



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:56 PM GMT

PDB ID : 3AXL
Title : Murine Valpha 10 Vbeta 8.1 T-cell receptor
Authors : Patel, O.; Rossjohn, J.
Deposited on : 2011-04-11
Resolution : 2.90 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

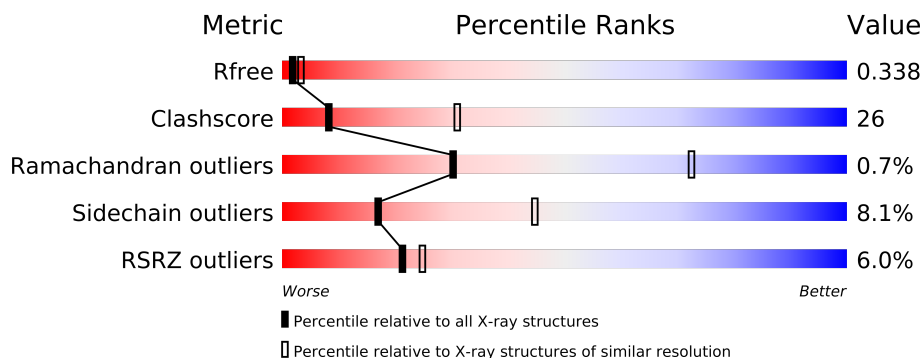
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.90 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}	66092	1053 (2.90-2.90)
Clashscore	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (2.90-2.90)
RSRZ outliers	66119	1054 (2.90-2.90)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	198	
1	G	198	
2	B	238	
2	H	238	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6728 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Valpha 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1393	866	232	287	8			
1	G	190	Total	C	N	O	S	0	0	0
			1437	890	240	299	8			

- Molecule 2 is a protein called Vbeta 8.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1887	1193	330	358	6			
2	H	238	Total	C	N	O	S	0	0	0
			1883	1189	330	358	6			

- Molecule 3 is water.

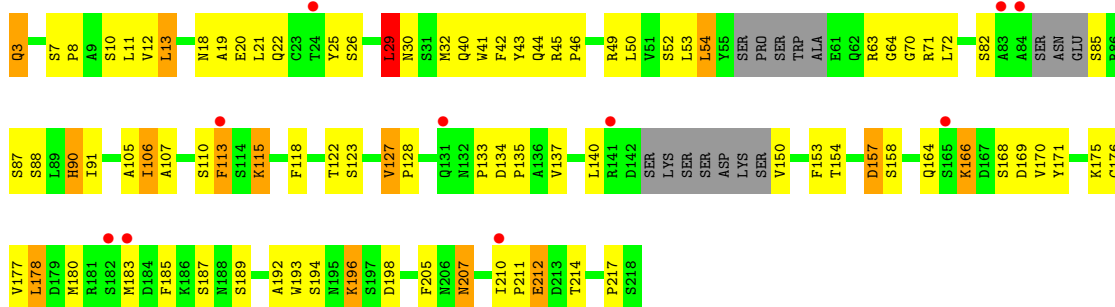
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	28	Total	O	0	0
			28	28		
3	B	43	Total	O	0	0
			43	43		
3	G	27	Total	O	0	0
			27	27		
3	H	30	Total	O	0	0
			30	30		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

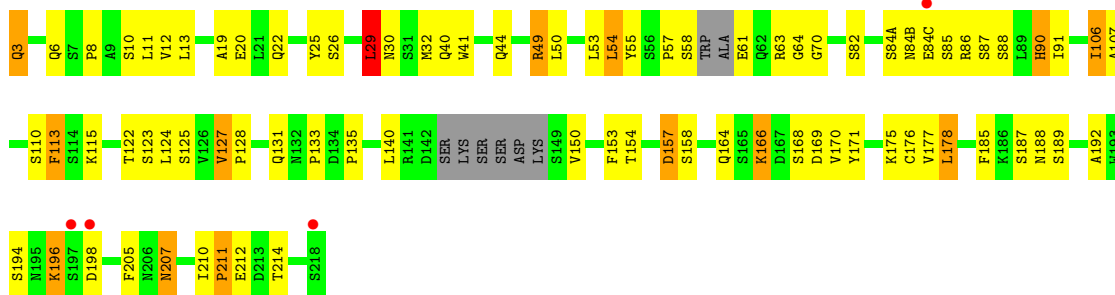
• Molecule 1: Valpha 10

Chain A: 



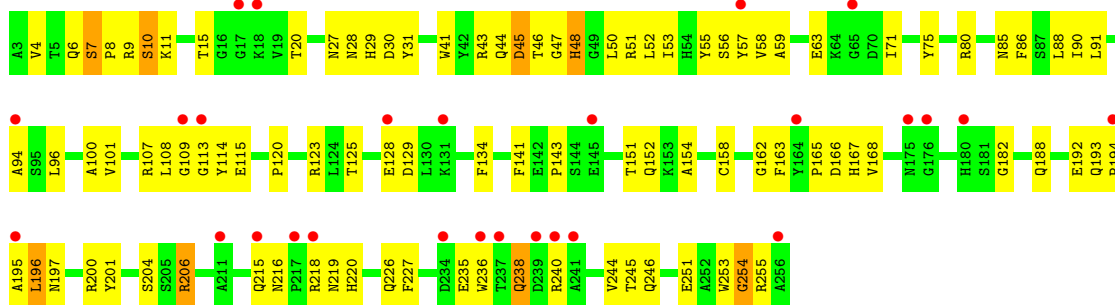
• Molecule 1: Valpha 10

Chain G: 

