



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 02:37 PM GMT

PDB ID : 1NGW
Title : Chimeric Affinity Matured Fab 7g12 complexed with mesoporphyrin
Authors : Yin, J.; Andryski, S.E.; Beuscher IV, A.E.; Stevens, R.C.; Schultz, P.G.
Deposited on : 2002-12-18
Resolution : 2.60 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>

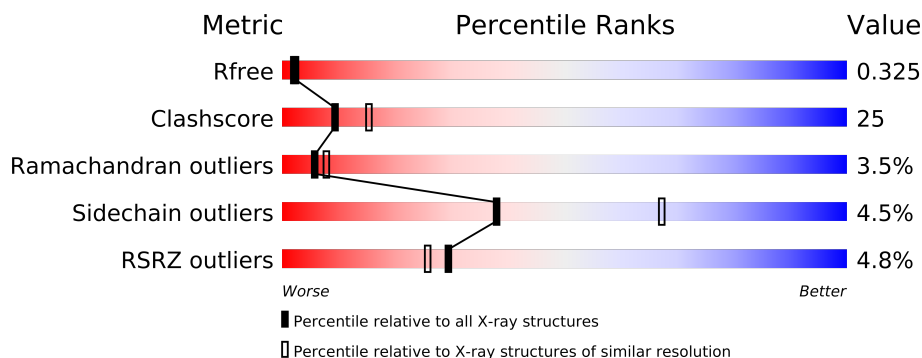
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.60 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	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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	213	
1	L	213	
2	B	216	
2	H	216	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	MMP	H	217	-	X

2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 6645 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.

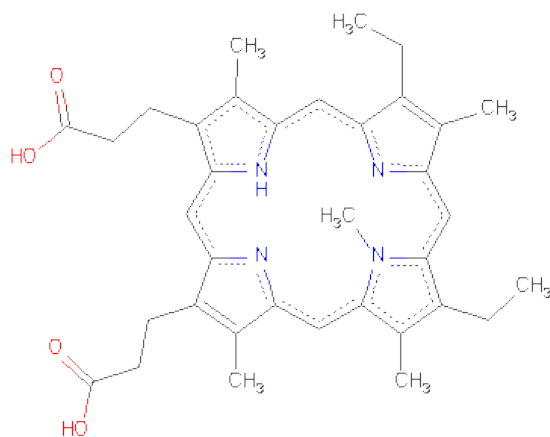
- Molecule 1 is a protein called Mature Metal Chelatase Catalytic Antibody, Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	0	0
			1646	1029	274	336	7			
1	A	213	Total	C	N	O	S	0	0	0
			1646	1029	274	336	7			

- Molecule 2 is a protein called Mature Metal Chelatase Catalytic Antibody, Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1622	1025	269	320	8			
2	B	216	Total	C	N	O	S	0	0	0
			1622	1025	269	320	8			

- Molecule 3 is N-METHYLMESOPORPHYRIN (three-letter code: MMP) (formula: $C_{35}H_{40}N_4O_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			42	34	4	4		
3	H	1	Total	C	N	O	0	0
			42	34	4	4		

- Molecule 4 is water.

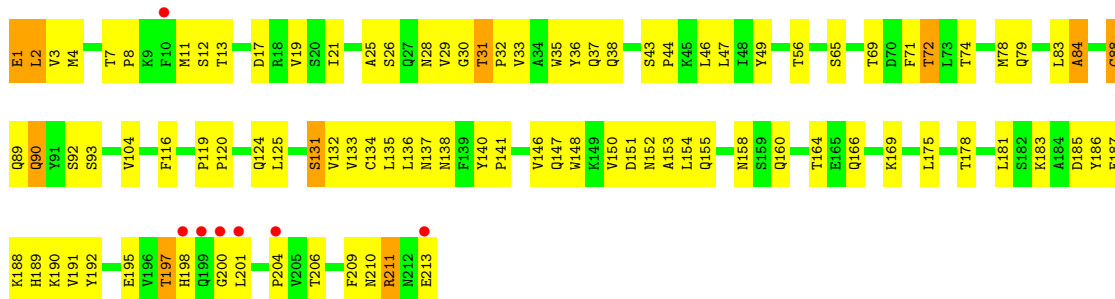
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		
4	B	7	Total	O	0	0
			7	7		
4	H	5	Total	O	0	0
			5	5		
4	L	6	Total	O	0	0
			6	6		

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.

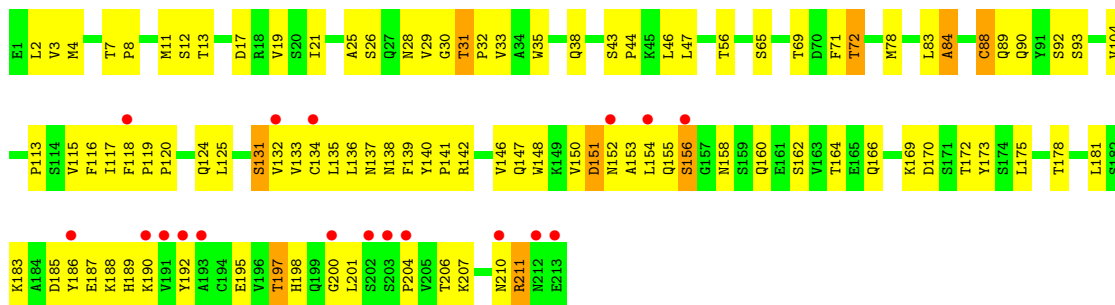
- Molecule 1: Mature Metal Chelatase Catalytic Antibody, Light chain

Chain L: 



