



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:14 AM UTC

PDB ID : 1P7K / pdb_00001p7k
Title : Crystal structure of an anti-ssDNA antigen-binding fragment (Fab) bound to 4-(2-Hydroxyethyl)piperazine-1-ethanesulfonic acid (HEPES)
Authors : Schuermann, J.P.; Henzl, M.T.; Deutscher, S.L.; Tanner, J.J.
Deposited on : 2003-05-02
Resolution : 1.75 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

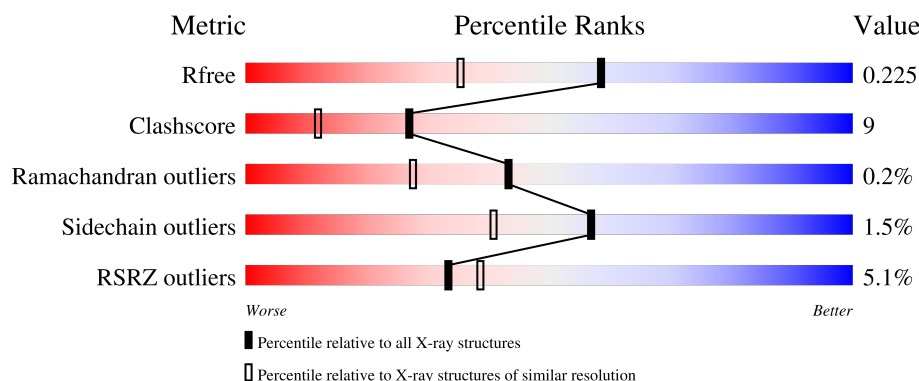
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	7% 76% 23% .
1	L	214	3% 84% 15%
2	B	224	5% 83% 13% ..
2	H	224	5% 80% 17% .

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
3	SO4	H	801	-	-	X	-
3	SO4	H	802	-	-	X	-
7	GOL	A	1116	-	-	X	-
7	GOL	H	1106	-	-	X	-
7	GOL	L	1103	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7565 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 antibody light chain FAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	213	Total	C	N	O	S	0	1	0
			1633	1021	273	333	6			
1	A	214	Total	C	N	O	S	0	1	0
			1623	1018	272	327	6			

- Molecule 2 is a protein called antibody heavy chain FAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	2	0
			1653	1055	263	324	11			
2	B	217	Total	C	N	O	S	0	3	0
			1636	1045	259	321	11			

There are 22 discrepancies between the modelled and reference sequences:

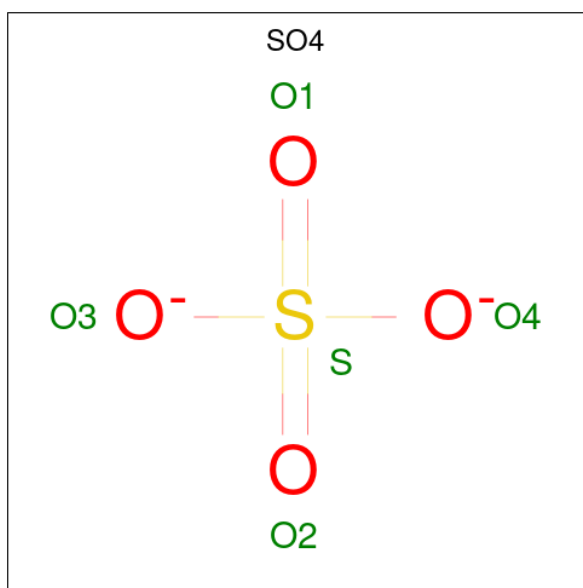
Chain	Residue	Modelled	Actual	Comment	Reference
H	1	GLN	-	cloning artifact	UNP Q9JL75
H	2	VAL	-	cloning artifact	UNP Q9JL75
H	3	LYS	-	cloning artifact	UNP Q9JL75
H	4	LEU	-	cloning artifact	UNP Q9JL75
H	52A	PRO	-	insertion	UNP Q9JL75
H	82A	SER	-	insertion	UNP Q9JL75
H	82B	SER	-	insertion	UNP Q9JL75
H	82C	LEU	-	insertion	UNP Q9JL75
H	100A	TYR	-	insertion	UNP Q9JL75
H	100B	ALA	-	insertion	UNP Q9JL75
H	100C	MET	-	insertion	UNP Q9JL75
B	1	GLN	-	cloning artifact	UNP Q9JL75
B	2	VAL	-	cloning artifact	UNP Q9JL75
B	3	LYS	-	cloning artifact	UNP Q9JL75
B	4	LEU	-	cloning artifact	UNP Q9JL75
B	52A	PRO	-	insertion	UNP Q9JL75

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Chain	Residue	Modelled	Actual	Comment	Reference
B	82A	SER	-	insertion	UNP Q9JL75
B	82B	SER	-	insertion	UNP Q9JL75
B	82C	LEU	-	insertion	UNP Q9JL75
B	100A	TYR	-	insertion	UNP Q9JL75
B	100B	ALA	-	insertion	UNP Q9JL75
B	100C	MET	-	insertion	UNP Q9JL75

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



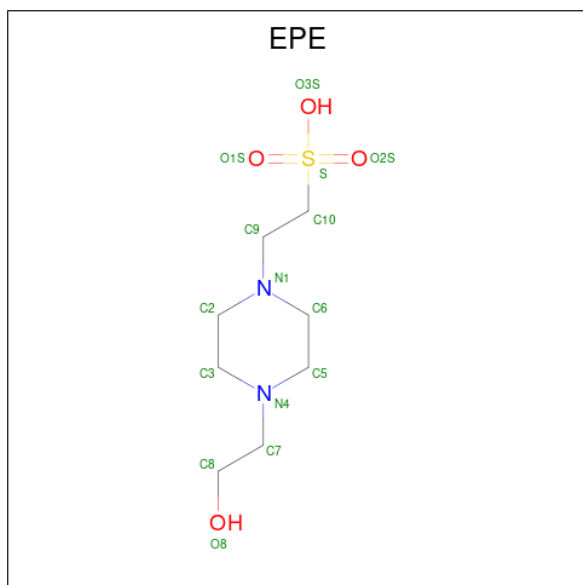
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	L	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	H	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

