:131:ARG:NH1	2:G:131:ARG:HB3	2.13	0.64
4:I:9:DA:C2'	4:I:10:DT:H5'	2.20	0.63
2:B:106:VAL:O	2:B:108:VAL:HG13	1.99	0.63
1:A:70:PHE:CE2	1:A:152:VAL:HG21	2.34	0.63
3:C:384:MET:SD	3:C:388:LYS:HB2	2.39	0.63
5:J:9:DT:H2'	5:J:10:DG:C8	2.34	0.63
3:C:326:LEU:HB3	3:C:421:LEU:O	2.00	0.62
1:A:109:ASN:C	1:A:109:ASN:HD22	2.02	0.62
4:I:11:DC:H2''	4:I:12:DC:C5'	2.23	0.62
2:G:90:ARG:HD2	2:G:96:TYR:CE2	2.35	0.62
2:B:41:GLN:HE21	2:B:117:HIS:HA	1.65	0.62
4:D:13:DT:H2''	4:D:14:DC:H5'	1.81	0.61
4:I:6:DC:H6	4:I:6:DC:H5'	1.64	0.60
2:B:63:ASN:N	2:B:63:ASN:HD22	1.98	0.60
1:F:149:THR:HG21	2:G:63:ASN:O	2.01	0.60
4:I:13:DT:H5'	4:I:13:DT:C6	2.36	0.60
2:B:113:TRP:C	2:B:114:ILE:HD12	2.22	0.60
1:A:145:SER:HB3	1:A:164:ARG:HA	1.84	0.59
1:F:63:VAL:HG22	1:F:64:ARG:N	2.17	0.59
1:F:157:PRO:HG2	2:G:67:GLN:HG3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:345:LEU:HB3	3:H:356:TRP:NE1	2.17	0.59
4:D:11:DC:H1'	4:D:12:DC:H5''	1.85	0.59
1:A:161:THR:H	2:B:104:ASN:HD21	1.51	0.59
1:A:139:ARG:HG3	1:A:139:ARG:NH1	2.18	0.58
3:H:419:GLN:HE21	3:H:426:PRO:HD3	1.68	0.58
1:F:174:ARG:NH2	5:J:10:DG:N7	2.51	0.58
3:H:335:ILE:HD12	3:H:339:GLN:HE21	1.67	0.58
4:D:11:DC:H2''	4:D:12:DC:H5'	1.86	0.58
1:A:71:LEU:HD11	1:A:94:LYS:HE3	1.85	0.58
4:I:8:DC:H4'	4:I:8:DC:OP1	2.03	0.57
1:A:149:THR:HG21	2:B:63:ASN:O	2.04	0.57
4:D:11:DC:H1'	4:D:12:DC:C5'	2.35	0.57
2:G:108:VAL:HG12	2:G:109:ILE:N	2.20	0.57
1:A:148:LEU:HB2	1:A:162:TYR:HB3	1.87	0.57
3:H:336:GLN:H	3:H:339:GLN:NE2	2.03	0.57
3:H:340:PHE:HZ	3:H:374:ARG:HB3	1.70	0.57
2:G:68:PHE:CZ	2:G:97:LEU:HD13	2.40	0.56
3:H:385:ASN:OD1	3:H:387:GLU:HB2	2.06	0.56
2:G:16:GLU:OE2	2:G:20:LYS:HB2	2.06	0.56
1:A:87:ILE:HD12	1:A:87:ILE:C	2.27	0.55
3:C:417:ASP:O	3:C:421:LEU:HD13	2.07	0.55
2:G:88:LEU:HD23	2:G:95:VAL:HG22	1.87	0.55
2:G:108:VAL:HG11	2:G:125:LEU:HB3	1.87	0.55
3:H:354:ILE:HD12	3:H:371:VAL:HG11	1.87	0.55
3:H:368:PRO:HG2	3:H:369:ASP:H	1.72	0.55
3:H:348:LYS:HE2	3:H:351:GLN:HE22	1.72	0.55
1:F:113:TYR:CE2	2:G:28:LYS:HD2	2.42	0.54
2:B:29:TYR:HA	2:B:55:ILE:HD13	1.89	0.54
1:A:90:LYS:HE3	1:A:130:ARG:HG2	1.90	0.54
