



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:55 AM UTC

PDB ID : 4ZZB / pdb_00004zzb
Title : The GLIC pentameric Ligand-Gated Ion Channel Locally-closed form complexed to xenon
Authors : Sauguet, L.; Fourati, Z.; Prange, T.; Delarue, M.; Colloc'h, N.
Deposited on : 2015-05-22
Resolution : 3.40 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

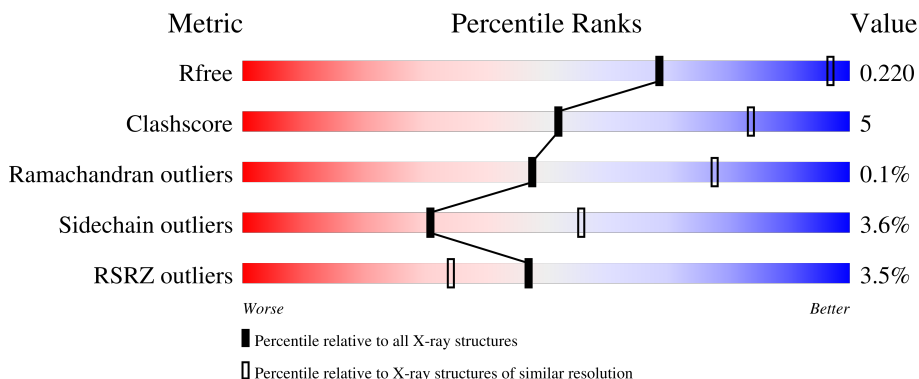
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	3% 87% 11% ..
1	B	317	3% 89% 8% ..
1	C	317	3% 82% 14% ..
1	D	317	3% 86% 11% .
1	E	317	4% 87% 10% ..

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
6	XE	A	405	-	-	X	-
6	XE	A	406	-	-	X	-
6	XE	A	407	-	-	X	-
6	XE	A	409	-	-	X	-
6	XE	B	404	-	-	X	-
6	XE	B	406	-	-	X	-
6	XE	C	407	-	-	X	-
6	XE	D	404	-	-	X	-
6	XE	D	406	-	-	X	-
6	XE	E	405	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12659 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
			2513	1654	401	453	5			
1	B	311	Total	C	N	O	S	0	1	0
			2522	1662	401	454	5			
1	C	311	Total	C	N	O	S	0	1	0
			2522	1662	401	454	5			
1	D	311	Total	C	N	O	S	0	0	0
			2513	1654	401	453	5			
1	E	311	Total	C	N	O	S	0	1	0
			2522	1662	401	454	5			

There are 20 discrepancies between the modelled and reference sequences:

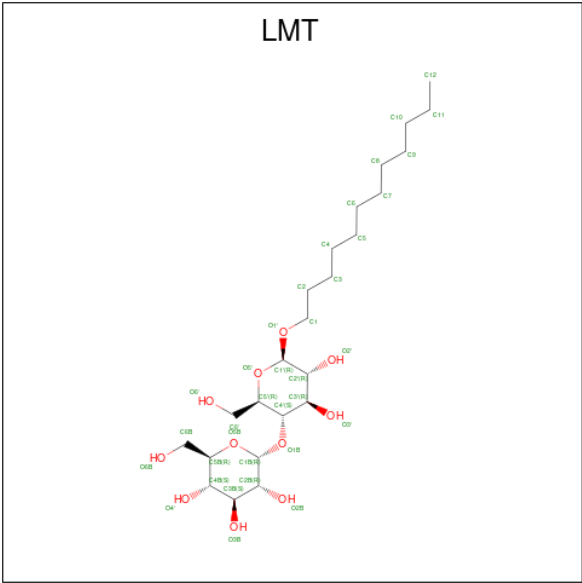
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	ALA	-	expression tag	UNP Q7NDN8
A	27	SER	CYS	engineered mutation	UNP Q7NDN8
A	33	CYS	LYS	engineered mutation	UNP Q7NDN8
A	246	CYS	LEU	engineered mutation	UNP Q7NDN8
B	1	ALA	-	expression tag	UNP Q7NDN8
B	27	SER	CYS	engineered mutation	UNP Q7NDN8
B	33	CYS	LYS	engineered mutation	UNP Q7NDN8
B	246	CYS	LEU	engineered mutation	UNP Q7NDN8
C	1	ALA	-	expression tag	UNP Q7NDN8
C	27	SER	CYS	engineered mutation	UNP Q7NDN8
C	33	CYS	LYS	engineered mutation	UNP Q7NDN8
C	246	CYS	LEU	engineered mutation	UNP Q7NDN8
D	1	ALA	-	expression tag	UNP Q7NDN8
D	27	SER	CYS	engineered mutation	UNP Q7NDN8
D	33	CYS	LYS	engineered mutation	UNP Q7NDN8
D	246	CYS	LEU	engineered mutation	UNP Q7NDN8
E	1	ALA	-	expression tag	UNP Q7NDN8
E	27	SER	CYS	engineered mutation	UNP Q7NDN8
E	33	CYS	LYS	engineered mutation	UNP Q7NDN8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	246	CYS	LEU	engineered mutation	UNP Q7NDN8

