



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:15 PM UTC

PDB ID : 4ZR6 / pdb_00004zr6
Title : Lymnaea Stagnalis Acetylcholine Binding Protein in Complex with 3-[(4E)-4-[(3-methylimidazol-4-yl)methylene]-2,3-dihydropyrrol-5-yl]pyridine
Authors : Bobango, J.; Bryant, D.; Denton, T.T.; Talley, T.T.
Deposited on : 2015-05-11
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

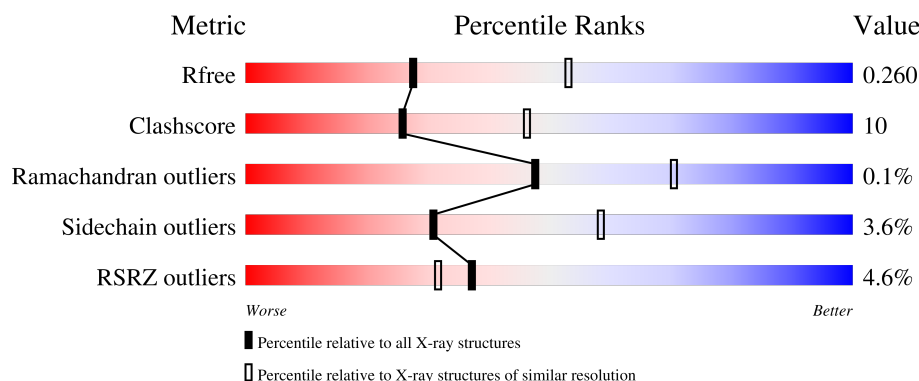
1 Overall quality at a glance

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	218	6% 72% 17% 11%
1	B	218	5% 74% 14% 12%
1	C	218	0% 70% 19% • 9%
1	D	218	3% 73% 17% • 9%
1	E	218	5% 70% 19% • 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7363 atoms, of which 70 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 Acetylcholine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1415	902	233	275	5			
1	B	192	Total	C	N	O	S	0	0	0
			1383	886	222	270	5			
1	C	198	Total	C	N	O	S	0	0	0
			1478	938	245	290	5			
1	D	199	Total	C	N	O	S	0	0	0
			1462	928	239	290	5			
1	E	199	Total	C	N	O	S	0	0	0
			1458	923	239	291	5			

There are 40 discrepancies between the modelled and reference sequences:

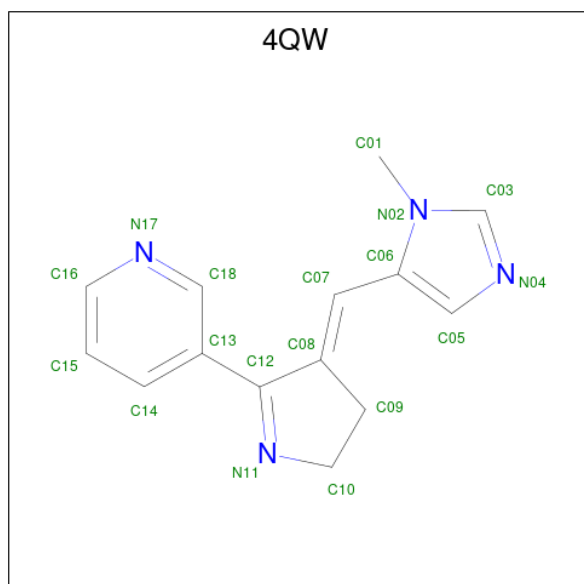
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	ASP	-	expression tag	UNP P58154
A	-6	TYR	-	expression tag	UNP P58154
A	-5	LYS	-	expression tag	UNP P58154
A	-4	ASP	-	expression tag	UNP P58154
A	-3	ASP	-	expression tag	UNP P58154
A	-2	ASP	-	expression tag	UNP P58154
A	-1	ASP	-	expression tag	UNP P58154
A	0	LYS	-	expression tag	UNP P58154
B	-7	ASP	-	expression tag	UNP P58154
B	-6	TYR	-	expression tag	UNP P58154
B	-5	LYS	-	expression tag	UNP P58154
B	-4	ASP	-	expression tag	UNP P58154
B	-3	ASP	-	expression tag	UNP P58154
B	-2	ASP	-	expression tag	UNP P58154
B	-1	ASP	-	expression tag	UNP P58154
B	0	LYS	-	expression tag	UNP P58154
C	-7	ASP	-	expression tag	UNP P58154
C	-6	TYR	-	expression tag	UNP P58154
C	-5	LYS	-	expression tag	UNP P58154