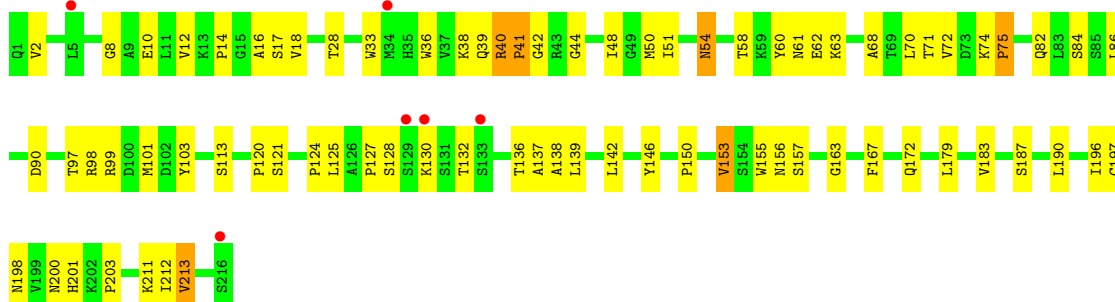
- Molecule 1: Mature Metal Chelatase Catalytic Antibody, Light chain

Chain A: 



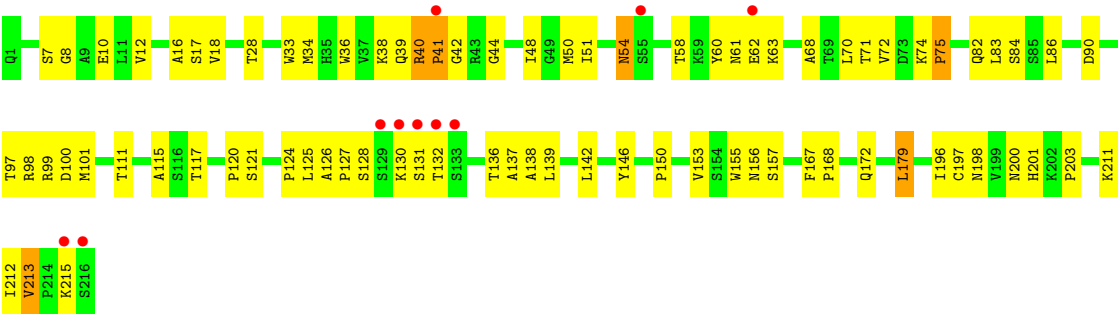
- Molecule 2: Mature Metal Chelatase Catalytic Antibody, Heavy chain

Chain H: 



● Molecule 2: Mature Metal Chelatase Catalytic Antibody, Heavy chain

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	100.72Å 134.22Å 72.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.60 20.15 – 2.40	Depositor EDS
% Data completeness (in resolution range)	89.0 (19.92-2.60) 84.2 (20.15-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.41Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.275 , 0.318 0.272 , 0.325	Depositor DCC
R_{free} test set	1379 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	53.0	Xtriage
Anisotropy	0.199	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 25.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	9 of 38586 reflections (0.023%)	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6645	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MMP

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	0.40	0/1681	0.65	0/2281
1	L	0.41	0/1681	0.66	0/2281
2	B	0.38	0/1662	0.68	0/2261
2	H	0.39	0/1662	0.68	0/2261
All	All	0.40	0/6686	0.67	0/9084

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1646	0	1596	108	1
1	L	1646	0	1596	98	0
2	B	1622	0	1609	63	0
2	H	1622	0	1609	65	0
3	B	42	0	37	1	0
3	H	42	0	37	2	0
4	A	7	0	0	6	0
4	B	7	0	0	3	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	5	0	0	3	0
4	L	6	0	0	1	0
All	All	6645	0	6484	324	1

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 25.