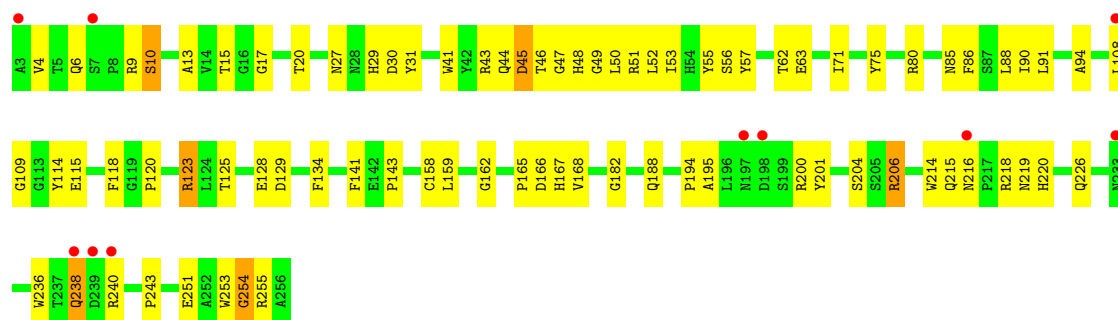


• Molecule 2: Vbeta 8.1

Chain B: 



● Molecule 2: Vbeta 8.1

Chain H: 

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.84Å 65.80Å 97.38Å 90.00° 97.14° 90.00°	Depositor
Resolution (Å)	42.95 – 2.90 42.95 – 2.77	Depositor EDS
% Data completeness (in resolution range)	96.4 (42.95-2.90) 97.2 (42.95-2.77)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.254 , 0.327 0.265 , 0.338	Depositor DCC
R_{free} test set	1185 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	39.7	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 24.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	1 of 23479 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	6728	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.53	0/1417	0.73	1/1921 (0.1%)
1	G	0.54	0/1463	0.71	1/1986 (0.1%)
2	B	0.53	0/1940	0.75	1/2644 (0.0%)
2	H	0.59	0/1936	0.98	3/2638 (0.1%)
All	All	0.55	0/6756	0.81	6/9189 (0.1%)

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
2	B	0	1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	29	LEU	CA-CB-CG	6.96	131.30	115.30
1	A	29	LEU	CA-CB-CG	6.44	130.12	115.30
2	B	7	SER	C-N-CD	-6.05	107.28	120.60
2	H	123	ARG	CA-CB-CG	5.28	125.02	113.40
2	H	159	LEU	CA-CB-CG	5.20	127.25	115.30
2	H	49	GLY	N-CA-C	-5.04	100.51	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	7	SER	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1393	0	1312	85	0
1	G	1437	0	1348	90	0
2	B	1887	0	1788	114	0
2	H	1883	0	1778	89	0
3	A	28	0	0	2	0
3	B	43	0	0	18	0
3	G	27	0	0	6	0
3	H	30	0	0	5	0
All	All	6728	0	6226	338	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 26.

