



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:08 AM UTC

PDB ID : 6HZ3 / pdb_00006hz3
Title : THE GLIC PENTAMERIC LIGAND-GATED ION CHANNEL MUTANT E243G
Authors : Hu, H.D.; Delarue, M.
Deposited on : 2018-10-22
Resolution : 3.15 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
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

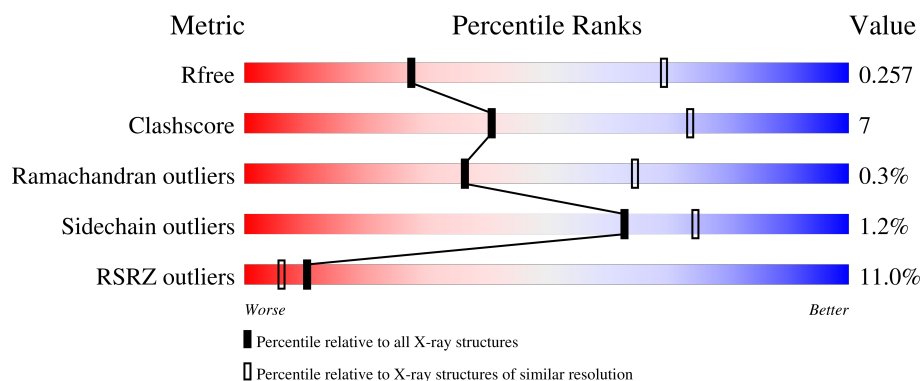
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.15 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	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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	317	9% 83% 14% ..
1	B	317	10% 81% 15% ..
1	C	317	9% 82% 15% ..
1	D	317	11% 80% 17% ..
1	E	317	14% 83% 15% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12637 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 Proton-gated ion channel.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	0	0	0
			2520	1661	404	451	4			
1	B	311	Total	C	N	O	S	0	0	0
			2520	1661	404	451	4			
1	C	311	Total	C	N	O	S	0	0	0
			2520	1661	404	451	4			
1	D	311	Total	C	N	O	S	0	0	0
			2520	1661	404	451	4			
1	E	311	Total	C	N	O	S	0	0	0
			2520	1661	404	451	4			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	243	GLY	GLU	engineered mutation	UNP Q7NDN8
B	243	GLY	GLU	engineered mutation	UNP Q7NDN8
C	243	GLY	GLU	engineered mutation	UNP Q7NDN8
D	243	GLY	GLU	engineered mutation	UNP Q7NDN8
E	243	GLY	GLU	engineered mutation	UNP Q7NDN8

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		
2	C	1	Total	Cl	0	0
			1	1		
2	D	1	Total	Cl	0	0
			1	1		
2	E	1	Total	Cl	0	0
			1	1		

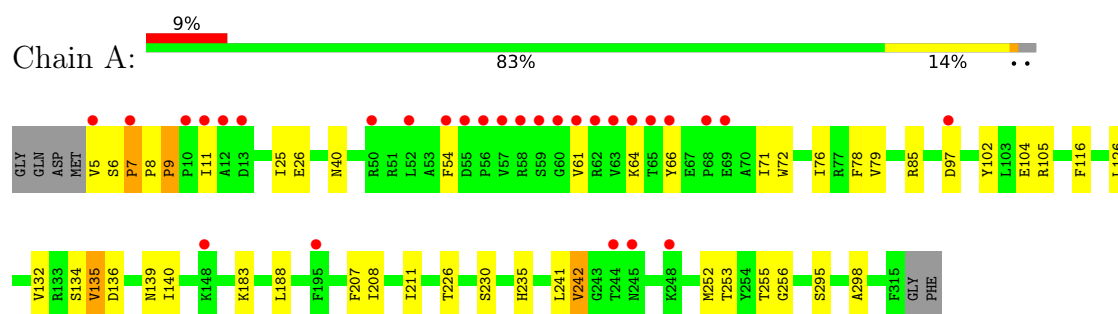
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total 7	O 7	0	0
3	B	8	Total 8	O 8	0	0
3	C	9	Total 9	O 9	0	0
3	D	3	Total 3	O 3	0	0
3	E	5	Total 5	O 5	0	0