- Molecule 2 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: C₂₄H₄₆O₁₁).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C 12 12	0	0

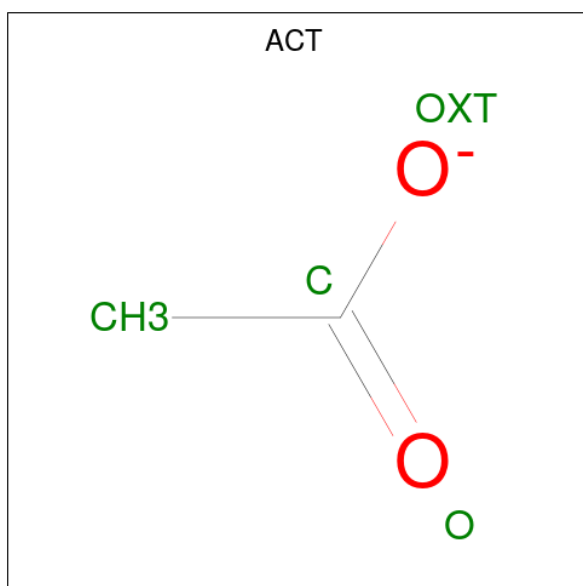
- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Na 1 1	0	0
4	B	1	Total Na 1 1	0	0
4	C	1	Total Na 1 1	0	0
4	D	1	Total Na 1 1	0	0
4	E	1	Total Na 1 1	0	0

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	B	1	Total C O 4 2 2	0	0
5	C	1	Total C O 4 2 2	0	0
5	C	1	Total C O 4 2 2	0	0
5	D	1	Total C O 4 2 2	0	0
5	E	1	Total C O 4 2 2	0	0

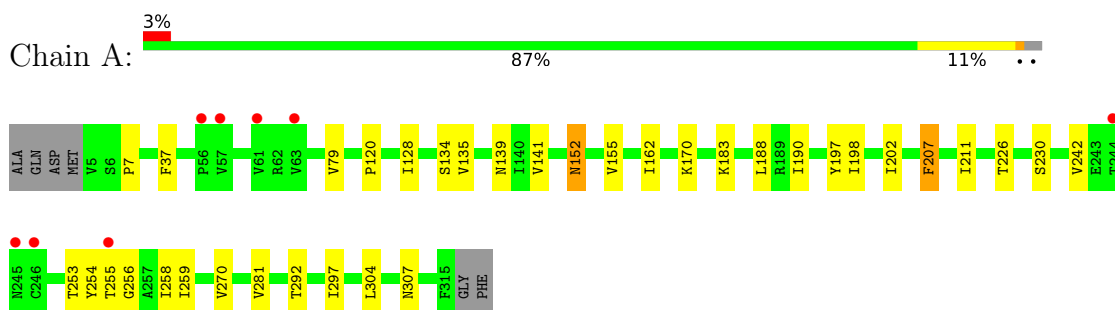
- Molecule 6 is XENON (CCD ID: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	6	Total 6	Xe 6	0	0
6	B	4	Total 4	Xe 4	0	0
6	C	4	Total 4	Xe 4	0	0
6	D	4	Total 4	Xe 4	0	0
6	E	3	Total 3	Xe 3	0	0

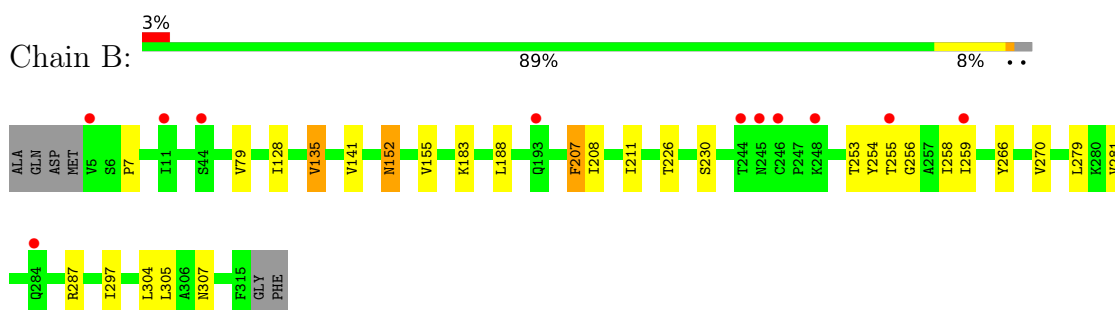
3 Residue-property plots [i](#)

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.