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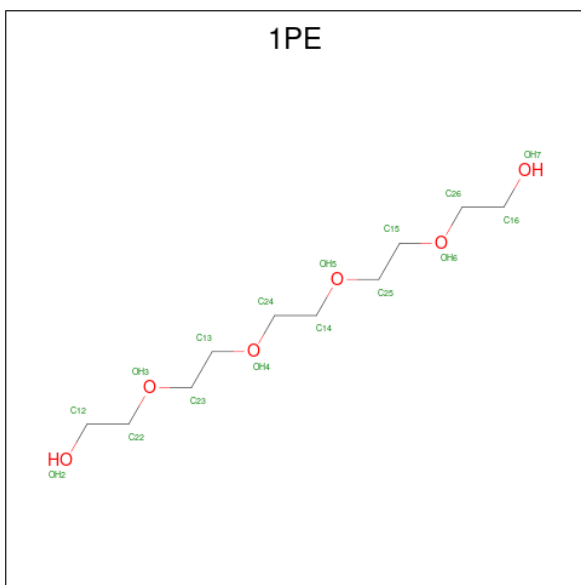
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



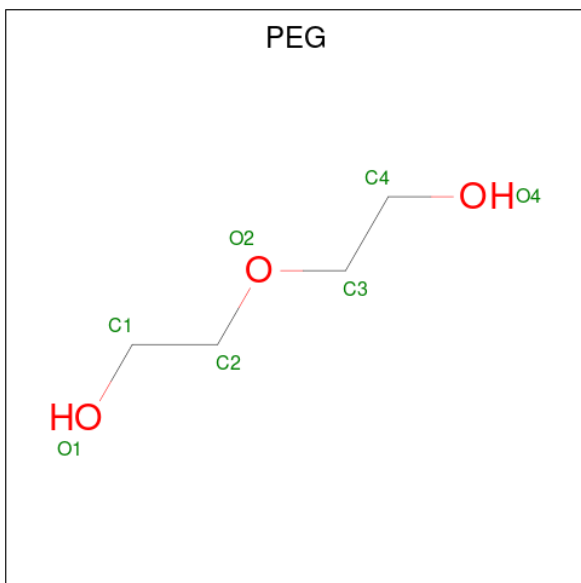
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	L	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
4	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 5 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



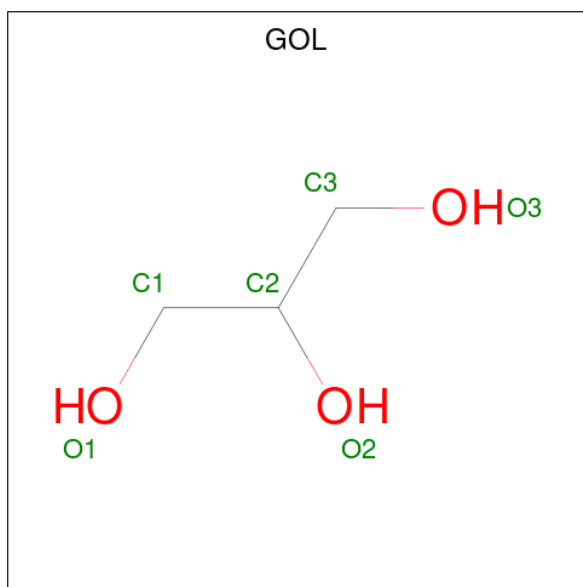
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	L	1	Total	C	O	0	0
			16	10	6		
5	H	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		
5	B	1	Total	C	O	0	0
			16	10	6		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total	C	O	0	0
			7	4	3		
6	H	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	L	1	Total	C	O	0	0
			6	3	3		
7	L	1	Total	C	O	0	0
			6	3	3		
7	L	1	Total	C	O	0	0
			6	3	3		
7	L	1	Total	C	O	0	0
			6	3	3		
7	L	1	Total	C	O	0	0
			6	3	3		
7	H	1	Total	C	O	0	0
			6	3	3		
7	H	1	Total	C	O	0	0
			6	3	3		
7	H	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		
7	A	1	Total	C	O	0	0
			6	3	3		

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

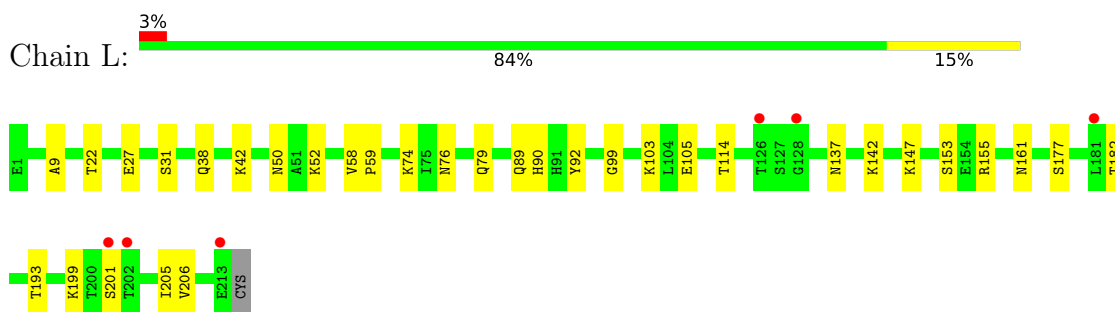
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	237	Total	O	0	0
			237	237		
8	H	195	Total	O	0	0
			195	195		
8	A	162	Total	O	0	0
			162	162		
8	B	177	Total	O	0	0
			177	177		

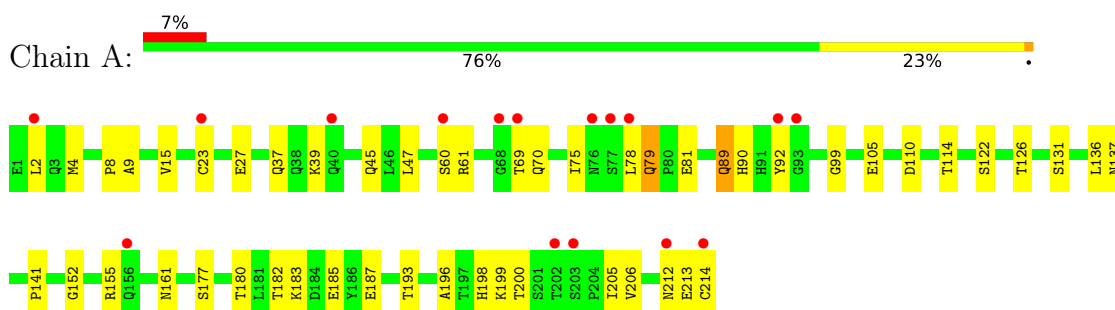
3 Residue-property plots [i](#)