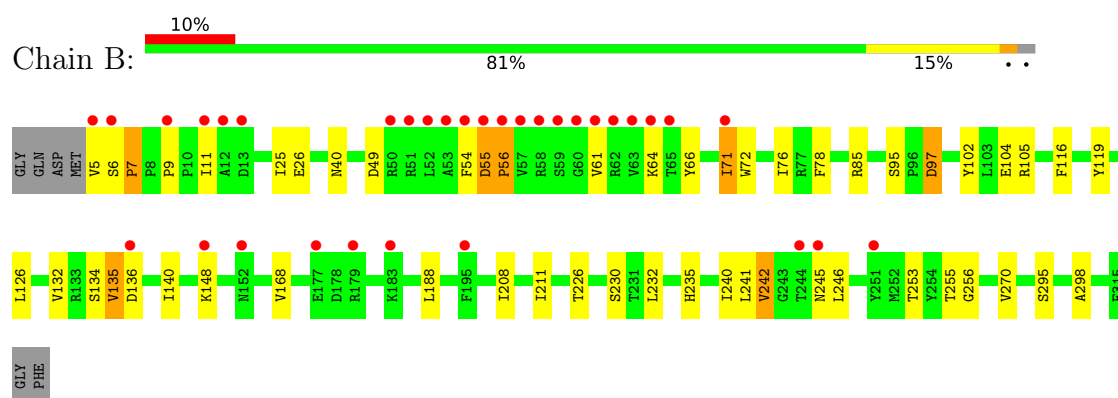
3 Residue-property plots

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.