All (324) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:83:LEU:HD11	1:L:166:GLN:HB3	1.35	1.04
2:B:132:THR:CG2	2:B:138:ALA:H	1.74	0.99
2:H:132:THR:CG2	2:H:138:ALA:H	1.76	0.98
2:B:156:ASN:HD21	2:B:196:ILE:H	1.12	0.96
1:A:150:VAL:HG12	1:A:155:GLN:NE2	1.83	0.94
1:A:83:LEU:HD11	1:A:166:GLN:HB3	1.48	0.93
2:H:156:ASN:HD21	2:H:196:ILE:H	1.14	0.92
1:L:150:VAL:HG12	1:L:155:GLN:NE2	1.85	0.92
2:B:132:THR:HG21	2:B:138:ALA:H	1.35	0.90
2:H:132:THR:HG21	2:H:138:ALA:H	1.38	0.88
2:B:156:ASN:ND2	2:B:196:ILE:H	1.71	0.87
2:H:156:ASN:ND2	2:H:196:ILE:H	1.72	0.86
2:B:34:MET:HE3	4:B:219:HOH:O	1.75	0.85
1:A:132:VAL:HA	4:A:219:HOH:O	1.77	0.84
1:L:90:GLN:HE21	1:L:92:SER:H	1.22	0.83
1:A:90:GLN:HE21	1:A:92:SER:H	1.23	0.82
2:H:153:VAL:HG23	4:H:219:HOH:O	1.80	0.82
1:A:181:LEU:HD21	4:A:218:HOH:O	1.80	0.80
1:A:113:PRO:HB3	1:A:139:PHE:HB3	1.65	0.79
2:H:120:PRO:HB3	2:H:146:TYR:HB3	1.66	0.78
2:B:120:PRO:HB3	2:B:146:TYR:HB3	1.66	0.78
2:B:34:MET:HB3	4:B:219:HOH:O	1.85	0.75
1:L:133:VAL:HG22	1:L:178:THR:HG22	1.68	0.75
1:L:124:GLN:HE22	1:L:131:SER:HB2	1.52	0.74
1:A:31:THR:H	1:A:32:PRO:HD3	1.52	0.74
2:H:197:CYS:HB2	4:H:219:HOH:O	1.86	0.74
1:A:28:ASN:HA	1:A:69:THR:HG22	1.69	0.74
2:B:157:SER:H	2:B:198:ASN:HD21	1.35	0.74
1:A:124:GLN:HE22	1:A:131:SER:HB2	1.53	0.73
1:L:31:THR:H	1:L:32:PRO:HD3	1.53	0.73
2:B:51:ILE:HG13	2:B:58:THR:HG22	1.71	0.72
1:L:74:THR:HG22	4:L:215:HOH:O	1.88	0.72
2:H:157:SER:H	2:H:198:ASN:HD21	1.37	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:133:VAL:HG22	1:A:178:THR:HG22	1.71	0.72
1:L:28:ASN:HA	1:L:69:THR:HG22	1.70	0.71
2:H:51:ILE:HG13	2:H:58:THR:HG22	1.70	0.70
3:H:217:MMP:H182	3:H:217:MMP:H161	1.74	0.70
1:L:135:LEU:HD11	1:L:137:ASN:ND2	2.08	0.69
2:B:201:HIS:CD2	2:B:203:PRO:HD2	2.28	0.69
1:A:181:LEU:HD11	1:A:186:TYR:HB2	1.75	0.68
1:A:181:LEU:HD11	4:A:218:HOH:O	1.92	0.68
2:H:201:HIS:CD2	2:H:203:PRO:HD2	2.29	0.67
1:L:181:LEU:HD11	1:L:186:TYR:HB2	1.76	0.67
1:A:31:THR:H	1:A:32:PRO:CD	2.08	0.67
1:A:78:MET:HE1	1:A:104:VAL:HG11	1.75	0.67
1:L:31:THR:H	1:L:32:PRO:CD	2.08	0.67
1:L:38:GLN:HE22	2:H:39:GLN:HE22	1.43	0.66
2:H:54:ASN:HA	2:H:74:LYS:HE2	1.77	0.66
1:A:38:GLN:HE22	2:B:39:GLN:HE22	1.42	0.66
1:L:78:MET:HE1	1:L:104:VAL:HG11	1.78	0.66
1:A:90:GLN:NE2	1:A:93:SER:H	1.94	0.66
1:L:2:LEU:HD12	1:L:2:LEU:H	1.61	0.66
1:L:190:LYS:HZ3	1:L:211:ARG:HH21	1.44	0.66
2:B:54:ASN:HA	2:B:74:LYS:HE2	1.77	0.65
1:L:164:THR:HG23	2:H:167:PHE:CE1	2.31	0.64
1:A:135:LEU:HD11	1:A:137:ASN:ND2	2.13	0.64
1:A:3:VAL:HG23	1:A:26:SER:HB3	1.80	0.63
1:A:150:VAL:HG12	1:A:155:GLN:HE22	1.64	0.63
2:B:86:LEU:HA	2:B:90:ASP:OD1	1.97	0.63
1:L:13:THR:HG21	1:L:19:VAL:HG22	1.81	0.63
1:A:31:THR:N	1:A:32:PRO:CD	2.62	0.63
2:H:86:LEU:HA	2:H:90:ASP:OD1	1.99	0.63
1:L:3:VAL:HG23	1:L:26:SER:HB3	1.80	0.62
1:L:90:GLN:NE2	1:L:93:SER:H	1.96	0.62
1:A:183:LYS:O	1:A:187:GLU:HG3	2.00	0.62
2:H:156:ASN:HD21	2:H:196:ILE:N	1.93	0.62
2:B:61:ASN:O	2:B:63:LYS:N	2.31	0.62
2:B:127:PRO:HG3	2:B:139:LEU:HB3	1.80	0.62
1:A:190:LYS:HZ3	1:A:211:ARG:HH21	1.47	0.61
2:H:132:THR:HG22	2:H:137:ALA:HA	1.82	0.61
1:L:31:THR:N	1:L:32:PRO:CD	2.63	0.61
2:B:86:LEU:HD12	2:B:90:ASP:HB3	1.83	0.61
2:H:127:PRO:HG3	2:H:139:LEU:CB	2.31	0.61
1:L:90:GLN:HE21	1:L:92:SER:N	1.97	0.61
2:B:127:PRO:HG3	2:B:139:LEU:CB	2.30	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:13:THR:HG21	1:A:19:VAL:HG22	1.82	0.61
2:H:58:THR:HG1	2:H:60:TYR:HE1	1.49	0.60
1:L:19:VAL:HG21	1:L:78:MET:HE2	1.83	0.60
2:B:132:THR:HG22	2:B:137:ALA:HA	1.83	0.60