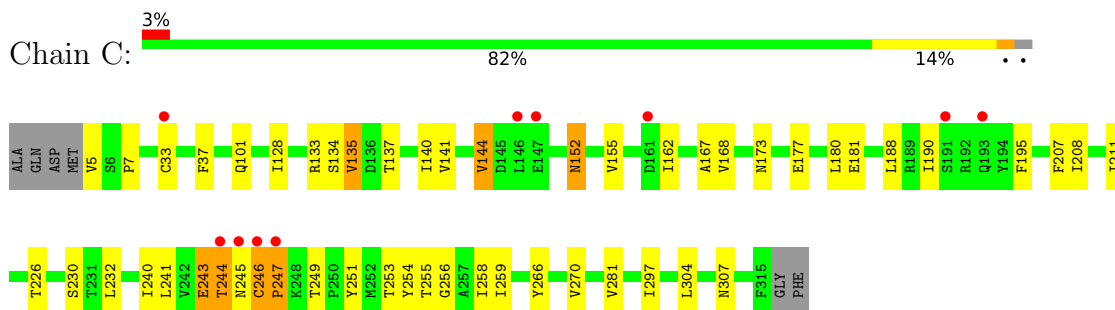
- Molecule 1: Proton-gated ion channel



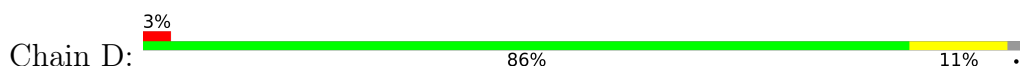
- Molecule 1: Proton-gated ion channel

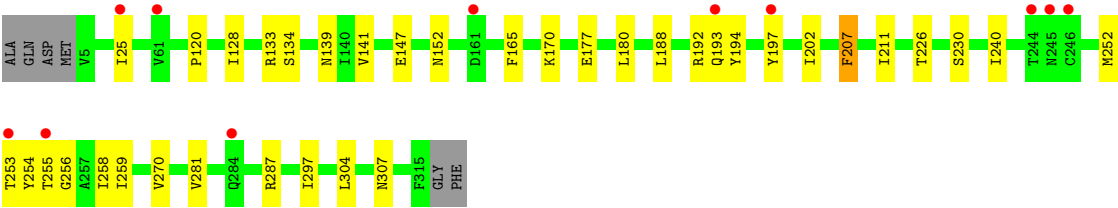


- Molecule 1: Proton-gated ion channel

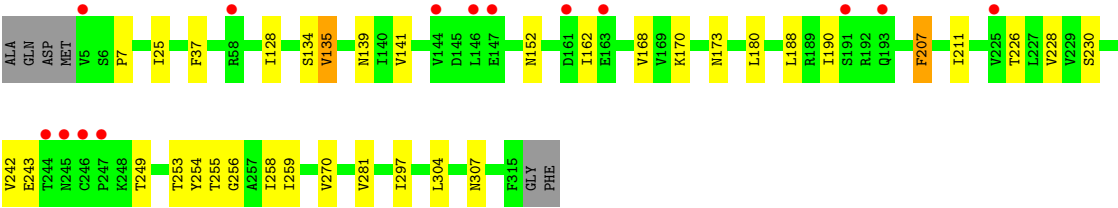
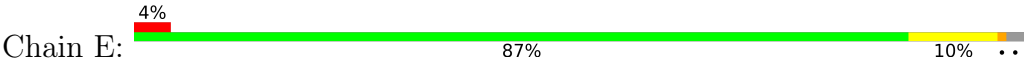


