



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:25 AM UTC

PDB ID : 5UTZ / pdb_00005utz
Title : Human IL-2/Fab complex
Authors : Jude, K.M.; Garcia, K.C.
Deposited on : 2017-02-15
Resolution : 2.75 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

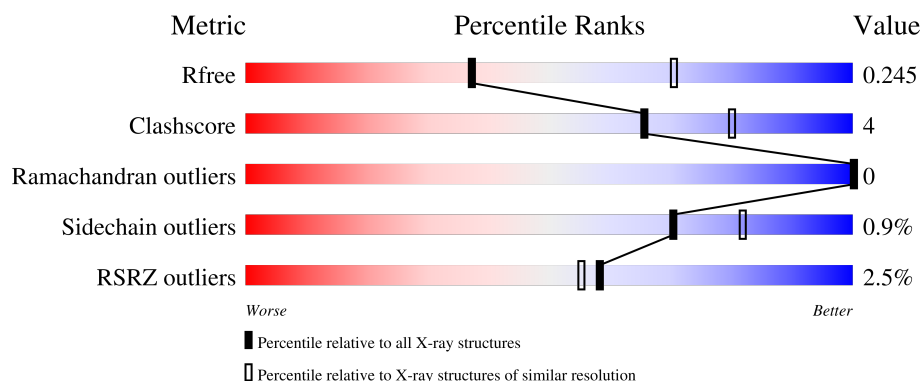
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	B	224	4% 86% 12% .
1	F	224	% 88% 12%
1	H	224	3% 86% 14%
1	J	224	2% 79% 18% .
2	C	216	91% 6% ..

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Mol	Chain	Length	Quality of chain
2	G	216	% 88% 10%
2	K	216	% 94% 6%
2	L	216	92% 7%
3	A	133	5% 80% 18%
3	D	133	2% 85% 10% 5%
3	E	133	3% 88% 7% 5%
3	I	133	10% 83% 11% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17641 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 Fab 5111 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	224	Total	C	N	O	S	0	1	0
			1683	1069	276	333	5			
1	B	220	Total	C	N	O	S	0	1	0
			1660	1057	272	326	5			
1	F	223	Total	C	N	O	S	0	1	0
			1677	1066	275	331	5			
1	J	219	Total	C	N	O	S	0	0	0
			1642	1046	268	323	5			

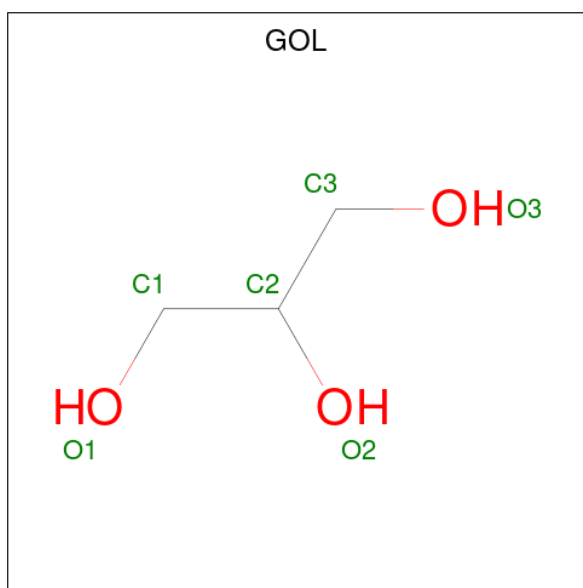
- Molecule 2 is a protein called Fab 5111 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	215	Total	C	N	O	S	0	1	0
			1626	1007	272	341	6			
2	C	213	Total	C	N	O	S	0	1	0
			1611	999	270	337	5			
2	G	213	Total	C	N	O	S	0	1	0
			1611	999	270	337	5			
2	K	215	Total	C	N	O	S	0	1	0
			1626	1007	272	341	6			

- Molecule 3 is a protein called Interleukin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	126	Total	C	N	O	S	0	0	0
			1034	666	171	190	7			
3	A	131	Total	C	N	O	S	0	0	0
			1065	684	176	198	7			
3	E	126	Total	C	N	O	S	0	0	0
			1034	666	171	190	7			
3	I	126	Total	C	N	O	S	0	0	0
			1033	665	171	190	7			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	G	1	Total	C	O	0	0
			6	3	3		
4	J	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	37	Total	O	0	0
			37	37		
5	L	39	Total	O	0	0
			39	39		
5	D	11	Total	O	0	0
			11	11		
5	B	31	Total	O	0	0
			31	31		
5	C	25	Total	O	0	0
			25	25		
5	A	14	Total	O	0	0
			14	14		