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	ASP	-	expression tag	UNP P58154
C	-3	ASP	-	expression tag	UNP P58154
C	-2	ASP	-	expression tag	UNP P58154
C	-1	ASP	-	expression tag	UNP P58154
C	0	LYS	-	expression tag	UNP P58154
D	-7	ASP	-	expression tag	UNP P58154
D	-6	TYR	-	expression tag	UNP P58154
D	-5	LYS	-	expression tag	UNP P58154
D	-4	ASP	-	expression tag	UNP P58154
D	-3	ASP	-	expression tag	UNP P58154
D	-2	ASP	-	expression tag	UNP P58154
D	-1	ASP	-	expression tag	UNP P58154
D	0	LYS	-	expression tag	UNP P58154
E	-7	ASP	-	expression tag	UNP P58154
E	-6	TYR	-	expression tag	UNP P58154
E	-5	LYS	-	expression tag	UNP P58154
E	-4	ASP	-	expression tag	UNP P58154
E	-3	ASP	-	expression tag	UNP P58154
E	-2	ASP	-	expression tag	UNP P58154
E	-1	ASP	-	expression tag	UNP P58154
E	0	LYS	-	expression tag	UNP P58154

- Molecule 2 is 3-{(4E)-4-[(1-methyl-1H-imidazol-5-yl)methylidene]-3,4-dihydro-2H-pyrrol-5-yl}pyridine (CCD ID: 4QW) (formula: C₁₄H₁₄N₄).



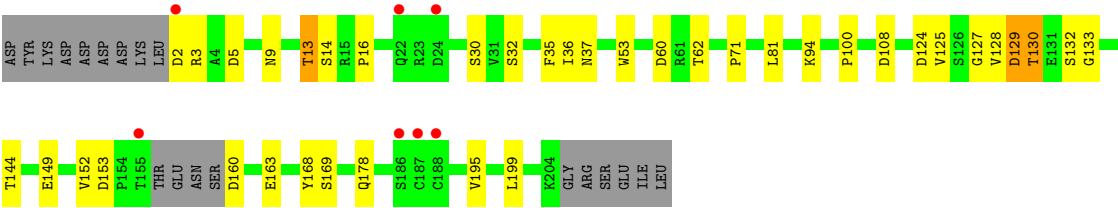
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 32	C 14	H 14	N 4	0	0
2	B	1	Total 32	C 14	H 14	N 4	0	0
2	C	1	Total 32	C 14	H 14	N 4	0	0
2	D	1	Total 32	C 14	H 14	N 4	0	0
2	E	1	Total 32	C 14	H 14	N 4	0	0

- Molecule 3 is water.