All (338) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:17:GLY:HA3	3:H:275:HOH:O	1.69	0.93
2:B:58:VAL:HA	3:B:271:HOH:O	1.72	0.88
2:B:216:ASN:HB3	2:B:219:ASN:HD22	1.35	0.88
2:B:216:ASN:HB3	2:B:219:ASN:ND2	1.89	0.86
2:H:6:GLN:HG2	2:H:120:PRO:HD2	1.57	0.84
2:H:48:HIS:HD2	2:H:51:ARG:CZ	1.90	0.84
2:B:6:GLN:HG2	2:B:120:PRO:HD2	1.61	0.83
2:H:216:ASN:HB3	2:H:219:ASN:ND2	1.95	0.81
1:A:153:PHE:HB2	1:A:205:PHE:CZ	2.14	0.81
1:A:166:LYS:O	1:G:166:LYS:O	1.99	0.81
2:B:162:GLY:O	2:B:200:ARG:HD3	1.81	0.80
2:H:43:ARG:HG2	2:H:53:ILE:HD11	1.65	0.78
2:B:236:TRP:CH2	2:B:238:GLN:HG3	2.17	0.78
1:G:153:PHE:HB2	1:G:205:PHE:CZ	2.18	0.78
2:H:236:TRP:CH2	2:H:238:GLN:HG3	2.19	0.78
2:B:194:PRO:HG2	1:G:166:LYS:HG3	1.64	0.77
2:B:43:ARG:HG2	2:B:53:ILE:HD11	1.66	0.77
2:H:216:ASN:HB3	2:H:219:ASN:HD22	1.49	0.77
2:H:13:ALA:HA	3:H:270:HOH:O	1.84	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:185:PHE:HE2	1:A:187:SER:HB2	1.48	0.76
1:G:185:PHE:HE2	1:G:187:SER:HB2	1.51	0.75
2:B:6:GLN:HB2	3:B:279:HOH:O	1.86	0.74
2:B:31:TYR:CE1	2:B:57:TYR:CE1	2.77	0.72
2:H:167:HIS:HE1	3:H:69:HOH:O	1.71	0.72
2:H:134:PHE:HZ	2:H:240:ARG:HH22	1.34	0.72
2:H:31:TYR:CE1	2:H:57:TYR:CE1	2.78	0.71
2:B:123:ARG:NH1	3:B:289:HOH:O	2.23	0.71
2:H:162:GLY:O	2:H:200:ARG:HD3	1.90	0.71
2:B:134:PHE:HZ	2:B:240:ARG:HH22	1.35	0.70
2:B:162:GLY:C	2:B:200:ARG:HD3	2.13	0.70
1:G:84(A):SER:C	1:G:86:ARG:NH2	2.45	0.69
1:A:185:PHE:CE2	1:A:187:SER:HB2	2.28	0.68
1:G:185:PHE:CE2	1:G:187:SER:HB2	2.28	0.68
1:A:110:SER:HB2	2:B:109:GLY:O	1.93	0.68
1:A:113:PHE:HB2	2:B:31:TYR:OH	1.95	0.67
2:B:196:LEU:HD13	2:B:196:LEU:H	1.57	0.67
1:G:169:ASP:OD2	1:G:196:LYS:HD3	1.95	0.67
1:G:128:PRO:HG3	1:G:177:VAL:HG21	1.77	0.66
1:A:169:ASP:O	1:A:194:SER:HB2	1.95	0.66
2:H:134:PHE:CZ	2:H:240:ARG:NH2	2.60	0.66
2:B:193:GLN:HB3	2:B:196:LEU:HD11	1.77	0.66
2:B:196:LEU:HD13	2:B:196:LEU:N	2.11	0.66
2:H:15:THR:HA	2:H:94:ALA:O	1.96	0.66
1:G:110:SER:HB3	3:G:239:HOH:O	1.95	0.66
1:G:55:TYR:HE2	1:G:57:PRO:HG3	1.61	0.66
2:H:215:GLN:HA	2:H:255:ARG:O	1.95	0.66
2:H:162:GLY:C	2:H:200:ARG:HD3	2.16	0.65
2:B:31:TYR:HE1	2:B:57:TYR:CE1	2.15	0.65
1:G:55:TYR:CE2	1:G:57:PRO:HG3	2.31	0.65
1:A:29:LEU:HB3	1:A:106:ILE:HD13	1.78	0.65
2:B:220:HIS:HE1	2:B:251:GLU:HB2	1.62	0.65
1:A:44:GLN:HB2	1:A:50:LEU:HD23	1.78	0.65
1:A:153:PHE:HB2	1:A:205:PHE:CE2	2.32	0.64
1:A:169:ASP:OD2	1:A:196:LYS:HD3	1.98	0.64
1:A:20:GLU:OE2	1:A:90:HIS:CG	2.50	0.64
2:B:193:GLN:CB	2:B:196:LEU:HD11	2.26	0.64
1:A:18:ASN:HB3	3:A:73:HOH:O	1.98	0.64
2:B:196:LEU:HD23	2:B:197:ASN:O	1.97	0.64
2:B:215:GLN:HA	2:B:255:ARG:O	1.98	0.64
2:B:48:HIS:HD2	2:B:51:ARG:CZ	2.11	0.64
1:A:113:PHE:HB2	2:B:31:TYR:CZ	2.33	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:31:TYR:HE1	2:H:57:TYR:CE1	2.16	0.64
1:A:25:TYR:CD1	1:A:26:SER:N	2.66	0.63
1:A:40:GLN:NE2	2:B:115:GLU:HG3	2.13	0.63
1:G:169:ASP:O	1:G:194:SER:HB2	1.97	0.63
1:G:44:GLN:HB2	1:G:50:LEU:HD23	1.80	0.63
1:G:110:SER:HB2	2:H:109:GLY:O	1.98	0.63
1:G:140:LEU:HB3	2:H:141:PHE:HB3	1.81	0.63
1:A:207:ASN:ND2	2:H:195:ALA:HB2	2.14	0.62
2:H:91:LEU:N	2:H:91:LEU:HD12	2.14	0.62
2:B:59:ALA:N	3:B:271:HOH:O	2.31	0.62
2:H:48:HIS:CD2	2:H:51:ARG:CZ	2.79	0.62
1:A:166:LYS:HG3	2:H:194:PRO:HG2	1.81	0.62
2:H:10:SER:HB2	2:H:167:HIS:HD2	1.65	0.61
2:B:158:CYS:O	2:B:204:SER:HA	2.00	0.61
1:G:19:ALA:HB3	1:G:91:ILE:HB	1.83	0.61
1:A:212:GLU:HG2	3:A:233:HOH:O	2.01	0.61
2:B:15:THR:HA	2:B:94:ALA:O	1.99	0.61
1:A:135:PRO:HB2	1:A:214:THR:HA	1.82	0.60
1:G:170:VAL:HA	1:G:194:SER:HB3	1.83	0.60
1:G:128:PRO:HG3	1:G:177:VAL:CG2	2.31	0.60
1:A:113:PHE:H	1:A:113:PHE:HD2	1.48	0.60
2:H:134:PHE:CE1	2:H:240:ARG:NH2	2.70	0.60
2:H:31:TYR:HA	2:H:56:SER:O	2.02	0.60
1:G:25:TYR:CD1	1:G:26:SER:N	2.70	0.60
1:G:113:PHE:HD2	1:G:113:PHE:H	1.50	0.59
1:G:135:PRO:HB2	1:G:214:THR:HA	1.85	0.59
2:B:236:TRP:N	3:B:273:HOH:O	2.36	0.59
1:A:3:GLN:HB3	1:A:25:TYR:HE1	1.66	0.59
1:G:29:LEU:HB3	1:G:106:ILE:HD13	1.83	0.58
1:A:170:VAL:HA	1:A:194:SER:HB3	1.85	0.58
1:G:178:LEU:HD23	3:G:240:HOH:O	2.03	0.58
2:B:91:LEU:HD12	2:B:91:LEU:N	2.18	0.58
1:G:153:PHE:HB2	1:G:205:PHE:CE2	2.39	0.58
1:G:40:GLN:NE2	2:H:115:GLU:HG3	2.18	0.58
2:H:220:HIS:HE1	2:H:251:GLU:HB2	1.67	0.58
2:B:128:GLU:OE1	2:B:128:GLU:HA	2.03	0.57
2:B:236:TRP:CZ2	2:B:238:GLN:HB2	2.39	0.57
2:B:134:PHE:CZ	2:B:240:ARG:NH2	2.64	0.57
2:B:216:ASN:HB2	3:B:112:HOH:O	2.04	0.57
2:B:48:HIS:CD2	2:B:51:ARG:CZ	2.88	0.57
2:B:100:ALA:C	3:B:289:HOH:O	2.43	0.56
1:A:19:ALA:HB3	1:A:91:ILE:HB	1.87	0.56

