



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 10:55 PM UTC

PDB ID : 3DRQ / pdb_00003drq
Title : Crystal structure of the HIV-1 broadly neutralizing antibody 2F5 in complex with the gp41 FP-MPER Hyb3K construct 514GIGALFLGFLGAAGS528KK-Ahx-655KNEQELLELDKWLWN671 soaked in PEG/2-propanol solution
Authors : Bryson, S.; Julien, J.P.; Pai, E.F.
Deposited on : 2008-07-11
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

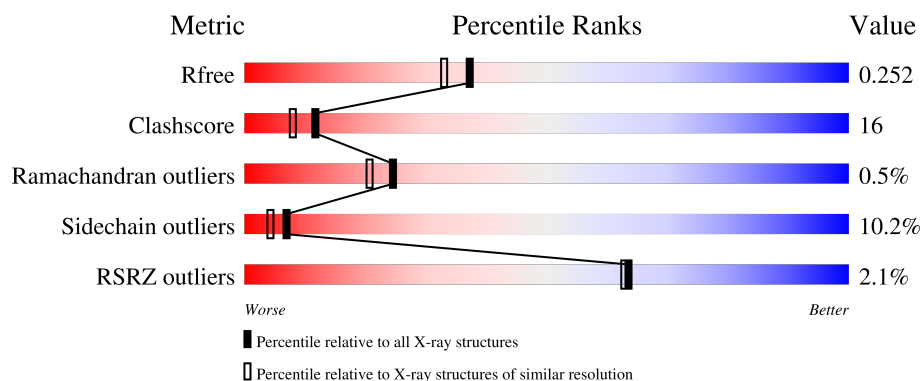
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	214	% 67% 26% 6%
2	B	235	3% 60% 27% • 8%
3	C	35	17% • • 77%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3567 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 Fab' light 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 Fab' heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	216	Total	C	N	O	S	0	0	0
			1623	1030	274	313	6			

- Molecule 3 is a protein called gp41 peptide epitope.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			67	44	10	13			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	140	Total	O	0	0
			140	140		
5	B	92	Total	O	0	0
			92	92		
5	C	2	Total	O	0	0
			2	2		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.50Å 76.60Å 94.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.90 – 2.00 19.90 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.7 (19.90-2.00) 97.7 (19.90-2.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.01Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.227 , 0.254 0.225 , 0.252	Depositor DCC
R_{free} test set	1583 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3567	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.46	0/1674	1.03	8/2276 (0.4%)
2	B	0.45	1/1662 (0.1%)	1.06	9/2273 (0.4%)
3	C	0.77	0/68	1.20	0/91
All	All	0.46	1/3404 (0.0%)	1.05	17/4640 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	100(N)	MET	SD-CE	-5.26	1.66	1.79

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	127	SER	N-CA-C	-9.67	100.74	112.54
2	B	64	ASN	N-CA-C	8.43	123.13	113.01
2	B	100(K)	VAL	N-CA-C	-7.45	102.94	113.07
2	B	143	LYS	N-CA-C	7.06	121.17	109.95
1	A	182	SER	N-CA-C	-6.95	101.17	110.55
2	B	101	ASP	N-CA-C	6.76	120.75	112.23
1	A	94	TYR	N-CA-C	-6.68	101.09	110.31
1	A	114	SER	N-CA-C	-6.61	98.46	109.24
2	B	120	SER	N-CA-C	-6.40	99.26	109.76
2	B	100(N)	MET	N-CA-C	-6.30	97.43	108.20
1	A	137	ASN	N-CA-C	6.22	119.39	109.24
2	B	65	THR	N-CA-C	-5.50	106.24	113.17
1	A	212	GLY	CA-C-O	5.39	125.71	119.13
1	A	205	VAL	CB-CA-C	-5.33	103.46	110.98
2	B	104	GLY	N-CA-C	-5.25	104.15	112.66
2	B	92	CYS	N-CA-C	-5.24	101.61	109.95
1	A	191	VAL	CB-CA-C	-5.09	104.60	110.91

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	54	1
2	B	1623	0	1630	58	1
3	C	67	0	64	5	0
4	A	6	0	8	1	0
5	A	140	0	0	3	0
5	B	92	0	0	3	0
5	C	2	0	0	0	0
All	All	3567	0	3285	105	1

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 16.