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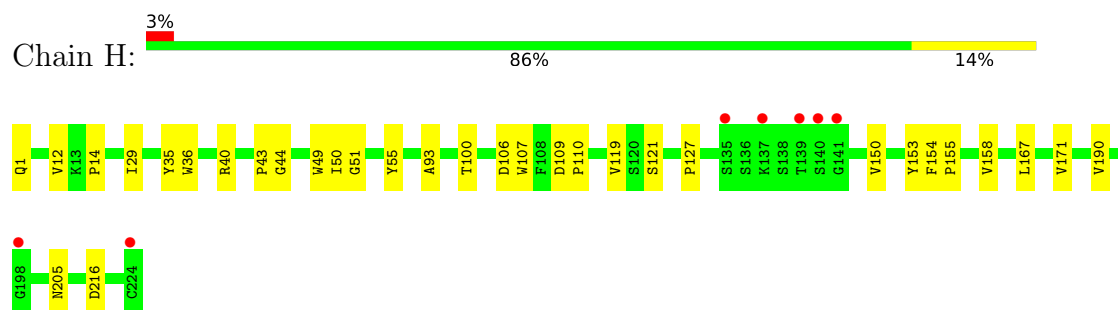
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	34	Total 34	O 34	0	0
5	G	46	Total 46	O 46	0	0
5	E	9	Total 9	O 9	0	0
5	J	29	Total 29	O 29	0	0
5	K	28	Total 28	O 28	0	0
5	I	6	Total 6	O 6	0	0

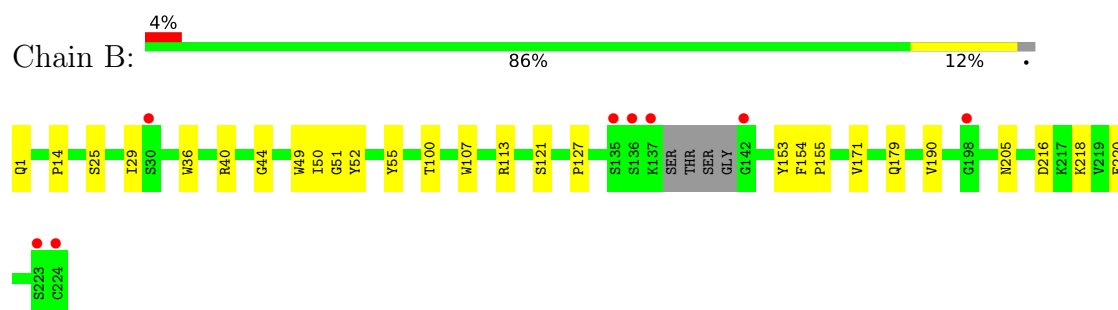
3 Residue-property plots [i](#)

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.