- Molecule 1: Proton-gated ion channel





● 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	176.97Å 130.53Å 157.53Å 90.00° 100.33° 90.00°	Depositor
Resolution (Å)	20.00 – 3.40 20.00 – 3.40	Depositor EDS
% Data completeness (in resolution range)	98.4 (20.00-3.40) 97.8 (20.00-3.40)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 3.40Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.216 , 0.223 (Not available) , 0.220	Depositor DCC
R_{free} test set	2425 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	107.6	Xtriage
Anisotropy	0.370	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 65.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12659	wwPDB-VP
Average B, all atoms (Å ²)	127.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.06% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: XE, NA, ACT, LMT, 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	0.70	0/2581	1.21	2/3531 (0.1%)
1	B	0.71	0/2594	1.21	4/3549 (0.1%)
1	C	0.70	0/2594	1.20	4/3549 (0.1%)
1	D	0.70	0/2581	1.19	3/3531 (0.1%)
1	E	0.70	0/2594	1.21	7/3549 (0.2%)
All	All	0.70	0/12944	1.20	20/17709 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	181	GLU	CB-CG-CD	6.39	123.46	112.60
1	E	152	ASN	CA-CB-CG	6.39	118.99	112.60
1	E	135	VAL	CA-C-N	6.35	128.70	120.44
1	E	135	VAL	C-N-CA	6.35	128.70	120.44
1	D	152	ASN	CA-CB-CG	6.20	118.80	112.60
1	C	152	ASN	CA-CB-CG	6.17	118.77	112.60
1	A	152	ASN	CA-CB-CG	6.03	118.63	112.60
1	B	152	ASN	CA-CB-CG	5.99	118.59	112.60
1	B	135	VAL	CA-C-N	5.74	128.29	120.54
1	B	135	VAL	C-N-CA	5.74	128.29	120.54
1	B	207	PHE	CA-CB-CG	5.62	119.42	113.80
1	E	207	PHE	CA-CB-CG	5.50	119.31	113.80
1	A	207	PHE	CA-CB-CG	5.48	119.28	113.80
1	D	207	PHE	CA-CB-CG	5.33	119.13	113.80
1	E	168	VAL	CA-C-N	5.08	127.70	120.53
1	E	168	VAL	C-N-CA	5.08	127.70	120.53
1	E	25	ILE	N-CA-C	-5.04	106.76	111.45
1	D	25	ILE	N-CA-C	-5.03	106.77	111.45
1	C	168	VAL	CA-C-N	5.03	127.62	120.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	168	VAL	C-N-CA	5.03	127.62	120.53

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	2513	0	2514	29	0
1	B	2522	0	2523	20	0
1	C	2522	0	2523	33	0
1	D	2513	0	2514	30	0
1	E	2522	0	2523	23	0
2	A	12	0	23	4	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
5	A	4	0	3	0	0
5	B	4	0	3	0	0
5	C	8	0	6	0	0
5	D	4	0	3	0	0
5	E	4	0	3	0	0
6	A	6	0	0	17	0
6	B	4	0	0	9	0
6	C	4	0	0	9	0
6	D	4	0	0	12	0
6	E	3	0	0	9	0
All	All	12659	0	12638	129	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 5.

