



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2018 – 11:19 am GMT

PDB ID : 3D0L
Title : Crystal structure of the HIV-1 broadly neutralizing antibody 2F5 in complex with the gp41 FP-MPER Hyb3K construct 514GIGALFLGFLGAAGS528KK-Ahx-655KNEQELLELDKWASLWN671
Authors : Bryson, S.; Julien, J.P.; Pai, E.F.
Deposited on : 2008-05-01
Resolution : 2.35 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

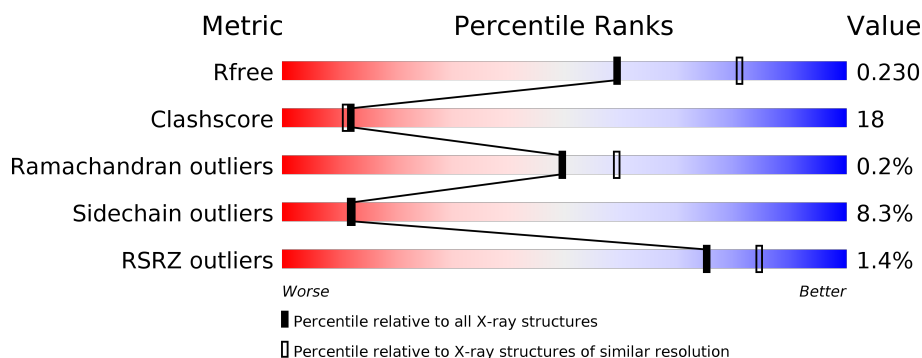
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	111664	1015 (2.36-2.36)
Clashscore	122126	1081 (2.36-2.36)
Ramachandran outliers	120053	1066 (2.36-2.36)
Sidechain outliers	120020	1067 (2.36-2.36)
RSRZ outliers	108989	1002 (2.36-2.36)

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. 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	214	
2	B	237	
3	C	35	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	B	6576	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3559 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 2F5 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1637	1022	280	331	4			

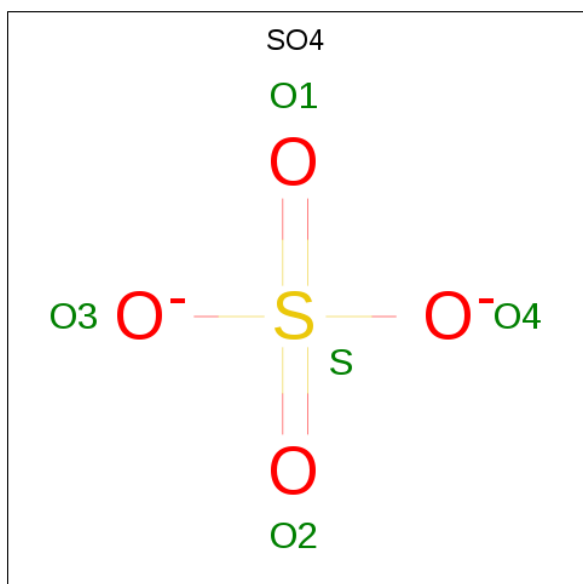
- Molecule 2 is a protein called 2F5 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	218	Total	C	N	O	S	0	0	0
			1639	1039	279	315	6			

- Molecule 3 is a protein called Hyb3k peptide construct.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			89	61	13	15			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 6 3 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	101	Total O 101 101	0	0
6	B	77	Total O 77 77	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.30Å 76.70Å 94.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.82 – 2.35 59.43 – 2.35	Depositor EDS
% Data completeness (in resolution range)	85.9 (48.82-2.35) 86.0 (59.43-2.35)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.34Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.211 , 0.229 0.211 , 0.230	Depositor DCC
R_{free} test set	2464 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.1	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3559	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, SO4

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	0.44	1/1674 (0.1%)	0.80	4/2276 (0.2%)
2	B	0.37	0/1678	0.73	1/2294 (0.0%)
3	C	0.79	0/92	1.80	2/125 (1.6%)
All	All	0.42	1/3444 (0.0%)	0.81	7/4695 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	81	GLU	CG-CD	5.66	1.60	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	10	LEU	CA-CB-CG	15.72	151.45	115.30
1	A	78	LEU	CA-CB-CG	11.12	140.88	115.30
1	A	213	GLU	N-CA-C	6.66	128.99	111.00
1	A	143	GLU	CA-CB-CG	6.55	127.80	113.40
3	C	11	TRP	N-CA-C	-5.58	95.94	111.00
2	B	178	LEU	CA-CB-CG	5.22	127.30	115.30
1	A	11	LEU	CA-CB-CG	5.08	126.97	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within

the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1637	0	1583	52	0
2	B	1639	0	1648	72	0
3	C	89	0	85	5	0
4	A	5	0	0	0	0
4	B	5	0	0	4	0
5	B	6	0	8	1	0
6	A	101	0	0	3	0
6	B	77	0	0	5	0
All	All	3559	0	3324	123	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 18.