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

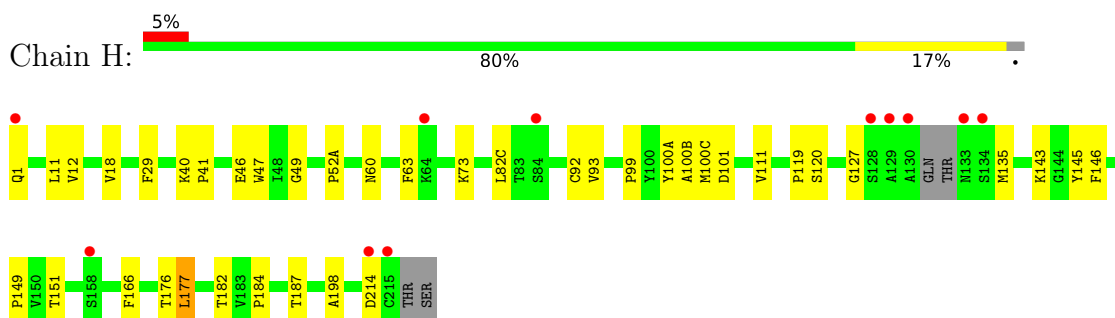
- Molecule 1: antibody light chain FAB



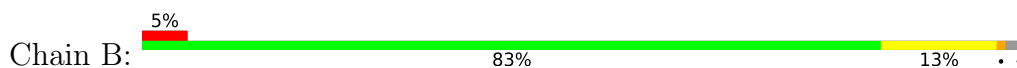
- Molecule 1: antibody light chain FAB

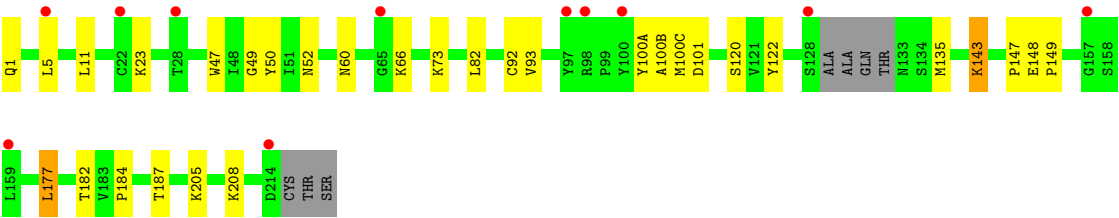


- Molecule 2: antibody heavy chain FAB



- Molecule 2: antibody heavy chain FAB





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.67Å 156.70Å 61.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.14 – 1.75 48.14 – 1.75	Depositor EDS
% Data completeness (in resolution range)	93.4 (48.14-1.75) 93.5 (48.14-1.75)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.99 (at 1.75Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.197 , 0.230 0.194 , 0.225	Depositor DCC
R_{free} test set	9796 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	23.2	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.3	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	0.96	EDS
Total number of atoms	7565	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: 1PE, SO4, GOL, EPE, PEG

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	0.50	0/1660	0.95	5/2257 (0.2%)
1	L	0.55	0/1671	1.00	8/2270 (0.4%)
2	B	0.49	0/1682	0.97	6/2298 (0.3%)
2	H	0.56	0/1699	0.99	8/2320 (0.3%)
All	All	0.53	0/6712	0.98	27/9145 (0.3%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	143	LYS	N-CA-C	7.50	121.62	109.40
2	B	92	CYS	N-CA-C	-7.33	97.64	109.96
2	H	92	CYS	N-CA-C	-7.24	97.79	109.96
2	H	143	LYS	N-CA-C	6.77	120.49	109.59
2	B	120	SER	N-CA-C	-6.49	98.69	109.46
2	H	101	ASP	N-CA-C	6.43	121.12	113.28
2	B	60	ASN	N-CA-C	-6.41	100.35	109.96
2	H	120	SER	N-CA-C	-6.31	98.95	109.24
1	L	114	THR	N-CA-C	-6.30	98.94	108.96
1	A	182	THR	N-CA-C	-6.26	101.26	110.52
2	B	100(B)	ALA	N-CA-C	-6.23	100.14	110.17
1	L	182	THR	N-CA-C	-6.01	101.62	110.52
1	A	137	ASN	N-CA-C	5.99	119.47	109.95
1	L	58	VAL	CA-C-N	5.73	125.69	119.78
1	L	58	VAL	C-N-CA	5.73	125.69	119.78
1	L	137	ASN	N-CA-C	5.68	118.99	109.95
1	A	99	GLY	N-CA-C	-5.57	103.99	112.51
2	B	101	ASP	N-CA-C	5.56	120.12	113.23
1	A	114	THR	N-CA-C	-5.55	100.14	108.96
1	L	142	LYS	N-CA-C	5.48	119.22	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	60	ASN	N-CA-C	-5.36	101.92	109.96
1	L	205	ILE	N-CA-C	-5.36	100.87	108.58
2	H	100(B)	ALA	N-CA-C	-5.35	100.97	109.96
1	L	99	GLY	N-CA-C	-5.33	104.35	112.51
2	H	166	PHE	N-CA-C	5.30	117.61	110.29
2	H	29	PHE	N-CA-C	5.19	117.33	111.11
1	A	205	ILE	N-CA-C	-5.11	100.41	108.23

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	1623	0	1529	40	0
1	L	1633	0	1541	27	0
2	B	1636	0	1569	24	0
2	H	1653	0	1595	25	0
3	A	5	0	0	0	0
3	B	10	0	0	0	0
3	H	25	0	0	0	8
3	L	5	0	0	0	0
4	B	15	0	18	1	0
4	L	15	0	17	2	0
5	B	32	0	44	8	0
5	H	16	0	22	3	0
5	L	16	0	22	2	0
6	H	7	0	10	1	0
6	L	7	0	10	3	0
7	A	24	0	31	11	0
7	B	24	0	32	4	0
7	H	18	0	24	7	0
7	L	30	0	40	13	0
8	A	162	0	0	8	0
8	B	177	0	0	3	0
8	H	195	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	L	237	0	0	8	0
All	All	7565	0	6504	123	8

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 9.

