



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:34 AM GMT

PDB ID : 2P8L
Title : Crystal structure of the HIV-1 Cross Neutralizing Monoclonal Antibody 2F5
in complex with gp41 Peptide ELLELDKWASLWN
Authors : Julien, J.P.; Bryson, S.; Pai, E.F.
Deposited on : 2007-03-22
Resolution : 2.44 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>

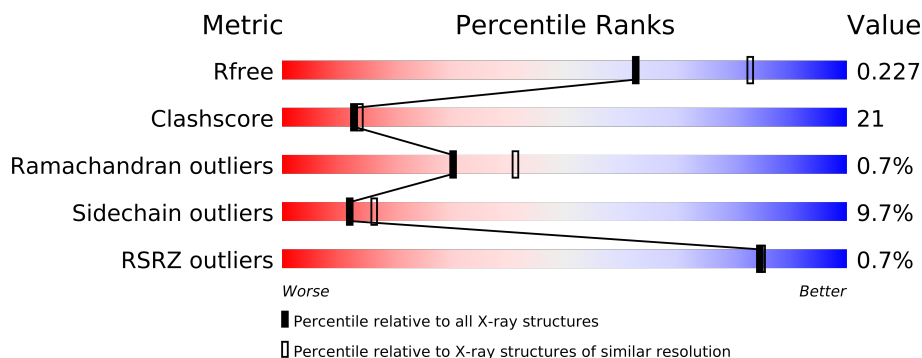
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.44 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	2989 (2.48-2.40)
Clashscore	79885	3698 (2.48-2.40)
Ramachandran outliers	78287	3639 (2.48-2.40)
Sidechain outliers	78261	3640 (2.48-2.40)
RSRZ outliers	66119	2993 (2.48-2.40)

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	214	
2	B	235	
3	C	13	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3677 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 nmAb 2F5, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1644	1025	281	333	5			

- Molecule 2 is a protein called nmAb 2F5, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	222	Total	C	N	O	S	0	0	0
			1674	1067	283	318	6			

- Molecule 3 is a protein called gp41 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	11	Total	C	N	O	0	0	0
			92	61	13	18			

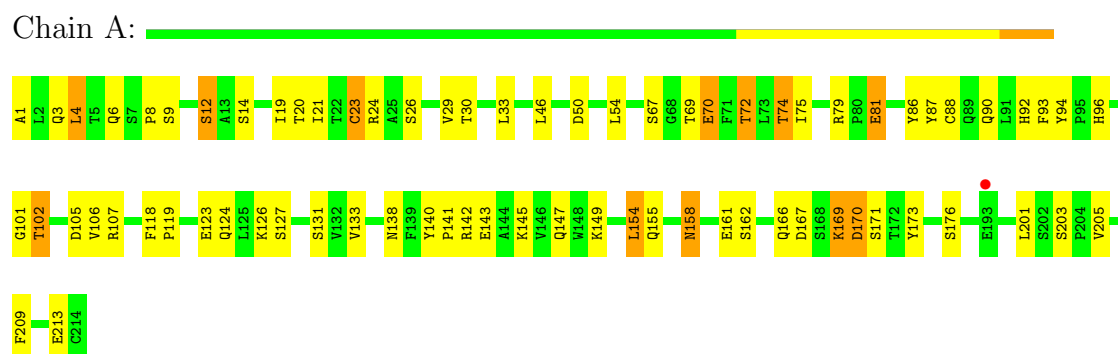
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	131	Total	O	0	0
			131	131		
4	C	6	Total	O	0	0
			6	6		