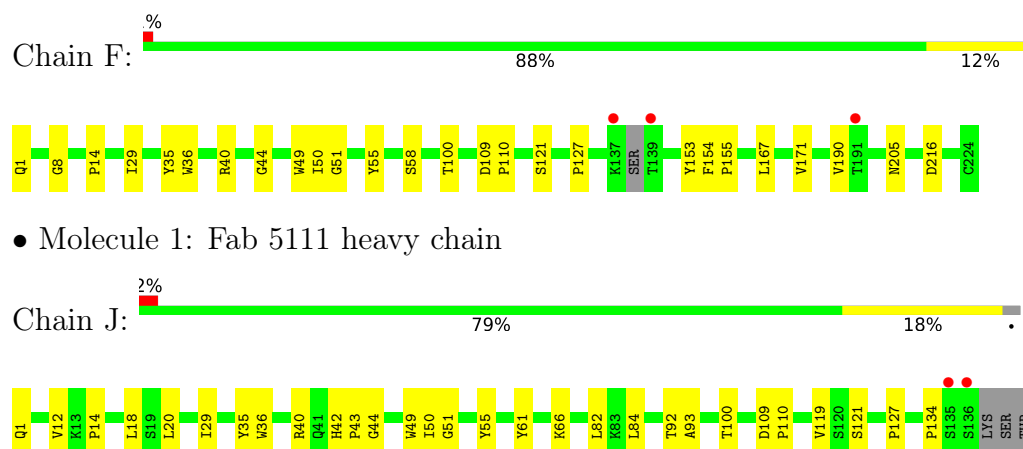
- Molecule 1: Fab 5111 heavy chain



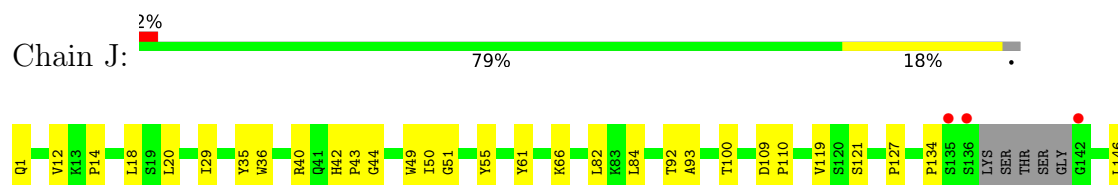
- Molecule 1: Fab 5111 heavy chain



- Molecule 1: Fab 5111 heavy chain



- Molecule 1: Fab 5111 heavy chain

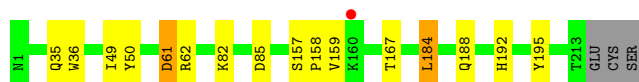




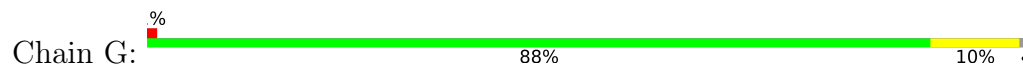
- Molecule 2: Fab 5111 light chain



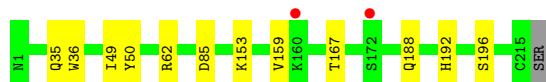
- Molecule 2: Fab 5111 light chain



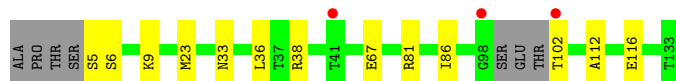
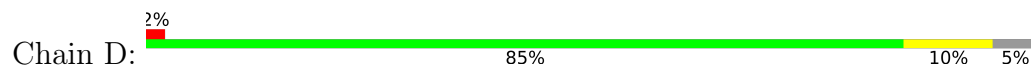
- Molecule 2: Fab 5111 light chain



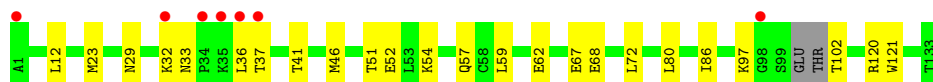
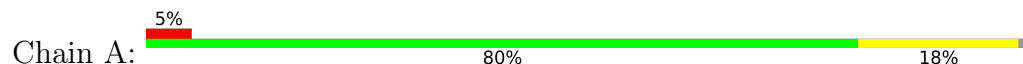
- Molecule 2: Fab 5111 light chain



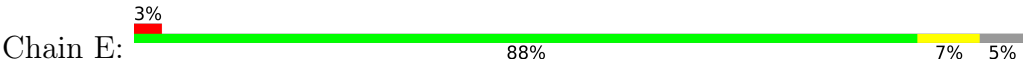
- Molecule 3: Interleukin-2



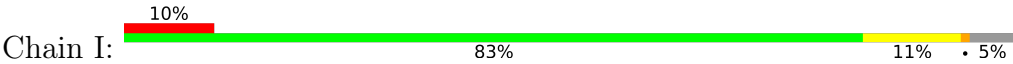
- Molecule 3: Interleukin-2



- Molecule 3: Interleukin-2



● Molecule 3: Interleukin-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	86.01Å 145.73Å 107.32Å 90.00° 95.38° 90.00°	Depositor
Resolution (Å)	86.17 – 2.75 86.17 – 2.75	Depositor EDS
% Data completeness (in resolution range)	98.8 (86.17-2.75) 98.9 (86.17-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.193 , 0.247 0.194 , 0.245	Depositor DCC
R_{free} test set	1973 reflections (2.87%)	wwPDB-VP
Wilson B-factor (Å ²)	51.5	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	17641	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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, PCA

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	B	0.25	0/1698	0.71	0/2323
1	F	0.26	0/1715	0.72	2/2346 (0.1%)
1	H	0.25	0/1722	0.71	0/2357
1	J	0.26	0/1680	0.72	0/2300
2	C	0.26	0/1650	0.69	0/2253
2	G	0.26	0/1650	0.69	0/2253
2	K	0.25	0/1665	0.69	0/2273
2	L	0.25	0/1665	0.71	0/2273
3	A	0.28	0/1082	0.70	0/1459
3	D	0.27	0/1050	0.68	0/1414
3	E	0.26	0/1050	0.68	0/1414
3	I	0.27	0/1049	0.74	0/1412
All	All	0.26	0/17676	0.70	2/24077 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	8	GLY	CA-C-N	5.00	124.69	119.64
1	F	8	GLY	C-N-CA	5.00	124.69	119.64