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:10:SER:HB2	2:B:167:HIS:HD2	1.69	0.56
2:B:31:TYR:CE1	2:B:57:TYR:CD1	2.93	0.56
1:G:20:GLU:OE2	1:G:90:HIS:CG	2.59	0.56
1:A:32:MET:HE1	1:A:41:TRP:HE1	1.70	0.56
1:A:175:LYS:HA	1:A:189:SER:O	2.05	0.56
2:B:134:PHE:CE1	2:B:240:ARG:NH2	2.73	0.56
2:H:158:CYS:O	2:H:204:SER:HA	2.05	0.56
1:G:178:LEU:HD13	2:H:182:GLY:HA2	1.86	0.55
1:G:63:ARG:C	1:G:70:GLY:H	2.10	0.55
2:H:31:TYR:CE1	2:H:57:TYR:CD1	2.93	0.55
2:B:108:LEU:HB3	2:B:114:TYR:HB2	1.89	0.55
2:B:143:PRO:HB3	2:B:154:ALA:HB1	1.88	0.55
1:G:53:LEU:O	1:G:54:LEU:CD1	2.54	0.55
1:G:22:GLN:OE1	1:G:88:SER:CB	2.53	0.55
2:B:27:ASN:HB3	2:B:29:HIS:CD2	2.42	0.55
2:H:30:ASP:OD2	2:H:31:TYR:CE2	2.60	0.55
1:G:3:GLN:HB3	1:G:25:TYR:HE1	1.73	0.54
1:A:110:SER:O	2:B:109:GLY:C	2.46	0.54
1:A:128:PRO:HG3	1:A:177:VAL:HG21	1.90	0.54
1:G:127:VAL:CG1	1:G:158:SER:HB3	2.39	0.53
1:A:63:ARG:C	1:A:70:GLY:H	2.12	0.53
2:B:216:ASN:CB	2:B:219:ASN:ND2	2.66	0.53
2:B:41:TRP:HH2	3:B:279:HOH:O	1.91	0.53
1:G:175:LYS:HA	1:G:189:SER:O	2.08	0.53
2:H:243:PRO:HA	3:H:282:HOH:O	2.06	0.53
1:G:113:PHE:HB2	2:H:31:TYR:OH	2.09	0.53
2:H:55:TYR:CE1	2:H:63:GLU:HG3	2.43	0.53
2:B:30:ASP:OD2	2:B:31:TYR:CE2	2.61	0.53
1:G:49:ARG:NH1	2:H:118:PHE:O	2.42	0.53
2:H:13:ALA:CA	3:H:270:HOH:O	2.51	0.52
2:B:48:HIS:HD2	2:B:51:ARG:NE	2.07	0.52
1:A:110:SER:HB2	2:B:113:GLY:CA	2.40	0.52
1:G:53:LEU:O	1:G:54:LEU:HD13	2.09	0.52
1:A:128:PRO:HG3	1:A:177:VAL:CG2	2.39	0.52
2:B:55:TYR:CE1	2:B:63:GLU:HG3	2.43	0.52
2:H:20:THR:HG23	2:H:90:ILE:CD1	2.40	0.52
2:B:254:GLY:O	2:B:255:ARG:HG3	2.10	0.52
1:G:188:ASN:N	3:G:220:HOH:O	2.20	0.52
2:B:41:TRP:HD1	2:B:86:PHE:CE2	2.28	0.52
1:G:189:SER:OG	2:H:206:ARG:HD2	2.10	0.52
2:B:166:ASP:HB3	2:B:201:TYR:CG	2.45	0.52
2:B:27:ASN:CB	2:B:29:HIS:CD2	2.94	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:236:TRP:CZ2	2:H:238:GLN:HB2	2.46	0.51
2:H:128:GLU:HA	2:H:128:GLU:OE1	2.10	0.51
2:B:162:GLY:HA2	2:B:200:ARG:CD	2.41	0.51
1:G:40:GLN:HE22	2:H:115:GLU:HG3	1.76	0.51
1:A:166:LYS:CE	1:G:168:SER:HA	2.41	0.51
2:B:27:ASN:HB3	2:B:29:HIS:NE2	2.26	0.51
1:A:22:GLN:OE1	1:A:88:SER:CB	2.60	0.50
1:G:30:ASN:HB3	1:G:107:ALA:O	2.12	0.50
1:G:61:GLU:N	3:G:222:HOH:O	2.43	0.50
1:A:110:SER:HA	2:B:113:GLY:HA2	1.92	0.50
1:G:11:LEU:N	1:G:11:LEU:HD23	2.26	0.50
2:B:20:THR:HG23	2:B:90:ILE:CD1	2.42	0.50
1:G:8:PRO:O	1:G:122:THR:HG23	2.12	0.50
1:A:178:LEU:HD12	1:A:178:LEU:C	2.32	0.50
1:G:22:GLN:OE1	1:G:88:SER:HB2	2.12	0.50
1:G:32:MET:HE1	1:G:41:TRP:HE1	1.77	0.50
2:B:31:TYR:HA	2:B:56:SER:O	2.13	0.49
1:A:110:SER:O	2:B:113:GLY:N	2.45	0.49
