



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:46 AM UTC

PDB ID : 6IVJ / pdb_00006ivj
Title : Crystal structure of a membrane protein G18A
Authors : Kittredge, A.; Fukuda, F.; Zhang, Y.; Yang, T.
Deposited on : 2018-12-04
Resolution : 2.77 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

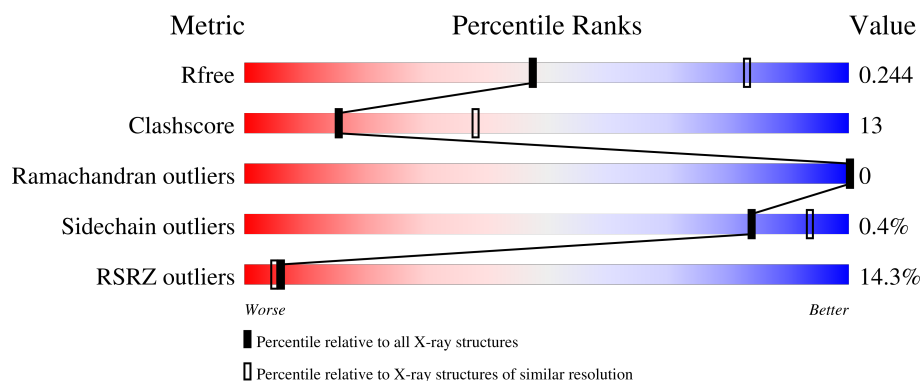
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	297	15% 68% 22% 9%
1	B	297	13% 62% 29% 9%
1	C	297	11% 72% 18% 9%
1	D	297	11% 69% 22% 10%
1	E	297	15% 67% 23% 9%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	CL	C	305	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10808 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 Ibestrophin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	269	Total	C	N	O	S	0	0	0
			2143	1389	362	383	9			
1	B	270	Total	C	N	O	S	0	0	0
			2152	1394	365	384	9			
1	C	269	Total	C	N	O	S	0	0	0
			2137	1385	363	380	9			
1	D	268	Total	C	N	O	S	0	0	0
			2124	1376	360	379	9			
1	E	270	Total	C	N	O	S	0	0	0
			2152	1396	364	383	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP W9BH30
A	-1	ASN	-	expression tag	UNP W9BH30
A	0	ALA	-	expression tag	UNP W9BH30
A	18	ALA	GLY	engineered mutation	UNP W9BH30
B	-2	SER	-	expression tag	UNP W9BH30
B	-1	ASN	-	expression tag	UNP W9BH30
B	0	ALA	-	expression tag	UNP W9BH30
B	18	ALA	GLY	engineered mutation	UNP W9BH30
C	-2	SER	-	expression tag	UNP W9BH30
C	-1	ASN	-	expression tag	UNP W9BH30
C	0	ALA	-	expression tag	UNP W9BH30
C	18	ALA	GLY	engineered mutation	UNP W9BH30
D	-2	SER	-	expression tag	UNP W9BH30
D	-1	ASN	-	expression tag	UNP W9BH30
D	0	ALA	-	expression tag	UNP W9BH30
D	18	ALA	GLY	engineered mutation	UNP W9BH30
E	-2	SER	-	expression tag	UNP W9BH30
E	-1	ASN	-	expression tag	UNP W9BH30
E	0	ALA	-	expression tag	UNP W9BH30

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Chain	Residue	Modelled	Actual	Comment	Reference
E	18	ALA	GLY	engineered mutation	UNP W9BH30

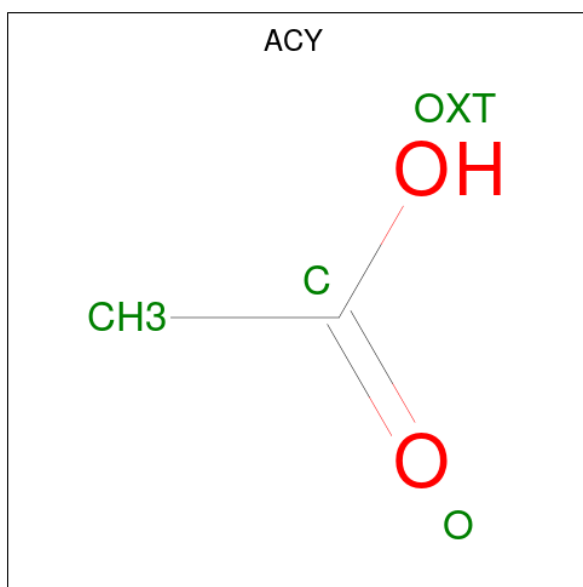
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Zn 3 3	0	0
2	B	5	Total Zn 5 5	0	0
2	C	3	Total Zn 3 3	0	0
2	D	3	Total Zn 3 3	0	0
2	E	2	Total Zn 2 2	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Cl 2 2	0	0
3	B	1	Total Cl 1 1	0	0
3	C	3	Total Cl 3 3	0	0
3	D	3	Total Cl 3 3	0	0
3	E	2	Total Cl 2 2	0	0

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



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

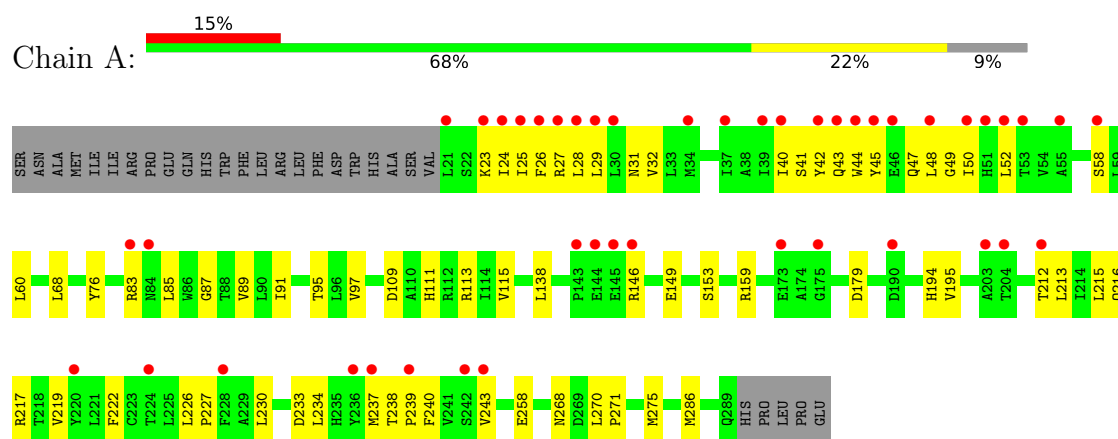
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	19	Total	O	0	0
			19	19		
5	B	15	Total	O	0	0
			15	15		
5	C	12	Total	O	0	0
			12	12		
5	D	12	Total	O	0	0
			12	12		
5	E	11	Total	O	0	0
			11	11		