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	B	1660	0	1630	14	0
1	F	1677	0	1645	13	0
1	H	1683	0	1650	18	0
1	J	1642	0	1609	20	0
2	C	1611	0	1541	9	0
2	G	1611	0	1541	15	0
2	K	1626	0	1551	7	0
2	L	1626	0	1551	10	0
3	A	1065	0	1104	13	0
3	D	1034	0	1072	9	0
3	E	1034	0	1072	6	0
3	I	1033	0	1070	9	0
4	B	6	0	8	1	0
4	C	6	0	8	1	0
4	G	6	0	8	0	0
4	J	6	0	8	0	0
4	L	6	0	8	1	0
5	A	14	0	0	0	0
5	B	31	0	0	1	0
5	C	25	0	0	0	0
5	D	11	0	0	0	0
5	E	9	0	0	1	0
5	F	34	0	0	0	0
5	G	46	0	0	2	0
5	H	37	0	0	0	0
5	I	6	0	0	0	0
5	J	29	0	0	0	0
5	K	28	0	0	1	0
5	L	39	0	0	0	0
All	All	17641	0	17076	132	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 4.

All (132) 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)
3:I:33:ASN:HB3	3:I:36:LEU:HB2	1.65	0.77
1:J:134:PRO:HG3	1:J:146:LEU:HB3	1.69	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:44:GLY:HA3	2:L:167:THR:HG21	1.74	0.69
1:J:127:PRO:HB3	1:J:153:TYR:HB3	1.79	0.64
1:F:14:PRO:HD2	1:F:121:SER:HB3	1.79	0.63
1:H:127:PRO:HB3	1:H:153:TYR:HB3	1.81	0.63
2:C:61:ASP:N	2:C:61:ASP:OD1	2.32	0.62
1:B:127:PRO:HB3	1:B:153:TYR:HB3	1.82	0.62
2:L:118:SER:HA	4:L:301:GOL:H32	1.83	0.61
1:B:14:PRO:HD2	1:B:121:SER:HB3	1.82	0.61
3:A:33:ASN:HB3	3:A:36:LEU:HB2	1.82	0.61
3:I:31:TYR:HH	3:I:75:SER:HG	1.49	0.60
3:E:67:GLU:HB2	3:E:86:ILE:HG13	1.83	0.60
3:E:48:LYS:HE2	3:E:106:GLU:HB3	1.84	0.60
1:F:127:PRO:HB3	1:F:153:TYR:HB3	1.83	0.59
1:J:14:PRO:HD2	1:J:121:SER:HB3	1.84	0.59
1:B:100:THR:HG23	1:B:107:TRP:H	1.67	0.58
3:D:5:SER:OG	3:D:6:SER:N	2.32	0.57
2:K:62:ARG:NH1	2:K:85:ASP:OD2	2.24	0.57
1:J:194:SER:HA	1:J:197:LEU:HD13	1.86	0.57
1:H:14:PRO:HD2	1:H:121:SER:HB3	1.86	0.56
3:A:67:GLU:HB2	3:A:86:ILE:HG13	1.87	0.56
2:L:154:ALA:HB1	2:L:192:HIS:CD2	2.40	0.56
2:K:188:GLN:O	2:K:192:HIS:ND1	2.39	0.56
1:B:205:ASN:ND2	1:B:216:ASP:OD1	2.38	0.55
3:D:67:GLU:HB2	3:D:86:ILE:HG13	1.88	0.55
2:C:188:GLN:O	2:C:192:HIS:ND1	2.40	0.55
3:D:33:ASN:HB3	3:D:36:LEU:HB2	1.90	0.54
3:I:106:GLU:N	3:I:106:GLU:OE1	2.38	0.53
1:F:35:TYR:HB2	1:F:100:THR:HG23	1.91	0.53
2:G:8[B]:HIS:ND1	5:G:402:HOH:O	2.34	0.52
1:J:40:ARG:HB3	1:J:50:ILE:HD11	1.90	0.52
1:F:29:ILE:HG23	1:F:36:TRP:NE1	2.25	0.52
1:H:35:TYR:OH	3:D:23:MET:HE1	2.10	0.51
1:J:205:ASN:ND2	1:J:216:ASP:OD1	2.33	0.51
1:F:40:ARG:HB3	1:F:50:ILE:HD11	1.93	0.51
3:D:38:ARG:NH2	3:A:62:GLU:OE1	2.44	0.51
1:B:29:ILE:HG23	1:B:36:TRP:NE1	2.26	0.51
1:B:171:VAL:HG22	1:B:190:VAL:HG22	1.91	0.51
3:A:37:THR:O	3:A:41:THR:N	2.43	0.51
2:C:62:ARG:NH1	2:C:85:ASP:OD2	2.38	0.51
2:G:154:ALA:HB1	2:G:192:HIS:CD2	2.45	0.51
2:G:62:ARG:NH1	2:G:85:ASP:OD2	2.30	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:29:ILE:HG23	1:J:36:TRP:NE1	2.26	0.50
1:H:35:TYR:HB2	1:H:100:THR:HG23	1.93	0.50
3:I:112:ALA:HB1	3:I:116:GLU:HB2	1.94	0.49
1:F:205:ASN:ND2	1:F:216:ASP:OD1	2.38	0.49
1:F:171:VAL:HG22	1:F:190:VAL:HG22	1.94	0.49
2:L:153:LYS:HD3	2:L:158:PRO:HA	1.94	0.49
2:G:148:VAL:HG12	2:G:201:HIS:HB2	1.95	0.49
3:A:51:THR:HG23	3:A:52:GLU:HG3	1.93	0.49
3:A:29:ASN:O	3:A:32:LYS:NZ	2.46	0.48
1:J:35:TYR:HB2	1:J:100:THR:HG23	1.95	0.48
1:H:40:ARG:HB3	1:H:50:ILE:HD11	1.95	0.48
1:H:29:ILE:HG23	1:H:36:TRP:NE1	2.28	0.48
2:C:36:TRP:HB2	2:C:49:ILE:HB	1.96	0.48
3:E:76:LYS:NZ	5:E:201:HOH:O	2.45	0.48