All (129) 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:244:THR:OG1	1:C:247:PRO:HD3	1.72	0.88
1:E:128:ILE:HG13	6:E:405:XE:XE	2.51	0.88
1:B:128:ILE:HG13	6:B:406:XE:XE	2.53	0.87
1:B:7:PRO:HG3	1:B:135:VAL:HG21	1.57	0.85
1:C:128:ILE:HG13	6:C:407:XE:XE	2.57	0.82
1:D:128:ILE:HG13	6:D:406:XE:XE	2.58	0.81
1:A:255:THR:HG23	6:A:408:XE:XE	2.61	0.79
1:B:255:THR:HG23	6:B:405:XE:XE	2.63	0.77
1:A:128:ILE:HG13	6:A:409:XE:XE	2.64	0.76
1:A:188:LEU:HB2	6:A:409:XE:XE	2.65	0.75
1:D:255:THR:HG23	6:D:405:XE:XE	2.67	0.72
1:C:188:LEU:HB2	6:C:407:XE:XE	2.68	0.71
1:D:188:LEU:HB2	6:D:406:XE:XE	2.68	0.71
1:C:253:THR:HG23	1:C:256:GLY:H	1.56	0.70
2:A:401:LMT:H123	6:A:405:XE:XE	2.71	0.69
1:A:253:THR:HG23	1:A:256:GLY:H	1.58	0.69
1:A:197:TYR:CD2	1:A:202:ILE:HD11	2.29	0.68
1:B:253:THR:HG23	1:B:256:GLY:H	1.58	0.68
1:C:255:THR:HG23	6:C:406:XE:XE	2.72	0.68
1:E:253:THR:HG23	1:E:256:GLY:H	1.58	0.68
1:D:253:THR:HG23	1:D:256:GLY:H	1.58	0.68
1:B:188:LEU:HB2	6:B:406:XE:XE	2.72	0.67
1:C:243:GLU:HG2	1:C:245:ASN:H	1.60	0.67
1:E:188:LEU:HB2	6:E:405:XE:XE	2.74	0.65
1:A:197:TYR:CE2	1:A:202:ILE:HD11	2.31	0.65
1:D:165:PHE:HE1	6:D:406:XE:XE	2.58	0.64
1:A:120:PRO:HG3	1:A:197:TYR:CD2	2.33	0.63
1:A:128:ILE:CD1	6:A:409:XE:XE	3.25	0.63
1:C:243:GLU:HG2	1:C:245:ASN:N	2.15	0.61
1:E:255:THR:HG23	6:E:404:XE:XE	2.79	0.61
1:C:208:ILE:HG22	1:C:266:TYR:OH	2.03	0.58
1:B:128:ILE:CG1	6:B:406:XE:XE	3.30	0.57
1:E:128:ILE:CG1	6:E:405:XE:XE	3.29	0.56
1:E:37:PHE:HZ	6:E:405:XE:XE	2.66	0.56
1:C:249:THR:HG22	1:C:251:TYR:O	2.06	0.56
1:B:207:PHE:HB2	6:B:404:XE:XE	2.83	0.56
1:C:128:ILE:CD1	6:C:407:XE:XE	3.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:128:ILE:CD1	6:E:405:XE:XE	3.32	0.55
1:E:37:PHE:CZ	6:E:405:XE:XE	3.38	0.55
1:C:128:ILE:CG1	6:C:407:XE:XE	3.32	0.55
1:D:128:ILE:CG1	6:D:406:XE:XE	3.32	0.55
1:D:194:TYR:CD2	1:D:197:TYR:HB2	2.42	0.54
1:B:207:PHE:CB	6:B:404:XE:XE	3.33	0.54
1:C:211:ILE:HG23	1:D:270:VAL:HG11	1.89	0.54
6:A:407:XE:XE	1:E:207:PHE:HB2	2.87	0.53
6:A:407:XE:XE	1:E:207:PHE:CB	3.35	0.52
1:A:37:PHE:CZ	6:A:409:XE:XE	3.40	0.52
1:C:188:LEU:CB	6:C:407:XE:XE	3.36	0.52
1:E:173:ASN:HB3	1:E:180:LEU:HD11	1.90	0.52
1:D:193:GLN:O	1:D:193:GLN:HG2	2.10	0.52
1:B:188:LEU:CB	6:B:406:XE:XE	3.36	0.52
1:C:7:PRO:HG3	1:C:135:VAL:HG21	1.91	0.52
1:A:188:LEU:CB	6:A:409:XE:XE	3.35	0.51
1:C:207:PHE:O	1:C:211:ILE:HG13	2.11	0.51
1:D:120:PRO:HG3	1:D:197:TYR:CD2	2.45	0.51
2:A:401:LMT:H52	1:D:240:ILE:HG21	1.92	0.51
1:D:193:GLN:O	1:D:193:GLN:CG	2.58	0.51
1:D:197:TYR:CE2	1:D:202:ILE:HD11	2.46	0.51
1:A:128:ILE:CG1	6:A:409:XE:XE	3.36	0.50
1:A:128:ILE:HD12	6:A:409:XE:XE	2.90	0.50
2:A:401:LMT:C12	6:A:405:XE:XE	3.38	0.50
1:D:128:ILE:CD1	6:D:406:XE:XE	3.38	0.50
1:D:165:PHE:CE1	6:D:406:XE:XE	3.42	0.50
1:D:188:LEU:CB	6:D:406:XE:XE	3.39	0.49
1:B:128:ILE:CD1	6:B:406:XE:XE	3.38	0.49
1:B:207:PHE:HB3	6:B:404:XE:XE	2.90	0.49
1:D:120:PRO:HG3	1:D:197:TYR:CG	2.47	0.49
1:A:37:PHE:HZ	6:A:409:XE:XE	2.74	0.49
6:A:407:XE:XE	1:E:207:PHE:HB3	2.90	0.49
1:E:188:LEU:CB	6:E:405:XE:XE	3.38	0.49
