



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:35 PM UTC

PDB ID : 8TQA / pdb_00008tqa
Title : Crystal structure of Fab.28.14.8 in complex with MHC-I (H2-Db)
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Deposited on : 2023-08-06
Resolution : 2.60 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

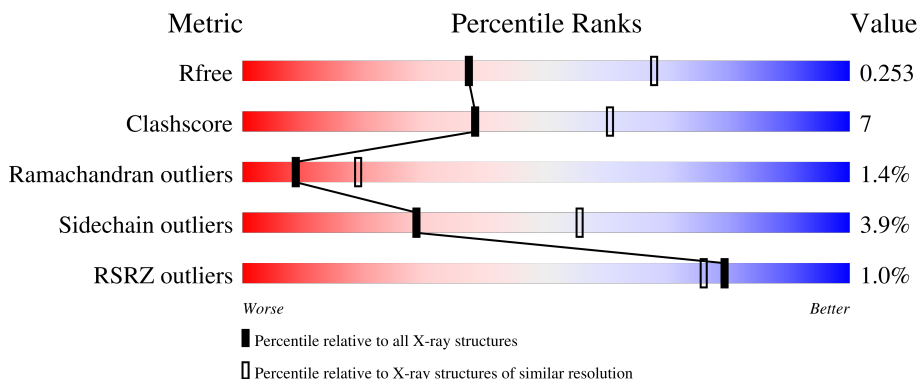
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	273	2% 82% 15% ..
2	B	99	% 81% 17% ..
3	H	217	% 76% 19% ..
4	L	214	79% 20%
5	P	9	89% 11%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6430 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 H-2 class I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2182	1381	383	409	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			772	494	126	145	7			

- Molecule 3 is a protein called Fab.28.14S Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	214	Total	C	N	O	S	0	0	0
			1620	1032	261	320	7			

- Molecule 4 is a protein called Fab.28.14S Light Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	213	Total	C	N	O	S	0	0	0
			1660	1034	280	340	6			

- Molecule 5 is a protein called influenza peptide.

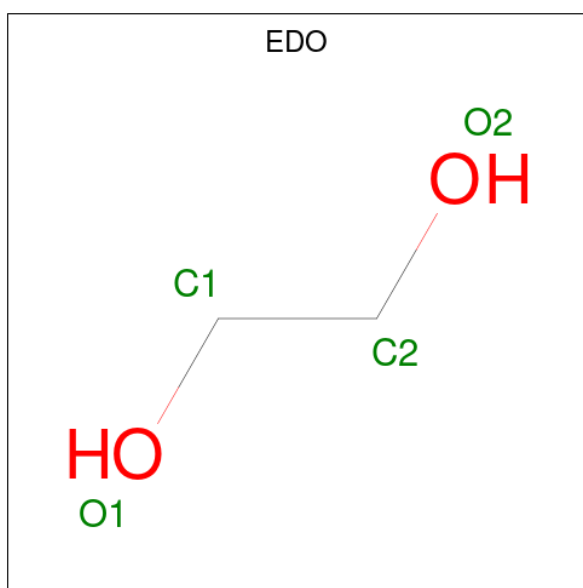
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	P	9	Total	C	N	O	S	0	0	0
			68	38	11	17	2			