2:H:17:SER:OG	2:H:84:SER:HA	2.02	0.59
2:H:74:LYS:HB2	2:H:75:PRO:HD3	1.82	0.59
2:B:74:LYS:HB2	2:B:75:PRO:HD3	1.82	0.59
1:A:125:LEU:O	1:A:183:LYS:HD2	2.02	0.59
2:H:61:ASN:O	2:H:63:LYS:N	2.31	0.59
2:B:17:SER:OG	2:B:84:SER:HA	2.02	0.59
2:H:127:PRO:HG3	2:H:139:LEU:HB3	1.84	0.59
1:L:2:LEU:N	1:L:2:LEU:HD12	2.16	0.59
1:A:120:PRO:N	4:A:219:HOH:O	2.36	0.58
1:A:175:LEU:HD23	1:A:175:LEU:C	2.23	0.58
2:H:211:LYS:HE2	2:H:213:VAL:HG13	1.84	0.58
2:H:86:LEU:HD12	2:H:90:ASP:HB3	1.85	0.58
1:L:183:LYS:O	1:L:187:GLU:HG3	2.04	0.58
1:L:11:MET:HG3	1:L:12:SER:H	1.68	0.58
1:L:198:HIS:HB3	1:L:201:LEU:HD23	1.84	0.58
2:B:8:GLY:HA2	4:B:220:HOH:O	2.03	0.58
1:A:198:HIS:HB3	1:A:201:LEU:HD23	1.85	0.58
2:H:132:THR:HG21	2:H:138:ALA:N	2.15	0.57
1:L:175:LEU:HD23	1:L:175:LEU:C	2.24	0.57
1:A:90:GLN:HE21	1:A:92:SER:N	1.98	0.57
1:L:125:LEU:O	1:L:183:LYS:HD2	2.04	0.57
1:A:11:MET:HG3	1:A:12:SER:H	1.70	0.57
1:A:2:LEU:HD12	1:A:2:LEU:N	2.19	0.57
2:B:156:ASN:HD21	2:B:196:ILE:N	1.92	0.56
1:L:150:VAL:HG12	1:L:155:GLN:HE22	1.67	0.56
1:A:113:PRO:CB	1:A:139:PHE:HB3	2.35	0.56
1:A:150:VAL:HG12	1:A:155:GLN:CD	2.25	0.56
1:A:29:VAL:HG23	1:A:30:GLY:N	2.20	0.56
1:L:7:THR:HG23	1:L:8:PRO:HA	1.87	0.56
2:B:124:PRO:HB3	2:B:212:ILE:HD12	1.88	0.56
1:A:89:GLN:HG2	1:A:90:GLN:N	2.20	0.56
1:L:197:THR:HG23	1:L:204:PRO:HB3	1.87	0.56
1:A:190:LYS:NZ	1:A:211:ARG:HH21	2.03	0.56
1:A:190:LYS:O	1:A:210:ASN:HA	2.06	0.56
2:H:51:ILE:HD13	2:H:72:VAL:HG13	1.87	0.56
1:A:197:THR:HG23	1:A:204:PRO:HB3	1.87	0.56
1:L:29:VAL:HG23	1:L:30:GLY:N	2.21	0.56
1:A:185:ASP:HA	1:A:188:LYS:HE3	1.88	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:12:VAL:HG21	2:B:18:VAL:HG13	1.88	0.55
1:A:117:ILE:HG22	2:B:131:SER:HB3	1.88	0.55
1:L:185:ASP:HA	1:L:188:LYS:HE3	1.88	0.55
1:L:155:GLN:HB3	1:L:158:ASN:HD21	1.72	0.55
1:L:83:LEU:O	1:L:84:ALA:HB2	2.06	0.55
2:H:120:PRO:HD3	2:H:201:HIS:ND1	2.22	0.55
1:L:190:LYS:O	1:L:210:ASN:HA	2.06	0.55
1:A:113:PRO:HB3	1:A:139:PHE:CB	2.35	0.55
2:H:40:ARG:HG3	2:H:41:PRO:HD2	1.89	0.55
1:A:11:MET:CE	1:A:12:SER:H	2.20	0.55
2:B:211:LYS:HE2	2:B:213:VAL:HG13	1.87	0.55
2:B:120:PRO:HD3	2:B:201:HIS:ND1	2.22	0.55
2:B:40:ARG:HG3	2:B:41:PRO:HD2	1.88	0.55
1:A:19:VAL:HG21	1:A:78:MET:HE2	1.89	0.54
1:L:150:VAL:HG12	1:L:155:GLN:CD	2.28	0.54
1:L:190:LYS:NZ	1:L:211:ARG:HH21	2.05	0.54
2:B:58:THR:HG1	2:B:60:TYR:HE1	1.54	0.54
2:B:132:THR:HG21	2:B:138:ALA:N	2.12	0.54
1:L:195:GLU:HG2	1:L:206:THR:HB	1.89	0.54
1:L:186:TYR:HA	1:L:192:TYR:OH	2.08	0.54
1:A:113:PRO:HB3	1:A:139:PHE:CD1	2.43	0.53
1:A:7:THR:HG23	1:A:8:PRO:HA	1.91	0.53
1:A:195:GLU:HG2	1:A:206:THR:HB	1.89	0.53
1:L:89:GLN:HG2	1:L:90:GLN:N	2.23	0.53
2:H:124:PRO:HB3	2:H:212:ILE:HD12	1.89	0.53
1:L:11:MET:CE	1:L:12:SER:H	2.21	0.53
1:A:186:TYR:HD2	4:A:218:HOH:O	1.92	0.52
1:A:116:PHE:HD2	2:B:132:THR:HA	1.75	0.52
1:A:2:LEU:H	1:A:2:LEU:HD12	1.75	0.52
2:H:14:PRO:HD3	2:H:113:SER:C	2.30	0.52
1:A:186:TYR:HA	1:A:192:TYR:OH	2.09	0.52
1:L:181:LEU:HD12	1:L:181:LEU:C	2.30	0.52
2:B:51:ILE:HD13	2:B:72:VAL:HG13	1.92	0.52
2:H:125:LEU:HD21	2:H:142:LEU:HB2	1.92	0.52
1:A:78:MET:HE1	1:A:104:VAL:CG1	2.39	0.51
1:L:7:THR:CG2	1:L:8:PRO:HA	2.40	0.51
1:A:83:LEU:O	1:A:84:ALA:HB2	2.09	0.51
2:B:124:PRO:CB	2:B:212:ILE:HD12	2.40	0.51
2:H:157:SER:N	2:H:198:ASN:HD21	2.07	0.51
1:A:19:VAL:HG21	1:A:78:MET:CE	2.40	0.51
1:A:155:GLN:HB3	1:A:158:ASN:HD21	1.73	0.50
1:L:119:PRO:HG3	1:L:209:PHE:CD2	2.47	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:38:LYS:HB2	2:B:48:ILE:HD11	1.92	0.50