All (123) 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)
2:B:1:ARG:HD3	2:B:2:ILE:H	1.22	1.05
1:A:37:ARG:NH1	1:A:81:GLU:HG3	1.84	0.93
2:B:1:ARG:CD	2:B:2:ILE:H	1.84	0.91
1:A:37:ARG:HH12	1:A:81:GLU:HG3	1.37	0.90
1:A:37:ARG:HH12	1:A:81:GLU:CG	1.90	0.83
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.59	0.83
2:B:100(H):ARG:HB2	4:B:6576:SO4:O2	1.79	0.82
2:B:99:THR:HB	2:B:100(G):ALA:CB	2.09	0.82
1:A:145:LYS:NZ	1:A:147:GLN:HG3	1.96	0.81
2:B:83:SER:OG	2:B:85:VAL:HG22	1.82	0.80
2:B:58:ARG:NE	3:C:3:GLU:OE2	2.12	0.80
1:A:81:GLU:HB3	6:A:6663:HOH:O	1.81	0.80
2:B:114:THR:HG21	2:B:175:LEU:CD2	2.14	0.76
1:A:145:LYS:HZ1	1:A:147:GLN:HG3	1.49	0.76
2:B:114:THR:HG21	2:B:175:LEU:HD23	1.71	0.73
2:B:1:ARG:CG	2:B:2:ILE:H	2.04	0.71
2:B:99:THR:HB	2:B:100(G):ALA:HB3	1.72	0.70
2:B:184:VAL:HB	2:B:185:PRO:HD2	1.74	0.68
2:B:56:ASP:OD2	3:C:6:LYS:HE3	1.94	0.68
2:B:99:THR:HB	2:B:100(G):ALA:HB1	1.74	0.68
1:A:22:THR:HG22	1:A:72:THR:HG22	1.75	0.68
2:B:99:THR:CB	2:B:100(G):ALA:HB1	2.23	0.68
2:B:54:ASP:OD1	3:C:6:LYS:NZ	2.25	0.68
2:B:105:GLN:HG3	6:B:6648:HOH:O	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:ARG:HD2	2:B:48:LEU:HD21	1.76	0.67
1:A:138:ASN:HD21	2:B:164:HIS:CE1	2.14	0.66
2:B:1:ARG:CG	2:B:2:ILE:N	2.57	0.65
1:A:2:LEU:HD22	1:A:90:GLN:HG2	1.79	0.65
1:A:54:LEU:CD2	1:A:58:VAL:HB	2.27	0.64
1:A:37:ARG:NH1	1:A:81:GLU:CG	2.56	0.64
2:B:1:ARG:HG2	2:B:2:ILE:N	2.12	0.63
1:A:132:VAL:HG13	1:A:179:LEU:HB3	1.81	0.62
2:B:123:PRO:HD3	2:B:209:LYS:HE2	1.81	0.62
1:A:118:PHE:HB2	1:A:133:VAL:HG22	1.81	0.61
2:B:40:PRO:HB2	2:B:43:LYS:HG3	1.82	0.61
1:A:191:VAL:HG22	6:A:6629:HOH:O	2.01	0.60
2:B:192:GLN:HB2	6:B:6629:HOH:O	2.01	0.60
2:B:146:PHE:CE1	2:B:147:PRO:HB3	2.36	0.60
2:B:38:ARG:HD3	2:B:46:GLU:OE1	2.02	0.60
1:A:77:THR:HG23	1:A:77:THR:O	2.02	0.60
1:A:149:LYS:HB2	1:A:193:GLU:HG3	1.85	0.59
2:B:100(L):ASN:O	2:B:100(M):ALA:HB2	2.04	0.58
2:B:13:LYS:O	2:B:16:GLN:HB2	2.04	0.57
2:B:212:GLU:HB2	2:B:213:PRO:HD2	1.87	0.57
1:A:28:GLY:O	6:A:6664:HOH:O	2.17	0.57
1:A:2:LEU:CD2	1:A:90:GLN:HG2	2.35	0.57
1:A:190:LYS:HE3	1:A:210:ASN:HB3	1.87	0.56
2:B:100(H):ARG:NH1	4:B:6576:SO4:O2	2.39	0.55
2:B:100(I):GLY:HA2	3:C:11:TRP:HZ2	1.70	0.55
2:B:1:ARG:HD3	2:B:2:ILE:N	2.06	0.55
2:B:114:THR:HG22	2:B:115:SER:N	2.21	0.54
1:A:124:GLN:HG2	1:A:129:THR:O	2.08	0.54
2:B:114:THR:HG21	2:B:175:LEU:HD21	1.88	0.54
2:B:100(I):GLY:N	2:B:100(J):PRO:HD2	2.23	0.54
2:B:192:GLN:NE2	6:B:6629:HOH:O	2.31	0.54
2:B:40:PRO:HB2	2:B:43:LYS:CG	2.38	0.54
2:B:99:THR:CB	2:B:100(G):ALA:CB	2.82	0.53
1:A:183:LYS:O	1:A:187:GLU:HG2	2.09	0.53
2:B:196:CYS:SG	2:B:209:LYS:HB3	2.48	0.53
1:A:159:SER:HA	1:A:178:THR:O	2.10	0.52
1:A:12:SER:HB3	1:A:140:TYR:OH	2.08	0.52
2:B:13:LYS:H	2:B:16:GLN:NE2	2.07	0.52
2:B:99:THR:OG1	2:B:100(G):ALA:HB1	2.10	0.52
2:B:178:LEU:HD12	2:B:178:LEU:C	2.30	0.51