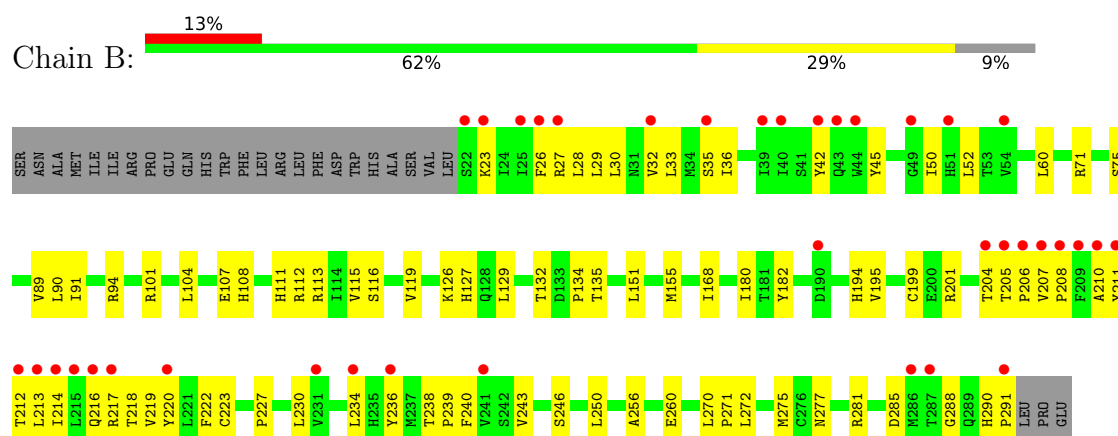
3 Residue-property plots [i](#)

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.

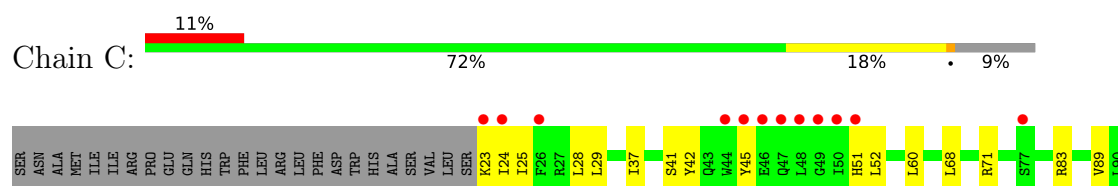
• Molecule 1: Ibestrophin

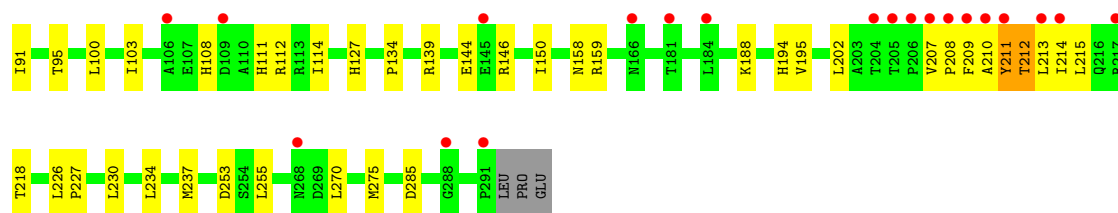


• Molecule 1: Ibestrophin

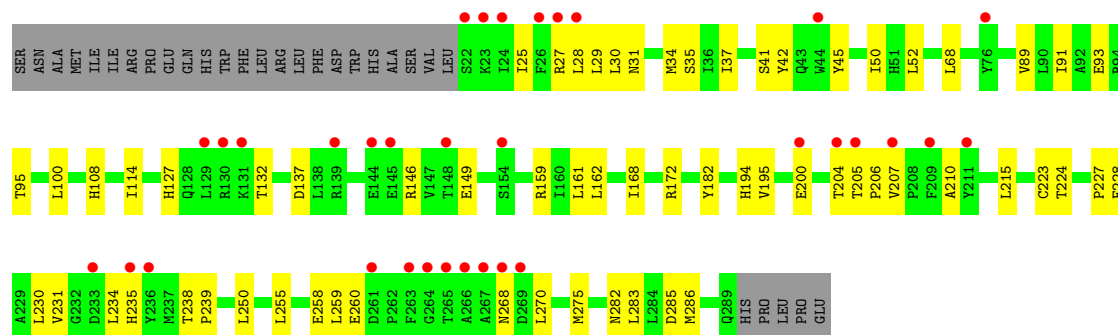


• Molecule 1: Ibestrophin

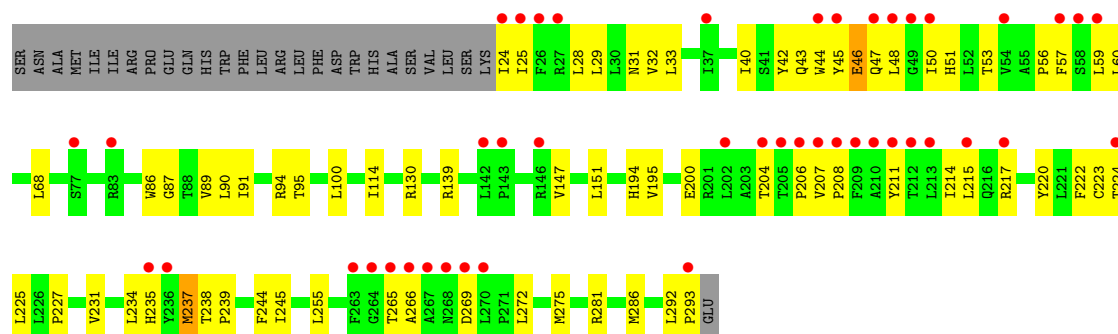




● Molecule 1: Ibestrophin



● Molecule 1: Ibestrophin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.12Å 160.03Å 161.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.76 – 2.77 48.76 – 2.77	Depositor EDS
% Data completeness (in resolution range)	98.7 (48.76-2.77) 99.1 (48.76-2.77)	Depositor EDS
R_{merge}	0.01	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0238, PHENIX	Depositor
R, R_{free}	0.215 , 0.244 0.215 , 0.244	Depositor DCC
R_{free} test set	3783 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	77.4	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 38.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10808	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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, ZN, ACY