2:B:125:LEU:HD21	2:B:142:LEU:HB2	1.92	0.50
1:L:147:GLN:CG	1:L:154:LEU:HD13	2.42	0.50
1:A:124:GLN:NE2	1:A:131:SER:HB2	2.26	0.50
1:L:3:VAL:CG2	1:L:26:SER:HB3	2.41	0.50
2:H:12:VAL:HG21	2:H:18:VAL:HG13	1.93	0.50
1:A:147:GLN:CG	1:A:154:LEU:HD13	2.42	0.50
2:H:38:LYS:HB2	2:H:48:ILE:HD11	1.94	0.50
1:L:49:TYR:CZ	3:H:217:MMP:H183	2.47	0.50
1:L:43:SER:HB2	1:L:44:PRO:CD	2.42	0.50
2:H:51:ILE:CD1	2:H:72:VAL:HG13	2.42	0.49
1:A:7:THR:CG2	1:A:8:PRO:HA	2.42	0.49
1:A:43:SER:HB2	1:A:44:PRO:CD	2.41	0.49
1:L:150:VAL:HG13	1:L:150:VAL:O	2.11	0.49
1:L:2:LEU:HD23	1:L:29:VAL:HG12	1.95	0.49
2:H:124:PRO:CB	2:H:212:ILE:HD12	2.42	0.49
2:H:33:TRP:CD1	2:H:33:TRP:N	2.79	0.49
2:B:120:PRO:CB	2:B:146:TYR:HB3	2.41	0.49
1:L:19:VAL:HG21	1:L:78:MET:CE	2.42	0.49
1:A:3:VAL:CG2	1:A:26:SER:HB3	2.41	0.49
1:A:7:THR:N	4:A:216:HOH:O	2.46	0.49
2:B:157:SER:N	2:B:198:ASN:HD21	2.06	0.49
2:B:33:TRP:CD1	2:B:33:TRP:N	2.81	0.48
1:A:150:VAL:HG13	1:A:150:VAL:O	2.13	0.48
2:H:127:PRO:HG3	2:H:139:LEU:HB2	1.96	0.48
1:L:11:MET:HG3	1:L:12:SER:N	2.28	0.48
1:A:190:LYS:NZ	1:A:211:ARG:NH2	2.61	0.48
1:A:13:THR:HB	1:A:17:ASP:HB3	1.96	0.48
2:B:17:SER:HA	2:B:86:LEU:HD23	1.96	0.48
1:L:124:GLN:NE2	1:L:131:SER:H	2.11	0.48
1:L:190:LYS:NZ	1:L:211:ARG:NH2	2.62	0.48
1:L:136:LEU:N	1:L:136:LEU:HD12	2.29	0.47
2:B:128:SER:C	2:B:130:LYS:H	2.17	0.47
1:A:115:VAL:HG23	1:A:207:LYS:HG3	1.95	0.47
2:H:39:GLN:O	2:H:40:ARG:HB3	2.14	0.47
1:A:146:VAL:O	1:A:146:VAL:HG23	2.14	0.47
1:L:13:THR:CG2	1:L:19:VAL:HG22	2.45	0.47
1:A:198:HIS:C	1:A:200:GLY:H	2.17	0.47
1:L:116:PHE:HD2	2:H:132:THR:HA	1.79	0.47
1:L:147:GLN:HG2	1:L:154:LEU:HD13	1.95	0.47
1:L:124:GLN:NE2	1:L:131:SER:HB2	2.25	0.47
2:B:39:GLN:O	2:B:40:ARG:HB3	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:H:201:HIS:HD2	2:H:203:PRO:HD2	1.80	0.47
1:L:198:HIS:C	1:L:200:GLY:H	2.16	0.47
1:L:78:MET:HE1	1:L:104:VAL:CG1	2.45	0.47
1:L:188:LYS:NZ	1:L:189:HIS:CE1	2.83	0.47
2:H:124:PRO:HB3	2:H:212:ILE:CD1	2.45	0.46
1:A:113:PRO:HB3	1:A:139:PHE:CG	2.51	0.46
1:A:181:LEU:HD12	1:A:181:LEU:C	2.35	0.46
1:L:188:LYS:HZ2	1:L:189:HIS:CE1	2.33	0.46
2:H:17:SER:HA	2:H:86:LEU:HD23	1.98	0.46
2:B:97:THR:OG1	2:B:101:MET:HG2	2.16	0.46
1:A:160:GLN:HE22	2:B:172:GLN:HA	1.81	0.46
2:B:51:ILE:CD1	2:B:72:VAL:HG13	2.46	0.46
1:A:140:TYR:CG	1:A:141:PRO:HA	2.51	0.46
1:A:186:TYR:O	1:A:211:ARG:HD3	2.15	0.46
2:H:120:PRO:CB	2:H:146:TYR:HB3	2.42	0.46
1:L:140:TYR:CG	1:L:141:PRO:HA	2.51	0.46
1:L:78:MET:CE	1:L:104:VAL:HG11	2.44	0.46
2:B:124:PRO:HB3	2:B:212:ILE:CD1	2.46	0.46
2:H:14:PRO:HD3	2:H:113:SER:O	2.16	0.46
1:L:35:TRP:CZ3	1:L:88:CYS:HB2	2.51	0.46
1:A:2:LEU:HD23	1:A:29:VAL:HG12	1.98	0.46
2:H:128:SER:C	2:H:130:LYS:H	2.18	0.46
1:A:33:VAL:HG21	1:A:71:PHE:CZ	2.50	0.46
2:B:127:PRO:HG3	2:B:139:LEU:HB2	1.98	0.45
2:B:201:HIS:HD2	2:B:203:PRO:HD2	1.77	0.45
1:L:146:VAL:HG23	1:L:146:VAL:O	2.16	0.45
1:A:136:LEU:HD12	1:A:136:LEU:N	2.31	0.45
1:L:79:GLN:HE21	1:L:79:GLN:HB3	1.56	0.45
1:A:124:GLN:NE2	1:A:131:SER:H	2.14	0.45
1:A:197:THR:CG2	1:A:204:PRO:HB3	2.47	0.45
1:A:33:VAL:HG21	1:A:71:PHE:CE2	2.51	0.45
1:L:36:TYR:OH	1:L:89:GLN:OE1	2.26	0.45
2:B:8:GLY:O	2:B:10:GLU:HG3	2.17	0.45
1:A:134:CYS:HB2	1:A:148:TRP:CZ2	2.52	0.45
1:L:33:VAL:HG21	1:L:71:PHE:CZ	2.51	0.45
1:A:153:ALA:O	1:A:155:GLN:NE2	2.49	0.45
2:B:86:LEU:HD12	2:B:90:ASP:CB	2.46	0.45
1:A:78:MET:CE	1:A:104:VAL:HG11	2.46	0.45
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.98	0.45
2:H:86:LEU:HD12	2:H:90:ASP:CB	2.47	0.44