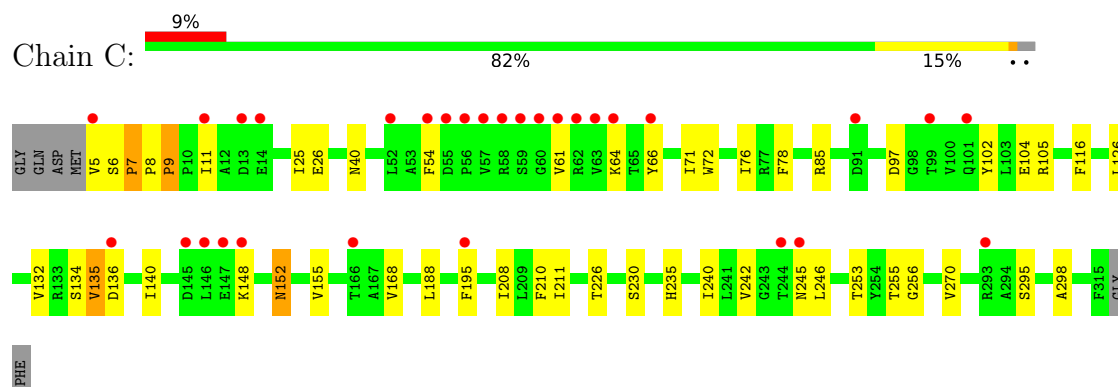
• Molecule 1: Proton-gated ion channel



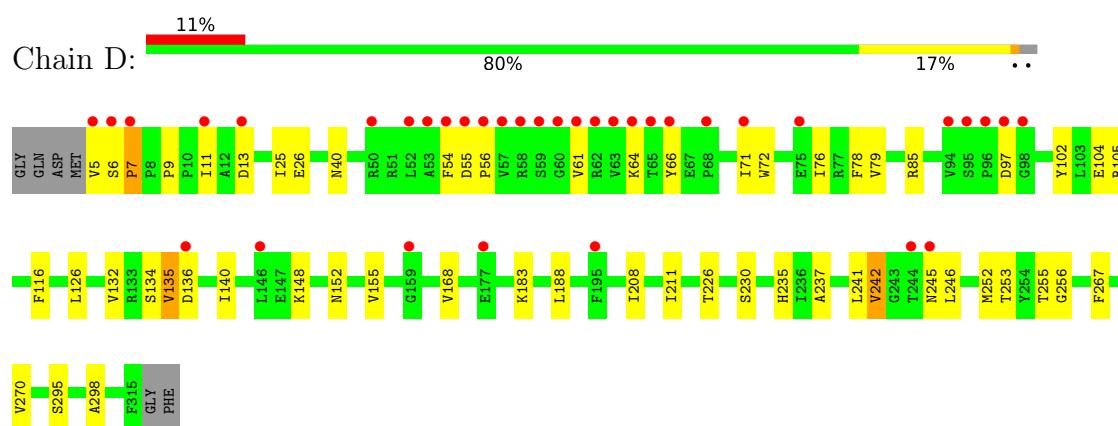
• Molecule 1: Proton-gated ion channel



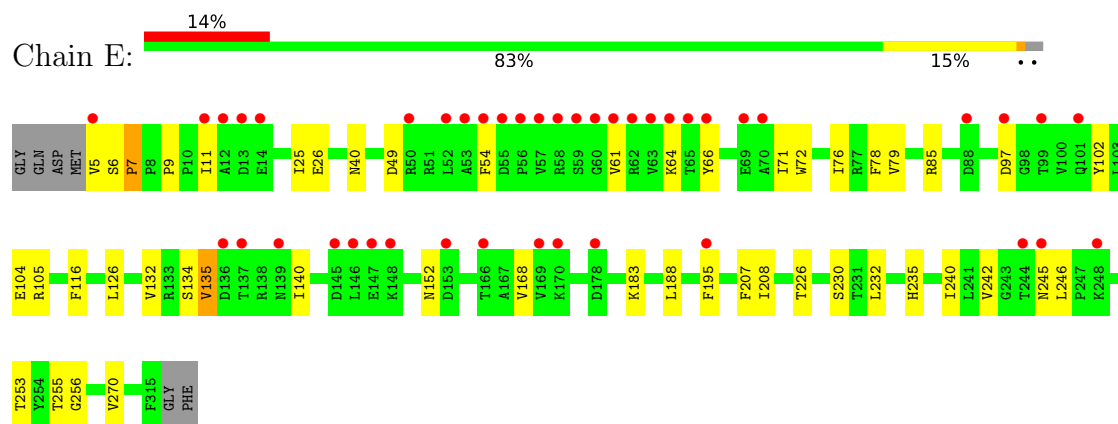
• Molecule 1: Proton-gated ion channel



• Molecule 1: Proton-gated ion channel



• Molecule 1: Proton-gated ion channel



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.45Å 129.07Å 159.61Å 90.00° 101.03° 90.00°	Depositor
Resolution (Å)	20.00 – 3.15 20.00 – 3.15	Depositor EDS
% Data completeness (in resolution range)	73.3 (20.00-3.15) 75.2 (20.00-3.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 3.09Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.230 , 0.249 0.241 , 0.257	Depositor DCC
R_{free} test set	2300 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.2	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	0.87	EDS
Total number of atoms	12637	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% 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:
CL

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	0.81	0/2588	1.29	6/3538 (0.2%)
1	B	0.86	1/2588 (0.0%)	1.32	11/3538 (0.3%)
1	C	0.80	0/2588	1.30	9/3538 (0.3%)
1	D	0.80	0/2588	1.28	9/3538 (0.3%)
1	E	0.81	0/2588	1.30	10/3538 (0.3%)
All	All	0.82	1/12940 (0.0%)	1.30	45/17690 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	97	ASP	CA-C	-8.19	1.41	1.52

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	55	ASP	CA-CB-CG	10.60	123.20	112.60
1	E	97	ASP	CA-CB-CG	9.05	121.65	112.60
1	C	97	ASP	CA-CB-CG	8.77	121.37	112.60
1	B	97	ASP	CA-CB-CG	8.58	121.18	112.60
1	A	97	ASP	CA-CB-CG	7.96	120.56	112.60
1	D	97	ASP	CA-CB-CG	7.75	120.35	112.60
1	A	9	PRO	O-C-N	-6.55	118.07	121.15
1	E	9	PRO	O-C-N	-6.28	118.20	121.15
1	C	9	PRO	O-C-N	-6.17	118.25	121.15
1	E	7	PRO	N-CA-C	6.16	118.22	110.70
1	B	9	PRO	O-C-N	-6.16	118.25	121.15
1	C	7	PRO	N-CA-C	6.06	118.09	110.70
1	A	7	PRO	N-CA-C	6.05	118.09	110.70
1	B	55	ASP	CB-CA-C	-5.93	100.25	109.08
1	D	7	PRO	N-CA-C	5.89	117.88	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	9	PRO	O-C-N	-5.88	118.39	121.15
1	B	7	PRO	N-CA-C	5.77	117.74	110.70
1	B	56	PRO	CB-CG-CD	-5.66	87.98	106.10
1	C	245	ASN	CA-CB-CG	5.55	118.15	112.60
1	E	152	ASN	CA-CB-CG	5.48	118.08	112.60
1	B	245	ASN	CA-CB-CG	5.45	118.05	112.60
1	D	245	ASN	CA-CB-CG	5.44	118.04	112.60
1	D	168	VAL	CA-C-N	5.29	127.99	120.53
1	D	168	VAL	C-N-CA	5.29	127.99	120.53
1	E	245	ASN	CA-CB-CG	5.28	117.88	112.60
1	E	135	VAL	CA-C-N	5.26	129.55	121.19
1	E	135	VAL	C-N-CA	5.26	129.55	121.19
1	C	135	VAL	CA-C-N	5.25	129.54	121.19
1	C	135	VAL	C-N-CA	5.25	129.54	121.19
1	A	135	VAL	CA-C-N	5.21	129.47	121.19
1	A	135	VAL	C-N-CA	5.21	129.47	121.19
1	C	168	VAL	CA-C-N	5.18	127.84	120.53
1	C	168	VAL	C-N-CA	5.18	127.84	120.53
1	A	207	PHE	CA-CB-CG	5.15	118.95	113.80
1	E	168	VAL	CA-C-N	5.12	127.75	120.53
1	E	168	VAL	C-N-CA	5.12	127.75	120.53
1	D	13	ASP	CA-CB-CG	5.09	117.69	112.60
1	B	135	VAL	CA-C-N	5.08	129.27	121.19
1	B	135	VAL	C-N-CA	5.08	129.27	121.19
1	B	168	VAL	CA-C-N	5.07	127.68	120.53
1	B	168	VAL	C-N-CA	5.07	127.68	120.53
1	C	152	ASN	CA-CB-CG	5.07	117.67	112.60
1	D	135	VAL	CA-C-N	5.06	129.23	121.19
1	D	135	VAL	C-N-CA	5.06	129.23	121.19
1	E	207	PHE	CA-CB-CG	5.01	118.81	113.80