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



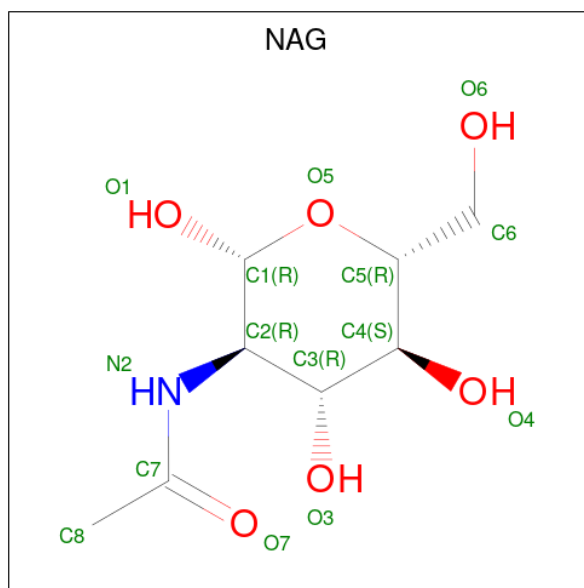
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	B	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			4	2	2		
7	H	1	Total	C	O	0	0
			4	2	2		
7	L	1	Total	C	O	0	0
			4	2	2		
7	L	1	Total	C	O	0	0
			4	2	2		
7	L	1	Total	C	O	0	0
			4	2	2		
7	P	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	H	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	24	Total	O	0	0
			24	24		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	3	Total 3	O 3	0	0
9	H	22	Total 22	O 22	0	0
9	L	17	Total 17	O 17	0	0

- Molecule 5: influenza peptide

Chain P:  89% 11%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.39Å 70.19Å 94.34Å 90.00° 98.00° 90.00°	Depositor
Resolution (Å)	60.57 – 2.60 60.57 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (60.57-2.60) 98.9 (60.57-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.181 , 0.253 0.181 , 0.253	Depositor DCC
R_{free} test set	1425 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.651	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 45.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6430	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.06% 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, EDO, NAG

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	0.43	0/2247	0.58	0/3054
2	B	0.35	0/798	0.56	0/1090
3	H	0.46	0/1665	0.67	1/2276 (0.0%)
4	L	0.41	0/1698	0.64	0/2307
5	P	0.42	0/67	0.68	0/88
All	All	0.42	0/6475	0.62	1/8815 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	H	61	ASN	N-CA-C	-5.40	101.86	109.96

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	A	2182	0	2023	31	0
2	B	772	0	702	10	0
3	H	1620	0	1559	23	0
4	L	1660	0	1586	23	0
5	P	68	0	61	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	12	0	16	4	0
7	A	4	0	6	3	0
7	B	8	0	12	0	0
7	H	4	0	6	0	0
7	L	16	0	24	2	0
7	P	4	0	6	0	0
8	H	14	0	13	1	0
9	A	24	0	0	1	0
9	B	3	0	0	0	0
9	H	22	0	0	0	0
9	L	17	0	0	0	0
All	All	6430	0	6014	85	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 7.