1:A:147:GLN:HG2	1:A:154:LEU:HD13	1.98	0.44
1:L:134:CYS:HB2	1:L:148:TRP:CZ2	2.53	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:L:197:THR:CG2	1:L:204:PRO:HB3	2.47	0.44
2:B:98:ARG:HG2	2:B:99:ARG:H	1.83	0.44
1:L:4:MET:HE1	1:L:25:ALA:HA	1.99	0.44
1:A:2:LEU:H	1:A:2:LEU:CD1	2.30	0.44
1:A:11:MET:HG3	1:A:12:SER:N	2.31	0.44
2:H:33:TRP:H	2:H:33:TRP:HD1	1.66	0.44
1:L:33:VAL:HG21	1:L:71:PHE:CE2	2.52	0.44
2:H:68:ALA:HA	2:H:82:GLN:O	2.18	0.44
2:H:187:SER:O	2:H:190:LEU:HD23	2.18	0.44
2:H:97:THR:OG1	2:H:101:MET:HG2	2.18	0.43
2:H:153:VAL:CG2	4:H:219:HOH:O	2.52	0.43
1:A:35:TRP:CZ3	1:A:88:CYS:HB2	2.53	0.43
2:H:8:GLY:O	2:H:10:GLU:HG3	2.18	0.43
1:L:21:ILE:O	1:L:72:THR:HA	2.18	0.43
1:A:162:SER:OG	2:B:168:PRO:HD2	2.18	0.43
1:A:170:ASP:HB3	1:A:172:THR:HG23	1.99	0.43
1:L:191:VAL:HG12	1:L:191:VAL:O	2.19	0.43
1:L:89:GLN:CG	1:L:90:GLN:N	2.82	0.43
2:B:68:ALA:HA	2:B:82:GLN:O	2.18	0.43
2:H:36:TRP:CD1	2:H:70:LEU:HD22	2.53	0.43
1:A:188:LYS:NZ	1:A:189:HIS:CE1	2.86	0.43
1:L:13:THR:HB	1:L:17:ASP:HB3	2.00	0.42
1:L:2:LEU:N	1:L:2:LEU:CD1	2.82	0.42
1:A:4:MET:HE1	1:A:25:ALA:HA	1.99	0.42
1:A:89:GLN:CG	1:A:90:GLN:N	2.82	0.42
2:B:33:TRP:CE3	2:B:50:MET:SD	3.12	0.42
1:L:135:LEU:CD1	1:L:137:ASN:ND2	2.80	0.42
1:L:1:GLU:O	1:L:2:LEU:C	2.58	0.42
1:A:188:LYS:HZ2	1:A:189:HIS:CE1	2.37	0.42
1:L:116:PHE:CD2	2:H:132:THR:HA	2.55	0.42
1:A:210:ASN:O	1:A:211:ARG:C	2.58	0.42
1:L:46:LEU:HD12	1:L:47:LEU:H	1.84	0.42
2:H:163:GLY:O	2:H:183:VAL:HA	2.19	0.42
2:B:17:SER:HA	2:B:86:LEU:CD2	2.50	0.42
1:L:11:MET:CG	1:L:12:SER:H	2.33	0.42
2:H:33:TRP:CE3	2:H:50:MET:SD	3.13	0.42
2:H:98:ARG:HG2	2:H:99:ARG:H	1.84	0.42
1:A:164:THR:HG23	2:B:167:PHE:CE1	2.54	0.42
1:L:210:ASN:O	1:L:211:ARG:C	2.57	0.42
1:A:189:HIS:HD1	1:A:189:HIS:H	1.67	0.42
2:B:128:SER:C	2:B:130:LYS:N	2.73	0.42
1:L:120:PRO:HD3	1:L:132:VAL:HG22	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:36:TRP:CD1	2:B:70:LEU:HD22	2.54	0.42
2:H:50:MET:HE2	2:H:50:MET:HB2	1.89	0.41
1:A:21:ILE:O	1:A:72:THR:HA	2.19	0.41
1:L:186:TYR:O	1:L:211:ARG:HD3	2.19	0.41
1:L:154:LEU:HD12	1:L:155:GLN:N	2.35	0.41
1:A:150:VAL:O	1:A:151:ASP:C	2.58	0.41
1:L:154:LEU:HD12	1:L:155:GLN:H	1.86	0.41
1:A:11:MET:HE2	1:A:12:SER:H	1.84	0.41
1:A:46:LEU:HD12	1:A:47:LEU:H	1.84	0.41
2:H:155:TRP:O	2:H:156:ASN:C	2.59	0.41
2:H:16:ALA:O	2:H:86:LEU:HD23	2.21	0.41
1:A:29:VAL:CG2	1:A:30:GLY:N	2.84	0.41
1:A:118:PHE:HA	1:A:119:PRO:HD3	1.68	0.41
1:A:28:ASN:CA	1:A:69:THR:HG22	2.45	0.41
2:B:179:LEU:HD23	2:B:179:LEU:C	2.41	0.41
2:H:2:VAL:HB	2:H:103:TYR:CE1	2.55	0.41
1:A:2:LEU:N	1:A:2:LEU:CD1	2.84	0.41
1:L:153:ALA:O	1:L:155:GLN:NE2	2.54	0.41
1:A:120:PRO:HD3	1:A:132:VAL:HG22	2.03	0.41
1:A:13:THR:CG2	1:A:19:VAL:HG22	2.47	0.41
2:B:16:ALA:O	2:B:86:LEU:HD23	2.21	0.41
1:A:188:LYS:HB2	1:A:188:LYS:HZ2	1.85	0.41
1:A:140:TYR:CD1	1:A:141:PRO:HA	2.56	0.41
1:A:142:ARG:HB3	1:A:173:TYR:CD2	2.55	0.41
1:L:29:VAL:CG2	1:L:30:GLY:N	2.85	0.41
1:A:65:SER:OG	1:A:72:THR:HG23	2.21	0.40
1:L:188:LYS:HB2	1:L:188:LYS:HZ2	1.86	0.40
2:B:155:TRP:CZ3	2:B:197:CYS:HB3	2.56	0.40
1:A:11:MET:CG	1:A:12:SER:H	2.34	0.40
1:L:65:SER:OG	1:L:72:THR:HG23	2.22	0.40
2:B:126:ALA:HB1	2:B:215:LYS:HB2	2.04	0.40
2:B:100:ASP:OD1	3:B:217:MMP:N2	2.53	0.40
1:L:160:GLN:HE22	2:H:172:GLN:HA	1.86	0.40
2:B:83:LEU:HB3	2:B:86:LEU:HD11	2.03	0.40
1:L:213:GLU:HG2	1:L:213:GLU:OXT	2.22	0.40
2:H:156:ASN:N	2:H:156:ASN:HD22	2.18	0.40
1:A:135:LEU:CD1	1:A:137:ASN:ND2	2.83	0.40
1:L:140:TYR:CD1	1:L:141:PRO:HA	2.56	0.40
2:H:128:SER:C	2:H:130:LYS:N	2.74	0.40