1:G:29:LEU:HD13	1:G:29:LEU:H	1.78	0.49
1:A:176:CYS:SG	2:B:206:ARG:NH1	2.86	0.49
2:H:44:GLN:HB2	2:H:50:LEU:HD23	1.93	0.49
1:G:171:TYR:O	1:G:192:ALA:HA	2.13	0.49
1:G:113:PHE:HB2	2:H:31:TYR:CZ	2.48	0.49
1:A:11:LEU:N	1:A:11:LEU:HD23	2.28	0.49
1:A:150:VAL:HG22	1:A:193:TRP:HB3	1.95	0.49
1:A:113:PHE:CB	2:B:31:TYR:OH	2.59	0.49
1:A:115:LYS:HA	2:B:107:ARG:HH12	1.77	0.49
1:A:171:TYR:O	1:A:192:ALA:HA	2.13	0.49
2:B:4:VAL:O	2:B:120:PRO:HD3	2.12	0.48
1:G:44:GLN:HA	3:G:233:HOH:O	2.13	0.48
2:H:141:PHE:N	2:H:141:PHE:CD2	2.81	0.48
1:A:29:LEU:HB3	1:A:106:ILE:CD1	2.42	0.48
1:G:3:GLN:NE2	1:G:26:SER:HB2	2.28	0.48
1:A:207:ASN:HD22	2:H:195:ALA:HB2	1.77	0.48
2:B:129:ASP:C	2:B:129:ASP:OD1	2.52	0.48
2:B:52:LEU:O	2:B:71:ILE:HG12	2.13	0.48
1:G:84(B):ASN:N	1:G:86:ARG:HH21	2.12	0.48
2:B:193:GLN:HB2	2:B:196:LEU:HD11	1.95	0.48
1:A:127:VAL:CG1	1:A:158:SER:HB3	2.42	0.48
1:A:210:ILE:O	1:A:211:PRO:C	2.51	0.48
1:A:7:SER:HB3	1:A:22:GLN:HB2	1.95	0.48
1:A:13:LEU:HD23	1:G:8:PRO:HB3	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:41:TRP:CD1	2:H:86:PHE:CD2	3.02	0.48
2:H:41:TRP:HD1	2:H:86:PHE:CE2	2.32	0.48
1:A:40:GLN:HE22	2:B:115:GLU:HG3	1.77	0.48
2:H:75:TYR:HB3	2:H:88:LEU:HD11	1.96	0.48
2:B:88:LEU:C	2:B:90:ILE:HD13	2.34	0.48
2:H:114:TYR:N	2:H:114:TYR:CD2	2.81	0.48
2:H:27:ASN:HB3	2:H:29:HIS:CD2	2.49	0.48
1:A:29:LEU:HD13	1:A:29:LEU:H	1.79	0.47
1:G:84(A):SER:C	1:G:86:ARG:HH22	2.15	0.47
1:A:54:LEU:HD12	1:A:54:LEU:HA	1.59	0.47
2:H:45:ASP:OD1	2:H:45:ASP:N	2.47	0.47
2:H:48:HIS:HD2	2:H:51:ARG:NE	2.12	0.47
1:G:40:GLN:HG3	1:G:54:LEU:O	2.14	0.47
1:A:178:LEU:HD13	2:B:182:GLY:HA2	1.97	0.47
1:A:46:PRO:HG3	3:B:285:HOH:O	2.15	0.47
2:H:30:ASP:OD2	2:H:31:TYR:CD2	2.68	0.47
1:G:84(B):ASN:N	1:G:86:ARG:NH2	2.63	0.47
1:A:30:ASN:HB3	1:A:107:ALA:O	2.15	0.47
2:B:41:TRP:CD1	2:B:86:PHE:CD2	3.01	0.47
1:A:3:GLN:NE2	1:A:26:SER:HB2	2.29	0.47
1:G:29:LEU:H	1:G:29:LEU:CD1	2.27	0.47
1:G:22:GLN:OE1	1:G:88:SER:HB3	2.13	0.47
1:G:210:ILE:O	1:G:211:PRO:C	2.53	0.47
2:H:168:VAL:HA	2:H:226:GLN:O	2.15	0.47
2:H:4:VAL:O	2:H:120:PRO:HD3	2.14	0.47
1:A:3:GLN:HB3	1:A:25:TYR:CE1	2.49	0.47
2:B:43:ARG:NH1	3:B:268:HOH:O	2.47	0.47
1:G:57:PRO:O	1:G:58:SER:O	2.33	0.46
1:A:29:LEU:H	1:A:29:LEU:CD1	2.28	0.46
1:G:3:GLN:HB3	1:G:25:TYR:CE1	2.50	0.46
2:H:125:THR:HG21	2:H:165:PRO:CG	2.45	0.46
1:A:168:SER:HA	1:G:166:LYS:CE	2.46	0.46
1:A:22:GLN:OE1	1:A:88:SER:HB2	2.15	0.46
1:G:63:ARG:O	1:G:70:GLY:N	2.48	0.46
2:H:162:GLY:HA2	2:H:200:ARG:CD	2.46	0.46
2:H:46:THR:HA	2:H:47:GLY:HA2	1.73	0.46
1:G:113:PHE:CB	2:H:31:TYR:OH	2.64	0.46
1:A:8:PRO:O	1:A:122:THR:HG23	2.15	0.46