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	2520	0	2542	34	0
1	B	2520	0	2542	38	6
1	C	2520	0	2542	35	0
1	D	2520	0	2542	40	0
1	E	2520	0	2542	31	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
3	A	7	0	0	0	0
3	B	8	0	0	2	0
3	C	9	0	0	2	0
3	D	3	0	0	2	0
3	E	5	0	0	0	0
All	All	12637	0	12710	165	6

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 (165) 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:C:148:LYS:HE3	3:C:501:HOH:O	1.69	0.91
1:B:148:LYS:CE	3:B:503:HOH:O	2.33	0.77
1:B:148:LYS:HE3	3:B:503:HOH:O	1.84	0.75
1:A:61:VAL:HG21	1:A:64:LYS:HG3	1.68	0.74
1:C:61:VAL:HG21	1:C:64:LYS:HG3	1.71	0.72
1:E:61:VAL:HG21	1:E:64:LYS:HG3	1.71	0.72
1:D:61:VAL:HG21	1:D:64:LYS:HG3	1.70	0.72
1:E:71:ILE:HD12	1:E:72:TRP:H	1.55	0.70
1:A:6:SER:HB3	1:A:7:PRO:HD2	1.74	0.70
1:C:71:ILE:HD12	1:C:72:TRP:H	1.57	0.70
1:B:6:SER:HB3	1:B:7:PRO:HD2	1.73	0.69
1:A:71:ILE:HD12	1:A:72:TRP:H	1.57	0.69
1:E:6:SER:HB3	1:E:7:PRO:HD2	1.74	0.69
1:D:6:SER:HB3	1:D:7:PRO:HD2	1.73	0.69
1:B:61:VAL:HG21	1:B:64:LYS:HG3	1.73	0.68
1:B:71:ILE:HD12	1:B:72:TRP:H	1.56	0.68
1:D:148:LYS:HE3	3:D:502:HOH:O	1.93	0.68
1:C:6:SER:HB3	1:C:7:PRO:HD2	1.75	0.67
1:D:71:ILE:HD12	1:D:72:TRP:H	1.58	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:148:LYS:CE	3:D:502:HOH:O	2.43	0.66
1:E:54:PHE:HD2	1:E:64:LYS:HZ1	1.44	0.64
1:A:25:ILE:HD13	1:A:105:ARG:HE	1.62	0.64
1:D:25:ILE:HD13	1:D:105:ARG:HE	1.63	0.62
1:B:25:ILE:HD13	1:B:105:ARG:HE	1.64	0.62
1:D:54:PHE:HD2	1:D:64:LYS:HZ1	1.47	0.61
1:E:25:ILE:HD13	1:E:105:ARG:HE	1.64	0.61
1:C:25:ILE:HD13	1:C:105:ARG:HE	1.66	0.60
1:E:116:PHE:O	1:E:253:THR:HG23	2.02	0.59
1:C:116:PHE:O	1:C:253:THR:HG23	2.02	0.58
1:B:116:PHE:O	1:B:253:THR:HG23	2.04	0.57
1:A:116:PHE:O	1:A:253:THR:HG23	2.04	0.57
1:C:54:PHE:HD2	1:C:64:LYS:HZ1	1.53	0.57
1:D:116:PHE:O	1:D:253:THR:HG23	2.04	0.56
1:B:54:PHE:HD2	1:B:64:LYS:HZ1	1.53	0.56
1:E:54:PHE:HD2	1:E:64:LYS:NZ	2.05	0.55
1:A:54:PHE:HD2	1:A:64:LYS:HZ1	1.54	0.55
1:A:25:ILE:HG21	1:A:105:ARG:HH21	1.72	0.55
1:A:54:PHE:HD2	1:A:64:LYS:NZ	2.05	0.55
1:C:54:PHE:HD2	1:C:64:LYS:NZ	2.05	0.54
1:D:54:PHE:HD2	1:D:64:LYS:NZ	2.05	0.54
1:E:25:ILE:HG21	1:E:105:ARG:HH21	1.72	0.54
1:D:134:SER:HA	1:D:140:ILE:HD12	1.90	0.53
1:C:134:SER:HA	1:C:140:ILE:HD12	1.91	0.53
1:B:25:ILE:HG21	1:B:105:ARG:HH21	1.75	0.52
1:B:78:PHE:CE2	1:B:85:ARG:HD3	2.44	0.52
1:D:25:ILE:HG21	1:D:105:ARG:HH21	1.74	0.52
1:B:54:PHE:HD2	1:B:64:LYS:NZ	2.07	0.52
1:C:25:ILE:HG21	1:C:105:ARG:HH21	1.74	0.52
1:A:78:PHE:CE2	1:A:85:ARG:HD3	2.45	0.51