All (85) 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)
6:A:303:GOL:H2	5:P:3:ASN:HB3	1.74	0.69
3:H:18:VAL:HG22	3:H:86:LEU:HD11	1.76	0.67
1:A:6:ARG:HH11	6:A:301:GOL:H2	1.60	0.67
4:L:66:GLY:H	7:L:302:EDO:H11	1.62	0.65
4:L:4:MET:HE3	4:L:23:CYS:SG	2.37	0.65
1:A:13:SER:O	1:A:92:SER:HA	1.98	0.64
1:A:6:ARG:HD2	6:A:301:GOL:H2	1.80	0.64
1:A:6:ARG:NH2	1:A:102:ASP:OD1	2.30	0.64
4:L:193:THR:HG23	4:L:208:SER:HB2	1.81	0.63
3:H:28:ASN:ND2	8:H:301:NAG:O5	2.32	0.62
3:H:141:LEU:HD22	3:H:213:ILE:HG21	1.80	0.62
1:A:35:ARG:NH2	2:B:54:MET:O	2.33	0.61
4:L:115:VAL:O	4:L:207:LYS:HE3	2.01	0.60
1:A:133:TRP:HB2	1:A:144:ARG:HG3	1.82	0.59
3:H:10:GLU:HB3	3:H:112:LEU:HD23	1.83	0.59
1:A:202:ARG:HD3	1:A:244:TRP:CD2	2.40	0.56
3:H:11:LEU:HA	3:H:113:THR:O	2.04	0.56
4:L:54:LEU:HD11	4:L:60:SER:HA	1.87	0.56
1:A:97:GLN:HG2	1:A:116:PHE:CD2	2.40	0.56
3:H:151:GLU:OE2	3:H:171:ALA:HB3	2.06	0.55
4:L:163:TRP:NE1	4:L:175:MET:HE3	2.22	0.54
4:L:54:LEU:HD22	4:L:58:VAL:HB	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:LEU:C	1:A:174:ASN:H	2.15	0.54
1:A:192:HIS:CE1	2:B:98:ASP:HB3	2.42	0.53
3:H:97:ALA:HB1	3:H:103:PHE:HB3	1.91	0.53
1:A:195:SER:OG	1:A:196:LYS:N	2.39	0.52
1:A:97:GLN:HG3	9:A:403:HOH:O	2.10	0.51
1:A:58:GLU:O	1:A:62:ARG:HG3	2.11	0.51
3:H:63:LYS:HD2	3:H:64:PHE:CE2	2.46	0.51
3:H:180:LEU:C	3:H:180:LEU:HD12	2.36	0.51
1:A:195:SER:HB3	1:A:198:GLU:HB2	1.92	0.50
4:L:25:ALA:O	4:L:69:THR:OG1	2.30	0.50
4:L:163:TRP:CE2	4:L:175:MET:HE3	2.46	0.50
4:L:19:VAL:HG21	4:L:78:LEU:HD22	1.94	0.49
3:H:166:VAL:HG22	3:H:184:VAL:HG23	1.95	0.48
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.95	0.48
4:L:190:ASN:O	4:L:210:ASN:HA	2.14	0.47
4:L:105:GLU:OE1	4:L:173:TYR:OH	2.29	0.47
4:L:184:ASP:O	4:L:188:ARG:HG2	2.15	0.47
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.50	0.46
4:L:21:ILE:HG12	4:L:102:THR:HG21	1.96	0.46
4:L:190:ASN:HA	4:L:211:ARG:HB3	1.98	0.46
1:A:264:GLU:N	7:A:302:EDO:O2	2.44	0.45
4:L:135:PHE:C	4:L:136:LEU:HD12	2.41	0.45
4:L:86:TYR:O	4:L:101:GLY:O	2.35	0.45
2:B:9:VAL:HA	2:B:24:ASN:O	2.16	0.45
3:H:140:THR:HB	3:H:185:THR:OG1	2.17	0.45
6:A:303:GOL:H12	5:P:3:ASN:H	1.83	0.44
7:A:302:EDO:H22	3:H:57:TYR:CZ	2.52	0.44
3:H:62:GLN:HA	3:H:65:LYS:HD3	1.98	0.44
4:L:117:ILE:HD13	4:L:209:PHE:HD2	1.82	0.44
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.98	0.44
1:A:255:GLN:H	1:A:255:GLN:CD	2.25	0.44
3:H:98:ARG:O	3:H:99:SER:HB2	2.16	0.44
2:B:12:ARG:CZ	2:B:22:ILE:HD13	2.47	0.44
1:A:203:CYS:HB2	1:A:217:TRP:CZ2	2.52	0.44
1:A:264:GLU:HG3	7:A:302:EDO:O2	2.17	0.43
3:H:33:TRP:HB2	3:H:99:SER:O	2.17	0.43
1:A:97:GLN:HG2	1:A:116:PHE:CE2	2.53	0.43
2:B:11:SER:OG	2:B:13:HIS:O	2.31	0.43
1:A:172:LEU:O	1:A:174:ASN:N	2.45	0.43
2:B:5:PRO:HB3	2:B:30:PHE:HB3	2.01	0.43
4:L:189:HIS:HB2	4:L:192:TYR:OH	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:GLU:HG2	1:A:233:THR:O	2.18	0.43
3:H:129:PRO:HD3	3:H:141:LEU:HD23	2.00	0.43
3:H:187:THR:O	3:H:190:THR:OG1	2.36	0.43
3:H:39:GLN:HG3	3:H:44:GLY:O	2.18	0.42
3:H:127:LEU:HD12	3:H:142:GLY:HA3	2.01	0.42
1:A:97:GLN:NE2	1:A:114:LEU:HD12	2.34	0.42
4:L:163:TRP:CZ2	4:L:175:MET:HE3	2.54	0.42
1:A:48:ARG:NH2	2:B:53:ASP:OD2	2.52	0.42
1:A:160:LEU:HD12	1:A:160:LEU:HA	1.92	0.42
3:H:50:TYR:CE1	3:H:59:GLU:HB2	2.55	0.41
4:L:7:THR:HA	4:L:8:PRO:HA	1.84	0.41
4:L:154:GLU:HG3	4:L:155:ARG:N	2.34	0.41
3:H:99:SER:HB3	3:H:100:PHE:H	1.68	0.41
4:L:101:GLY:HA3	7:L:304:EDO:O1	2.20	0.41
1:A:202:ARG:HG3	1:A:246:SER:HB3	2.03	0.41
2:B:46:ILE:HA	2:B:47:PRO:HD3	1.92	0.41
1:A:20:PRO:HD3	1:A:75:ARG:HG3	2.02	0.41
3:H:141:LEU:HD22	3:H:213:ILE:CG2	2.49	0.41
1:A:45:TYR:CE2	1:A:67:ALA:HB2	2.56	0.41
2:B:41:LYS:HD2	2:B:78:TYR:CE1	2.57	0.41
3:H:191:TRP:CD1	3:H:196:ILE:HD12	2.56	0.40
1:A:227:ASP:OD1	1:A:227:ASP:N	2.39	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	A	265/273 (97%)	254 (96%)	8 (3%)	3 (1%)	11	25
2	B	96/99 (97%)	85 (88%)	10 (10%)	1 (1%)	12	28
3	H	210/217 (97%)	200 (95%)	5 (2%)	5 (2%)	4	9