2:B:53:ILE:N	2:B:53:ILE:HD12	2.31	0.46
1:G:29:LEU:HB3	1:G:106:ILE:CD1	2.43	0.46
2:H:108:LEU:HB3	2:H:114:TYR:HB2	1.98	0.46
2:B:236:TRP:CZ2	2:B:238:GLN:HG3	2.51	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:101:VAL:HG23	3:B:289:HOH:O	2.15	0.46
1:G:196:LYS:N	3:G:228:HOH:O	2.48	0.46
1:A:133:PRO:HB3	1:A:157:ASP:HB3	1.98	0.45
1:G:54:LEU:HD12	1:G:54:LEU:HA	1.67	0.45
2:B:28:ASN:ND2	3:B:278:HOH:O	2.49	0.45
2:B:235:GLU:HA	3:B:273:HOH:O	2.16	0.45
2:H:51:ARG:HB3	2:H:71:ILE:CD1	2.46	0.45
2:H:134:PHE:CE1	2:H:240:ARG:NH1	2.85	0.45
2:H:254:GLY:O	2:H:255:ARG:HG3	2.16	0.45
1:A:63:ARG:O	1:A:70:GLY:N	2.49	0.45
1:G:178:LEU:CD1	2:H:182:GLY:HA2	2.47	0.45
2:B:195:ALA:HB2	1:G:207:ASN:ND2	2.32	0.45
1:G:127:VAL:HG11	1:G:158:SER:HB3	1.99	0.45
2:H:88:LEU:C	2:H:90:ILE:HD13	2.37	0.45
1:G:84(C):GLU:O	1:G:86:ARG:N	2.49	0.45
2:B:226:GLN:NE2	3:B:266:HOH:O	2.46	0.45
1:G:176:CYS:SG	2:H:206:ARG:NH1	2.89	0.45
1:A:82:SER:O	1:A:87:SER:HA	2.17	0.45
2:B:216:ASN:ND2	2:B:219:ASN:ND2	2.65	0.45
2:B:192:GLU:HG3	2:B:200:ARG:O	2.17	0.45
2:H:218:ARG:O	2:H:253:TRP:HZ3	2.00	0.44
1:G:84(A):SER:C	1:G:86:ARG:HH21	2.18	0.44
2:H:236:TRP:CZ2	2:H:238:GLN:HG3	2.51	0.44
2:H:10:SER:HA	2:H:123:ARG:O	2.17	0.44
2:B:11:LYS:NZ	3:B:288:HOH:O	2.39	0.44
2:B:216:ASN:O	2:B:254:GLY:HA3	2.16	0.44
2:B:162:GLY:CA	2:B:200:ARG:HD3	2.48	0.44
1:G:113:PHE:HB2	2:H:31:TYR:CE1	2.53	0.44
2:B:75:TYR:HB3	2:B:88:LEU:HD11	1.99	0.44
2:B:151:THR:O	2:B:152:GLN:HB2	2.17	0.44
1:A:8:PRO:HB2	1:G:12:VAL:O	2.16	0.44
2:B:143:PRO:CB	2:B:154:ALA:HB1	2.48	0.44
2:B:163:PHE:CE2	2:B:201:TYR:HB2	2.51	0.44
2:B:80:ARG:HD2	2:B:85:ASN:O	2.17	0.44
2:B:134:PHE:CE1	2:B:240:ARG:NH1	2.86	0.44
1:G:169:ASP:N	1:G:169:ASP:OD1	2.49	0.44
2:H:129:ASP:OD1	2:H:129:ASP:C	2.56	0.44
2:B:141:PHE:N	2:B:141:PHE:CD2	2.85	0.43
2:H:10:SER:HB2	2:H:167:HIS:CD2	2.49	0.43
1:A:53:LEU:O	1:A:54:LEU:CD1	2.66	0.43
1:A:140:LEU:HB3	2:B:141:PHE:HB3	2.00	0.43
2:B:41:TRP:CD1	2:B:86:PHE:CE2	3.06	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:216:ASN:CB	2:H:219:ASN:ND2	2.76	0.43
2:H:166:ASP:HB3	2:H:201:TYR:CG	2.54	0.43
2:B:125:THR:HG21	2:B:165:PRO:CG	2.49	0.43
2:H:219:ASN:O	2:H:253:TRP:HA	2.19	0.43
2:H:53:ILE:N	2:H:53:ILE:HD12	2.33	0.43
1:A:113:PHE:HB2	2:B:31:TYR:CE1	2.53	0.43
1:G:178:LEU:HD12	1:G:178:LEU:C	2.38	0.43
2:B:96:LEU:HD11	2:B:128:GLU:OE2	2.19	0.43
1:A:12:VAL:O	1:G:8:PRO:HB2	2.18	0.43
2:H:166:ASP:HB3	2:H:201:TYR:CD2	2.54	0.43
1:A:180:MET:HE1	1:A:183:MET:SD	2.59	0.43
2:H:52:LEU:O	2:H:71:ILE:HG12	2.19	0.43
2:H:27:ASN:CB	2:H:29:HIS:CD2	3.02	0.43
2:B:101:VAL:CG2	3:B:289:HOH:O	2.66	0.43