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	0.42	0/2189	0.59	0/2979
1	B	0.44	0/2200	0.71	0/2995
1	C	0.40	0/2185	0.63	0/2975
1	D	0.36	0/2170	0.56	0/2954
1	E	0.43	0/2201	0.69	0/2999
All	All	0.41	0/10945	0.64	0/14902

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	46	GLU	Peptide

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	2143	0	2192	68	0
1	B	2152	0	2195	75	0
1	C	2137	0	2170	64	0
1	D	2124	0	2162	63	0
1	E	2152	0	2195	55	0
2	A	3	0	0	0	0
2	B	5	0	0	1	0
2	C	3	0	0	1	0
2	D	3	0	0	0	0
2	E	2	0	0	0	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	3	0	0	3	0
3	D	3	0	0	1	0
3	E	2	0	0	0	0
4	C	4	0	3	1	0
5	A	19	0	0	0	0
5	B	15	0	0	4	0
5	C	12	0	0	1	0
5	D	12	0	0	0	0
5	E	11	0	0	0	0
All	All	10808	0	10917	289	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 13.

All (289) 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:A:42:TYR:HA	1:A:45:TYR:CD1	1.82	1.13
1:C:207:VAL:HG22	1:C:208:PRO:HD2	1.20	1.10
1:C:188:LYS:HE3	3:C:305:CL:CL	1.91	1.08
1:B:71:ARG:NH1	1:B:260:GLU:OE2	1.89	1.06
1:C:207:VAL:HG22	1:C:208:PRO:CD	1.93	0.98
1:B:45:TYR:CD2	1:B:50:ILE:HG13	2.01	0.94
1:A:42:TYR:HA	1:A:45:TYR:CE1	2.01	0.94
1:E:231:VAL:HA	1:E:238:THR:HG21	1.52	0.91
1:C:211:TYR:HE1	1:E:255:LEU:HD21	1.36	0.91
1:C:42:TYR:HA	1:C:45:TYR:CD1	2.08	0.88
1:B:111:HIS:ND1	5:B:401:HOH:O	2.07	0.87
1:E:206:PRO:HB2	1:E:208:PRO:HD2	1.57	0.86
1:A:42:TYR:CA	1:A:45:TYR:CD1	2.59	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:TYR:CE1	1:E:255:LEU:HD21	2.12	0.84
1:A:58:SER:HA	1:E:59:LEU:HD12	1.59	0.84
1:B:45:TYR:HD2	1:B:50:ILE:HG13	1.42	0.81
1:B:210:ALA:O	1:B:214:ILE:HG12	1.79	0.81
1:C:210:ALA:O	1:C:214:ILE:HG13	1.83	0.79
1:B:236:TYR:O	1:B:239:PRO:HD2	1.83	0.78
1:A:68:LEU:HD23	1:A:215:LEU:HD13	1.64	0.77
1:A:41:SER:C	1:A:45:TYR:CE1	2.63	0.76
1:A:41:SER:C	1:A:45:TYR:HE1	1.93	0.75
1:A:42:TYR:C	1:A:45:TYR:HD1	1.95	0.74
1:A:29:LEU:HA	1:A:32:VAL:HG23	1.70	0.74
1:B:211:TYR:CD2	1:D:255:LEU:HD21	2.23	0.74
1:C:194:HIS:HA	1:E:91:ILE:HD13	1.71	0.73
1:E:46:GLU:HG2	1:E:51:HIS:ND1	2.04	0.73
1:B:42:TYR:O	1:B:45:TYR:N	2.18	0.72
1:C:255:LEU:CD1	1:D:210:ALA:HB3	2.19	0.72
1:C:207:VAL:CG2	1:C:208:PRO:HD2	2.12	0.72
2:B:304:ZN:ZN	5:B:401:HOH:O	1.39	0.72
1:A:42:TYR:CA	1:A:45:TYR:HD1	2.02	0.70
1:D:234:LEU:HB2	1:D:238:THR:HG22	1.73	0.70
2:C:303:ZN:ZN	3:C:307:CL:CL	1.77	0.70
1:A:42:TYR:O	1:A:45:TYR:HD1	1.74	0.70
1:A:237:MET:C	1:A:239:PRO:HD2	2.16	0.69
1:C:91:ILE:HD13	1:D:194:HIS:HA	1.75	0.69
1:E:292:LEU:HD22	1:E:293:PRO:HD2	1.74	0.69
1:C:207:VAL:CG2	1:C:208:PRO:CD	2.70	0.67
1:A:41:SER:O	1:A:45:TYR:CE1	2.48	0.67
1:B:213:LEU:HA	1:B:216:GLN:HG2	1.75	0.66
1:E:234:LEU:HB2	1:E:238:THR:HG22	1.77	0.66
1:B:71:ARG:HH22	1:B:212:THR:HG22	1.60	0.66
1:C:68:LEU:HD23	1:C:215:LEU:HD13	1.79	0.65
1:D:231:VAL:HA	1:D:238:THR:HG21	1.77	0.65
1:A:41:SER:O	1:A:45:TYR:CD1	2.49	0.65
1:D:258:GLU:O	1:D:268:ASN:ND2	2.25	0.64
1:C:214:ILE:O	1:C:218:THR:HG23	1.98	0.64
1:A:83:ARG:HH21	1:A:87:GLY:HA2	1.62	0.64
1:C:42:TYR:HA	1:C:45:TYR:HD1	1.58	0.64
1:A:194:HIS:HA	1:B:91:ILE:HD13	1.80	0.63
1:A:258:GLU:OE2	1:A:268:ASN:ND2	2.31	0.63
1:B:208:PRO:HG2	1:D:259:LEU:HD21	1.81	0.62
1:C:255:LEU:HD13	1:D:210:ALA:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:204:THR:O	1:D:206:PRO:HD3	1.98	0.62
1:C:188:LYS:CE	3:C:305:CL:CL	2.77	0.62
1:B:214:ILE:O	1:B:218:THR:HG23	2.00	0.61
1:B:104:LEU:HB3	1:B:107:GLU:HG3	1.82	0.61
1:E:47:GLN:HB2	1:E:48:LEU:HD12	1.82	0.61
1:C:71:ARG:NH2	1:C:253:ASP:OD1	2.33	0.61
1:E:60:LEU:HD23	1:E:222:PHE:HD2	1.64	0.61