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	L	211/214 (99%)	200 (95%)	9 (4%)	2 (1%)	14	30
5	P	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	789/812 (97%)	745 (94%)	33 (4%)	11 (1%)	9	19

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	H	99	SER
2	B	47	PRO
4	L	68	GLY
4	L	102	THR
1	A	17	LEU
1	A	173	LYS
1	A	174	ASN
3	H	11	LEU
3	H	121	ALA
3	H	130	VAL
3	H	216	VAL

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	221/232 (95%)	214 (97%)	7 (3%)	34	62
2	B	83/93 (89%)	81 (98%)	2 (2%)	43	70
3	H	181/185 (98%)	169 (93%)	12 (7%)	15	34
4	L	191/192 (100%)	185 (97%)	6 (3%)	35	63
5	P	8/8 (100%)	8 (100%)	0	100	100
All	All	684/710 (96%)	657 (96%)	27 (4%)	28	55

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	88	SER
1	A	110	LEU
1	A	160	LEU
1	A	183	ASP
1	A	227	ASP
1	A	272	LEU
2	B	77	THR
2	B	92	THR
3	H	62	GLN
3	H	69	THR
3	H	112	LEU
3	H	130	VAL
3	H	140	THR
3	H	162	LEU
3	H	173	LEU
3	H	175	SER
3	H	176	ASP
3	H	180	LEU
3	H	197	THR
3	H	206	SER
4	L	1	LEU
4	L	22	SER
4	L	77	ASN
4	L	165	ASP
4	L	181	LEU
4	L	202	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	188	HIS
1	A	218	GLN
1	A	220	ASN
2	B	6	GLN
3	H	62	GLN
3	H	82	GLN
4	L	27	GLN
4	L	31	ASN
4	L	92	ASN
4	L	156	GLN
5	P	3	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 ⓘ