2:G:36:TRP:HB2	2:G:49:ILE:HB	1.96	0.47
2:L:35:GLN:HG3	2:L:50:TYR:HA	1.94	0.47
2:G:35:GLN:HG3	2:G:50:TYR:HA	1.96	0.47
1:J:43:PRO:HD3	1:J:93:ALA:HA	1.95	0.47
2:L:40:ARG:NH1	2:L:84:GLU:O	2.45	0.47
2:G:108:THR:HG21	2:G:145:PRO:HB3	1.97	0.47
2:G:55:ARG:NH1	5:G:407:HOH:O	2.48	0.47
1:J:20:LEU:HD12	1:J:82:LEU:HD23	1.97	0.47
1:B:113:ARG:NH2	5:B:403:HOH:O	2.47	0.47
3:D:112:ALA:HB1	3:D:116:GLU:HB3	1.97	0.46
1:F:167:LEU:HD21	1:F:190:VAL:HG11	1.97	0.46
2:G:22:CYS:N	2:G:74:ALA:O	2.48	0.46
2:C:184:LEU:HD21	2:C:195:TYR:CZ	2.50	0.46
1:H:106:ASP:OD2	3:D:81:ARG:NH2	2.48	0.46
2:C:35:GLN:HG3	2:C:50:TYR:HA	1.97	0.46
2:L:10:VAL:HG21	2:L:20:ILE:HG12	1.98	0.46
2:K:153:LYS:HB2	2:K:196:SER:HB2	1.96	0.46
2:C:82:LYS:HG2	4:C:301:GOL:H2	1.98	0.46
2:K:36:TRP:HB2	2:K:49:ILE:HB	1.97	0.46
1:B:218:LYS:HE2	1:B:220:GLU:HG2	1.98	0.45
1:F:49:TRP:CZ2	1:F:51:GLY:HA2	2.52	0.45
1:F:58:SER:OG	3:E:20:ASP:OD2	2.34	0.45
3:A:52:GLU:OE1	3:A:54:LYS:HD3	2.16	0.45
2:K:35:GLN:HG3	2:K:50:TYR:HA	1.98	0.45
1:B:40:ARG:HB3	1:B:50:ILE:HD11	1.97	0.45
2:G:157:SER:HA	2:G:158:PRO:HD3	1.87	0.45
2:G:36:TRP:CD2	2:G:76:LEU:HB2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:171:VAL:HG22	1:H:190:VAL:HG22	1.98	0.45
3:I:44:PHE:CE2	3:I:114:ILE:HG12	2.52	0.45
2:G:33:TYR:CE1	3:E:76:LYS:HD3	2.53	0.44
1:B:49:TRP:CZ2	1:B:51:GLY:HA2	2.53	0.43
1:H:109:ASP:HA	1:H:110:PRO:HA	1.79	0.43
1:H:205:ASN:ND2	1:H:216:ASP:OD1	2.37	0.43
3:D:102:THR:O	3:D:102:THR:OG1	2.34	0.43
1:B:154:PHE:HA	1:B:155:PRO:HA	1.78	0.43
1:J:44:GLY:HA3	2:K:167:THR:HG21	1.99	0.43
3:A:59:LEU:HD13	3:A:121:TRP:CD1	2.53	0.43
1:J:18:LEU:HB3	1:J:84:LEU:HB3	2.01	0.43
1:J:49:TRP:CZ2	1:J:51:GLY:HA2	2.54	0.43
2:L:62:ARG:NH1	2:L:85:ASP:OD1	2.50	0.43
1:J:134:PRO:HD2	1:J:221:PRO:HA	2.01	0.43
1:J:42:HIS:ND1	1:J:92:THR:O	2.52	0.43
1:H:154:PHE:HA	1:H:155:PRO:HA	1.79	0.43
3:A:23:MET:HE3	3:A:80:LEU:HD22	2.01	0.43
1:J:167:LEU:HD21	1:J:190:VAL:HG11	2.01	0.43
3:A:37:THR:O	3:A:41:THR:OG1	2.19	0.42
2:K:192:HIS:NE2	5:K:302:HOH:O	2.37	0.42
1:H:49:TRP:CZ2	1:H:51:GLY:HA2	2.54	0.42
1:H:43:PRO:HD3	1:H:93:ALA:HA	2.01	0.42
3:A:68:GLU:O	3:A:72:LEU:HD13	2.20	0.42
3:I:37:THR:O	3:I:41:THR:N	2.52	0.42
3:I:48:LYS:HE2	3:I:106:GLU:HB2	2.02	0.42
1:F:154:PHE:HA	1:F:155:PRO:HA	1.79	0.42
3:I:15:GLU:HA	3:I:18:LEU:HB3	2.02	0.42
2:L:22:CYS:N	2:L:74:ALA:O	2.53	0.41
1:J:150:VAL:HG11	1:J:158:VAL:HG11	2.03	0.41
2:C:157:SER:HA	2:C:158:PRO:HD3	1.91	0.41
3:I:106:GLU:H	3:I:106:GLU:CD	2.26	0.41
3:A:46:MET:HE1	3:A:120:ARG:HB2	2.01	0.41
3:A:57:GLN:OE1	3:A:97:LYS:NZ	2.48	0.41
3:D:5:SER:HB3	3:D:9:LYS:HD2	2.03	0.41
1:F:109:ASP:HA	1:F:110:PRO:HA	1.79	0.41
2:G:36:TRP:CZ3	2:G:91:CYS:HB3	2.55	0.41
2:G:95:ASP:O	3:E:81:ARG:NH1	2.53	0.41
1:J:12:VAL:HG13	1:J:119:VAL:HG22	2.03	0.41
1:J:109:ASP:HA	1:J:110:PRO:HA	1.82	0.41
1:H:167:LEU:HD21	1:H:190:VAL:HG11	2.03	0.41
1:B:25:SER:HA	4:B:301:GOL:H2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:12:VAL:O	1:H:119:VAL:HA	2.20	0.41
1:B:179[B]:GLN:HE21	1:B:179[B]:GLN:HB2	1.66	0.41
1:H:107:TRP:CD1	2:L:35:GLN:HG2	2.56	0.40
1:H:150:VAL:HG11	1:H:158:VAL:HG11	2.02	0.40
1:B:44:GLY:HA3	2:C:167:THR:HG21	2.03	0.40
1:J:61:TYR:HD2	1:J:66:LYS:HD2	1.86	0.40
1:F:44:GLY:HA3	2:G:167:THR:HG21	2.03	0.40