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)
1:L:155:ARG:HE	6:L:1201:PEG:H11	1.27	0.99
1:A:8:PRO:HB2	7:A:1116:GOL:H12	1.48	0.94
7:L:1103:GOL:H2	1:A:15:VAL:HG12	1.52	0.91
1:L:9:ALA:H	7:L:1107:GOL:H11	1.38	0.89
5:H:1001:1PE:H251	8:H:1331:HOH:O	1.74	0.86
1:A:9:ALA:H	7:A:1116:GOL:C3	1.92	0.82
1:A:9:ALA:H	7:A:1116:GOL:H31	1.43	0.82
2:B:93:VAL:HG11	2:B:100(C):MET:HB3	1.61	0.81
1:L:31:SER:HB2	7:L:1103:GOL:O1	1.81	0.79
2:H:1:GLN:HE21	2:H:1:GLN:CA	1.95	0.78
1:A:212:ASN:C	1:A:214:CYS:H	1.94	0.76
1:L:50:ASN:OD1	7:L:1103:GOL:H11	1.85	0.75
2:B:73:LYS:HZ1	7:B:1112:GOL:H31	1.50	0.75
2:H:18:VAL:HG22	2:H:82(C):LEU:HD11	1.68	0.75
2:H:176:THR:HG22	7:H:1106:GOL:H32	1.70	0.73
8:A:1169:HOH:O	7:B:1117:GOL:H32	1.87	0.73
2:H:52(A):PRO:O	2:H:73:LYS:HE2	1.88	0.73
2:H:93:VAL:HG11	2:H:100(C):MET:HB3	1.71	0.72
1:L:42:LYS:O	7:L:1115:GOL:H12	1.90	0.72
1:A:185:GLU:HG2	8:A:1185:HOH:O	1.91	0.71
1:A:9:ALA:HB3	7:A:1116:GOL:H2	1.72	0.71
2:B:93:VAL:CG1	2:B:100(C):MET:HB3	2.23	0.69
7:L:1103:GOL:H32	8:A:1198:HOH:O	1.93	0.68
1:A:45:GLN:HG3	8:A:1206:HOH:O	1.93	0.68
2:B:73:LYS:NZ	7:B:1112:GOL:H31	2.08	0.68
1:L:52:LYS:HB3	5:L:1003:1PE:H261	1.74	0.68
1:L:50:ASN:CG	7:L:1103:GOL:H11	2.19	0.67
1:L:79:GLN:NE2	8:L:1360:HOH:O	2.27	0.66
1:A:90:HIS:HD2	1:A:92:TYR:H	1.42	0.66
2:B:11:LEU:CD2	5:B:1004:1PE:H152	2.26	0.66
1:A:89:GLN:NE2	2:B:100(C):MET:HE3	2.11	0.65
2:B:11:LEU:HD21	5:B:1004:1PE:H152	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:147:LYS:HE2	8:L:1320:HOH:O	1.98	0.63
1:L:9:ALA:N	7:L:1107:GOL:H11	2.11	0.63
1:A:198:HIS:HD2	1:A:200:THR:OG1	1.82	0.62
1:A:79[B]:GLN:HG2	8:A:1198:HOH:O	1.98	0.61
1:L:59:PRO:HD3	7:L:1119:GOL:H31	1.83	0.61
1:A:131:SER:OG	1:A:180:THR:HG22	2.01	0.60
7:A:1118:GOL:H11	2:B:122:TYR:OH	2.01	0.60
5:B:1002:1PE:H142	7:B:1111:GOL:H2	1.84	0.60
7:L:1110:GOL:H32	2:H:127:GLY:HA2	1.84	0.59
1:A:2:LEU:HG	1:A:27:GLU:HG2	1.85	0.58
1:L:90:HIS:HD2	1:L:92:TYR:H	1.51	0.58
2:B:135:MET:HE3	2:B:182:THR:HG22	1.84	0.58
2:H:135:MET:HE3	2:H:182:THR:HG22	1.86	0.58
1:A:9:ALA:N	7:A:1116:GOL:H31	2.15	0.57
1:A:61:ARG:NH2	1:A:79[A]:GLN:HG2	2.20	0.57
2:B:52:ASN:HD22	5:B:1002:1PE:H261	1.70	0.57
1:L:155:ARG:NE	6:L:1201:PEG:H11	2.10	0.55
2:B:148:GLU:O	5:B:1004:1PE:H231	2.06	0.55
2:H:1:GLN:CA	2:H:1:GLN:NE2	2.66	0.55
2:H:177:LEU:HA	7:H:1106:GOL:C1	2.36	0.55
1:A:39:LYS:HE2	1:A:81:GLU:O	2.06	0.55
1:A:212:ASN:C	1:A:214:CYS:N	2.61	0.55
1:A:69:THR:HG23	1:A:70:GLN:HG2	1.90	0.53
1:L:22:THR:HG23	8:L:1364:HOH:O	2.08	0.53
2:B:100(A):TYR:HB2	4:B:902:EPE:O1S	2.07	0.53
1:L:9:ALA:H	7:L:1107:GOL:C1	2.18	0.53
1:A:193:THR:CG2	1:A:206:VAL:HG13	2.39	0.53
2:H:176:THR:HG22	7:H:1106:GOL:C3	2.36	0.52
1:A:155:ARG:NH2	8:A:1185:HOH:O	2.37	0.52
1:L:59:PRO:CD	7:L:1119:GOL:H31	2.38	0.52