1:E:91:ILE:O	1:E:95:THR:HG23	2.00	0.61
1:A:42:TYR:CA	1:A:45:TYR:CE1	2.79	0.61
1:B:211:TYR:CE2	1:D:255:LEU:HD21	2.35	0.61
1:D:235:HIS:O	1:D:238:THR:HG23	2.00	0.60
1:B:213:LEU:O	1:B:217:ARG:HG3	2.01	0.60
1:C:209:PHE:O	1:C:210:ALA:C	2.45	0.60
1:C:255:LEU:HD13	1:D:210:ALA:CB	2.31	0.59
1:C:52:LEU:HD21	1:E:237:MET:SD	2.42	0.59
1:A:24:ILE:HG13	1:A:27:ARG:HB2	1.84	0.58
1:B:29:LEU:O	1:B:32:VAL:HG12	2.02	0.58
1:A:271:PRO:O	1:A:275:MET:HG3	2.03	0.58
1:A:32:VAL:HG22	1:A:243:VAL:HG21	1.85	0.58
1:B:135:THR:HG23	1:B:151:LEU:HD11	1.84	0.58
1:D:172:ARG:HB3	1:D:172:ARG:HH21	1.68	0.58
1:E:130:ARG:NH1	1:E:269:ASP:OD2	2.36	0.58
1:E:217:ARG:HA	1:E:220:TYR:HB2	1.85	0.58
1:C:60:LEU:HD21	1:E:244:PHE:CE2	2.39	0.57
1:A:42:TYR:O	1:A:45:TYR:CD1	2.57	0.57
1:A:239:PRO:HG2	1:A:240:PHE:HD1	1.70	0.57
1:B:127:HIS:CD2	1:B:134:PRO:HA	2.40	0.57
1:B:168:ILE:HG22	1:B:182:TYR:CE1	2.39	0.57
1:B:207:VAL:HG21	1:B:260:GLU:CD	2.30	0.56
1:B:155:MET:HE1	1:D:285:ASP:HB3	1.86	0.56
1:E:40:ILE:O	1:E:43:GLN:HG3	2.05	0.56
1:E:100:LEU:HD11	1:E:114:ILE:HD13	1.87	0.56
1:B:246:SER:O	1:B:250:LEU:HD12	2.06	0.56
1:A:238:THR:N	1:A:239:PRO:HD2	2.21	0.56
1:E:24:ILE:HG23	1:E:25:ILE:H	1.70	0.56
1:C:91:ILE:O	1:C:95:THR:HG23	2.07	0.55
1:B:129:LEU:HD21	1:B:199:CYS:HB3	1.87	0.55
1:A:230:LEU:HB3	1:A:234:LEU:HD12	1.89	0.55
1:B:71:ARG:HD3	1:B:256:ALA:HB1	1.89	0.55
1:B:155:MET:HE1	1:D:285:ASP:CB	2.37	0.55
1:D:27:ARG:HH12	1:D:224:THR:CG2	2.19	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:270:LEU:HD22	1:D:275:MET:HE1	1.89	0.55
1:C:37:ILE:O	1:C:41:SER:OG	2.19	0.54
1:A:25:ILE:O	1:A:29:LEU:HG	2.06	0.54
1:D:34:MET:HE3	1:D:228:PHE:HE1	1.72	0.54
1:B:108:HIS:HA	5:B:401:HOH:O	2.07	0.54
1:D:91:ILE:O	1:D:95:THR:HG23	2.06	0.54
1:A:40:ILE:O	1:A:44:TRP:HZ3	1.91	0.53
1:C:100:LEU:HD11	1:C:114:ILE:HD13	1.89	0.53
1:C:150:ILE:HA	1:C:159:ARG:HG2	1.90	0.53
1:D:146:ARG:HA	1:D:149:GLU:HG3	1.90	0.53
1:B:23:LYS:O	1:B:27:ARG:NH2	2.42	0.53
1:C:42:TYR:HD1	1:C:45:TYR:CD1	2.26	0.53
1:B:277:ASN:O	1:B:281:ARG:HG3	2.09	0.53
1:C:42:TYR:HA	1:C:45:TYR:CE1	2.43	0.53
1:D:224:THR:O	1:D:227:PRO:HD2	2.09	0.53
1:B:71:ARG:CD	1:B:256:ALA:HB1	2.38	0.53
1:D:168:ILE:HG22	1:D:182:TYR:CD2	2.43	0.53
1:C:52:LEU:HA	1:E:234:LEU:HD22	1.91	0.53
4:C:304:ACY:H3	1:E:281:ARG:HD3	1.91	0.52
1:E:89:VAL:HG23	1:E:195:VAL:HG11	1.91	0.52
1:B:271:PRO:HG2	1:B:275:MET:HE3	1.91	0.52
1:C:23:LYS:HB3	1:C:24:ILE:HD12	1.91	0.52
1:D:231:VAL:CA	1:D:238:THR:HG21	2.39	0.52
1:E:46:GLU:HG2	1:E:51:HIS:CE1	2.44	0.52
1:A:91:ILE:O	1:A:95:THR:HG23	2.10	0.52
1:C:237:MET:HE2	1:D:52:LEU:HD21	1.92	0.52
1:A:43:GLN:O	1:A:47:GLN:NE2	2.42	0.52
1:C:270:LEU:HD22	1:C:275:MET:HE1	1.91	0.52
1:B:101:ARG:HD3	1:B:290:HIS:HB3	1.91	0.52
1:E:200:GLU:O	1:E:204:THR:HG23	2.09	0.52
1:B:216:GLN:HA	1:B:219:VAL:HG12	1.91	0.52
1:B:201:ARG:O	1:B:205:THR:HB	2.10	0.51
1:C:111:HIS:CE1	5:C:401:HOH:O	2.64	0.51
1:B:42:TYR:O	1:B:45:TYR:HD1	1.93	0.51
1:B:108:HIS:ND1	5:B:401:HOH:O	2.34	0.51
1:A:89:VAL:HG23	1:A:195:VAL:HG11	1.92	0.51
1:B:28:LEU:HG	1:B:243:VAL:HG22	1.92	0.51
1:A:58:SER:HA	1:E:59:LEU:CD1	2.36	0.51
1:C:210:ALA:O	1:C:214:ILE:CG1	2.55	0.50
1:D:35:SER:HB3	1:D:227:PRO:HB3	1.93	0.50
1:C:208:PRO:C	1:C:210:ALA:N	2.67	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:LEU:HD21	1:A:217:ARG:CZ	2.41	0.50
1:B:126:LYS:HG3	1:B:272:LEU:HB3	1.93	0.50
1:C:83:ARG:HG2	1:D:205:THR:HG21	1.94	0.50
1:C:108:HIS:O	1:C:112:ARG:HG3	2.11	0.50
1:B:52:LEU:HD23	1:D:234:LEU:HD22	1.94	0.50
1:D:108:HIS:HB3	3:D:306:CL:CL	2.49	0.49