1:A:159:VAL:HG13	2:B:103:LEU:HA	1.89	0.54
5:J:8:DG:H2''	5:J:9:DT:O5'	2.08	0.54
3:H:348:LYS:HA	3:H:351:GLN:HE21	1.73	0.53
3:H:353:PHE:C	3:H:366:SER:HB2	2.29	0.53
3:H:354:ILE:HG12	3:H:354:ILE:O	2.07	0.53
2:G:109:ILE:HG13	2:G:128:ASP:HB2	1.91	0.53
3:H:419:GLN:NE2	3:H:425:THR:HA	2.23	0.53
3:H:334:PRO:HG2	3:H:335:ILE:H	1.74	0.53
2:B:98:LYS:HE3	2:B:109:ILE:HG21	1.91	0.53
3:C:321:ILE:HD13	3:C:321:ILE:N	2.24	0.53
2:G:63:ASN:N	2:G:63:ASN:ND2	2.58	0.53
3:H:393:LEU:HA	3:H:396:TYR:HD2	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:2:DA:C2'	4:I:3:DA:C5'	2.85	0.53
2:B:3:ARG:HH21	2:B:135:GLU:CD	2.12	0.52
3:C:425:THR:OG1	3:C:428:GLU:HG3	2.09	0.52
1:F:115:ALA:HB2	1:F:146:PHE:CZ	2.45	0.52
2:B:114:ILE:HG23	2:B:120:ASP:N	2.25	0.52
3:H:365:LEU:HD21	3:H:368:PRO:HA	1.90	0.52
1:F:177:ARG:NH1	5:J:11:DG:N7	2.57	0.52
2:G:137:ALA:O	2:G:138:LEU:HD23	2.10	0.52
3:C:321:ILE:HD13	3:C:321:ILE:H	1.73	0.52
2:G:55:ILE:HG12	2:G:114:ILE:HD11	1.92	0.52
2:B:5:VAL:HG12	2:B:105:GLY:O	2.09	0.52
4:D:14:DC:H2''	4:D:15:DT:H5'	1.92	0.52
2:G:45:GLN:NE2	2:G:49:ARG:HH21	2.07	0.51
3:C:345:LEU:HD21	3:C:354:ILE:HG12	1.93	0.51
1:F:70:PHE:CE2	1:F:152:VAL:HG21	2.45	0.51
1:A:70:PHE:O	1:A:71:LEU:HD12	2.11	0.51
4:D:9:DA:H1'	4:D:10:DT:H5''	1.92	0.51
2:G:63:ASN:HD22	2:G:63:ASN:H	1.56	0.51
2:G:87:ASP:OD2	2:G:90:ARG:HB2	2.09	0.51
3:H:379:LYS:HG3	4:I:9:DA:OP1	2.11	0.51
2:G:3:ARG:HD3	2:G:135:GLU:OE1	2.10	0.51
3:H:419:GLN:HE22	3:H:425:THR:HG22	1.75	0.51
3:H:421:LEU:O	3:H:422:LEU:HG	2.11	0.51
3:H:347:ASP:OD2	3:H:349:SER:HB3	2.11	0.50
3:H:344:LEU:HD23	3:H:371:VAL:HG12	1.93	0.50
3:C:320:VAL:HG12	3:C:321:ILE:HG23	1.93	0.50
4:I:13:DT:H1'	4:I:14:DC:H5'	1.93	0.50
1:A:109:ASN:HD22	1:A:112:ASN:H	1.59	0.50
4:D:6:DC:H6	4:D:6:DC:H5'	1.76	0.50
4:D:8:DC:H2''	4:D:9:DA:C8	2.47	0.50
1:A:109:ASN:ND2	1:A:109:ASN:C	2.66	0.49
3:H:344:LEU:CD2	3:H:371:VAL:HG12	2.43	0.49
2:B:70:PRO:HD3	2:B:85:TYR:HE2	1.77	0.49
2:G:89:GLU:C	2:G:91:GLU:H	2.15	0.49
4:D:13:DT:H1'	4:D:14:DC:H5''	1.95	0.49
1:A:111:GLU:OE1	1:A:144:LYS:HD2	2.13	0.49