1:A:122:SER:O	1:A:126:THR:HG23	2.09	0.52
2:H:151:THR:OG1	2:H:198:ALA:HB3	2.09	0.52
2:H:184:PRO:HB2	2:H:187:THR:HG23	1.93	0.51
1:A:141:PRO:O	1:A:198:HIS:HE1	1.94	0.51
1:L:153:SER:HB2	6:L:1201:PEG:H32	1.92	0.51
2:B:147:PRO:HD3	5:B:1004:1PE:H151	1.93	0.51
1:L:76:ASN:ND2	8:L:1312:HOH:O	2.43	0.51
2:H:40:LYS:N	7:H:1105:GOL:O3	2.33	0.51
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.46	0.51
2:B:184:PRO:HB2	2:B:187:THR:HG23	1.93	0.50
2:B:50:TYR:C	2:B:50:TYR:CD1	2.90	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:1001:1PE:H252	8:H:1339:HOH:O	2.12	0.49
2:H:93:VAL:CG1	2:H:100(C):MET:HB3	2.42	0.49
2:B:205:LYS:HD3	8:B:1273:HOH:O	2.13	0.48
1:A:39:LYS:NZ	1:A:81:GLU:CG	2.77	0.48
1:A:136:LEU:CD2	1:A:196:ALA:HB2	2.44	0.48
1:A:75:ILE:HG21	1:A:78:LEU:HD23	1.95	0.48
2:B:177:LEU:C	2:B:177:LEU:HD12	2.38	0.48
2:H:63:PHE:CZ	7:H:1101:GOL:H2	2.49	0.48
1:L:74:LYS:NZ	8:L:1340:HOH:O	2.46	0.48
1:A:4:MET:HE3	1:A:23:CYS:SG	2.53	0.48
5:B:1002:1PE:H152	8:B:1223:HOH:O	2.13	0.48
1:A:206:VAL:O	7:A:1108:GOL:H2	2.14	0.47
5:H:1001:1PE:H162	6:H:1202:PEG:H41	1.97	0.47
1:A:8:PRO:HA	7:A:1116:GOL:H31	1.97	0.46
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.97	0.46
2:H:119:PRO:HB3	2:H:145:TYR:HB3	1.98	0.45
1:A:110:ASP:OD2	1:A:199:LYS:NZ	2.49	0.45
1:L:201:SER:HB2	8:L:1285:HOH:O	2.16	0.45
1:A:9:ALA:CB	7:A:1116:GOL:H2	2.45	0.45
2:B:1:GLN:HG2	2:B:1:GLN:O	2.17	0.45
1:L:193:THR:CG2	1:L:206:VAL:HG13	2.47	0.45
2:H:12:VAL:O	2:H:111:VAL:HA	2.17	0.45
2:H:46:GLU:HB2	7:H:1105:GOL:H11	1.98	0.45
1:A:161:ASN:HD22	1:A:177:SER:HA	1.81	0.45
2:H:73:LYS:NZ	8:H:1314:HOH:O	2.42	0.45
1:A:9:ALA:H	7:A:1116:GOL:C2	2.30	0.44
1:A:79[A]:GLN:CD	1:A:81:GLU:CB	2.91	0.44
2:B:208:LYS:HE3	8:B:1234:HOH:O	2.17	0.44
2:B:5:LEU:HB3	2:B:23:LYS:HB3	1.99	0.44
2:B:47:TRP:CZ2	2:B:49:GLY:HA2	2.52	0.44
1:A:69:THR:HG22	8:A:1203:HOH:O	2.18	0.44
5:L:1003:1PE:H151	5:L:1003:1PE:H161	1.71	0.44
2:H:11:LEU:HD11	2:H:146:PHE:CZ	2.53	0.44
1:A:183:LYS:O	1:A:187:GLU:HG3	2.17	0.44
1:A:193:THR:HG23	1:A:206:VAL:HG13	2.00	0.43
4:L:901:EPE:O2S	2:H:99:PRO:HA	2.19	0.43
1:L:161:ASN:HD22	1:L:177:SER:HA	1.83	0.43
1:A:152:GLY:HA2	8:A:1214:HOH:O	2.19	0.43
1:L:38:GLN:NE2	7:L:1115:GOL:H11	2.33	0.43
1:A:79[A]:GLN:CD	1:A:81:GLU:HB2	2.43	0.43
2:H:11:LEU:HD11	2:H:146:PHE:HZ	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:1118:GOL:H12	2:B:143:LYS:HD3	2.02	0.42
2:B:66:LYS:O	2:B:82:LEU:HA	2.19	0.42
1:L:27:GLU:HG3	8:L:1321:HOH:O	2.20	0.41
1:L:199:LYS:HB2	1:L:199:LYS:HE3	1.86	0.41
4:L:901:EPE:O1S	2:H:100(A):TYR:HB2	2.19	0.41
2:H:177:LEU:HA	7:H:1106:GOL:H12	2.02	0.41
1:L:193:THR:HG23	1:L:206:VAL:HG13	2.02	0.41
2:B:11:LEU:HD23	5:B:1004:1PE:H152	2.00	0.40
1:L:103:LYS:CG	8:L:1392:HOH:O	2.69	0.40