2:B:100(H):ARG:CB	4:B:6576:SO4:O2	2.55	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:PHE:HD2	1:A:133:VAL:HG23	1.76	0.51
1:A:126:LYS:C	1:A:128:GLY:N	2.64	0.51
2:B:37:ILE:N	2:B:37:ILE:HD12	2.26	0.51
1:A:149:LYS:HB2	1:A:193:GLU:CG	2.40	0.50
2:B:114:THR:HG22	2:B:115:SER:O	2.12	0.50
1:A:145:LYS:HB2	1:A:197:THR:HB	1.95	0.49
2:B:13:LYS:HB2	2:B:16:GLN:NE2	2.28	0.49
2:B:31:ASP:HB2	2:B:34:VAL:HG21	1.95	0.49
2:B:43:LYS:O	6:B:6641:HOH:O	2.19	0.48
1:A:145:LYS:HZ3	1:A:147:GLN:HG3	1.78	0.48
1:A:34:ALA:O	1:A:88:CYS:HA	2.14	0.48
1:A:184:ALA:O	1:A:188:LYS:HG3	2.13	0.48
1:A:4:LEU:HA	1:A:24:ARG:O	2.12	0.48
2:B:100(G):ALA:N	4:B:6576:SO4:O4	2.47	0.48
1:A:54:LEU:HD21	1:A:58:VAL:HB	1.96	0.47
2:B:13:LYS:HB2	2:B:16:GLN:CD	2.34	0.47
2:B:114:THR:CG2	2:B:115:SER:N	2.77	0.47
2:B:82(B):ARG:O	2:B:82(B):ARG:HG3	2.15	0.47
1:A:211:ARG:HG2	1:A:211:ARG:HH11	1.80	0.46
1:A:138:ASN:HD21	2:B:164:HIS:HE1	1.58	0.46
1:A:193:GLU:HB3	1:A:208:SER:HB3	1.98	0.46
1:A:32:ALA:HB3	1:A:92:HIS:HB2	1.96	0.46
2:B:82(B):ARG:HG2	2:B:82(B):ARG:HH11	1.81	0.46
2:B:178:LEU:HA	5:B:3731:GOL:O3	2.15	0.45
1:A:115:VAL:HA	1:A:135:LEU:O	2.16	0.45
1:A:201:LEU:HD13	1:A:205:VAL:HG12	1.98	0.45
2:B:136:ALA:CB	2:B:189:LEU:HD11	2.47	0.45
1:A:118:PHE:HD2	1:A:133:VAL:CG2	2.30	0.45
1:A:132:VAL:HG22	1:A:148:TRP:CH2	2.52	0.44
1:A:145:LYS:HB3	1:A:145:LYS:HE3	1.41	0.44
2:B:114:THR:CG2	2:B:175:LEU:HD21	2.48	0.43
2:B:116:THR:HA	2:B:146:PHE:O	2.17	0.43
3:C:7:TRP:HB3	3:C:10:LEU:HD22	2.01	0.43
2:B:84:PRO:HA	2:B:111:ILE:HB	2.00	0.43
2:B:82(B):ARG:HG2	2:B:82(B):ARG:NH1	2.34	0.43
2:B:114:THR:HG23	6:B:6625:HOH:O	2.19	0.43
1:A:8:PRO:O	1:A:102:THR:HG23	2.19	0.42
1:A:140:TYR:CD1	1:A:141:PRO:HA	2.55	0.42
1:A:32:ALA:O	1:A:90:GLN:HA	2.19	0.42
2:B:151:THR:OG1	2:B:199:ASN:HB3	2.20	0.42
2:B:31:ASP:HB2	2:B:34:VAL:CG2	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:212:GLU:HB2	2:B:213:PRO:CD	2.49	0.42
2:B:51:ILE:HB	2:B:69:ILE:HD13	2.02	0.42
1:A:61:ARG:HD2	1:A:77:THR:O	2.20	0.41
1:A:79:ARG:HB3	1:A:80:PRO:HD2	2.02	0.41
1:A:167:ASP:HB3	1:A:170:ASP:OD2	2.21	0.41
2:B:24:PHE:CZ	2:B:29:LEU:HD22	2.54	0.41
2:B:96:ARG:O	2:B:100(K):VAL:HG12	2.20	0.41
2:B:126:PRO:HD3	2:B:138:LEU:HB3	2.01	0.41
2:B:144:ASP:HB3	2:B:175:LEU:HD13	2.02	0.41
1:A:79:ARG:HB3	1:A:80:PRO:CD	2.50	0.41
2:B:100(L):ASN:O	2:B:100(M):ALA:CB	2.69	0.41
1:A:21:ILE:HD11	1:A:73:LEU:HD23	2.03	0.41
2:B:184:VAL:HB	2:B:185:PRO:CD	2.47	0.41
2:B:54:ASP:O	2:B:55:ASP:CB	2.69	0.41
1:A:125:LEU:O	1:A:183:LYS:HD2	2.21	0.40
1:A:2:LEU:HD12	1:A:27:GLN:HB2	2.04	0.40
1:A:58:VAL:HA	1:A:59:PRO:HD3	1.99	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	211/214 (99%)	202 (96%)	9 (4%)	0	100	100
2	B	212/237 (90%)	199 (94%)	12 (6%)	1 (0%)	31	34
3	C	8/35 (23%)	8 (100%)	0	0	100	100
All	All	431/486 (89%)	409 (95%)	21 (5%)	1 (0%)	49	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	100(M)	ALA