1:A:153:SER:OG	1:B:285:ASP:OD2	2.25	0.49
1:D:50:ILE:O	1:D:50:ILE:HG13	2.11	0.49
1:C:25:ILE:HA	1:C:28:LEU:HB2	1.95	0.49
1:C:208:PRO:O	1:C:210:ALA:N	2.45	0.49
1:A:47:GLN:H	1:A:47:GLN:CD	2.15	0.49
1:A:109:ASP:HB3	1:A:113:ARG:HH12	1.76	0.49
1:B:127:HIS:HD1	1:B:132:THR:HG1	1.54	0.49
1:B:227:PRO:HA	1:B:230:LEU:HB2	1.95	0.49
1:D:100:LEU:HD11	1:D:114:ILE:HD13	1.93	0.49
1:B:194:HIS:HA	1:D:91:ILE:HD13	1.94	0.49
1:E:31:ASN:ND2	1:E:223:CYS:O	2.37	0.49
1:A:212:THR:O	1:A:216:GLN:HG2	2.12	0.48
1:D:238:THR:N	1:D:239:PRO:HD2	2.28	0.48
1:A:47:GLN:OE1	1:A:47:GLN:N	2.29	0.48
1:A:76:TYR:CE2	1:E:207:VAL:HG21	2.48	0.48
1:D:127:HIS:NE2	1:D:137:ASP:OD2	2.32	0.48
1:D:200:GLU:O	1:D:204:THR:HG23	2.14	0.48
1:A:44:TRP:O	1:A:48:LEU:HD12	2.12	0.48
1:A:83:ARG:NH1	1:A:275:MET:SD	2.86	0.48
1:C:211:TYR:O	1:C:215:LEU:HB2	2.13	0.48
1:E:139:ARG:HG3	1:E:139:ARG:HH11	1.79	0.48
1:B:112:ARG:NH2	1:B:288:GLY:O	2.46	0.48
1:E:238:THR:N	1:E:239:PRO:HD2	2.29	0.47
1:B:126:LYS:CG	1:B:272:LEU:HB3	2.44	0.47
1:C:158:ASN:OD1	1:E:94:ARG:HD3	2.14	0.47
1:C:230:LEU:HB3	1:C:234:LEU:HD12	1.96	0.47
1:B:90:LEU:HG	1:B:94:ARG:HD2	1.96	0.47
1:C:255:LEU:CD1	1:D:210:ALA:CB	2.89	0.47
1:E:207:VAL:N	1:E:208:PRO:HD2	2.29	0.47
1:C:89:VAL:HG23	1:C:195:VAL:HG11	1.95	0.47
1:A:97:VAL:HG12	1:A:286:MET:HE2	1.96	0.47
1:A:233:ASP:HB3	1:E:53:THR:OG1	2.15	0.47
1:B:127:HIS:ND1	1:B:132:THR:OG1	2.41	0.47
1:C:146:ARG:HH11	1:C:146:ARG:HA	1.80	0.47
1:D:27:ARG:HH12	1:D:224:THR:HG23	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:VAL:HA	1:D:238:THR:CG2	2.44	0.47
1:E:211:TYR:O	1:E:214:ILE:HG13	2.15	0.47
1:A:111:HIS:O	1:A:115:VAL:HG23	2.15	0.47
1:A:215:LEU:HD23	1:A:215:LEU:HA	1.68	0.47
1:A:83:ARG:HH21	1:A:87:GLY:CA	2.28	0.47
1:D:146:ARG:HH11	1:D:149:GLU:CD	2.22	0.47
1:C:209:PHE:O	1:C:212:THR:N	2.48	0.47
1:D:42:TYR:O	1:D:45:TYR:HB2	2.15	0.46
1:B:155:MET:HE3	1:D:282:ASN:HA	1.98	0.46
1:D:100:LEU:HD11	1:D:114:ILE:CD1	2.45	0.46
1:A:270:LEU:HB3	1:A:275:MET:HE1	1.98	0.46
1:B:155:MET:HE1	1:D:285:ASP:CG	2.40	0.46
1:E:53:THR:HB	1:E:56:PRO:HD3	1.98	0.46
1:B:35:SER:OG	1:B:238:THR:HG22	2.16	0.46
1:A:91:ILE:HD13	1:E:194:HIS:HA	1.97	0.46
1:B:42:TYR:HA	1:B:45:TYR:CD1	2.51	0.46
1:B:204:THR:O	1:B:206:PRO:HD3	2.16	0.46
1:A:52:LEU:HD23	1:B:234:LEU:HD22	1.98	0.46
1:E:68:LEU:HD23	1:E:215:LEU:HD21	1.97	0.46
1:B:33:LEU:HA	1:B:36:ILE:HB	1.98	0.45
1:B:223:CYS:SG	1:B:246:SER:HB3	2.57	0.45
1:C:103:ILE:HG23	1:D:172:ARG:NH1	2.31	0.45
1:C:255:LEU:HD11	1:D:210:ALA:HB3	1.96	0.45
1:D:172:ARG:HB3	1:D:172:ARG:NH2	2.31	0.45
1:E:28:LEU:O	1:E:32:VAL:HG23	2.17	0.45
1:A:42:TYR:N	1:A:45:TYR:HE1	2.15	0.45
1:A:270:LEU:HD22	1:A:275:MET:HE1	1.98	0.45
1:B:45:TYR:CD2	1:B:50:ILE:CG1	2.88	0.45
1:C:285:ASP:O	1:D:159:ARG:NH2	2.48	0.45
1:A:85:LEU:HD22	1:A:195:VAL:HA	1.99	0.45
1:D:25:ILE:HA	1:D:28:LEU:HG	1.99	0.44
1:D:93:GLU:HB3	1:D:283:LEU:HD11	1.99	0.44
1:B:60:LEU:HD23	1:B:222:PHE:HD2	1.83	0.44
1:D:29:LEU:HD23	1:D:29:LEU:HA	1.71	0.44
1:B:89:VAL:CG2	1:B:195:VAL:HG11	2.47	0.44
1:A:24:ILE:C	1:A:26:PHE:N	2.75	0.44
1:C:208:PRO:C	1:C:210:ALA:H	2.24	0.44
1:C:214:ILE:O	1:C:218:THR:CG2	2.66	0.44
1:B:75:SER:OG	1:B:260:GLU:HG3	2.18	0.44
1:D:31:ASN:ND2	1:D:223:CYS:O	2.50	0.44
1:E:272:LEU:HA	1:E:272:LEU:HD12	1.69	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:LEU:O	1:A:31:ASN:HB3	2.18	0.44
1:D:250:LEU:HD23	1:D:250:LEU:HA	1.82	0.44
1:B:236:TYR:C	1:B:238:THR:H	2.25	0.43
1:E:87:GLY:O	1:E:91:ILE:HG13	2.18	0.43
1:D:37:ILE:O	1:D:41:SER:HB3	2.18	0.43
1:C:51:HIS:C	1:C:52:LEU:HD23	2.44	0.43