There are no symmetry-related clashes.

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	B	217/224 (97%)	212 (98%)	5 (2%)	0	100	100
1	F	220/224 (98%)	215 (98%)	5 (2%)	0	100	100
1	H	223/224 (100%)	216 (97%)	7 (3%)	0	100	100
1	J	215/224 (96%)	209 (97%)	6 (3%)	0	100	100
2	C	212/216 (98%)	206 (97%)	6 (3%)	0	100	100
2	G	212/216 (98%)	207 (98%)	5 (2%)	0	100	100
2	K	214/216 (99%)	210 (98%)	4 (2%)	0	100	100
2	L	214/216 (99%)	208 (97%)	6 (3%)	0	100	100
3	A	127/133 (96%)	124 (98%)	3 (2%)	0	100	100
3	D	122/133 (92%)	122 (100%)	0	0	100	100
3	E	122/133 (92%)	121 (99%)	1 (1%)	0	100	100
3	I	122/133 (92%)	120 (98%)	2 (2%)	0	100	100
All	All	2220/2292 (97%)	2170 (98%)	50 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	B	190/192 (99%)	188 (99%)	2 (1%)	65	79
1	F	192/192 (100%)	191 (100%)	1 (0%)	81	88
1	H	193/192 (100%)	192 (100%)	1 (0%)	81	88
1	J	188/192 (98%)	186 (99%)	2 (1%)	65	79
2	C	187/189 (99%)	184 (98%)	3 (2%)	55	72
2	G	187/189 (99%)	186 (100%)	1 (0%)	81	88
2	K	189/189 (100%)	188 (100%)	1 (0%)	81	88
2	L	189/189 (100%)	189 (100%)	0	100	100
3	A	124/126 (98%)	122 (98%)	2 (2%)	55	72
3	D	120/126 (95%)	120 (100%)	0	100	100
3	E	120/126 (95%)	118 (98%)	2 (2%)	53	71
3	I	120/126 (95%)	118 (98%)	2 (2%)	53	71
All	All	1999/2028 (99%)	1982 (99%)	17 (1%)	70	82