1:G:84(B):ASN:O	1:G:84(C):GLU:C	2.56	0.43
2:H:91:LEU:N	2:H:91:LEU:CD1	2.82	0.42
2:H:41:TRP:CD1	2:H:86:PHE:CE2	3.07	0.42
1:A:72:LEU:N	1:A:72:LEU:HD12	2.34	0.42
2:B:218:ARG:O	2:B:253:TRP:HZ3	2.02	0.42
1:A:29:LEU:CB	1:A:106:ILE:HD13	2.47	0.42
1:G:53:LEU:O	1:G:54:LEU:HD12	2.18	0.42
1:A:212:GLU:O	1:A:212:GLU:HG2	2.19	0.42
1:A:53:LEU:O	1:A:54:LEU:HD13	2.18	0.42
1:G:6:GLN:HA	1:G:22:GLN:O	2.20	0.42
2:B:168:VAL:HA	2:B:226:GLN:O	2.20	0.42
1:A:105:ALA:HB2	1:A:118:PHE:CD1	2.55	0.42
2:H:27:ASN:HB3	2:H:29:HIS:NE2	2.35	0.42
1:A:21:LEU:HD22	1:A:122:THR:HG21	2.02	0.42
2:B:216:ASN:HD22	2:B:219:ASN:HD21	1.68	0.42
2:B:10:SER:HB2	2:B:167:HIS:CD2	2.51	0.42
2:B:43:ARG:HH21	2:B:100:ALA:HB2	1.85	0.41
1:A:45:ARG:HG3	1:A:46:PRO:O	2.19	0.41
2:B:45:ASP:N	2:B:45:ASP:OD1	2.52	0.41
2:B:196:LEU:N	2:B:196:LEU:CD1	2.81	0.41
1:A:113:PHE:CD2	1:A:113:PHE:N	2.88	0.41
2:B:30:ASP:OD2	2:B:31:TYR:CD2	2.73	0.41
1:G:210:ILE:O	1:G:210:ILE:HG13	2.18	0.41
1:A:71:ARG:C	1:A:72:LEU:HD12	2.40	0.41
2:B:255:ARG:HB3	3:B:290:HOH:O	2.20	0.41
2:H:236:TRP:CZ2	2:H:243:PRO:HD3	2.56	0.41
2:B:43:ARG:NH2	2:B:100:ALA:HB2	2.36	0.41
2:B:255:ARG:NH1	3:B:270:HOH:O	2.53	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:227:PHE:O	2:B:245:THR:HG23	2.20	0.41
1:G:127:VAL:HG13	1:G:158:SER:OG	2.20	0.41
1:G:82:SER:O	1:G:87:SER:HA	2.20	0.41
1:G:29:LEU:N	1:G:29:LEU:CD1	2.84	0.41
1:G:11:LEU:HD21	1:G:124:LEU:HD13	2.03	0.41
2:B:114:TYR:HB3	2:B:115:GLU:H	1.61	0.41
1:G:12:VAL:HA	1:G:125:SER:O	2.21	0.41
1:G:131:GLN:O	1:G:133:PRO:HD3	2.21	0.41
2:H:162:GLY:CA	2:H:200:ARG:HD3	2.51	0.41
1:G:133:PRO:HB3	1:G:157:ASP:HB3	2.03	0.41
2:H:143:PRO:HD2	2:H:214:TRP:CE2	2.56	0.41
2:B:244:VAL:O	2:B:246:GLN:HG2	2.20	0.41
1:A:42:PHE:CD1	1:A:52:SER:HA	2.56	0.41
2:B:46:THR:HA	2:B:47:GLY:HA2	1.74	0.41
2:H:80:ARG:HD2	2:H:85:ASN:O	2.21	0.41
1:A:166:LYS:HE2	1:G:168:SER:HA	2.02	0.40
1:A:134:ASP:O	1:A:135:PRO:C	2.58	0.40
1:A:169:ASP:OD1	1:A:169:ASP:N	2.53	0.40
1:A:3:GLN:HE21	1:A:3:GLN:HB3	1.69	0.40
2:B:44:GLN:HB2	2:B:50:LEU:HD23	2.04	0.40
2:B:219:ASN:O	2:B:253:TRP:HA	2.22	0.40
2:H:51:ARG:CB	2:H:71:ILE:HD11	2.52	0.40
2:H:62:THR:HG22	2:H:63:GLU:N	2.36	0.40
1:A:43:TYR:C	1:A:43:TYR:CD2	2.94	0.40
1:A:137:VAL:O	1:A:217:PRO:HD2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	175/198 (88%)	162 (93%)	12 (7%)	1 (1%)	33	76
1	G	184/198 (93%)	170 (92%)	12 (6%)	2 (1%)	21	60
2	B	236/238 (99%)	222 (94%)	12 (5%)	2 (1%)	27	68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	236/238 (99%)	225 (95%)	10 (4%)	1 (0%)	43	82
All	All	831/872 (95%)	779 (94%)	46 (6%)	6 (1%)	30	72