1:A:120:PRO:HG3	1:A:197:TYR:CE2	2.48	0.48
1:A:207:PHE:CB	6:A:406:XE:XE	3.40	0.48
1:D:207:PHE:CB	6:D:404:XE:XE	3.40	0.48
1:C:37:PHE:HZ	6:C:407:XE:XE	2.75	0.48
1:C:133:ARG:NH1	1:C:177:GLU:HB2	2.29	0.47
1:B:208:ILE:HG22	1:B:266:TYR:OH	2.14	0.47
1:A:207:PHE:HB2	6:A:406:XE:XE	2.93	0.47
1:C:37:PHE:CZ	6:C:407:XE:XE	3.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:162:ILE:HG13	1:E:190:ILE:HG22	1.97	0.47
1:D:287:ARG:HH11	1:D:287:ARG:HA	1.79	0.47
1:D:207:PHE:HB3	6:D:404:XE:XE	2.94	0.46
2:A:401:LMT:H51	1:C:240:ILE:HG21	1.97	0.46
1:D:128:ILE:HD12	6:D:406:XE:XE	2.93	0.46
1:A:207:PHE:HB3	6:A:406:XE:XE	2.94	0.46
1:A:211:ILE:HG23	1:B:270:VAL:HG11	1.98	0.46
1:B:279:LEU:HD13	1:B:287:ARG:HB3	1.98	0.46
1:C:244:THR:HG1	1:C:247:PRO:HD3	1.75	0.46
1:D:207:PHE:HB2	6:D:404:XE:XE	2.94	0.45
1:A:197:TYR:CD2	1:A:202:ILE:CD1	2.98	0.45
1:E:226:THR:O	1:E:230:SER:HB2	2.17	0.45
1:A:152:ASN:HB3	1:A:155:VAL:HG23	1.99	0.45
1:B:152:ASN:HB3	1:B:155:VAL:HG23	1.99	0.45
1:D:211:ILE:HG23	1:E:270:VAL:HG11	1.97	0.44
1:C:144:VAL:HG11	1:C:167:ALA:HB3	1.99	0.44
1:A:198:ILE:HA	1:A:202:ILE:HD12	1.99	0.44
1:C:135:VAL:HG23	1:C:137:THR:H	1.82	0.44
1:B:208:ILE:HG12	1:C:232:LEU:HD21	1.99	0.44
1:C:128:ILE:HD12	6:C:407:XE:XE	2.96	0.44
1:D:226:THR:O	1:D:230:SER:HB2	2.18	0.44
1:C:195:PHE:HE1	1:D:252:MET:HB2	1.83	0.44
1:D:133:ARG:NH1	1:D:177:GLU:HB2	2.32	0.44
1:A:226:THR:O	1:A:230:SER:HB2	2.17	0.44
1:B:226:THR:O	1:B:230:SER:HB2	2.17	0.44
1:A:7:PRO:HG3	1:A:135:VAL:HG21	2.00	0.43
1:A:253:THR:HG23	1:A:256:GLY:N	2.31	0.43
1:C:226:THR:O	1:C:230:SER:HB2	2.19	0.43
1:D:134:SER:HB3	1:D:139:ASN:HA	2.01	0.43
1:A:79:VAL:HG21	1:A:183:LYS:HD2	2.00	0.43
1:C:152:ASN:HB3	1:C:155:VAL:HG23	1.99	0.43
1:D:253:THR:HG23	1:D:256:GLY:N	2.31	0.43
1:A:270:VAL:HG11	1:E:211:ILE:HG12	2.00	0.43
1:B:79:VAL:HG21	1:B:183:LYS:HD2	2.01	0.43
1:C:253:THR:HG23	1:C:256:GLY:N	2.29	0.42
1:A:134:SER:HB3	1:A:139:ASN:HA	2.01	0.42
1:E:134:SER:HB3	1:E:139:ASN:HA	2.00	0.42
1:E:254:TYR:CZ	1:E:258:ILE:HD11	2.54	0.42
1:E:242:VAL:O	1:E:243:GLU:C	2.63	0.42
1:A:254:TYR:CZ	1:A:258:ILE:HD11	2.55	0.41
1:C:134:SER:HA	1:C:140:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:254:TYR:CZ	1:C:258:ILE:HD11	2.55	0.41
1:B:211:ILE:HG23	1:C:270:VAL:HG11	2.01	0.41
1:D:254:TYR:CZ	1:D:258:ILE:HD11	2.55	0.41
1:B:254:TYR:CZ	1:B:258:ILE:HD11	2.56	0.41
1:E:7:PRO:HG3	1:E:135:VAL:HG21	2.01	0.41
1:C:173:ASN:HB3	1:C:180:LEU:HD11	2.02	0.41
1:C:162:ILE:HG13	1:C:190:ILE:HG22	2.03	0.41
1:E:128:ILE:HD12	6:E:405:XE:XE	2.99	0.41
1:D:211:ILE:HG12	1:E:270:VAL:HG11	2.01	0.40
1:A:162:ILE:HG13	1:A:190:ILE:HG22	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	309/317 (98%)	302 (98%)	7 (2%)	0	100	100
1	B	310/317 (98%)	305 (98%)	5 (2%)	0	100	100
1	C	310/317 (98%)	304 (98%)	4 (1%)	2 (1%)	21	50
1	D	309/317 (98%)	302 (98%)	7 (2%)	0	100	100
1	E	310/317 (98%)	301 (97%)	9 (3%)	0	100	100
All	All	1548/1585 (98%)	1514 (98%)	32 (2%)	2 (0%)	48	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	247	PRO
1	C	246	CYS