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	55	TYR
1	B	52	TYR
1	B	55	TYR
2	C	61	ASP
2	C	159	VAL
2	C	184	LEU
3	A	12	LEU
3	A	102	THR
1	F	55	TYR
2	G	76	LEU
3	E	102	THR
3	E	103	PHE
1	J	55	TYR
1	J	179	GLN

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Mol	Chain	Res	Type
2	K	159	VAL
3	I	38	ARG
3	I	41	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	H	200	GLN
2	L	38	GLN
1	B	79	GLN
1	B	207	ASN
2	C	54	GLN
3	A	79	HIS
1	F	16	GLN
1	F	200	GLN
2	G	192	HIS
3	E	88	ASN
1	J	16	GLN
1	J	207	ASN
2	K	38	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	PCA	J	1	1	7,8,9	1.92	1 (14%)	9,10,12	2.30	5 (55%)
1	PCA	H	1	1	7,8,9	1.92	1 (14%)	9,10,12	2.38	5 (55%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	PCA	B	1	1	7,8,9	1.92	1 (14%)	9,10,12	2.24	5 (55%)
1	PCA	F	1	1	7,8,9	1.91	1 (14%)	9,10,12	2.29	5 (55%)

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
1	PCA	J	1	1	-	0/0/11/13	0/1/1/1
1	PCA	H	1	1	-	0/0/11/13	0/1/1/1
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1
1	PCA	F	1	1	-	0/0/11/13	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	1	PCA	CD-N	4.97	1.46	1.34
1	J	1	PCA	CD-N	4.96	1.46	1.34
1	B	1	PCA	CD-N	4.95	1.46	1.34
1	F	1	PCA	CD-N	4.93	1.46	1.34

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	1	PCA	CB-CA-C	-3.77	107.49	112.66
1	J	1	PCA	CB-CA-C	-3.48	107.89	112.66
1	F	1	PCA	CB-CA-C	-3.25	108.20	112.66
1	B	1	PCA	OE-CD-CG	-2.98	121.41	126.72
1	H	1	PCA	CA-N-CD	-2.97	103.42	113.58
1	B	1	PCA	CB-CA-C	-2.96	108.60	112.66
1	F	1	PCA	OE-CD-CG	-2.95	121.45	126.72
1	H	1	PCA	OE-CD-CG	-2.94	121.47	126.72
1	J	1	PCA	CA-N-CD	-2.93	103.56	113.58
1	F	1	PCA	CA-N-CD	-2.91	103.60	113.58
1	J	1	PCA	OE-CD-CG	-2.90	121.54	126.72
1	B	1	PCA	CA-N-CD	-2.89	103.67	113.58
1	B	1	PCA	CB-CA-N	2.78	110.89	103.24
1	J	1	PCA	CB-CA-N	2.74	110.77	103.24
1	H	1	PCA	CB-CA-N	2.72	110.72	103.24
1	F	1	PCA	CB-CA-N	2.72	110.71	103.24
1	B	1	PCA	CG-CD-N	2.59	114.73	108.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	1	PCA	CG-CD-N	2.56	114.64	108.39
1	F	1	PCA	CG-CD-N	2.55	114.63	108.39
1	J	1	PCA	CG-CD-N	2.53	114.59	108.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	GOL	G	301	-	5,5,5	0.36	0	5,5,5	0.34	0
4	GOL	L	301	-	5,5,5	0.37	0	5,5,5	0.29	0
4	GOL	J	301	-	5,5,5	0.38	0	5,5,5	0.33	0
4	GOL	B	301	-	5,5,5	0.37	0	5,5,5	0.32	0
4	GOL	C	301	-	5,5,5	0.37	0	5,5,5	0.35	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	G	301	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	L	301	-	-	2/4/4/4	-
4	GOL	J	301	-	-	0/4/4/4	-
4	GOL	B	301	-	-	4/4/4/4	-
4	GOL	C	301	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	301	GOL	O1-C1-C2-C3
4	B	301	GOL	O1-C1-C2-O2
4	B	301	GOL	O1-C1-C2-C3
4	B	301	GOL	C1-C2-C3-O3
4	B	301	GOL	O2-C2-C3-O3
4	C	301	GOL	O1-C1-C2-C3
4	G	301	GOL	O1-C1-C2-C3
4	L	301	GOL	O1-C1-C2-O2
4	G	301	GOL	O1-C1-C2-O2
4	C	301	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	301	GOL	1	0
4	B	301	GOL	1	0
4	C	301	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	B	219/224 (97%)	-0.14	8 (3%) 45 43	27, 45, 86, 158	1 (0%)
1	F	222/224 (99%)	-0.11	3 (1%) 73 72	25, 47, 77, 133	1 (0%)
1	H	223/224 (99%)	-0.15	7 (3%) 51 49	25, 43, 78, 161	1 (0%)
1	J	218/224 (97%)	-0.04	5 (2%) 61 58	37, 53, 102, 180	0
2	C	213/216 (98%)	-0.07	1 (0%) 87 87	34, 58, 78, 113	1 (0%)
2	G	213/216 (98%)	-0.11	2 (0%) 81 81	28, 51, 73, 111	1 (0%)
2	K	215/216 (99%)	-0.03	2 (0%) 81 81	28, 57, 85, 176	1 (0%)
2	L	215/216 (99%)	-0.20	1 (0%) 87 87	24, 48, 72, 127	1 (0%)
3	A	131/133 (98%)	0.19	7 (5%) 32 31	36, 56, 105, 151	0
3	D	126/133 (94%)	0.36	3 (2%) 59 56	43, 67, 116, 144	0
3	E	126/133 (94%)	0.35	4 (3%) 50 48	38, 66, 102, 137	0
3	I	126/133 (94%)	0.73	13 (10%) 12 12	47, 79, 136, 167	0
All	All	2247/2292 (98%)	0.01	56 (2%) 58 55	24, 53, 101, 180	7 (0%)