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	GLY
2	B	254	GLY
2	H	254	GLY
2	B	8	PRO
1	G	64	GLY
1	G	211	PRO

5.3.2 Protein sidechains ⓘ

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	160/177 (90%)	138 (86%)	22 (14%)	5	14
1	G	166/177 (94%)	143 (86%)	23 (14%)	5	14
2	B	202/205 (98%)	194 (96%)	8 (4%)	42	81
2	H	201/205 (98%)	195 (97%)	6 (3%)	53	89
All	All	729/764 (95%)	670 (92%)	59 (8%)	17	45

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	10	SER
1	A	13	LEU
1	A	29	LEU
1	A	49	ARG
1	A	54	LEU
1	A	85	SER
1	A	90	HIS
1	A	106	ILE

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Mol	Chain	Res	Type
1	A	113	PHE
1	A	115	LYS
1	A	123	SER
1	A	127	VAL
1	A	154	THR
1	A	157	ASP
1	A	164	GLN
1	A	166	LYS
1	A	178	LEU
1	A	196	LYS
1	A	198	ASP
1	A	207	ASN
1	A	212	GLU
2	B	9	ARG
2	B	10	SER
2	B	45	ASP
2	B	48	HIS
2	B	188	GLN
2	B	196	LEU
2	B	206	ARG
2	B	238	GLN
1	G	3	GLN
1	G	10	SER
1	G	13	LEU
1	G	29	LEU
1	G	49	ARG
1	G	54	LEU
1	G	85	SER
1	G	90	HIS
1	G	106	ILE
1	G	113	PHE
1	G	115	LYS
1	G	123	SER
1	G	127	VAL
1	G	150	VAL
1	G	154	THR
1	G	157	ASP
1	G	164	GLN
1	G	166	LYS
1	G	178	LEU
1	G	196	LYS
1	G	198	ASP

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Mol	Chain	Res	Type
1	G	207	ASN
1	G	212	GLU
2	H	9	ARG
2	H	10	SER
2	H	45	ASP
2	H	188	GLN
2	H	206	ARG
2	H	238	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	159	GLN
1	A	195	ASN
2	B	44	GLN
2	B	48	HIS
2	B	132	ASN
2	B	167	HIS
2	B	219	ASN
2	B	220	HIS
1	G	3	GLN
1	G	44	GLN
1	G	159	GLN
1	G	195	ASN
2	H	44	GLN
2	H	48	HIS
2	H	167	HIS
2	H	219	ASN
2	H	220	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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 ⓘ

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, 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	183/198 (92%)	0.27	10 (5%) 24 29	10, 29, 55, 82	0
1	G	190/198 (95%)	0.12	4 (2%) 60 69	11, 28, 56, 83	0
2	B	238/238 (100%)	0.61	27 (11%) 6 7	20, 35, 63, 98	0
2	H	238/238 (100%)	0.26	10 (4%) 35 41	17, 33, 61, 98	0
All	All	849/872 (97%)	0.33	51 (6%) 21 25	10, 32, 58, 98	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	182	SER	4.5
2	H	240	ARG	4.4
2	B	240	ARG	4.3
2	B	239	ASP	4.1
1	G	218	SER	3.6
1	A	165	SER	3.5
1	G	198	ASP	3.5
2	B	241	ALA	3.4
2	B	57	TYR	3.4
2	H	239	ASP	3.2
1	A	113	PHE	3.1
1	A	210	ILE	3.1
1	A	183	MET	3.0
2	H	233	ASN	2.9
2	B	215	GLN	2.9
2	B	237	THR	2.9
1	G	84(C)	GLU	2.8
2	B	109	GLY	2.8
2	B	195	ALA	2.8
2	H	197	ASN	2.7
2	B	175	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
2	B	17	GLY	2.6
2	B	236	TRP	2.6
2	B	145	GLU	2.6
1	A	131	GLN	2.6
2	B	113	GLY	2.5
2	H	3	ALA	2.5
1	A	84	ALA	2.5
2	B	234	ASP	2.5
1	G	197	SER	2.5
1	A	141	ARG	2.5
2	B	128	GLU	2.4
2	B	217	PRO	2.4
2	B	94	ALA	2.3
2	B	176	GLY	2.3
2	B	194	PRO	2.3
2	H	108	LEU	2.3
1	A	24	THR	2.3
2	B	211	ALA	2.3
1	A	83	ALA	2.2
2	B	218	ARG	2.2
2	B	65	GLY	2.2
2	H	7	SER	2.2
2	H	198	ASP	2.2
2	B	256	ALA	2.1
2	B	131	LYS	2.1
2	B	18	LYS	2.1
2	B	164	TYR	2.1
2	H	238	GLN	2.1
2	H	216	ASN	2.1
2	B	180	HIS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.