All (8) 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)
3:H:801:SO4:S	3:H:801:SO4:O2[2_655]	1.46	0.74
3:H:801:SO4:S	3:H:801:SO4:O1[2_655]	1.47	0.73
3:H:802:SO4:S	3:H:802:SO4:O4[2_655]	1.47	0.73
3:H:801:SO4:S	3:H:801:SO4:O4[2_655]	1.48	0.72
3:H:802:SO4:S	3:H:802:SO4:O2[2_655]	1.48	0.72
3:H:802:SO4:S	3:H:802:SO4:O1[2_655]	1.48	0.72
3:H:802:SO4:S	3:H:802:SO4:O3[2_655]	1.48	0.72
3:H:801:SO4:S	3:H:801:SO4:O3[2_655]	1.49	0.71

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	213/214 (100%)	206 (97%)	6 (3%)	1 (0%)	24	11
1	L	212/214 (99%)	208 (98%)	4 (2%)	0	100	100
2	B	216/224 (96%)	212 (98%)	4 (2%)	0	100	100
2	H	218/224 (97%)	210 (96%)	7 (3%)	1 (0%)	24	11

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	859/876 (98%)	836 (97%)	21 (2%)	2 (0%)	43	27

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	213	GLU
2	H	214	ASP

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	177/188 (94%)	172 (97%)	5 (3%)	38	18
1	L	181/188 (96%)	179 (99%)	2 (1%)	65	52
2	B	183/194 (94%)	181 (99%)	2 (1%)	65	52
2	H	185/194 (95%)	182 (98%)	3 (2%)	55	38
All	All	726/764 (95%)	714 (98%)	12 (2%)	57	36

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	89	GLN
1	L	105	GLU
2	H	41	PRO
2	H	149	PRO
2	H	177	LEU
1	A	60	SER
1	A	79[A]	GLN
1	A	79[B]	GLN
1	A	89	GLN
1	A	105	GLU
2	B	149	PRO
2	B	177	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	GLN
1	L	40	GLN
1	L	70	GLN
1	L	76	ASN
1	L	89	GLN
1	L	90	HIS
1	L	137	ASN
1	L	145	ASN
1	L	161	ASN
1	A	38	GLN
1	A	40	GLN
1	A	89	GLN
1	A	90	HIS
1	A	137	ASN
1	A	161	ASN
1	A	198	HIS
1	A	210	ASN
2	B	39	GLN
2	B	105	GLN
2	B	171	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

33 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	$\# Z > 2$
7	GOL	H	1105	-	5,5,5	0.33	0	5,5,5	1.42	1 (20%)
7	GOL	B	1117	-	5,5,5	0.26	0	5,5,5	1.32	1 (20%)
7	GOL	A	1118	-	5,5,5	0.29	0	5,5,5	1.79	2 (40%)
7	GOL	L	1103	-	5,5,5	0.27	0	5,5,5	1.85	2 (40%)
3	SO4	H	802	-	4,4,4	0.37	0	6,6,6	0.12	0
3	SO4	L	808	-	4,4,4	0.42	0	6,6,6	0.12	0
3	SO4	H	805	-	4,4,4	0.40	0	6,6,6	0.07	0
7	GOL	L	1107	-	5,5,5	0.35	0	5,5,5	1.18	1 (20%)
3	SO4	B	809	-	4,4,4	0.39	0	6,6,6	0.05	0
3	SO4	H	804	-	4,4,4	0.36	0	6,6,6	0.13	0
7	GOL	L	1115	-	5,5,5	0.40	0	5,5,5	1.27	1 (20%)
7	GOL	A	1108	-	5,5,5	0.36	0	5,5,5	0.48	0
3	SO4	H	803	-	4,4,4	0.39	0	6,6,6	0.19	0
3	SO4	B	807	-	4,4,4	0.38	0	6,6,6	0.14	0
4	EPE	L	901	-	15,15,15	0.82	0	19,20,20	1.43	3 (15%)
7	GOL	A	1116	-	5,5,5	0.24	0	5,5,5	0.80	0
5	1PE	H	1001	-	15,15,15	0.71	0	14,14,14	0.79	0
7	GOL	A	1120	-	5,5,5	0.33	0	5,5,5	1.24	1 (20%)
7	GOL	L	1119	-	5,5,5	0.31	0	5,5,5	0.37	0
3	SO4	H	801	-	4,4,4	0.37	0	6,6,6	0.11	0
6	PEG	H	1202	-	6,6,6	0.52	0	5,5,5	1.58	1 (20%)
5	1PE	L	1003	-	15,15,15	0.73	0	14,14,14	0.76	0
5	1PE	B	1002	-	15,15,15	0.71	0	14,14,14	0.77	0
6	PEG	L	1201	-	6,6,6	0.59	0	5,5,5	1.68	1 (20%)
7	GOL	B	1102	-	5,5,5	0.30	0	5,5,5	1.16	1 (20%)
7	GOL	H	1101	-	5,5,5	0.31	0	5,5,5	1.90	2 (40%)
7	GOL	B	1112	-	5,5,5	0.25	0	5,5,5	0.37	0
7	GOL	H	1106	-	5,5,5	0.27	0	5,5,5	1.58	1 (20%)
3	SO4	A	806	-	4,4,4	0.42	0	6,6,6	0.12	0
7	GOL	L	1110	-	5,5,5	0.35	0	5,5,5	1.29	1 (20%)
4	EPE	B	902	-	15,15,15	0.86	0	19,20,20	1.37	2 (10%)
5	1PE	B	1004	-	15,15,15	0.78	0	14,14,14	0.89	0
7	GOL	B	1111	-	5,5,5	0.31	0	5,5,5	0.75	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
7	GOL	H	1105	-	-	4/4/4/4	-
7	GOL	B	1117	-	-	4/4/4/4	-
7	GOL	A	1118	-	-	3/4/4/4	-
7	GOL	L	1103	-	-	2/4/4/4	-
7	GOL	L	1107	-	-	4/4/4/4	-
7	GOL	A	1108	-	-	4/4/4/4	-
7	GOL	L	1115	-	-	2/4/4/4	-
4	EPE	L	901	-	-	3/9/19/19	0/1/1/1
7	GOL	A	1116	-	-	4/4/4/4	-
5	1PE	H	1001	-	-	7/13/13/13	-
7	GOL	A	1120	-	-	4/4/4/4	-
7	GOL	L	1119	-	-	2/4/4/4	-
6	PEG	H	1202	-	-	0/4/4/4	-
5	1PE	B	1002	-	-	6/13/13/13	-
5	1PE	L	1003	-	-	10/13/13/13	-
6	PEG	L	1201	-	-	2/4/4/4	-
7	GOL	B	1102	-	-	1/4/4/4	-
7	GOL	H	1101	-	-	2/4/4/4	-
7	GOL	B	1112	-	-	4/4/4/4	-
7	GOL	H	1106	-	-	2/4/4/4	-
7	GOL	L	1110	-	-	1/4/4/4	-
4	EPE	B	902	-	-	5/9/19/19	0/1/1/1
5	1PE	B	1004	-	-	10/13/13/13	-
7	GOL	B	1111	-	-	4/4/4/4	-