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	278/284 (98%)	269 (97%)	9 (3%)	34	58
1	B	279/284 (98%)	272 (98%)	7 (2%)	42	62
1	C	279/284 (98%)	264 (95%)	15 (5%)	20	47
1	D	278/284 (98%)	268 (96%)	10 (4%)	31	56
1	E	279/284 (98%)	270 (97%)	9 (3%)	34	58
All	All	1393/1420 (98%)	1343 (96%)	50 (4%)	31	56

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	141	VAL
1	A	170	LYS
1	A	242	VAL
1	A	259	ILE
1	A	281	VAL
1	A	292	THR
1	A	297	ILE
1	A	304	LEU
1	A	307	ASN
1	B	141	VAL
1	B	259	ILE
1	B	281	VAL
1	B	297	ILE
1	B	304	LEU
1	B	305	LEU
1	B	307	ASN
1	C	5	VAL
1	C	33	CYS
1	C	101	GLN
1	C	135	VAL
1	C	141	VAL
1	C	144	VAL
1	C	241	LEU

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Mol	Chain	Res	Type
1	C	243	GLU
1	C	244	THR
1	C	246	CYS
1	C	259	ILE
1	C	281	VAL
1	C	297	ILE
1	C	304	LEU
1	C	307	ASN
1	D	141	VAL
1	D	147	GLU
1	D	170	LYS
1	D	180	LEU
1	D	192	ARG
1	D	259	ILE
1	D	281	VAL
1	D	297	ILE
1	D	304	LEU
1	D	307	ASN
1	E	141	VAL
1	E	170	LYS
1	E	228	VAL
1	E	249	THR
1	E	259	ILE
1	E	281	VAL
1	E	297	ILE
1	E	304	LEU
1	E	307	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	GLN
1	A	239	ASN
1	A	277	HIS
1	A	284	GLN
1	A	307	ASN
1	B	239	ASN
1	B	277	HIS
1	B	307	ASN
1	C	193	GLN
1	C	277	HIS
1	C	284	GLN

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Mol	Chain	Res	Type
1	D	277	HIS
1	D	307	ASN
1	E	239	ASN
1	E	276	GLN
1	E	277	HIS
1	E	307	ASN

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](#)

Of 38 ligands modelled in this entry, 31 are monoatomic - leaving 7 for Mogul analysis.

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	$\# Z > 2$
5	ACT	C	404	-	3,3,3	1.20	0	3,3,3	0.87	0
5	ACT	D	403	-	3,3,3	1.26	0	3,3,3	0.71	0
5	ACT	E	403	-	3,3,3	1.03	0	3,3,3	1.02	0
5	ACT	B	403	-	3,3,3	1.22	0	3,3,3	0.78	0
5	ACT	C	401	-	3,3,3	1.19	0	3,3,3	0.87	0
5	ACT	A	404	-	3,3,3	1.14	0	3,3,3	0.88	0
2	LMT	A	401	-	11,11,36	0.25	0	10,10,47	0.79	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
2	LMT	A	401	-	-	0/9/9/61	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	LMT	4	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	311/317 (98%)	0.16	8 (2%) 57 42	95, 125, 179, 210	0
1	B	311/317 (98%)	0.16	11 (3%) 47 34	59, 117, 145, 169	1 (0%)
1	C	311/317 (98%)	0.18	10 (3%) 50 37	58, 121, 158, 181	1 (0%)
1	D	311/317 (98%)	0.19	11 (3%) 47 34	93, 127, 179, 227	0
1	E	311/317 (98%)	0.33	14 (4%) 38 28	55, 126, 172, 196	1 (0%)
All	All	1555/1585 (98%)	0.20	54 (3%) 47 34	55, 123, 171, 227	3 (0%)