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	186/186 (100%)	172 (92%)	14 (8%)	15	15
2	B	190/207 (92%)	175 (92%)	15 (8%)	13	13
3	C	9/25 (36%)	6 (67%)	3 (33%)	0	0
All	All	385/418 (92%)	353 (92%)	32 (8%)	12	12

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	12	SER
1	A	21	ILE
1	A	33	LEU
1	A	54	LEU
1	A	78	LEU
1	A	81	GLU
1	A	97	THR
1	A	132	VAL
1	A	145	LYS
1	A	152	ASN
1	A	191	VAL
1	A	203	SER
1	A	213	GLU
2	B	1	ARG
2	B	5	LYS
2	B	11	LEU
2	B	29	LEU
2	B	38	ARG
2	B	63	LEU
2	B	100(H)	ARG
2	B	135	THR

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Mol	Chain	Res	Type
2	B	149	PRO
2	B	150	VAL
2	B	156	SER
2	B	159	LEU
2	B	178	LEU
2	B	183	THR
2	B	199	ASN
3	C	2	LEU
3	C	3	GLU
3	C	10	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	89	GLN
1	A	138	ASN
1	A	147	GLN
1	A	152	ASN
1	A	199	GLN
2	B	16	GLN
2	B	39	GLN
2	B	199	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
4	SO4	A	6577	-	4,4,4	1.07	0	6,6,6	0.67	0
5	GOL	B	3731	-	5,5,5	0.38	0	5,5,5	0.66	0
4	SO4	B	6576	-	4,4,4	0.97	0	6,6,6	0.72	0

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
4	SO4	A	6577	-	-	0/0/0/0	0/0/0/0
5	GOL	B	3731	-	-	0/4/4/4	0/0/0/0
4	SO4	B	6576	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	3731	GOL	1	0
4	B	6576	SO4	4	0

5.7 Other polymers [i](#)

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 [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	213/214 (99%)	-0.18	1 (0%) 90 95	7, 17, 30, 50	0
2	B	218/237 (91%)	0.03	4 (1%) 68 78	9, 22, 38, 49	0
3	C	10/35 (28%)	0.67	1 (10%) 7 11	32, 42, 50, 55	0
All	All	441/486 (90%)	-0.06	6 (1%) 75 84	7, 20, 38, 55	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	99	THR	3.9
2	B	114	THR	3.0
3	C	11	TRP	2.7
2	B	100(G)	ALA	2.7
1	A	212	GLY	2.5
2	B	1	ARG	2.4

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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	A	6577	5/5	0.85	0.17	79,79,80,80	0
5	GOL	B	3731	6/6	0.86	0.21	32,32,33,34	0
4	SO4	B	6576	5/5	0.90	0.21	80,80,81,81	0

6.5 Other polymers [i](#)

There are no such residues in this entry.