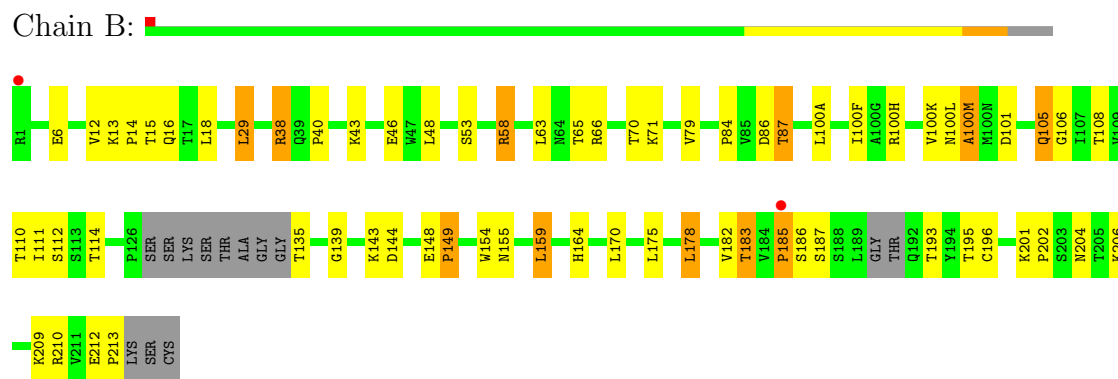
3 Residue-property plots

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: nmAb 2F5, light chain



- Molecule 2: nmAb 2F5, heavy chain



- Molecule 3: gp41 peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.00Å 64.70Å 177.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.69 – 2.44 48.57 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.5 (43.69-2.44) 68.9 (48.57-1.95)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.94 (at 1.95Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.201 , 0.225 0.202 , 0.227	Depositor DCC
R_{free} test set	1227 reflections (4.90%)	DCC
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.242	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 44.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 35271 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3677	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.43% 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	0.47	0/1681	0.76	0/2284
2	B	0.47	0/1715	0.81	0/2346
3	C	1.13	0/93	1.12	1/125 (0.8%)
All	All	0.50	0/3489	0.80	1/4755 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	LEU	CA-CB-CG	-7.70	97.59	115.30

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	1644	0	1588	77	1
2	B	1674	0	1687	65	0
3	C	92	0	95	7	0
4	A	130	0	0	20	1
4	B	131	0	0	16	2
4	C	6	0	0	0	0
All	All	3677	0	3370	142	2

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 21.