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	161	ASP	5.2
1	D	244	THR	5.0
1	B	244	THR	4.4
1	D	246	CYS	4.2
1	A	245	ASN	4.1
1	A	244	THR	4.0
1	E	244	THR	4.0
1	B	246	CYS	3.9
1	A	61	VAL	3.8
1	D	245	ASN	3.7
1	E	245	ASN	3.6
1	A	246	CYS	3.5
1	C	147	GLU	3.5
1	A	56	PRO	3.5
1	E	146	LEU	3.4
1	C	246	CYS	3.4
1	B	245	ASN	3.3
1	C	33	CYS	3.3
1	A	57	VAL	3.3
1	A	63	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	247	PRO	3.0
1	B	255	THR	3.0
1	C	244	THR	3.0
1	D	197	TYR	2.9
1	C	161	ASP	2.9
1	E	191	SER	2.6
1	E	5	VAL	2.6
1	E	246	CYS	2.5
1	B	259	ILE	2.5
1	E	193	GLN	2.5
1	E	163	GLU	2.5
1	B	193	GLN	2.4
1	D	193	GLN	2.4
1	E	147	GLU	2.4
1	D	61	VAL	2.4
1	C	193	GLN	2.3
1	C	245	ASN	2.3
1	A	255	THR	2.3
1	D	161	ASP	2.3
1	E	58	ARG	2.3
1	D	25	ILE	2.3
1	D	255	THR	2.2
1	E	144	VAL	2.2
1	E	225	VAL	2.2
1	D	253	THR	2.2
1	C	191	SER	2.1
1	B	11	ILE	2.1
1	E	247	PRO	2.1
1	D	284	GLN	2.1
1	B	5	VAL	2.1
1	B	248	LYS	2.1
1	B	284	GLN	2.1
1	C	146	LEU	2.0
1	B	44	SER	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(Å ²)	Q<0.9
4	NA	A	403	1/1	0.66	0.11	136,136,136,136	0
4	NA	B	402	1/1	0.70	0.34	132,132,132,132	0
4	NA	C	403	1/1	0.76	0.10	116,116,116,116	0
5	ACT	D	403	4/4	0.76	0.29	133,133,134,138	0
5	ACT	E	403	4/4	0.77	0.18	116,117,117,117	0
5	ACT	B	403	4/4	0.83	0.19	108,112,112,114	0
6	XE	D	406	1/1	0.83	0.20	90,90,90,90	1
6	XE	B	407	1/1	0.86	0.21	72,72,72,72	1
4	NA	D	402	1/1	0.87	0.15	146,146,146,146	0
3	CL	D	401	1/1	0.87	0.11	113,113,113,113	0
2	LMT	A	401	12/35	0.87	0.31	102,104,109,112	0
3	CL	B	401	1/1	0.89	0.12	103,103,103,103	0
6	XE	B	406	1/1	0.89	0.13	67,67,67,67	1
6	XE	E	405	1/1	0.89	0.17	68,68,68,68	1
5	ACT	C	404	4/4	0.90	0.14	105,107,107,109	0
3	CL	A	402	1/1	0.92	0.13	105,105,105,105	0
6	XE	E	406	1/1	0.92	0.13	74,74,74,74	1
5	ACT	A	404	4/4	0.93	0.15	125,127,128,128	0
6	XE	C	407	1/1	0.93	0.22	67,67,67,67	1
6	XE	A	409	1/1	0.93	0.16	72,72,72,72	1
6	XE	A	410	1/1	0.93	0.09	64,64,64,64	1
4	NA	E	402	1/1	0.93	0.07	111,111,111,111	0
6	XE	E	404	1/1	0.94	0.12	161,161,161,161	1
5	ACT	C	401	4/4	0.94	0.13	102,103,103,106	0
6	XE	B	405	1/1	0.94	0.19	139,139,139,139	1
6	XE	C	406	1/1	0.95	0.11	152,152,152,152	1
3	CL	C	402	1/1	0.95	0.08	109,109,109,109	0
3	CL	E	401	1/1	0.96	0.07	104,104,104,104	0
6	XE	D	405	1/1	0.97	0.11	153,153,153,153	1
6	XE	C	408	1/1	0.97	0.07	61,61,61,61	1
6	XE	D	407	1/1	0.97	0.15	70,70,70,70	1
6	XE	A	405	1/1	0.98	0.10	146,146,146,146	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	XE	A	408	1/1	0.99	0.13	144,144,144,144	1
6	XE	B	404	1/1	0.99	0.02	100,100,100,100	1
6	XE	C	405	1/1	0.99	0.02	107,107,107,107	1
6	XE	A	407	1/1	0.99	0.03	106,106,106,106	1
6	XE	A	406	1/1	1.00	0.01	110,110,110,110	1
6	XE	D	404	1/1	1.00	0.02	111,111,111,111	1

6.5 Other polymers [i](#)

There are no such residues in this entry.