All (1) 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	Distance(Å)	Clash(Å)
1:A:156:SER:O	1:A:156:SER:O[2.665]	1.76	0.44

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	211/213 (99%)	181 (86%)	23 (11%)	7 (3%)	6	9
1	L	211/213 (99%)	181 (86%)	23 (11%)	7 (3%)	6	9
2	B	214/216 (99%)	183 (86%)	22 (10%)	9 (4%)	4	5
2	H	214/216 (99%)	181 (85%)	26 (12%)	7 (3%)	6	9
All	All	850/858 (99%)	726 (85%)	94 (11%)	30 (4%)	6	8

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	2	LEU
2	H	54	ASN
2	H	62	GLU
1	A	152	ASN
2	B	42	GLY
2	B	54	ASN
2	B	62	GLU
1	L	138	ASN
1	L	152	ASN
1	L	211	ARG
2	H	42	GLY
2	H	75	PRO
1	A	138	ASN
1	A	211	ARG
2	B	75	PRO
1	L	84	ALA
1	L	151	ASP
2	H	40	ARG
2	H	41	PRO
2	H	44	GLY

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	84	ALA
1	A	151	ASP
2	B	40	ARG
2	B	44	GLY
2	B	115	ALA
1	A	156	SER
2	B	7	SER
2	B	41	PRO
1	L	31	THR
1	A	31	THR

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	189/189 (100%)	183 (97%)	6 (3%)	51	80
1	L	189/189 (100%)	181 (96%)	8 (4%)	40	71
2	B	185/185 (100%)	174 (94%)	11 (6%)	28	52
2	H	185/185 (100%)	176 (95%)	9 (5%)	35	62
All	All	748/748 (100%)	714 (96%)	34 (4%)	38	67

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	GLU
1	L	56	THR
1	L	72	THR
1	L	88	CYS
1	L	90	GLN
1	L	131	SER
1	L	169	LYS
1	L	197	THR
2	H	28	THR
2	H	71	THR
2	H	121	SER
2	H	136	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	150	PRO
2	H	153	VAL
2	H	179	LEU
2	H	200	ASN
2	H	213	VAL
1	A	56	THR
1	A	72	THR
1	A	88	CYS
1	A	131	SER
1	A	169	LYS
1	A	197	THR
2	B	28	THR
2	B	71	THR
2	B	111	THR
2	B	117	THR
2	B	121	SER
2	B	136	THR
2	B	150	PRO
2	B	153	VAL
2	B	179	LEU
2	B	200	ASN
2	B	213	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	42	GLN
1	L	79	GLN
1	L	90	GLN
1	L	124	GLN
1	L	137	ASN
1	L	147	GLN
1	L	152	ASN
1	L	155	GLN
1	L	189	HIS
1	L	212	ASN
2	H	39	GLN
2	H	156	ASN
2	H	198	ASN
2	H	205	ASN
1	A	42	GLN
1	A	79	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	90	GLN
1	A	124	GLN
1	A	137	ASN
1	A	147	GLN
1	A	152	ASN
1	A	155	GLN
1	A	189	HIS
1	A	212	ASN
2	B	39	GLN
2	B	156	ASN
2	B	198	ASN
2	B	205	ASN

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 ⓘ

2 ligands are modelled in this entry.