1:A:134:SER:HA	1:A:140:ILE:HD12	1.91	0.51
1:E:134:SER:HA	1:E:140:ILE:HD12	1.92	0.51
1:B:134:SER:HA	1:B:140:ILE:HD12	1.91	0.50
1:C:78:PHE:CE2	1:C:85:ARG:HD3	2.46	0.50
1:D:126:LEU:HD12	1:D:188:LEU:HD23	1.93	0.50
1:E:6:SER:HB3	1:E:7:PRO:CD	2.42	0.50
1:A:126:LEU:HD12	1:A:188:LEU:HD23	1.93	0.50
1:E:78:PHE:CE2	1:E:85:ARG:HD3	2.47	0.49
1:D:78:PHE:CE2	1:D:85:ARG:HD3	2.47	0.49
1:C:211:ILE:HG23	1:D:270:VAL:HG11	1.94	0.49
1:A:226:THR:O	1:A:230:SER:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:PRO:HG3	1:B:135:VAL:CG1	2.42	0.48
1:C:226:THR:O	1:C:230:SER:HB2	2.14	0.48
1:A:6:SER:HB3	1:A:7:PRO:CD	2.42	0.48
1:A:64:LYS:HD2	1:A:66:TYR:OH	2.14	0.48
1:B:7:PRO:HG3	1:B:135:VAL:HG11	1.95	0.48
1:D:64:LYS:HD2	1:D:66:TYR:OH	2.14	0.48
1:C:126:LEU:HD12	1:C:188:LEU:HD23	1.95	0.48
1:B:126:LEU:HD12	1:B:188:LEU:HD23	1.95	0.48
1:E:126:LEU:HD12	1:E:188:LEU:HD23	1.94	0.47
1:C:195:PHE:CZ	1:D:252:MET:HE2	2.49	0.47
1:D:226:THR:O	1:D:230:SER:HB2	2.14	0.47
1:C:6:SER:HB3	1:C:7:PRO:CD	2.43	0.47
1:D:7:PRO:HG3	1:D:135:VAL:CG1	2.45	0.47
1:C:7:PRO:HG3	1:C:135:VAL:CG1	2.45	0.47
1:E:226:THR:O	1:E:230:SER:HB2	2.15	0.47
1:A:252:MET:HE2	1:E:195:PHE:CZ	2.49	0.47
1:C:64:LYS:HD2	1:C:66:TYR:OH	2.15	0.47
1:C:7:PRO:HG3	1:C:135:VAL:HG11	1.97	0.47
1:D:211:ILE:HG23	1:E:270:VAL:HG11	1.97	0.47
1:B:226:THR:O	1:B:230:SER:HB2	2.15	0.46
1:E:7:PRO:HG3	1:E:135:VAL:CG1	2.45	0.46
1:E:208:ILE:HD13	1:E:235:HIS:HB2	1.98	0.46
1:B:6:SER:HB3	1:B:7:PRO:CD	2.42	0.46
1:D:54:PHE:CD2	1:D:64:LYS:NZ	2.82	0.46
1:E:7:PRO:HG3	1:E:135:VAL:HG11	1.98	0.46
1:E:54:PHE:CD2	1:E:64:LYS:NZ	2.82	0.46
1:A:7:PRO:HG3	1:A:135:VAL:CG1	2.45	0.46
1:A:208:ILE:HD13	1:A:235:HIS:HB2	1.98	0.46
1:D:26:GLU:HB2	1:D:40:ASN:HB3	1.98	0.46
1:E:64:LYS:HD2	1:E:66:TYR:OH	2.16	0.46
1:D:208:ILE:HD13	1:D:235:HIS:HB2	1.98	0.45
1:C:54:PHE:CD2	1:C:64:LYS:NZ	2.82	0.45
1:B:54:PHE:CD2	1:B:64:LYS:NZ	2.84	0.45
1:B:211:ILE:HG23	1:C:270:VAL:HG11	1.99	0.45
1:C:253:THR:HB	1:C:256:GLY:H	1.82	0.45
1:D:208:ILE:HG22	1:E:232:LEU:HD21	1.99	0.45
1:A:7:PRO:HG3	1:A:135:VAL:HG11	1.99	0.45
1:D:6:SER:HB3	1:D:7:PRO:CD	2.42	0.45
1:C:208:ILE:HD13	1:C:235:HIS:HB2	1.99	0.45
1:A:54:PHE:CD2	1:A:64:LYS:NZ	2.82	0.44
1:B:64:LYS:HD2	1:B:66:TYR:OH	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:GLU:HB2	1:C:40:ASN:HB3	1.99	0.44
1:E:26:GLU:HB2	1:E:40:ASN:HB3	1.99	0.44
1:E:253:THR:HB	1:E:256:GLY:H	1.82	0.44
1:D:7:PRO:HG3	1:D:135:VAL:HG11	1.98	0.44
1:B:241:LEU:HB2	1:C:240:ILE:HG13	2.00	0.44
1:B:208:ILE:HD13	1:B:235:HIS:HB2	1.99	0.44