1:C:188:LYS:HA	1:C:188:LYS:HD3	1.88	0.43
1:B:113:ARG:O	1:B:116:SER:OG	2.36	0.43
1:E:100:LEU:HD23	1:E:100:LEU:HA	1.87	0.43
1:A:89:VAL:CG2	1:A:195:VAL:HG11	2.48	0.43
1:B:26:PHE:CE1	1:B:30:LEU:HD12	2.53	0.43
1:E:50:ILE:HG13	1:E:50:ILE:O	2.19	0.43
1:A:215:LEU:O	1:A:219:VAL:HG23	2.18	0.43
1:B:240:PHE:O	1:B:243:VAL:HG12	2.19	0.43
1:A:159:ARG:NH1	1:B:291:PRO:HD3	2.34	0.43
1:A:49:GLY:HA3	1:B:236:TYR:CD2	2.54	0.42
1:C:60:LEU:HD13	1:E:245:ILE:HG12	2.01	0.42
1:A:42:TYR:O	1:A:45:TYR:HB2	2.18	0.42
1:E:224:THR:O	1:E:227:PRO:HD2	2.18	0.42
1:C:83:ARG:HD2	1:C:83:ARG:O	2.19	0.42
1:D:207:VAL:HG11	1:D:260:GLU:HG2	2.02	0.42
1:A:60:LEU:HD23	1:A:222:PHE:CD2	2.54	0.42
1:E:86:TRP:HB3	1:E:275:MET:SD	2.60	0.42
1:A:138:LEU:HD23	1:A:138:LEU:HA	1.81	0.42
1:C:202:LEU:HD23	1:C:202:LEU:HA	1.90	0.42
1:C:215:LEU:HA	1:C:215:LEU:HD23	1.74	0.42
1:A:23:LYS:NZ	1:A:23:LYS:HB3	2.35	0.42
1:E:265:THR:OG1	1:E:266:ALA:N	2.53	0.42
1:B:27:ARG:HG2	1:B:220:TYR:CZ	2.55	0.42
1:D:230:LEU:HD23	1:D:230:LEU:HA	1.89	0.42
1:B:271:PRO:O	1:B:275:MET:HG3	2.20	0.42
1:C:83:ARG:HB2	1:C:270:LEU:HD21	2.02	0.42
1:D:89:VAL:HG23	1:D:195:VAL:HG11	2.02	0.42
1:E:25:ILE:O	1:E:29:LEU:HG	2.19	0.42
1:A:146:ARG:NH2	1:A:149:GLU:OE2	2.53	0.41
1:B:28:LEU:HG	1:B:243:VAL:CG2	2.50	0.41
1:E:42:TYR:O	1:E:45:TYR:HB2	2.20	0.41
1:E:147:VAL:HG12	1:E:151:LEU:HD12	2.02	0.41
1:B:29:LEU:O	1:B:32:VAL:N	2.52	0.41
1:B:126:LYS:HE3	1:B:126:LYS:HB3	1.83	0.41
1:A:179:ASP:HB3	1:B:180:ILE:HG21	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:90:LEU:O	1:E:94:ARG:HG3	2.20	0.41
1:A:238:THR:N	1:A:239:PRO:CD	2.83	0.41
1:C:146:ARG:HA	1:C:146:ARG:HD3	1.92	0.41
1:E:29:LEU:O	1:E:33:LEU:HG	2.21	0.41
1:E:235:HIS:O	1:E:238:THR:HG23	2.20	0.41
1:C:51:HIS:O	1:C:52:LEU:HD23	2.21	0.41
1:C:139:ARG:HE	1:C:144:GLU:CD	2.28	0.41
1:E:130:ARG:NH1	1:E:130:ARG:HG3	2.36	0.41
1:A:60:LEU:HD23	1:A:222:PHE:HD2	1.86	0.41
1:A:226:LEU:N	1:A:227:PRO:HD2	2.36	0.41
1:B:270:LEU:HD22	1:B:275:MET:HE1	2.03	0.41
1:D:68:LEU:HD23	1:D:215:LEU:HD13	2.03	0.41
1:A:50:ILE:HG13	1:A:50:ILE:O	2.21	0.41
1:B:71:ARG:HH11	1:B:71:ARG:HG2	1.86	0.41
1:D:127:HIS:ND1	1:D:132:THR:OG1	2.53	0.41
1:D:286:MET:HE2	1:D:286:MET:HB3	1.86	0.41
1:E:57:PHE:HZ	1:E:225:LEU:HD23	1.84	0.41
1:E:207:VAL:N	1:E:208:PRO:CD	2.84	0.41
1:E:286:MET:HE2	1:E:286:MET:HB3	1.92	0.40
1:C:127:HIS:CD2	1:C:134:PRO:HA	2.56	0.40
1:C:226:LEU:N	1:C:227:PRO:HD2	2.36	0.40
1:D:30:LEU:O	1:D:34:MET:HG2	2.21	0.40
1:D:238:THR:HG23	1:D:238:THR:H	1.63	0.40
1:A:24:ILE:CG1	1:A:27:ARG:HB2	2.51	0.40
1:C:25:ILE:O	1:C:29:LEU:HD23	2.21	0.40
1:B:115:VAL:O	1:B:119:VAL:HG23	2.21	0.40
1:B:243:VAL:HA	1:B:246:SER:OG	2.21	0.40
1:D:159:ARG:HD3	1:D:162:LEU:HD23	2.04	0.40
1:D:161:LEU:HD23	1:D:161:LEU:HA	1.75	0.40
1:D:207:VAL:HG11	1:D:260:GLU:CD	2.46	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	267/297 (90%)	253 (95%)	14 (5%)	0	100	100
1	B	268/297 (90%)	250 (93%)	18 (7%)	0	100	100
1	C	267/297 (90%)	251 (94%)	16 (6%)	0	100	100
1	D	266/297 (90%)	255 (96%)	11 (4%)	0	100	100
1	E	268/297 (90%)	260 (97%)	8 (3%)	0	100	100
All	All	1336/1485 (90%)	1269 (95%)	67 (5%)	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	A	234/260 (90%)	234 (100%)	0	100	100
1	B	235/260 (90%)	235 (100%)	0	100	100
1	C	231/260 (89%)	228 (99%)	3 (1%)	61	83
1	D	230/260 (88%)	230 (100%)	0	100	100
1	E	235/260 (90%)	233 (99%)	2 (1%)	70	87
All	All	1165/1300 (90%)	1160 (100%)	5 (0%)	84	93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	211	TYR
1	C	212	THR
1	C	213	LEU
1	E	44	TRP
1	E	237	MET