3:C:385:ASN:OD1	3:C:388:LYS:HG2	2.13	0.49
3:H:389:LEU:O	3:H:389:LEU:HD23	2.13	0.49
1:F:161:THR:HG22	1:F:162:TYR:N	2.27	0.48
3:H:354:ILE:HD12	3:H:371:VAL:CG1	2.43	0.48
2:B:11:LYS:HE3	2:B:15:GLU:OE1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ASN:HD21	1:A:118:ARG:NH1	2.00	0.48
2:B:32:PHE:N	2:B:32:PHE:CD2	2.80	0.48
3:H:336:GLN:H	3:H:339:GLN:HE21	1.61	0.48
1:F:102:LEU:N	1:F:102:LEU:HD22	2.29	0.48
3:H:365:LEU:HB3	3:H:410:TYR:HB3	1.96	0.48
2:G:108:VAL:HG12	2:G:109:ILE:H	1.78	0.47
1:A:112:ASN:C	1:A:112:ASN:HD22	2.18	0.47
2:B:111:LYS:HD2	2:B:126:GLU:OE2	2.14	0.47
2:B:8:GLN:O	2:B:127:PHE:HE2	1.98	0.47
1:F:87:ILE:HD13	1:F:87:ILE:H	1.79	0.47
2:G:90:ARG:HG3	2:G:90:ARG:HH11	1.80	0.47
2:B:70:PRO:HD3	2:B:85:TYR:CE2	2.50	0.47
2:B:67:GLN:O	2:B:99:ALA:HB1	2.14	0.47
4:D:9:DA:H2''	4:D:10:DT:H5'	1.97	0.47
2:G:68:PHE:CE1	2:G:97:LEU:HB3	2.50	0.47
2:B:29:TYR:CE1	2:B:44:PHE:HB2	2.51	0.46
1:F:105:VAL:HG23	1:F:131:PHE:CZ	2.50	0.46
2:G:95:VAL:O	2:G:113:TRP:HA	2.16	0.46
3:H:361:TRP:CE3	3:H:361:TRP:HA	2.49	0.46
4:I:6:DC:H5'	4:I:6:DC:C6	2.47	0.46
4:D:7:DA:H2''	4:D:8:DC:O5'	2.15	0.46
4:D:10:DT:C2'	4:D:11:DC:H5'	2.43	0.46
3:H:361:TRP:NE1	3:H:426:PRO:HG3	2.30	0.46
2:G:86:VAL:O	2:G:86:VAL:HG13	2.15	0.46
2:B:98:LYS:CE	2:B:111:LYS:HE2	2.46	0.46
2:B:134:GLN:O	2:B:137:ALA:HB3	2.16	0.46
3:C:321:ILE:HB	3:C:322:PRO:HD2	1.97	0.46
2:G:71:ALA:HB1	2:G:131:ARG:HE	1.80	0.46
3:H:348:LYS:HA	3:H:351:GLN:HB2	1.97	0.46
2:G:27:ILE:HB	2:G:55:ILE:CG2	2.45	0.46
2:B:11:LYS:HD2	2:B:106:VAL:HG21	1.98	0.46
3:C:364:LYS:CB	3:C:411:VAL:HG22	2.45	0.46
1:A:97:VAL:H	1:A:127:GLN:NE2	2.03	0.46
3:C:385:ASN:HD21	3:C:387:GLU:HB2	1.81	0.46
5:E:15:DC:H2'	5:J:1:DA:O4'	2.16	0.46
1:A:115:ALA:HB2	1:A:146:PHE:CZ	2.51	0.46
2:B:95:VAL:O	2:B:113:TRP:HA	2.16	0.46
3:C:324:ALA:O	3:C:329:TYR:HB3	2.16	0.46
1:A:71:LEU:CD1	1:A:94:LYS:HE3	2.46	0.46
2:G:85:TYR:HA	2:G:98:LYS:HB3	1.98	0.46
3:H:362:GLU:HB2	3:H:413:ARG:HD2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:424:TYR:CD2	3:H:429:LEU:HD21	2.52	0.45
1:A:65:THR:HG23	1:A:71:LEU:HA	1.98	0.45