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	902	EPE	O3S-S-O2S	-3.61	102.36	111.40
4	L	901	EPE	O3S-S-O2S	-3.44	102.79	111.40
7	A	1118	GOL	C3-C2-C1	3.12	123.24	111.80
7	H	1101	GOL	C3-C2-C1	-3.06	100.57	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	H	1106	GOL	O2-C2-C3	3.05	121.80	109.18
7	L	1103	GOL	C3-C2-C1	-2.95	100.97	111.80
7	L	1103	GOL	O2-C2-C3	2.81	120.81	109.18
7	A	1120	GOL	O2-C2-C3	2.45	119.33	109.18
7	L	1107	GOL	C3-C2-C1	-2.45	102.80	111.80
4	B	902	EPE	O2S-S-O1S	2.45	121.78	113.82
6	L	1201	PEG	O2-C2-C1	2.43	120.83	110.11
7	H	1101	GOL	O2-C2-C1	2.42	119.19	109.18
6	H	1202	PEG	O2-C2-C1	2.41	120.75	110.11
7	H	1105	GOL	C3-C2-C1	-2.37	103.09	111.80
7	A	1118	GOL	O2-C2-C1	2.34	118.85	109.18
7	L	1115	GOL	O2-C2-C3	2.30	118.70	109.18
4	L	901	EPE	O2S-S-O1S	2.22	121.04	113.82
4	L	901	EPE	O1S-S-C10	2.17	110.00	106.73
7	B	1102	GOL	O2-C2-C3	2.15	118.10	109.18
7	B	1117	GOL	O2-C2-C3	2.12	117.96	109.18
7	L	1110	GOL	O2-C2-C1	2.07	117.76	109.18

There are no chirality outliers.

All (90) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	901	EPE	C10-C9-N1-C2
4	L	901	EPE	C10-C9-N1-C6
4	B	902	EPE	C10-C9-N1-C2
4	B	902	EPE	C10-C9-N1-C6
4	B	902	EPE	C9-C10-S-O1S
7	L	1103	GOL	C1-C2-C3-O3
7	L	1107	GOL	O1-C1-C2-O2
7	L	1107	GOL	O1-C1-C2-C3
7	L	1115	GOL	O1-C1-C2-C3
7	L	1119	GOL	O1-C1-C2-C3
7	H	1101	GOL	C1-C2-C3-O3
7	H	1105	GOL	O1-C1-C2-C3
7	H	1106	GOL	C1-C2-C3-O3
7	A	1108	GOL	O1-C1-C2-C3
7	A	1108	GOL	C1-C2-C3-O3
7	A	1116	GOL	O1-C1-C2-C3
7	A	1118	GOL	C1-C2-C3-O3
7	A	1120	GOL	O1-C1-C2-C3
7	B	1111	GOL	O1-C1-C2-C3
7	B	1112	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
7	B	1112	GOL	O1-C1-C2-C3
7	B	1117	GOL	O1-C1-C2-C3
5	L	1003	1PE	C16-C26-OH6-C15
5	H	1001	1PE	C16-C26-OH6-C15
5	B	1002	1PE	C16-C26-OH6-C15
5	B	1004	1PE	C16-C26-OH6-C15
5	L	1003	1PE	OH2-C12-C22-OH3
5	H	1001	1PE	OH6-C15-C25-OH5
5	H	1001	1PE	OH5-C14-C24-OH4
7	H	1105	GOL	O1-C1-C2-O2
7	B	1117	GOL	O1-C1-C2-O2
5	B	1004	1PE	OH6-C15-C25-OH5
4	B	902	EPE	C9-C10-S-O3S
5	B	1002	1PE	OH2-C12-C22-OH3
5	B	1004	1PE	OH2-C12-C22-OH3
5	L	1003	1PE	OH4-C13-C23-OH3
5	H	1001	1PE	OH4-C13-C23-OH3
5	B	1002	1PE	OH4-C13-C23-OH3
5	B	1004	1PE	OH4-C13-C23-OH3
7	L	1107	GOL	C1-C2-C3-O3
7	L	1110	GOL	C1-C2-C3-O3
7	L	1115	GOL	C1-C2-C3-O3
7	H	1105	GOL	C1-C2-C3-O3
7	A	1116	GOL	C1-C2-C3-O3
7	A	1120	GOL	C1-C2-C3-O3
7	B	1102	GOL	C1-C2-C3-O3
7	B	1111	GOL	C1-C2-C3-O3
7	B	1112	GOL	C1-C2-C3-O3
7	B	1117	GOL	C1-C2-C3-O3
7	L	1103	GOL	O2-C2-C3-O3
7	L	1119	GOL	O1-C1-C2-O2
7	A	1108	GOL	O1-C1-C2-O2
7	A	1120	GOL	O2-C2-C3-O3
6	L	1201	PEG	O1-C1-C2-O2
7	L	1107	GOL	O2-C2-C3-O3
7	H	1101	GOL	O2-C2-C3-O3
7	H	1105	GOL	O2-C2-C3-O3
7	A	1108	GOL	O2-C2-C3-O3
7	B	1111	GOL	O2-C2-C3-O3
7	B	1112	GOL	O2-C2-C3-O3
5	L	1003	1PE	OH6-C15-C25-OH5
7	A	1120	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
7	B	1117	GOL	O2-C2-C3-O3
7	H	1106	GOL	O2-C2-C3-O3
7	A	1116	GOL	O2-C2-C3-O3
7	B	1111	GOL	O1-C1-C2-O2
4	B	902	EPE	C9-C10-S-O2S
5	B	1004	1PE	C13-C23-OH3-C22
5	L	1003	1PE	C25-C15-OH6-C26
5	L	1003	1PE	C12-C22-OH3-C23
5	L	1003	1PE	C14-C24-OH4-C13
5	H	1001	1PE	OH2-C12-C22-OH3
5	H	1001	1PE	C24-C14-OH5-C25
5	H	1001	1PE	C12-C22-OH3-C23
5	B	1004	1PE	C24-C14-OH5-C25
5	B	1004	1PE	C15-C25-OH5-C14
7	A	1118	GOL	O1-C1-C2-O2
5	L	1003	1PE	OH7-C16-C26-OH6
7	A	1118	GOL	O2-C2-C3-O3
5	B	1004	1PE	C25-C15-OH6-C26
5	L	1003	1PE	C24-C14-OH5-C25
5	B	1002	1PE	OH5-C14-C24-OH4
5	L	1003	1PE	C15-C25-OH5-C14
5	B	1004	1PE	OH5-C14-C24-OH4
7	A	1116	GOL	O1-C1-C2-O2
5	B	1002	1PE	C25-C15-OH6-C26
5	B	1004	1PE	C14-C24-OH4-C13
5	B	1002	1PE	OH7-C16-C26-OH6
4	L	901	EPE	N4-C7-C8-O8
6	L	1201	PEG	C4-C3-O2-C2

There are no ring outliers.