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	216	GLN
1	B	99	GLN
1	B	102	ASN
1	B	216	GLN
1	B	268	ASN
1	B	289	GLN
1	C	72	ASN
1	C	289	GLN
1	D	102	ASN
1	E	158	ASN
1	E	268	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 28 ligands modelled in this entry, 27 are monoatomic - leaving 1 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	$\# Z > 2$
4	ACY	C	304	-	3,3,3	3.14	1 (33%)	3,3,3	1.88	1 (33%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	304	ACY	CH3-C	4.99	1.68	1.49

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	304	ACY	O-C-CH3	-2.53	112.14	122.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	304	ACY	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	269/297 (90%)	0.69	45 (16%) 4 3	45, 70, 133, 171	0
1	B	270/297 (90%)	0.63	38 (14%) 6 5	48, 79, 130, 146	0
1	C	269/297 (90%)	0.58	32 (11%) 9 7	50, 73, 137, 176	0
1	D	268/297 (90%)	0.70	33 (12%) 8 6	53, 83, 120, 144	0
1	E	270/297 (90%)	0.65	45 (16%) 4 3	48, 70, 129, 165	0
All	All	1346/1485 (90%)	0.65	193 (14%) 6 5	45, 75, 130, 176	0

All (193) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	204	THR	9.3
1	C	47	GLN	7.7
1	D	145	GLU	7.3
1	B	214	ILE	6.3
1	C	213	LEU	5.9
1	B	211	TYR	5.8
1	C	208	PRO	5.7
1	D	265	THR	5.7
1	B	39	ILE	5.3
1	D	267	ALA	5.2
1	E	24	ILE	5.1
1	E	58	SER	5.1
1	A	204	THR	5.0
1	A	29	LEU	5.0
1	D	264	GLY	5.0
1	B	207	VAL	4.9
1	A	28	LEU	4.7
1	D	268	ASN	4.7
1	D	204	THR	4.7
1	C	291	PRO	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	220	TYR	4.6
1	C	214	ILE	4.6
1	E	44	TRP	4.5
1	E	206	PRO	4.5
1	E	265	THR	4.5
1	C	51	HIS	4.4
1	E	208	PRO	4.4
1	E	209	PHE	4.4
1	A	51	HIS	4.3
1	B	210	ALA	4.3
1	D	130	ARG	4.3
1	C	209	PHE	4.3
1	D	129	LEU	4.3
1	A	26	PHE	4.3
1	B	44	TRP	4.2
1	C	207	VAL	4.2
1	A	175	GLY	4.2
1	D	207	VAL	4.1
1	E	57	PHE	4.0
1	C	23	LYS	4.0
1	C	206	PRO	4.0
1	A	21	LEU	4.0
1	B	51	HIS	4.0
1	D	131	LYS	4.0
1	D	205	THR	3.9
1	A	145	GLU	3.9
1	A	220	TYR	3.8
1	C	48	LEU	3.8
1	C	210	ALA	3.8
1	D	209	PHE	3.8
1	A	42	TYR	3.7
1	E	266	ALA	3.7
1	B	287	THR	3.7
1	E	263	PHE	3.7
1	B	204	THR	3.6
1	E	235	HIS	3.6
1	C	26	PHE	3.6
1	C	145	GLU	3.5
1	B	23	LYS	3.5
1	B	236	TYR	3.5
1	A	37	ILE	3.4
1	E	205	THR	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	77	SER	3.4
1	A	46	GLU	3.4
1	A	143	PRO	3.4
1	E	142	LEU	3.4
1	E	264	GLY	3.4
1	C	44	TRP	3.3
1	A	27	ARG	3.3
1	A	144	GLU	3.3
1	B	209	PHE	3.3
1	E	26	PHE	3.3
1	C	204	THR	3.3
1	C	24	ILE	3.3
1	D	24	ILE	3.3
1	D	154	SER	3.2
1	A	236	TYR	3.2
1	C	184	LEU	3.2
1	E	45	TYR	3.2
1	A	50	ILE	3.1
1	E	207	VAL	3.1
1	E	267	ALA	3.1
1	B	216	GLN	3.1
1	A	212	THR	3.1
1	A	25	ILE	3.1
1	D	266	ALA	3.1
1	A	23	LYS	3.1
1	A	203	ALA	3.1
1	A	83	ARG	3.1
1	B	32	VAL	3.1
1	A	43	GLN	3.0
1	E	50	ILE	3.0
1	D	44	TRP	3.0
1	D	269	ASP	3.0
1	B	54	VAL	3.0
1	D	23	LYS	3.0
1	E	212	THR	3.0
1	B	49	GLY	2.9
1	A	58	SER	2.9
1	A	24	ILE	2.9
1	A	34	MET	2.9
1	E	83	ARG	2.9
1	A	44	TRP	2.8