12 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	EDO	A	302	-	3,3,3	0.70	0	2,2,2	0.50	0
8	NAG	H	301	-	14,14,15	0.19	0	17,19,21	0.73	1 (5%)
6	GOL	A	301	-	5,5,5	1.35	1 (20%)	5,5,5	1.03	0
7	EDO	B	102	-	3,3,3	0.34	0	2,2,2	0.64	0
7	EDO	H	302	-	3,3,3	0.61	0	2,2,2	0.14	0
7	EDO	P	101	-	3,3,3	0.58	0	2,2,2	0.34	0
7	EDO	L	301	-	3,3,3	0.61	0	2,2,2	0.13	0
7	EDO	L	302	-	3,3,3	0.57	0	2,2,2	0.21	0
7	EDO	B	101	-	3,3,3	0.62	0	2,2,2	0.18	0
7	EDO	L	304	-	3,3,3	0.48	0	2,2,2	0.60	0
7	EDO	L	303	-	3,3,3	0.45	0	2,2,2	0.45	0
6	GOL	A	303	-	5,5,5	0.82	0	5,5,5	1.11	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	EDO	A	302	-	-	0/1/1/1	-
8	NAG	H	301	-	-	2/6/23/26	0/1/1/1
6	GOL	A	301	-	-	0/4/4/4	-
7	EDO	B	102	-	-	1/1/1/1	-
7	EDO	H	302	-	-	0/1/1/1	-
7	EDO	P	101	-	-	0/1/1/1	-
7	EDO	L	301	-	-	1/1/1/1	-
7	EDO	L	302	-	-	1/1/1/1	-
7	EDO	B	101	-	-	1/1/1/1	-
7	EDO	L	304	-	-	0/1/1/1	-
7	EDO	L	303	-	-	1/1/1/1	-
6	GOL	A	303	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	301	GOL	C3-C2	2.58	1.61	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	H	301	NAG	C1-O5-C5	2.34	115.33	112.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	H	301	NAG	C4-C5-C6-O6
8	H	301	NAG	O5-C5-C6-O6
6	A	303	GOL	O1-C1-C2-C3
7	L	302	EDO	O1-C1-C2-O2
7	L	301	EDO	O1-C1-C2-O2
7	L	303	EDO	O1-C1-C2-O2
6	A	303	GOL	O1-C1-C2-O2
7	B	101	EDO	O1-C1-C2-O2
7	B	102	EDO	O1-C1-C2-O2

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	302	EDO	3	0
8	H	301	NAG	1	0
6	A	301	GOL	2	0
7	L	302	EDO	1	0
7	L	304	EDO	1	0
6	A	303	GOL	2	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 [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	269/273 (98%)	-0.28	5 (1%) 66 61	20, 36, 61, 83	0
2	B	98/99 (98%)	0.22	1 (1%) 79 76	28, 59, 93, 101	0
3	H	214/217 (98%)	-0.31	2 (0%) 81 78	18, 37, 67, 87	0
4	L	213/214 (99%)	-0.28	0 100 100	25, 42, 59, 83	0
5	P	9/9 (100%)	-0.18	0 100 100	35, 39, 47, 47	0
All	All	803/812 (98%)	-0.23	8 (0%) 79 76	18, 40, 73, 101	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	H	216	VAL	3.5
1	A	180	LEU	3.3
1	A	17	LEU	3.2
1	A	181	ARG	3.1
1	A	16	GLY	2.8
1	A	174	ASN	2.4
2	B	75	THR	2.3
3	H	131	CYS	2.1

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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	NAG	H	301	14/15	0.71	0.19	70,78,88,90	0
7	EDO	L	301	4/4	0.73	0.20	51,51,52,53	0
7	EDO	L	302	4/4	0.75	0.17	51,54,60,63	0
7	EDO	B	102	4/4	0.78	0.20	51,53,54,54	0
7	EDO	L	304	4/4	0.79	0.17	41,41,44,44	0
7	EDO	H	302	4/4	0.80	0.16	45,46,50,50	0
6	GOL	A	303	6/6	0.83	0.14	37,42,43,46	0
6	GOL	A	301	6/6	0.89	0.12	36,38,38,41	0
7	EDO	A	302	4/4	0.89	0.14	37,38,39,40	0
7	EDO	P	101	4/4	0.90	0.12	36,39,40,43	0
7	EDO	L	303	4/4	0.92	0.13	44,46,46,48	0
7	EDO	B	101	4/4	0.94	0.10	41,41,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.