All (142) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:58:ARG:NH1	3:C:3:GLU:OE1	1.91	1.03
2:B:201:LYS:HE3	4:B:324:HOH:O	1.66	0.96
3:C:0:GLU:O	3:C:1:LEU:HB2	1.67	0.92
2:B:84:PRO:O	2:B:87:THR:HG23	1.72	0.90
1:A:90:GLN:HE22	1:A:93:PHE:H	0.95	0.89
1:A:90:GLN:NE2	1:A:93:PHE:H	1.71	0.87
2:B:66:ARG:HD3	4:B:309:HOH:O	1.75	0.86
1:A:90:GLN:HE22	1:A:93:PHE:N	1.77	0.83
1:A:67:SER:HB3	4:A:312:HOH:O	1.78	0.82
3:C:0:GLU:O	3:C:1:LEU:CB	2.29	0.81
1:A:24:ARG:NE	1:A:70:GLU:OE2	2.14	0.80
1:A:6:GLN:HG2	1:A:23:CYS:SG	2.20	0.80
2:B:108:THR:HG23	4:B:320:HOH:O	1.81	0.79
1:A:8:PRO:O	1:A:102:THR:HB	1.84	0.78
2:B:38:ARG:HD2	2:B:48:LEU:HD21	1.65	0.78
1:A:81:GLU:HG3	4:A:308:HOH:O	1.85	0.77
2:B:40:PRO:HB2	2:B:43:LYS:HG3	1.66	0.75
1:A:90:GLN:NE2	1:A:92:HIS:H	1.84	0.74
2:B:155:ASN:OD1	4:B:310:HOH:O	2.09	0.71
1:A:90:GLN:HE21	1:A:92:HIS:H	1.38	0.70
2:B:38:ARG:HB3	2:B:48:LEU:HD11	1.73	0.70
2:B:154:TRP:CH2	2:B:196:CYS:SG	2.85	0.69
1:A:201:LEU:HD13	1:A:205:VAL:HG12	1.73	0.69
1:A:6:GLN:CG	1:A:23:CYS:SG	2.81	0.68
1:A:19:ILE:HD11	1:A:75:ILE:HD12	1.75	0.68
2:B:209:LYS:HE2	4:B:315:HOH:O	1.92	0.67
1:A:20:THR:HG22	1:A:74:THR:HB	1.77	0.67
2:B:209:LYS:CG	4:B:315:HOH:O	2.42	0.67
1:A:155:GLN:HB3	1:A:158:ASN:HD21	1.60	0.66
1:A:107:ARG:HD2	4:A:305:HOH:O	1.94	0.66
2:B:209:LYS:HG3	4:B:315:HOH:O	1.95	0.66
1:A:203:SER:HB2	4:A:341:HOH:O	1.93	0.66
1:A:12:SER:HB2	1:A:107:ARG:HB2	1.78	0.66
1:A:4:LEU:HD13	1:A:88:CYS:SG	2.37	0.65
1:A:81:GLU:CD	1:A:81:GLU:H	1.99	0.64
1:A:94:TYR:CZ	2:B:58:ARG:HG3	2.32	0.64
2:B:193:THR:O	4:B:310:HOH:O	2.14	0.64
1:A:29:VAL:HG11	1:A:90:GLN:HG3	1.79	0.64
1:A:1:ALA:HA	3:C:0:GLU:OE2	1.98	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:206:LYS:HB2	4:B:337:HOH:O	1.98	0.62
1:A:123:GLU:HB2	4:A:302:HOH:O	1.98	0.61
1:A:20:THR:HB	4:A:343:HOH:O	2.00	0.61
2:B:135:THR:N	4:B:308:HOH:O	2.32	0.61
1:A:96:HIS:C	4:A:316:HOH:O	2.37	0.61
2:B:114:THR:HG21	2:B:175:LEU:HD23	1.83	0.60
3:C:0:GLU:O	3:C:0:GLU:HG3	2.01	0.59
2:B:100(H):ARG:O	2:B:100(K):VAL:HG22	2.02	0.59
2:B:87:THR:HB	2:B:110:THR:HA	1.84	0.58
2:B:13:LYS:NZ	4:B:321:HOH:O	2.37	0.58
1:A:90:GLN:NE2	1:A:92:HIS:N	2.52	0.57
1:A:81:GLU:CG	4:A:308:HOH:O	2.49	0.57
1:A:74:THR:HG22	4:A:281:HOH:O	2.03	0.57
1:A:29:VAL:HG11	1:A:90:GLN:CG	2.34	0.57
2:B:12:VAL:HG12	2:B:16:GLN:HE21	1.68	0.57
1:A:79:ARG:HB3	1:A:81:GLU:OE2	2.05	0.57
2:B:178:LEU:HD12	2:B:178:LEU:C	2.26	0.56
1:A:124:GLN:O	1:A:127:SER:OG	2.17	0.56
1:A:67:SER:CB	4:A:312:HOH:O	2.47	0.55
1:A:79:ARG:HD2	4:A:321:HOH:O	2.06	0.55
1:A:170:ASP:O	1:A:171:SER:HB2	2.07	0.55
1:A:21:ILE:HG23	1:A:102:THR:HG21	1.89	0.55
2:B:183:THR:HG22	4:B:343:HOH:O	2.07	0.55
1:A:169:LYS:HB3	1:A:169:LYS:NZ	2.22	0.54
1:A:94:TYR:CE2	2:B:58:ARG:HG3	2.42	0.54
1:A:106:VAL:H	1:A:166:GLN:HE22	1.56	0.54
1:A:24:ARG:CD	1:A:70:GLU:OE2	2.56	0.53
1:A:118:PHE:HD2	1:A:133:VAL:HG22	1.73	0.53
1:A:12:SER:HB3	1:A:140:TYR:OH	2.09	0.53
1:A:158:ASN:ND2	1:A:158:ASN:H	2.07	0.53
2:B:209:LYS:CB	4:B:315:HOH:O	2.56	0.52
2:B:65:THR:OG1	4:B:307:HOH:O	2.18	0.52