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	$\# Z > 2$
3	MMP	B	217	-	46,46,47	3.11	22 (47%)	60,68,70	1.11	4 (6%)
3	MMP	H	217	-	46,46,47	3.12	20 (43%)	60,68,70	1.34	5 (8%)

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
3	MMP	B	217	-	-	1/30/78/82	0/0/5/5
3	MMP	H	217	-	-	1/30/78/82	0/0/5/5

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	217	MMP	C25-C24	14.19	1.44	1.35
3	H	217	MMP	C25-C24	13.26	1.44	1.35
3	H	217	MMP	C35-C41	8.49	1.40	1.35
3	B	217	MMP	C35-C41	6.78	1.39	1.35
3	B	217	MMP	C37-C33	4.79	1.60	1.52
3	H	217	MMP	O42-C49	4.61	1.38	1.22
3	H	217	MMP	C37-C33	4.59	1.60	1.52
3	B	217	MMP	O32-C39	4.52	1.38	1.22
3	H	217	MMP	O32-C39	4.47	1.38	1.22
3	B	217	MMP	O42-C49	4.45	1.38	1.22
3	H	217	MMP	C45-C11	3.89	1.44	1.39
3	H	217	MMP	C47-C42	3.51	1.60	1.51
3	B	217	MMP	C45-C11	3.43	1.43	1.39
3	H	217	MMP	C15-C14	3.31	1.43	1.39
3	B	217	MMP	C47-C42	3.26	1.60	1.51
3	B	217	MMP	C46-C43	2.82	1.57	1.51
3	H	217	MMP	C46-C43	2.65	1.56	1.51
3	H	217	MMP	C16-C12	2.63	1.55	1.47
3	B	217	MMP	C31-N3	2.63	1.39	1.36
3	H	217	MMP	C48-C49	2.54	1.57	1.50
3	B	217	MMP	C15-C14	2.52	1.42	1.39
3	H	217	MMP	C48-C47	2.47	1.59	1.52
3	B	217	MMP	C34-C33	2.44	1.48	1.42
3	B	217	MMP	C31-C32	2.41	1.48	1.42
3	B	217	MMP	C38-C39	2.39	1.56	1.50
3	H	217	MMP	C34-C33	2.34	1.47	1.42
3	B	217	MMP	C48-C49	2.34	1.56	1.50
3	B	217	MMP	O41-C49	2.31	1.38	1.30
3	H	217	MMP	C31-C32	2.30	1.47	1.42
3	H	217	MMP	C36-C32	2.26	1.56	1.51
3	B	217	MMP	C16-C12	2.26	1.54	1.47
3	B	217	MMP	O31-C39	2.24	1.38	1.30
3	H	217	MMP	O41-C49	2.23	1.38	1.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	217	MMP	C36-C32	2.23	1.56	1.51
3	H	217	MMP	O31-C39	2.15	1.38	1.30
3	B	217	MMP	C27-C23	2.12	1.57	1.51
3	H	217	MMP	C17-C13	2.12	1.59	1.51
3	B	217	MMP	C48-C47	2.12	1.58	1.52
3	B	217	MMP	C13-C12	-2.08	1.40	1.43
3	H	217	MMP	C27-C23	2.04	1.56	1.51
3	B	217	MMP	C34-N3	2.04	1.38	1.36
3	H	217	MMP	C38-C39	2.01	1.55	1.50

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	217	MMP	C34-C35-C41	-5.26	119.66	129.92
3	B	217	MMP	C34-C35-C41	-3.47	123.15	129.92
3	H	217	MMP	C31-C25-C24	-3.43	123.22	129.92
3	H	217	MMP	C35-C41-N4	-3.03	123.00	128.59
3	B	217	MMP	C35-C41-N4	-2.95	123.14	128.59
3	H	217	MMP	C25-C24-N2	-2.76	123.49	128.59
3	B	217	MMP	C31-C25-C24	-2.62	124.82	129.92
3	H	217	MMP	C45-C44-N4	-2.39	120.45	124.99
3	B	217	MMP	C25-C24-N2	-2.21	124.52	128.59

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	217	MMP	C34-C35-C41-N4
3	B	217	MMP	C34-C35-C41-N4

There are no ring outliers.

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	213/213 (100%)	0.49	18 (8%) 11 8	25, 58, 109, 146	0
1	L	213/213 (100%)	0.25	7 (3%) 44 41	27, 53, 96, 140	0
2	B	216/216 (100%)	0.33	10 (4%) 31 27	30, 58, 93, 128	0
2	H	216/216 (100%)	0.07	6 (2%) 50 48	20, 50, 92, 125	0
All	All	858/858 (100%)	0.28	41 (4%) 29 26	20, 55, 99, 146	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	131	SER	6.5
1	A	213	GLU	6.3
2	B	216	SER	5.0
2	B	133	SER	4.4
1	L	213	GLU	4.2
2	H	133	SER	4.1
2	B	129	SER	4.0
1	L	200	GLY	4.0
1	A	152	ASN	3.7
1	A	191	VAL	3.7
1	A	202	SER	3.6
2	H	129	SER	3.3
2	B	130	LYS	3.3
2	H	216	SER	3.2
2	H	130	LYS	3.2
1	L	198	HIS	3.2
2	B	132	THR	3.1
1	A	190	LYS	3.1
1	A	203	SER	2.9
1	A	156	SER	2.9
1	A	192	TYR	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	193	ALA	2.8
1	A	134	CYS	2.6
1	L	201	LEU	2.6
1	A	186	TYR	2.6
1	A	154	LEU	2.6
1	L	204	PRO	2.6
1	A	204	PRO	2.5
1	L	10	PHE	2.4
2	B	41	PRO	2.4
1	A	210	ASN	2.3
1	A	200	GLY	2.3
1	A	118	PHE	2.3
1	L	199	GLN	2.2
2	B	55	SER	2.2
2	H	34	MET	2.1
2	H	5	LEU	2.1
1	A	132	VAL	2.1
2	B	62	GLU	2.0
1	A	212	ASN	2.0
2	B	215	LYS	2.0

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 ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MMP	H	217	42/43	0.29	2.42	48,96,130,135	0
3	MMP	B	217	42/43	0.26	1.38	29,81,109,115	0

6.5 Other polymers ⓘ

There are no such residues in this entry.