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	141	GLY	4.9
3	A	37	THR	4.6
1	F	137	LYS	4.0
3	I	103	PHE	3.7
1	H	140	SER	3.4
3	E	102	THR	3.2
2	G	8[A]	HIS	3.2
3	I	37	THR	3.2
3	A	98	GLY	3.1
3	I	34	PRO	3.0
2	G	213	THR	2.9

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Mol	Chain	Res	Type	RSRZ
3	E	6	SER	2.9
3	E	7	THR	2.9
3	D	98	GLY	2.8
1	H	135	SER	2.8
1	H	139	THR	2.8
1	J	136	SER	2.7
3	A	36	LEU	2.7
3	A	32	LYS	2.7
1	J	222	LYS	2.6
3	D	102	THR	2.6
3	I	7	THR	2.6
3	I	36	LEU	2.6
1	B	224	CYS	2.6
1	J	224	CYS	2.6
1	B	142	GLY	2.5
1	B	137	LYS	2.5
1	H	198	GLY	2.5
1	B	198	GLY	2.5
3	A	34	PRO	2.4
1	B	135	SER	2.4
1	B	136	SER	2.4
2	K	160	LYS	2.4
1	B	30	SER	2.3
3	I	4	SER	2.3
1	H	137	LYS	2.3
1	F	139	THR	2.3
3	D	41	THR	2.3
1	J	142	GLY	2.3
3	A	35	LYS	2.2
1	H	224	CYS	2.2
1	B	223	SER	2.2
3	I	32	LYS	2.2
3	I	104	MET	2.2
3	I	48	LYS	2.2
3	I	51	THR	2.1
3	I	35	LYS	2.1
2	C	160	LYS	2.1
3	I	49	LYS	2.1
2	K	172	SER	2.1
2	L	213	THR	2.1
3	I	10	THR	2.1
1	J	135	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	F	191	THR	2.0
3	A	1	ALA	2.0
3	E	98	GLY	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	PCA	B	1	8/9	0.87	0.13	39,73,77,78	0
1	PCA	H	1	8/9	0.88	0.11	57,64,75,78	0
1	PCA	F	1	8/9	0.89	0.14	59,73,83,87	0
1	PCA	J	1	8/9	0.92	0.10	47,61,76,77	0

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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	C	301	6/6	0.81	0.17	57,66,71,74	0
4	GOL	L	301	6/6	0.84	0.12	59,75,83,87	0
4	GOL	J	301	6/6	0.84	0.15	71,74,76,77	0
4	GOL	B	301	6/6	0.88	0.17	55,64,69,69	0
4	GOL	G	301	6/6	0.90	0.14	45,55,58,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.