2:B:13:LYS:O	2:B:16:GLN:HB2	2.09	0.52
1:A:167:ASP:HB3	1:A:170:ASP:HB2	1.92	0.52
1:A:147:GLN:NE2	1:A:154:LEU:HG	2.25	0.52
2:B:12:VAL:CG1	2:B:16:GLN:HE21	2.23	0.51
1:A:161:GLU:HA	1:A:176:SER:O	2.11	0.51
2:B:38:ARG:HD3	2:B:46:GLU:OE1	2.11	0.51
1:A:6:GLN:HE22	1:A:87:TYR:HA	1.75	0.51
2:B:66:ARG:NH2	2:B:86:ASP:OD2	2.43	0.51
2:B:13:LYS:H	2:B:16:GLN:NE2	2.08	0.50
1:A:105:ASP:HB2	1:A:166:GLN:NE2	2.25	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:158:ASN:H	1:A:158:ASN:HD22	1.58	0.50
2:B:70:THR:HG22	2:B:79:VAL:HB	1.94	0.50
2:B:100(L):ASN:O	2:B:100(M):ALA:HB2	2.11	0.50
2:B:100(A):LEU:HB2	2:B:100(F):ILE:HD13	1.93	0.49
1:A:72:THR:HB	4:A:343:HOH:O	2.12	0.49
1:A:140:TYR:CG	1:A:141:PRO:HA	2.47	0.49
1:A:213:GLU:HA	4:A:309:HOH:O	2.13	0.48
1:A:213:GLU:CA	4:A:309:HOH:O	2.59	0.48
2:B:212:GLU:HB2	2:B:213:PRO:HD2	1.95	0.48
1:A:1:ALA:HA	3:C:0:GLU:CD	2.34	0.48
2:B:84:PRO:HA	2:B:111:ILE:HB	1.96	0.48
2:B:29:LEU:O	2:B:53:SER:HB2	2.14	0.47
1:A:119:PRO:HB3	1:A:209:PHE:CZ	2.49	0.47
1:A:169:LYS:HG3	4:A:323:HOH:O	2.15	0.47
2:B:185:PRO:C	2:B:187:SER:N	2.67	0.47
2:B:84:PRO:O	2:B:87:THR:CG2	2.53	0.47
1:A:124:GLN:HE22	1:A:131:SER:HB2	1.80	0.47
1:A:21:ILE:CG2	1:A:102:THR:HG21	2.45	0.46
2:B:185:PRO:O	2:B:187:SER:N	2.49	0.46
1:A:3:GLN:HB2	1:A:26:SER:HB3	1.99	0.45
1:A:142:ARG:HD2	1:A:173:TYR:CE2	2.52	0.45
1:A:24:ARG:HA	1:A:69:THR:O	2.17	0.45
1:A:46:LEU:HD22	2:B:101:ASP:HA	1.99	0.45
2:B:38:ARG:NH2	2:B:86:ASP:OD1	2.50	0.44
1:A:158:ASN:N	1:A:158:ASN:ND2	2.64	0.44
1:A:6:GLN:HG3	1:A:23:CYS:SG	2.57	0.44
1:A:72:THR:CB	4:A:343:HOH:O	2.66	0.44
1:A:142:ARG:HB2	1:A:173:TYR:CE1	2.52	0.43
2:B:105:GLN:HG2	2:B:106:GLY:N	2.33	0.43
1:A:90:GLN:NE2	1:A:93:PHE:N	2.49	0.43
2:B:209:LYS:CE	4:B:315:HOH:O	2.59	0.43
2:B:159:LEU:HD13	2:B:182:VAL:HG21	1.99	0.43
2:B:66:ARG:HH22	2:B:86:ASP:CG	2.22	0.43
2:B:66:ARG:CD	4:B:309:HOH:O	2.49	0.43
1:A:166:GLN:NE2	4:A:337:HOH:O	2.52	0.43
2:B:201:LYS:HB2	2:B:202:PRO:HD3	2.01	0.42
2:B:185:PRO:C	2:B:187:SER:H	2.21	0.42
2:B:14:PRO:O	2:B:15:THR:OG1	2.24	0.42
1:A:86:TYR:O	1:A:101:GLY:HA2	2.19	0.42
1:A:1:ALA:HB3	4:A:304:HOH:O	2.18	0.42
2:B:210:ARG:NH1	2:B:212:GLU:OE1	2.52	0.42
2:B:29:LEU:HA	2:B:29:LEU:HD12	1.75	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:6:GLU:OE1	2:B:6:GLU:N	2.50	0.42
1:A:127:SER:N	4:A:328:HOH:O	2.53	0.42
2:B:212:GLU:HA	2:B:213:PRO:HD3	1.80	0.41
2:B:143:LYS:HE3	2:B:144:ASP:OD2	2.20	0.41
3:C:1:LEU:HD13	3:C:1:LEU:C	2.40	0.41
2:B:12:VAL:HG21	2:B:18:LEU:HD13	2.01	0.41
1:A:138:ASN:OD1	2:B:164:HIS:HE1	2.03	0.41
2:B:43:LYS:HE2	2:B:43:LYS:HB3	1.88	0.41
1:A:118:PHE:CD2	1:A:133:VAL:HG22	2.54	0.41
2:B:148:GLU:HA	2:B:149:PRO:HA	1.80	0.41
2:B:139:GLY:HA2	2:B:154:TRP:CH2	2.56	0.41
2:B:159:LEU:HD23	2:B:159:LEU:HA	1.78	0.41
2:B:143:LYS:HG2	2:B:144:ASP:N	2.36	0.41
1:A:105:ASP:OD1	1:A:105:ASP:C	2.60	0.41
1:A:29:VAL:O	1:A:30:THR:OG1	2.36	0.40
2:B:112:SER:C	2:B:114:THR:H	2.24	0.40
1:A:126:LYS:HB2	4:A:328:HOH:O	2.21	0.40
2:B:212:GLU:HB2	2:B:213:PRO:CD	2.51	0.40
1:A:201:LEU:CD1	1:A:205:VAL:HG12	2.45	0.40