24 monomers are involved in 61 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	H	1105	GOL	2	0
7	B	1117	GOL	1	0
7	A	1118	GOL	2	0
7	L	1103	GOL	5	0
3	H	802	SO4	0	4
7	L	1107	GOL	3	0
7	L	1115	GOL	2	0
7	A	1108	GOL	1	0
4	L	901	EPE	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1116	GOL	8	0
5	H	1001	1PE	3	0
7	L	1119	GOL	2	0
3	H	801	SO4	0	4
6	H	1202	PEG	1	0
5	L	1003	1PE	2	0
5	B	1002	1PE	3	0
6	L	1201	PEG	3	0
7	H	1101	GOL	1	0
7	B	1112	GOL	2	0
7	H	1106	GOL	4	0
7	L	1110	GOL	1	0
4	B	902	EPE	1	0
5	B	1004	1PE	5	0
7	B	1111	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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%)	0.56	16 (7%) 20 23	17, 28, 42, 50	1 (0%)
1	L	213/214 (99%)	0.14	6 (2%) 55 61	13, 23, 35, 49	1 (0%)
2	B	217/224 (96%)	0.50	11 (5%) 33 39	12, 26, 44, 48	3 (1%)
2	H	220/224 (98%)	0.22	11 (5%) 34 39	9, 22, 36, 56	3 (1%)
All	All	864/876 (98%)	0.36	44 (5%) 33 39	9, 25, 41, 56	8 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	CYS	5.2
2	H	64	LYS	4.6
2	B	128	SER	4.5
2	H	130	ALA	4.4
2	H	215	CYS	4.3
2	H	129	ALA	4.3
2	B	214	ASP	4.3
1	L	213	GLU	4.1
1	A	93	GLY	3.8
2	B	97	TYR	3.7
1	A	69	THR	3.4
1	L	126	THR	3.3
2	H	128	SER	3.1
2	B	65	GLY	3.0
1	A	203	SER	2.9
1	A	202	THR	2.8
2	H	1	GLN	2.7
1	A	2	LEU	2.7
2	H	133	ASN	2.6
1	A	212	ASN	2.6
1	L	128	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	40	GLN	2.5
2	B	98	ARG	2.4
1	A	76	ASN	2.4
2	B	28	THR	2.4
2	B	5	LEU	2.4
2	B	159	LEU	2.3
1	L	181	LEU	2.3
1	A	156	GLN	2.3
2	B	157	GLY	2.2
1	L	201	SER	2.2
2	B	22[A]	CYS	2.2
1	A	60	SER	2.2
2	H	134	SER	2.2
2	H	158	SER	2.2
1	A	23	CYS	2.1
1	A	68	GLY	2.1
2	H	84	SER	2.1
1	A	78	LEU	2.1
2	H	214	ASP	2.1
1	L	202	THR	2.1
2	B	100	TYR	2.1
1	A	77	SER	2.1
1	A	92	TYR	2.0

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 [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.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	GOL	A	1116	6/6	0.65	0.19	60,61,62,62	0
7	GOL	H	1101	6/6	0.69	0.17	40,41,42,42	0
7	GOL	B	1102	6/6	0.71	0.21	49,51,52,53	0
7	GOL	B	1112	6/6	0.72	0.20	66,66,67,68	0
3	SO4	A	806	5/5	0.73	0.15	80,80,80,81	0
7	GOL	A	1120	6/6	0.74	0.20	52,54,55,56	0
7	GOL	L	1103	6/6	0.75	0.21	65,65,66,67	0
7	GOL	H	1106	6/6	0.75	0.20	45,46,48,48	0
5	1PE	B	1002	16/16	0.76	0.23	63,64,68,69	0
5	1PE	B	1004	16/16	0.77	0.17	56,57,62,63	0
5	1PE	L	1003	16/16	0.77	0.25	68,69,70,71	0
7	GOL	L	1119	6/6	0.77	0.20	60,61,63,65	0
7	GOL	L	1107	6/6	0.79	0.16	41,45,46,48	0
7	GOL	L	1115	6/6	0.79	0.18	52,55,55,56	0
6	PEG	H	1202	7/7	0.79	0.21	45,45,46,48	0
7	GOL	H	1105	6/6	0.81	0.19	50,51,52,55	0
7	GOL	L	1110	6/6	0.82	0.16	61,62,62,62	0
7	GOL	B	1111	6/6	0.82	0.15	65,65,65,65	0
5	1PE	H	1001	16/16	0.82	0.20	62,64,67,68	0
7	GOL	B	1117	6/6	0.82	0.18	58,59,59,60	0
7	GOL	A	1118	6/6	0.83	0.14	40,42,44,45	0
7	GOL	A	1108	6/6	0.83	0.16	58,59,59,59	0
3	SO4	H	805	5/5	0.83	0.12	69,69,69,70	0
6	PEG	L	1201	7/7	0.84	0.18	54,56,56,57	0
3	SO4	B	809	5/5	0.84	0.12	64,65,65,66	0
4	EPE	B	902	15/15	0.87	0.15	45,52,57,57	0
3	SO4	H	802	5/5	0.88	0.19	32,32,34,36	4
3	SO4	L	808	5/5	0.89	0.10	59,60,60,61	0
3	SO4	B	807	5/5	0.89	0.10	52,52,53,53	0
3	SO4	H	804	5/5	0.93	0.09	52,52,54,54	0
4	EPE	L	901	15/15	0.95	0.12	23,40,48,48	0
3	SO4	H	803	5/5	0.96	0.07	41,43,44,44	0
3	SO4	H	801	5/5	0.99	0.04	22,22,23,23	4

6.5 Other polymers ⓘ

There are no such residues in this entry.