All (105) 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:CG	2:B:2:ILE:H	1.56	1.13
1:A:149:LYS:HB2	1:A:193:GLU:HG3	1.34	1.07
2:B:1:ARG:HG3	2:B:2:ILE:N	1.70	1.02
2:B:1:ARG:HG3	2:B:2:ILE:H	0.84	1.00
1:A:138:ASN:HD21	2:B:164:HIS:HE1	1.12	0.96
1:A:103:ARG:HH12	1:A:142:ARG:HD2	1.30	0.93
1:A:103:ARG:HH12	1:A:142:ARG:CD	1.84	0.90
1:A:138:ASN:HD21	2:B:164:HIS:CE1	1.99	0.78
1:A:198:HIS:CD2	1:A:200:GLY:H	2.02	0.77
1:A:115:VAL:HG12	1:A:207:LYS:HG3	1.66	0.77
1:A:103:ARG:NH1	1:A:142:ARG:HD2	1.99	0.76
1:A:90:GLN:HE22	1:A:93:PHE:H	1.30	0.76
1:A:125:LEU:O	1:A:183:LYS:HD3	1.86	0.75
1:A:149:LYS:HB2	1:A:193:GLU:CG	2.16	0.74
2:B:14:PRO:O	2:B:15:THR:HB	1.88	0.74
2:B:38:ARG:HD3	2:B:46:GLU:OE1	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:ARG:HH11	1:A:79:ARG:HG2	1.54	0.71
1:A:138:ASN:ND2	2:B:164:HIS:HE1	1.87	0.70
1:A:201:LEU:HD13	1:A:205:VAL:HG23	1.74	0.69
2:B:119:PRO:HB3	2:B:145:TYR:HB3	1.74	0.69
2:B:210:ARG:NH1	5:B:265:HOH:O	2.25	0.68
2:B:13:LYS:HG3	2:B:113:SER:HA	1.77	0.67
1:A:61:ARG:NE	1:A:79:ARG:HG3	2.09	0.67
2:B:83:SER:OG	2:B:85:VAL:HG22	1.94	0.67
1:A:201:LEU:HD13	1:A:205:VAL:CG2	2.24	0.66
1:A:54:LEU:CD2	1:A:58:VAL:HB	2.27	0.64
1:A:27:GLN:HE22	3:C:2:LEU:HD22	1.62	0.63
1:A:55:GLU:CD	2:B:96:ARG:HH12	2.08	0.61
1:A:61:ARG:CZ	1:A:79:ARG:HG3	2.30	0.61
1:A:90:GLN:NE2	1:A:92:HIS:H	1.99	0.60
2:B:38:ARG:HD2	2:B:48:LEU:HD21	1.82	0.60
2:B:40:PRO:HB2	2:B:43:LYS:HG3	1.84	0.60
1:A:115:VAL:O	1:A:207:LYS:HD2	2.04	0.58
1:A:2:LEU:HD22	1:A:90:GLN:HG2	1.84	0.58
2:B:186:SER:HA	2:B:189:LEU:HD12	1.84	0.58
1:A:184:ALA:O	1:A:188:LYS:HG2	2.02	0.57
2:B:95:ARG:HB2	2:B:100(N):MET:HE1	1.86	0.57
2:B:1:ARG:CG	2:B:2:ILE:N	2.31	0.56
2:B:95:ARG:HD2	2:B:100(N):MET:HE1	1.87	0.56
2:B:54:ASP:OD2	3:C:6:LYS:NZ	2.34	0.56
1:A:115:VAL:CG1	1:A:207:LYS:HG3	2.34	0.56
1:A:55:GLU:OE2	2:B:96:ARG:NH1	2.40	0.55
2:B:178:LEU:HD12	2:B:178:LEU:C	2.31	0.55
1:A:54:LEU:HD22	1:A:58:VAL:HB	1.88	0.54
2:B:24:PHE:CZ	2:B:29:LEU:HD22	2.42	0.54
1:A:49:TYR:CE1	2:B:100(J):PRO:HB3	2.42	0.54
1:A:132:VAL:HG22	1:A:148:TRP:CH2	2.44	0.53
2:B:100(L):ASN:O	2:B:100(M):ALA:HB2	2.10	0.52
1:A:77:THR:O	1:A:77:THR:HG23	2.10	0.50
1:A:103:ARG:HH12	1:A:142:ARG:HD3	1.71	0.49
2:B:146:PHE:CE1	2:B:147:PRO:HB3	2.48	0.49
1:A:79:ARG:HG2	1:A:79:ARG:NH1	2.25	0.49
1:A:78:LEU:HD22	1:A:82:ASP:HB2	1.94	0.48
2:B:38:ARG:HB3	2:B:48:LEU:HD11	1.94	0.48