1	B	26	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	22	SER	2.8
1	A	190	ASP	2.8
1	C	211	TYR	2.8
1	C	50	ILE	2.8
1	A	52	LEU	2.8
1	E	48	LEU	2.8
1	A	84	ASN	2.8
1	A	224	THR	2.8
1	B	286	MET	2.8
1	A	40	ILE	2.7
1	D	235	HIS	2.7
1	C	205	THR	2.7
1	D	144	GLU	2.7
1	A	45	TYR	2.7
1	B	43	GLN	2.7
1	E	211	TYR	2.6
1	E	25	ILE	2.6
1	C	46	GLU	2.6
1	D	26	PHE	2.6
1	D	211	TYR	2.6
1	C	106	ALA	2.6
1	B	205	THR	2.6
1	D	236	TYR	2.6
1	E	49	GLY	2.6
1	D	261	ASP	2.5
1	E	213	LEU	2.5
1	D	200	GLU	2.5
1	C	49	GLY	2.5
1	E	47	GLN	2.5
1	E	269	ASP	2.5
1	E	215	LEU	2.5
1	E	268	ASN	2.5
1	A	30	LEU	2.5
1	A	146	ARG	2.5
1	E	293	PRO	2.5
1	B	231	VAL	2.5
1	A	48	LEU	2.4
1	E	59	LEU	2.4
1	A	39	ILE	2.4
1	B	25	ILE	2.4
1	E	270	LEU	2.4
1	B	22	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	E	210	ALA	2.4
1	D	233	ASP	2.4
1	B	291	PRO	2.4
1	C	77	SER	2.3
1	B	27	ARG	2.3
1	D	28	LEU	2.3
1	A	237	MET	2.3
1	B	35	SER	2.3
1	B	208	PRO	2.3
1	E	236	TYR	2.3
1	A	173	GLU	2.3
1	C	288	GLY	2.3
1	B	213	LEU	2.3
1	B	190	ASP	2.3
1	C	217	ARG	2.3
1	E	54	VAL	2.3
1	E	143	PRO	2.2
1	E	224	THR	2.2
1	C	45	TYR	2.2
1	B	234	LEU	2.2
1	E	202	LEU	2.2
1	A	228	PHE	2.2
1	A	55	ALA	2.2
1	D	139	ARG	2.2
1	C	268	ASN	2.2
1	B	40	ILE	2.2
1	C	166	ASN	2.1
1	B	42	TYR	2.1
1	A	242	SER	2.1
1	A	243	VAL	2.1
1	B	241	VAL	2.1
1	B	212	THR	2.1
1	E	37	ILE	2.1
1	E	27	ARG	2.1
1	A	53	THR	2.1
1	D	76	TYR	2.1
1	D	263	PHE	2.1
1	B	217	ARG	2.1
1	E	146	ARG	2.1
1	C	109	ASP	2.0
1	E	217	ARG	2.0
1	C	181	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	239	PRO	2.0
1	B	206	PRO	2.0
1	D	27	ARG	2.0
1	D	148	THR	2.0
1	B	215	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ACY	C	304	4/4	0.70	0.31	57,58,66,74	0
2	ZN	B	302	1/1	0.91	0.12	153,153,153,153	0
2	ZN	D	302	1/1	0.92	0.11	110,110,110,110	0
2	ZN	D	303	1/1	0.96	0.07	139,139,139,139	0
3	CL	C	307	1/1	0.96	0.08	86,86,86,86	0
3	CL	D	305	1/1	0.96	0.08	92,92,92,92	0
2	ZN	B	306	1/1	0.96	0.16	154,154,154,154	0
3	CL	A	304	1/1	0.97	0.08	74,74,74,74	0
3	CL	C	305	1/1	0.98	0.06	67,67,67,67	0
3	CL	C	306	1/1	0.98	0.10	76,76,76,76	0
2	ZN	A	303	1/1	0.98	0.05	87,87,87,87	0
3	CL	D	304	1/1	0.98	0.06	75,75,75,75	0
2	ZN	B	304	1/1	0.98	0.07	93,93,93,93	0
3	CL	D	306	1/1	0.98	0.06	97,97,97,97	0
3	CL	E	303	1/1	0.98	0.06	74,74,74,74	0
3	CL	B	305	1/1	0.98	0.08	77,77,77,77	0
2	ZN	E	302	1/1	0.99	0.04	65,65,65,65	0
2	ZN	C	302	1/1	0.99	0.02	82,82,82,82	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	A	305	1/1	0.99	0.05	90,90,90,90	0
3	CL	E	304	1/1	0.99	0.07	79,79,79,79	0
2	ZN	C	303	1/1	0.99	0.04	92,92,92,92	0
2	ZN	E	301	1/1	1.00	0.03	62,62,62,62	0
2	ZN	C	301	1/1	1.00	0.02	61,61,61,61	0
2	ZN	A	302	1/1	1.00	0.04	78,78,78,78	0
2	ZN	B	303	1/1	1.00	0.02	79,79,79,79	0
2	ZN	D	301	1/1	1.00	0.02	90,90,90,90	0
2	ZN	A	301	1/1	1.00	0.02	72,72,72,72	0
2	ZN	B	301	1/1	1.00	0.02	76,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.