1:A:78:HIS:CD2	1:A:167:LYS:HE3	2.51	0.45
2:B:4:VAL:HA	2:B:105:GLY:HA2	1.98	0.45
3:H:353:PHE:HB3	3:H:370:GLU:HG2	1.98	0.45
5:J:9:DT:H4'	5:J:9:DT:OP1	2.16	0.45
4:I:11:DC:C2'	4:I:12:DC:H5'	2.28	0.45
2:G:29:TYR:CD2	2:G:44:PHE:HD2	2.34	0.45
2:B:21:LEU:HD21	2:B:60:THR:HG21	1.97	0.45
2:G:45:GLN:HG2	2:G:49:ARG:HE	1.82	0.45
2:B:114:ILE:HA	2:B:120:ASP:O	2.16	0.45
1:F:66:ASP:OD2	1:F:161:THR:HB	2.17	0.45
2:B:67:GLN:HE21	2:B:67:GLN:CA	2.26	0.44
2:B:98:LYS:HA	2:B:110:TRP:O	2.18	0.44
2:B:130:GLU:O	2:B:134:GLN:HG3	2.17	0.44
1:F:145:SER:OG	1:F:164:ARG:HA	2.18	0.44
2:G:68:PHE:CD1	2:G:99:ALA:HB2	2.52	0.44
3:H:418:LEU:O	3:H:422:LEU:HB2	2.17	0.44
3:H:338:TRP:NE1	3:H:339:GLN:HG3	2.32	0.44
3:H:348:LYS:HA	3:H:351:GLN:NE2	2.32	0.44
3:H:430:HIS:O	3:H:433:LEU:HG	2.18	0.44
2:B:5:VAL:CG2	2:B:6:PRO:HD2	2.47	0.44
1:F:78:HIS:CD2	1:F:167:LYS:HG2	2.52	0.44
4:I:7:DA:H2''	4:I:8:DC:O5'	2.17	0.44
5:E:9:DT:H2''	5:E:10:DG:C8	2.53	0.44
2:G:71:ALA:CB	2:G:131:ARG:HE	2.30	0.44
2:G:106:VAL:O	2:G:108:VAL:HG23	2.18	0.43
3:H:419:GLN:HE22	3:H:425:THR:HA	1.83	0.43
3:H:422:LEU:HD13	3:H:429:LEU:HD11	2.01	0.43
3:C:327:ALA:HA	3:C:421:LEU:HD23	1.99	0.43
3:H:373:ARG:NH1	3:H:373:ARG:HB2	2.32	0.43
2:G:11:LYS:HE3	2:G:15:GLU:OE2	2.18	0.43
3:H:350:CYS:O	3:H:354:ILE:HG22	2.19	0.43
3:H:381:LYS:O	3:H:384:MET:HB2	2.18	0.43
1:A:101:THR:HG23	1:A:153:PHE:O	2.18	0.43
1:F:157:PRO:CG	2:G:67:GLN:HG3	2.47	0.43
1:A:75:LEU:HD21	1:A:89:PHE:CD2	2.54	0.43
2:G:12:PHE:HB2	2:G:106:VAL:HG11	2.01	0.43
2:G:134:GLN:O	2:G:138:LEU:HG	2.19	0.43
1:A:144:LYS:CE	3:H:407:GLY:HA3	2.43	0.43
2:B:98:LYS:HE3	2:B:111:LYS:HE2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:44:PHE:CD1	2:G:44:PHE:C	2.92	0.43
3:H:393:LEU:HA	3:H:396:TYR:CD2	2.53	0.43
1:A:144:LYS:NZ	3:H:408:LYS:HE3	2.34	0.43
3:H:362:GLU:HB2	3:H:413:ARG:HH11	1.84	0.43
2:B:9:ARG:HA	2:B:127:PHE:CE2	2.54	0.42
2:G:54:GLU:C	2:G:55:ILE:HD12	2.40	0.42
2:G:131:ARG:HB3	2:G:131:ARG:HH11	1.82	0.42
1:A:126:ASN:O	1:A:127:GLN:HB2	2.20	0.42
2:B:86:VAL:HG13	2:B:86:VAL:O	2.19	0.42