2:B:95:ARG:CG	2:B:100(N):MET:HE1	2.42	0.48
2:B:71:LYS:HE3	5:B:263:HOH:O	2.13	0.48
1:A:7:SER:HA	1:A:8:PRO:C	2.37	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:201:LYS:N	2:B:202:PRO:CD	2.77	0.48
1:A:33:LEU:HD22	1:A:71:PHE:CG	2.49	0.47
1:A:138:ASN:ND2	2:B:164:HIS:CE1	2.72	0.47
2:B:2:ILE:HG23	2:B:27:PHE:CD2	2.50	0.47
1:A:79:ARG:NH1	1:A:79:ARG:CG	2.76	0.47
2:B:14:PRO:O	2:B:15:THR:CB	2.62	0.47
1:A:132:VAL:HG22	1:A:148:TRP:CZ3	2.50	0.46
2:B:2:ILE:HA	2:B:25:SER:O	2.15	0.46
2:B:24:PHE:HZ	2:B:29:LEU:HD22	1.80	0.46
1:A:38:GLN:O	1:A:84:ALA:HB1	2.16	0.46
2:B:201:LYS:H	2:B:201:LYS:HG2	1.58	0.46
2:B:39:GLN:C	2:B:88:ALA:HB1	2.40	0.46
2:B:186:SER:HA	2:B:189:LEU:CD1	2.46	0.45
1:A:141:PRO:O	1:A:198:HIS:HE1	1.99	0.45
2:B:37:ILE:HD12	2:B:37:ILE:N	2.32	0.45
1:A:124:GLN:NE2	1:A:131:SER:H	2.15	0.45
2:B:95:ARG:CD	2:B:100(N):MET:HE1	2.47	0.45
1:A:1:ALA:HB3	5:A:3762:HOH:O	2.15	0.45
2:B:125:ALA:HB1	2:B:126:PRO:HD2	1.99	0.44
2:B:87:THR:N	2:B:111:ILE:HD13	2.32	0.44
1:A:1:ALA:HB1	3:C:2:LEU:HD11	1.99	0.44
1:A:198:HIS:HD2	1:A:200:GLY:H	1.58	0.44
2:B:14:PRO:HD3	2:B:112:SER:C	2.42	0.44
2:B:2:ILE:HG23	2:B:27:PHE:HD2	1.82	0.43
1:A:145:LYS:HB3	1:A:197:THR:HB	2.01	0.43
1:A:159:SER:HA	1:A:178:THR:O	2.18	0.43
2:B:135:THR:O	2:B:135:THR:HG23	2.17	0.43
2:B:95:ARG:HB2	2:B:100(N):MET:CE	2.48	0.43
1:A:183:LYS:O	1:A:187:GLU:HG2	2.19	0.43
2:B:9:PRO:HA	2:B:10:PRO:HD3	1.82	0.43
2:B:95:ARG:CB	2:B:100(N):MET:HE1	2.49	0.43
1:A:176:SER:CB	4:A:3731:GOL:H11	2.49	0.42
2:B:56:ASP:OD1	3:C:6:LYS:HE3	2.19	0.42
2:B:107:ILE:O	2:B:107:ILE:HG23	2.20	0.42
1:A:124:GLN:HE22	1:A:131:SER:H	1.68	0.42
2:B:86:ASP:C	2:B:111:ILE:HD13	2.44	0.42
1:A:11:LEU:HD22	1:A:104:VAL:HG13	2.02	0.41
2:B:82(B):ARG:HB3	5:B:237:HOH:O	2.20	0.41
1:A:118:PHE:HD2	1:A:133:VAL:HG22	1.85	0.41
2:B:87:THR:HG23	2:B:110:THR:HA	2.03	0.41
5:A:3813:HOH:O	2:B:164:HIS:HD2	2.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:36:TRP:C	2:B:37:ILE:HD12	2.45	0.41
2:B:89:THR:HA	2:B:108:THR:HA	2.02	0.40
2:B:125:ALA:HA	2:B:126:PRO:HD3	1.89	0.40
1:A:115:VAL:HA	1:A:135:LEU:O	2.22	0.40
1:A:27:GLN:NE2	3:C:2:LEU:HD22	2.33	0.40
1:A:81:GLU:HG2	5:A:3812:HOH:O	2.21	0.40
2:B:14:PRO:HG3	2:B:111:ILE:HG22	2.03	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	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:GLU:OE1	2:B:17:THR:OG1[3_545]	2.12	0.08