1:D:241:LEU:HB2	1:E:240:ILE:HG13	1.98	0.44
1:A:26:GLU:HB2	1:A:40:ASN:HB3	2.00	0.44
1:A:76:ILE:HD13	1:A:132:VAL:HB	2.00	0.44
1:B:26:GLU:HB2	1:B:40:ASN:HB3	1.99	0.43
1:B:102:TYR:CE1	1:B:104:GLU:HG2	2.54	0.43
1:A:135:VAL:HG22	1:A:136:ASP:H	1.83	0.43
1:D:242:VAL:O	1:D:242:VAL:CG1	2.67	0.43
1:A:8:PRO:HA	1:A:9:PRO:HD3	1.90	0.43
1:B:49:ASP:OD2	1:B:71:ILE:HD11	2.19	0.43
1:B:253:THR:HB	1:B:256:GLY:H	1.82	0.43
1:D:76:ILE:HD13	1:D:132:VAL:HB	2.00	0.43
1:D:135:VAL:HG22	1:D:136:ASP:H	1.83	0.42
1:D:253:THR:HB	1:D:256:GLY:H	1.84	0.42
1:C:8:PRO:HA	1:C:9:PRO:HD3	1.89	0.42
1:C:148:LYS:CE	3:C:501:HOH:O	2.44	0.42
1:D:152:ASN:HB3	1:D:155:VAL:HG23	2.02	0.42
1:E:102:TYR:CE1	1:E:104:GLU:HG2	2.55	0.42
1:A:242:VAL:O	1:A:242:VAL:CG1	2.68	0.42
1:B:242:VAL:O	1:B:242:VAL:HG12	2.20	0.42
1:B:242:VAL:O	1:B:242:VAL:CG1	2.67	0.41
1:C:210:PHE:HB2	1:D:267:PHE:CE1	2.55	0.41
1:D:55:ASP:HA	1:D:56:PRO:HD3	1.94	0.41
1:A:253:THR:HB	1:A:256:GLY:H	1.84	0.41
1:C:102:TYR:CE1	1:C:104:GLU:HG2	2.55	0.41
1:E:76:ILE:HD13	1:E:132:VAL:HB	2.02	0.41
1:B:119:TYR:CD1	1:B:246:LEU:HD11	2.55	0.41
1:E:116:PHE:CD1	1:E:246:LEU:HD22	2.56	0.41
1:C:76:ILE:HD13	1:C:132:VAL:HB	2.01	0.41
1:A:211:ILE:HG23	1:B:270:VAL:HG11	2.02	0.41
1:C:116:PHE:CD1	1:C:246:LEU:HD22	2.55	0.41
1:D:116:PHE:CD1	1:D:246:LEU:HD22	2.56	0.41
1:A:208:ILE:HG22	1:B:232:LEU:HD21	2.03	0.41
1:B:78:PHE:HE2	1:B:85:ARG:HD3	1.86	0.41
1:B:135:VAL:HG22	1:B:136:ASP:H	1.86	0.41
1:D:102:TYR:CE1	1:D:104:GLU:HG2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ILE:HD13	1:B:132:VAL:HB	2.02	0.41
1:C:126:LEU:HB2	1:C:188:LEU:HB3	2.03	0.41
1:A:102:TYR:HE1	1:A:104:GLU:HG2	1.86	0.40
1:C:135:VAL:HG22	1:C:136:ASP:H	1.86	0.40
1:D:237:ALA:O	1:E:240:ILE:HD11	2.21	0.40
1:D:242:VAL:O	1:D:242:VAL:HG12	2.21	0.40
1:A:241:LEU:HB2	1:B:240:ILE:HG13	2.03	0.40
1:E:102:TYR:HE1	1:E:104:GLU:HG2	1.86	0.40
1:A:102:TYR:CE1	1:A:104:GLU:HG2	2.55	0.40
1:A:295:SER:HA	1:A:298:ALA:HB3	2.04	0.40
1:D:79:VAL:HG21	1:D:183:LYS:HD2	2.03	0.40
1:D:126:LEU:HB2	1:D:188:LEU:HB3	2.04	0.40
1:D:295:SER:HA	1:D:298:ALA:HB3	2.04	0.40
1:E:79:VAL:HG21	1:E:183:LYS:HD2	2.03	0.40
1:A:79:VAL:HG21	1:A:183:LYS:HD2	2.03	0.40
1:B:295:SER:HA	1:B:298:ALA:HB3	2.03	0.40
1:C:152:ASN:HB3	1:C:155:VAL:HG23	2.02	0.40
1:A:126:LEU:HB2	1:A:188:LEU:HB3	2.04	0.40
1:A:134:SER:HB3	1:A:139:ASN:HA	2.03	0.40
1:B:102:TYR:HE1	1:B:104:GLU:HG2	1.86	0.40
1:B:116:PHE:CD1	1:B:246:LEU:HD22	2.56	0.40
1:C:295:SER:HA	1:C:298:ALA:HB3	2.03	0.40
1:E:49:ASP:OD2	1:E:71:ILE:HD11	2.22	0.40