1:F:79:TRP:CZ2	1:F:86:PRO:HD3	2.54	0.42
3:H:345:LEU:HD22	3:H:356:TRP:NE1	2.35	0.42
1:F:82:ASN:OD1	1:F:118:ARG:NH1	2.45	0.42
4:I:5:DC:H2''	4:I:6:DC:H5'	2.01	0.42
1:F:111:GLU:OE2	1:F:141:GLY:HA3	2.19	0.42
3:C:405:THR:HG21	3:C:408:LYS:HD2	2.01	0.42
2:G:115:ASP:O	2:G:117:HIS:N	2.53	0.42
1:A:62:LEU:HD21	1:A:92:VAL:HG21	2.01	0.42
3:C:364:LYS:HB2	3:C:411:VAL:HG22	2.01	0.42
4:I:8:DC:H2''	4:I:9:DA:C8	2.55	0.42
3:C:330:THR:C	3:C:332:SER:H	2.23	0.42
2:G:118:ARG:HB3	2:G:120:ASP:OD2	2.20	0.42
1:A:80:ARG:HG3	1:A:170:GLY:O	2.20	0.41
2:B:29:TYR:CZ	2:B:31:GLY:HA3	2.55	0.41
3:H:383:LYS:O	3:H:383:LYS:HG3	2.19	0.41
1:A:79:TRP:CZ2	1:A:86:PRO:HD3	2.56	0.41
1:A:174:ARG:NH2	5:E:10:DG:N7	2.65	0.41
2:B:68:PHE:HA	2:B:99:ALA:HB2	2.02	0.41
2:B:130:GLU:H	2:B:130:GLU:HG2	1.57	0.41
4:I:12:DC:H2''	4:I:13:DT:OP2	2.21	0.41
3:C:418:LEU:HD23	3:C:418:LEU:HA	1.82	0.41
1:F:87:ILE:HD13	1:F:87:ILE:N	2.36	0.41
1:A:105:VAL:HG23	1:A:131:PHE:CZ	2.56	0.41
2:B:27:ILE:HD12	2:B:114:ILE:HD11	2.02	0.41
3:H:354:ILE:O	3:H:354:ILE:HG23	2.20	0.41
2:B:68:PHE:CZ	2:B:97:LEU:HD13	2.56	0.41
1:F:175:GLU:OE1	1:F:175:GLU:HA	2.20	0.41
3:C:354:ILE:HD12	3:C:365:LEU:HD22	2.01	0.40
1:F:75:LEU:HD11	1:F:148:LEU:HD11	2.03	0.40
2:G:34:ASP:O	2:G:35:ARG:HD2	2.20	0.40
3:H:361:TRP:HB3	3:H:414:PHE:HB2	2.03	0.40
2:B:115:ASP:O	2:B:117:HIS:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:322:PRO:HG2	3:C:326:LEU:HG	2.03	0.40
2:G:98:LYS:HA	2:G:110:TRP:O	2.20	0.40
2:B:37:HIS:O	2:B:41:GLN:HG3	2.21	0.40
2:G:136:ASP:C	2:G:138:LEU:H	2.25	0.40
3:C:341:LEU:HD11	3:C:393:LEU:HD21	2.04	0.40
3:C:385:ASN:ND2	3:C:387:GLU:HB2	2.36	0.40
1:A:99:ASP:OD1	1:A:125:LYS:HD2	2.22	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	
1	A	116/204 (57%)	109 (94%)	6 (5%)	1 (1%)	17	40
1	F	116/204 (57%)	109 (94%)	6 (5%)	1 (1%)	17	40
2	B	126/142 (89%)	98 (78%)	26 (21%)	2 (2%)	9	24
2	G	125/142 (88%)	103 (82%)	19 (15%)	3 (2%)	6	15
3	C	116/166 (70%)	112 (97%)	4 (3%)	0	100	100
3	H	99/166 (60%)	68 (69%)	22 (22%)	9 (9%)	1	0
All	All	698/1024 (68%)	599 (86%)	83 (12%)	16 (2%)	6	16