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

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:ASP:OD2	4:B:329:HOH:O[4_566]	1.96	0.24
4:A:299:HOH:O	4:B:310:HOH:O[3_555]	2.05	0.15

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	212/214 (99%)	202 (95%)	10 (5%)	0	100	100
2	B	216/235 (92%)	207 (96%)	7 (3%)	2 (1%)	25	33
3	C	9/13 (69%)	8 (89%)	0	1 (11%)	1	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	437/462 (95%)	417 (95%)	17 (4%)	3 (1%)	30	41

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	1	LEU
2	B	100(M)	ALA
2	B	186	SER

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	187/187 (100%)	167 (89%)	20 (11%)	10	12
2	B	194/204 (95%)	179 (92%)	15 (8%)	18	27
3	C	10/12 (83%)	7 (70%)	3 (30%)	0	0
All	All	391/403 (97%)	353 (90%)	38 (10%)	12	16

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	9	SER
1	A	12	SER
1	A	14	SER
1	A	23	CYS
1	A	33	LEU
1	A	54	LEU
1	A	70	GLU
1	A	72	THR
1	A	74	THR
1	A	81	GLU
1	A	102	THR
1	A	143	GLU
1	A	145	LYS
1	A	149	LYS

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Mol	Chain	Res	Type
1	A	154	LEU
1	A	158	ASN
1	A	162	SER
1	A	169	LYS
1	A	170	ASP
2	B	29	LEU
2	B	38	ARG
2	B	58	ARG
2	B	63	LEU
2	B	71	LYS
2	B	87	THR
2	B	105	GLN
2	B	149	PRO
2	B	159	LEU
2	B	170	LEU
2	B	178	LEU
2	B	183	THR
2	B	185	PRO
2	B	195	THR
2	B	204	ASN
3	C	0	GLU
3	C	1	LEU
3	C	2	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	27	GLN
1	A	38	GLN
1	A	90	GLN
1	A	124	GLN
1	A	137	ASN
1	A	147	GLN
1	A	155	GLN
1	A	158	ASN
1	A	166	GLN
1	A	199	GLN
2	B	16	GLN
2	B	39	GLN
2	B	76	ASN
2	B	164	HIS

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Mol	Chain	Res	Type
2	B	192	GLN
2	B	199	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 ⓘ

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	214/214 (100%)	-0.39	1 (0%) 88 89	30, 42, 57, 65	0
2	B	222/235 (94%)	-0.29	2 (0%) 81 81	27, 40, 59, 82	0
3	C	11/13 (84%)	0.02	0 100 100	34, 50, 79, 80	0
All	All	447/462 (96%)	-0.33	3 (0%) 84 85	27, 41, 58, 82	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	185	PRO	4.5
2	B	1	ARG	2.2
1	A	193	GLU	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.