All (6) 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:B:55:ASP:OD2	1:B:97:ASP:CG[2_555]	1.47	0.73
1:B:55:ASP:OD2	1:B:97:ASP:OD1[2_555]	1.47	0.73
1:B:55:ASP:OD2	1:B:97:ASP:CB[2_555]	1.60	0.60
1:B:55:ASP:OD1	1:B:95:SER:OG[2_555]	1.99	0.21
1:B:55:ASP:CG	1:B:97:ASP:OD1[2_555]	2.05	0.15
1:B:55:ASP:OD1	1:B:97:ASP:OD1[2_555]	2.15	0.05

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	309/317 (98%)	291 (94%)	17 (6%)	1 (0%)	36	64
1	B	309/317 (98%)	290 (94%)	18 (6%)	1 (0%)	36	64
1	C	309/317 (98%)	289 (94%)	19 (6%)	1 (0%)	36	64
1	D	309/317 (98%)	291 (94%)	17 (6%)	1 (0%)	36	64
1	E	309/317 (98%)	290 (94%)	18 (6%)	1 (0%)	36	64
All	All	1545/1585 (98%)	1451 (94%)	89 (6%)	5 (0%)	36	64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	VAL
1	B	242	VAL
1	D	242	VAL
1	E	242	VAL
1	C	242	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	279/283 (99%)	276 (99%)	3 (1%)	65	75
1	B	279/283 (99%)	274 (98%)	5 (2%)	51	70
1	C	279/283 (99%)	276 (99%)	3 (1%)	65	75
1	D	279/283 (99%)	276 (99%)	3 (1%)	65	75

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	279/283 (99%)	276 (99%)	3 (1%)	65	75
All	All	1395/1415 (99%)	1378 (99%)	17 (1%)	63	74

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

Mol	Chain	Res	Type
1	A	5	VAL
1	A	11	ILE
1	A	255	THR
1	B	5	VAL
1	B	11	ILE
1	B	56	PRO
1	B	71	ILE
1	B	255	THR
1	C	5	VAL
1	C	11	ILE
1	C	255	THR
1	D	5	VAL
1	D	11	ILE
1	D	255	THR
1	E	5	VAL
1	E	11	ILE
1	E	255	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	152	ASN
1	A	239	ASN
1	A	277	HIS
1	A	284	GLN
1	A	307	ASN
1	B	152	ASN
1	B	239	ASN
1	B	277	HIS
1	B	284	GLN
1	B	307	ASN
1	C	83	ASN
1	C	152	ASN
1	C	239	ASN
1	C	277	HIS

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Mol	Chain	Res	Type
1	C	284	GLN
1	C	307	ASN
1	D	152	ASN
1	D	239	ASN
1	D	277	HIS
1	D	284	GLN
1	D	307	ASN
1	E	152	ASN
1	E	239	ASN
1	E	277	HIS
1	E	307	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 ⓘ

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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 ⓘ

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	311/317 (98%)	0.38	29 (9%) 14 8	8, 35, 98, 171	0
1	B	311/317 (98%)	0.59	33 (10%) 11 6	9, 33, 73, 146	0
1	C	311/317 (98%)	0.42	30 (9%) 13 8	8, 34, 86, 145	0
1	D	311/317 (98%)	0.60	36 (11%) 9 6	6, 37, 96, 169	0
1	E	311/317 (98%)	0.58	43 (13%) 6 4	11, 36, 91, 154	0
All	All	1555/1585 (98%)	0.51	171 (10%) 10 6	6, 35, 93, 171	0

All (171) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	57	VAL	10.8
1	B	56	PRO	10.4
1	D	61	VAL	8.6
1	E	60	GLY	8.3
1	E	61	VAL	7.9
1	A	60	GLY	7.6
1	B	61	VAL	7.5
1	D	57	VAL	7.2
1	A	61	VAL	7.0
1	C	61	VAL	6.6
1	B	11	ILE	6.5
1	B	53	ALA	6.4
1	B	58	ARG	6.4
1	B	60	GLY	6.3
1	B	59	SER	6.3
1	C	60	GLY	6.2
1	E	54	PHE	6.2
1	D	64	LYS	6.2
1	B	54	PHE	6.1
1	D	63	VAL	6.0