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%)	207 (98%)	4 (2%)	0	100	100
2	B	210/235 (89%)	199 (95%)	9 (4%)	2 (1%)	12	8
3	C	6/35 (17%)	6 (100%)	0	0	100	100
All	All	427/484 (88%)	412 (96%)	13 (3%)	2 (0%)	24	21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	15	THR
2	B	100(M)	ALA

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	186/187 (100%)	165 (89%)	21 (11%)	5	3
2	B	189/204 (93%)	172 (91%)	17 (9%)	9	6
3	C	7/25 (28%)	6 (86%)	1 (14%)	3	2
All	All	382/416 (92%)	343 (90%)	39 (10%)	7	4

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	54	LEU
1	A	70	GLU
1	A	73	LEU
1	A	78	LEU
1	A	79	ARG
1	A	90	GLN
1	A	97	THR
1	A	103	ARG
1	A	126	LYS
1	A	129	THR
1	A	132	VAL
1	A	133	VAL
1	A	145	LYS
1	A	152	ASN
1	A	165	GLU
1	A	187	GLU
1	A	190	LYS
1	A	191	VAL
1	A	203	SER
1	A	205	VAL
2	B	3	THR
2	B	13	LYS
2	B	29	LEU
2	B	30	SER
2	B	34	VAL

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Mol	Chain	Res	Type
2	B	38	ARG
2	B	53	SER
2	B	63	LEU
2	B	105	GLN
2	B	108	THR
2	B	126	PRO
2	B	149	PRO
2	B	156	SER
2	B	159	LEU
2	B	178	LEU
2	B	183	THR
2	B	210	ARG
3	C	2	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	38	GLN
1	A	90	GLN
1	A	124	GLN
1	A	138	ASN
1	A	147	GLN
1	A	152	ASN
1	A	198	HIS
1	A	199	GLN
2	B	16	GLN
2	B	39	GLN
2	B	164	HIS
2	B	192	GLN
2	B	199	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry

1 ligand is 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	GOL	A	3731	-	5,5,5	0.49	0	5,5,5	0.63	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	GOL	A	3731	-	-	0/4/4/4	-

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	3731	GOL	1	0

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 [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%)	-0.19	2 (0%) 81 80	14, 22, 33, 53	0
2	B	216/235 (91%)	0.13	7 (3%) 50 49	15, 26, 43, 52	0
3	C	8/35 (22%)	0.62	0 100 100	32, 36, 42, 43	0
All	All	437/484 (90%)	-0.02	9 (2%) 63 63	14, 24, 42, 53	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	193	GLU	3.6
2	B	213	PRO	3.5
2	B	100(J)	PRO	3.2
2	B	30	SER	2.8
1	A	1	ALA	2.6
2	B	190	GLY	2.5
2	B	114	THR	2.4
2	B	100(I)	GLY	2.4
2	B	189	LEU	2.3

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 oligosaccharides 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. 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($\text{\AA}^2$)	Q<0.9
4	GOL	A	3731	6/6	0.81	0.13	34,37,39,39	0

6.5 Other polymers

There are no such residues in this entry.