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	G	23	ARG
3	H	354	ILE
3	H	359	ASP
3	H	361	TRP
2	B	70	PRO

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Mol	Chain	Res	Type
1	F	111	GLU
3	H	370	GLU
2	B	116	LEU
2	G	116	LEU
3	H	355	SER
3	H	368	PRO
3	H	422	LEU
1	A	95	GLY
2	G	130	GLU
3	H	358	GLY
3	H	334	PRO

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	99/178 (56%)	94 (95%)	5 (5%)	24	50
1	F	99/178 (56%)	95 (96%)	4 (4%)	31	60
2	B	113/123 (92%)	108 (96%)	5 (4%)	28	56
2	G	112/123 (91%)	107 (96%)	5 (4%)	27	55
3	C	102/145 (70%)	100 (98%)	2 (2%)	55	81
3	H	91/145 (63%)	87 (96%)	4 (4%)	28	56
All	All	616/892 (69%)	591 (96%)	25 (4%)	30	59

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	77	THR
1	A	96	ASP
1	A	109	ASN
1	A	112	ASN
1	A	145	SER
2	B	9	ARG
2	B	63	ASN

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Mol	Chain	Res	Type
2	B	67	GLN
2	B	103	LEU
2	B	130	GLU
3	C	321	ILE
3	C	351	GLN
1	F	83	LYS
1	F	87	ILE
1	F	96	ASP
1	F	112	ASN
2	G	7	ASP
2	G	63	ASN
2	G	90	ARG
2	G	103	LEU
2	G	120	ASP
3	H	361	TRP
3	H	373	ARG
3	H	384	MET
3	H	398	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	ASN
1	A	109	ASN
1	A	112	ASN
1	A	119	ASN
1	A	127	GLN
1	A	132	ASN
1	A	155	ASN
2	B	8	GLN
2	B	41	GLN
2	B	63	ASN
2	B	67	GLN
2	B	104	ASN
2	B	134	GLN
3	C	380	ASN
1	F	112	ASN
1	F	132	ASN
1	F	163	HIS
2	G	8	GLN
2	G	46	ASN
2	G	63	ASN

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Mol	Chain	Res	Type
2	G	67	GLN
3	H	339	GLN
3	H	351	GLN
3	H	380	ASN
3	H	419	GLN

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](#)

There are no ligands in this entry.

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 [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	A	118/204 (57%)	0.32	1 (0%) 86 87	50, 72, 95, 110	0
1	F	118/204 (57%)	0.25	1 (0%) 86 87	47, 62, 82, 103	0
2	B	130/142 (91%)	0.80	9 (6%) 16 15	77, 101, 128, 135	0
2	G	129/142 (90%)	0.67	7 (5%) 25 24	63, 88, 115, 120	0
3	C	118/166 (71%)	0.41	5 (4%) 36 35	46, 69, 100, 111	0
3	H	101/166 (60%)	1.95	45 (44%) 0 0	71, 131, 173, 179	0
4	D	15/15 (100%)	-0.17	0 100 100	57, 60, 76, 78	0
4	I	15/15 (100%)	-0.14	0 100 100	57, 69, 84, 85	0
5	E	15/15 (100%)	-0.22	0 100 100	50, 61, 73, 77	0
5	J	15/15 (100%)	-0.05	0 100 100	54, 66, 79, 87	0
All	All	774/1084 (71%)	0.64	68 (8%) 10 8	46, 80, 143, 179	0

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	426	PRO	6.2
3	H	421	LEU	5.8
3	H	346	THR	5.5
3	H	423	GLY	5.5
3	H	416	CYS	4.9
3	H	417	ASP	4.7
3	H	357	THR	4.7
3	H	424	TYR	4.7
3	H	431	ALA	4.7
3	H	335	ILE	4.6
3	H	361	TRP	4.3
3	C	332	SER	4.2
3	H	425	THR	4.1

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Mol	Chain	Res	Type	RSRZ
2	B	96	TYR	4.1
3	H	422	LEU	3.7
3	H	345	LEU	3.7
3	H	364	LYS	3.7
3	C	327	ALA	3.7
3	H	334	PRO	3.3
3	H	353	PHE	3.3
3	H	418	LEU	3.2
3	H	419	GLN	3.2
3	H	427	GLU	3.1
2	G	133	GLN	3.1
1	A	60	GLY	3.0
3	H	354	ILE	3.0
3	H	356	TRP	2.9
3	H	420	SER	2.9
3	H	342	LEU	2.9
3	H	333	GLY	2.8
3	H	379	LYS	2.8
3	H	371	VAL	2.7
3	H	429	LEU	2.7
3	H	363	PHE	2.7
2	B	82	SER	2.7
2	B	93	GLY	2.7
2	B	94	LYS	2.6
2	G	82	SER	2.6
3	H	414	PHE	2.6
3	H	355	SER	2.5
3	H	409	ARG	2.5
2	G	109	ILE	2.5
3	H	432	MET	2.5
3	H	359	ASP	2.5
2	B	98	LYS	2.5
3	H	383	LYS	2.4
3	H	348	LYS	2.4
3	H	415	VAL	2.4
3	H	341	LEU	2.3
1	F	177	ARG	2.3
2	B	86	VAL	2.3
3	C	421	LEU	2.3
2	G	127	PHE	2.2
3	H	428	GLU	2.2
3	C	333	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
2	G	93	GLY	2.2
2	B	116	LEU	2.2
3	C	320	VAL	2.1
3	H	338	TRP	2.1
3	H	401	ILE	2.1
2	B	5	VAL	2.1
2	B	95	VAL	2.1
3	H	366	SER	2.1
2	G	131	ARG	2.0
3	H	398	ASP	2.0
2	G	111	LYS	2.0
3	H	405	THR	2.0
3	H	337	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.