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Mol	Chain	Res	Type	RSRZ
1	C	62	ARG	6.0
1	E	56	PRO	5.7
1	E	59	SER	5.5
1	B	62	ARG	5.4
1	D	62	ARG	5.3
1	B	52	LEU	5.2
1	B	55	ASP	5.2
1	A	57	VAL	5.2
1	E	57	VAL	5.1
1	E	55	ASP	5.0
1	D	56	PRO	5.0
1	C	59	SER	5.0
1	A	63	VAL	5.0
1	C	57	VAL	4.9
1	C	147	GLU	4.9
1	C	56	PRO	4.8
1	D	60	GLY	4.8
1	E	5	VAL	4.8
1	C	54	PHE	4.7
1	E	58	ARG	4.7
1	D	54	PHE	4.7
1	E	62	ARG	4.6
1	C	58	ARG	4.6
1	A	5	VAL	4.6
1	C	5	VAL	4.6
1	D	5	VAL	4.5
1	E	136	ASP	4.5
1	D	52	LEU	4.4
1	E	148	LYS	4.4
1	A	64	LYS	4.4
1	A	56	PRO	4.4
1	D	11	ILE	4.3
1	B	50	ARG	4.2
1	A	11	ILE	4.2
1	E	147	GLU	4.1
1	A	65	THR	4.0
1	D	6	SER	4.0
1	B	5	VAL	4.0
1	A	59	SER	4.0
1	D	98	GLY	3.9
1	C	55	ASP	3.9
1	D	13	ASP	3.9

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Mol	Chain	Res	Type	RSRZ
1	E	11	ILE	3.9
1	D	65	THR	3.8
1	D	58	ARG	3.8
1	B	64	LYS	3.7
1	A	58	ARG	3.6
1	E	146	LEU	3.6
1	E	145	ASP	3.6
1	E	166	THR	3.6
1	D	59	SER	3.6
1	C	136	ASP	3.6
1	E	64	LYS	3.5
1	A	54	PHE	3.5
1	C	245	ASN	3.5
1	E	52	LEU	3.4
1	D	96	PRO	3.4
1	B	195	PHE	3.4
1	D	97	ASP	3.4
1	C	148	LYS	3.4
1	D	95	SER	3.3
1	E	13	ASP	3.3
1	D	245	ASN	3.3
1	A	66	TYR	3.2
1	C	52	LEU	3.1
1	E	153	ASP	3.1
1	C	14	GLU	3.1
1	E	63	VAL	3.1
1	E	139	ASN	3.1
1	A	13	ASP	3.1
1	C	64	LYS	3.1
1	A	69	GLU	3.1
1	C	13	ASP	3.1
1	E	97	ASP	3.1
1	D	53	ALA	3.0
1	C	63	VAL	3.0
1	D	66	TYR	3.0
1	B	6	SER	2.9
1	C	244	THR	2.9
1	E	245	ASN	2.9
1	C	146	LEU	2.9
1	A	245	ASN	2.9
1	A	68	PRO	2.8
1	D	146	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	71	ILE	2.8
1	B	179	ARG	2.8
1	D	71	ILE	2.8
1	B	245	ASN	2.7
1	B	148	LYS	2.7
1	E	99	THR	2.7
1	C	66	TYR	2.7
1	E	53	ALA	2.6
1	B	136	ASP	2.6
1	B	251	TYR	2.5
1	E	14	GLU	2.5
1	D	195	PHE	2.5
1	D	7	PRO	2.5
1	E	12	ALA	2.5
1	A	50	ARG	2.5
1	A	52	LEU	2.5
1	B	9	PRO	2.5
1	B	51	ARG	2.4
1	E	137	THR	2.4
1	B	244	THR	2.4
1	E	50	ARG	2.4
1	C	195	PHE	2.4
1	E	169	VAL	2.4
1	B	152	ASN	2.3
1	A	195	PHE	2.3
1	E	69	GLU	2.3
1	B	63	VAL	2.3
1	C	166	THR	2.3
1	A	97	ASP	2.3
1	C	145	ASP	2.3
1	D	75	GLU	2.3
1	D	159	GLY	2.3
1	E	70	ALA	2.3
1	E	101	GLN	2.3
1	B	183	LYS	2.3
1	B	13	ASP	2.3
1	D	94	VAL	2.3
1	A	62	ARG	2.2
1	D	50	ARG	2.2
1	A	10	PRO	2.2
1	E	66	TYR	2.2
1	B	65	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	244	THR	2.2
1	B	12	ALA	2.2
1	D	55	ASP	2.2
1	A	12	ALA	2.2
1	D	177	GLU	2.2
1	C	293	ARG	2.2
1	E	195	PHE	2.2
1	C	11	ILE	2.1
1	C	101	GLN	2.1
1	D	136	ASP	2.1
1	E	65	THR	2.1
1	E	88	ASP	2.1
1	A	148	LYS	2.1
1	E	170	LYS	2.1
1	E	178	ASP	2.1
1	A	7	PRO	2.1
1	D	68	PRO	2.1
1	C	99	THR	2.1
1	E	248	LYS	2.0
1	A	55	ASP	2.0
1	A	244	THR	2.0
1	B	177	GLU	2.0
1	D	244	THR	2.0
1	C	91	ASP	2.0
1	A	248	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

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
2	CL	B	401	1/1	0.91	0.12	37,37,37,37	0
2	CL	E	401	1/1	0.94	0.06	34,34,34,34	0
2	CL	A	401	1/1	0.97	0.06	32,32,32,32	0
2	CL	C	401	1/1	0.98	0.10	25,25,25,25	0
2	CL	D	401	1/1	0.99	0.03	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.