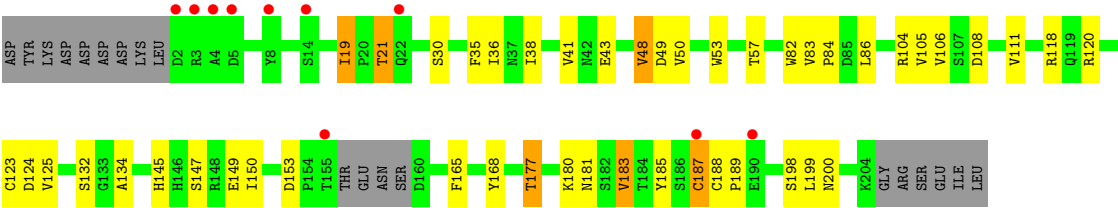
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total 2	O 2	0	0
3	C	3	Total 3	O 3	0	0
3	D	1	Total 1	O 1	0	0
3	E	1	Total 1	O 1	0	0

- Molecule 1: Acetylcholine-binding protein





● Molecule 1: Acetylcholine-binding protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.92Å 78.10Å 120.50Å 90.00° 102.32° 90.00°	Depositor
Resolution (Å)	44.26 – 2.60 44.26 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.26-2.60) 96.8 (44.26-2.60)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 2.51Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.210 , 0.266 (Not available) , 0.260	Depositor DCC
R_{free} test set	2001 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.526	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.5	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	0.92	EDS
Total number of atoms	7363	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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: 4QW

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	0.57	0/1447	0.94	2/1996 (0.1%)
1	B	0.63	0/1415	0.95	2/1952 (0.1%)
1	C	0.66	0/1513	0.95	2/2082 (0.1%)
1	D	0.63	0/1497	0.91	2/2066 (0.1%)
1	E	0.58	0/1491	0.90	1/2054 (0.0%)
All	All	0.62	0/7363	0.93	9/10150 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	188	CYS	CA-C-N	7.26	127.88	120.04
1	C	188	CYS	C-N-CA	7.26	127.88	120.04
1	D	94	LYS	CA-C-N	-6.21	113.11	120.13
1	D	94	LYS	C-N-CA	-6.21	113.11	120.13
1	A	25	ARG	CA-C-N	5.77	125.78	119.90
1	A	25	ARG	C-N-CA	5.77	125.78	119.90
1	E	21	THR	N-CA-C	5.69	118.42	111.82
1	B	23	ARG	CB-CA-C	-5.61	110.09	116.54
1	B	95	PRO	CA-C-O	-5.33	115.52	121.23

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	1415	0	1258	31	0
1	B	1383	0	1224	23	0
1	C	1478	0	1362	31	0
1	D	1462	0	1308	28	0
1	E	1458	0	1307	29	0
2	A	18	14	0	1	0
2	B	18	14	0	1	0
2	C	18	14	0	0	0
2	D	18	14	0	1	0
2	E	18	14	0	0	0
3	B	2	0	0	0	0
3	C	3	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	1	0
All	All	7293	70	6459	131	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 10.

All (131) 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:41:VAL:HG13	1:A:125:VAL:HG11	1.56	0.87
1:A:130:THR:HG22	1:A:132:SER:H	1.39	0.87
1:B:30:SER:HB2	1:B:57:THR:HG22	1.62	0.81
1:E:177:THR:HG22	1:E:198:SER:HB2	1.64	0.80
1:C:152:VAL:HG23	1:C:195:VAL:HG23	1.64	0.78
1:B:130:THR:HG22	1:B:132:SER:H	1.49	0.78
1:A:137:ARG:NH1	1:A:198:SER:OG	2.17	0.77
1:A:130:THR:HG22	1:A:132:SER:N	1.99	0.77
1:B:152:VAL:HG12	1:B:195:VAL:HG23	1.66	0.76
1:D:152:VAL:HG12	1:D:195:VAL:HG23	1.68	0.75
1:D:130:THR:HG22	1:D:132:SER:H	1.53	0.73
1:C:130:THR:HG22	1:C:132:SER:H	1.54	0.73
1:D:130:THR:CG2	1:D:132:SER:H	2.05	0.70
1:C:19:ILE:HG22	1:C:82:TRP:CZ2	2.26	0.69
1:D:130:THR:HG22	1:D:132:SER:N	2.09	0.67
1:C:124:ASP:HB2	1:D:168:TYR:CE1	2.31	0.66
1:B:152:VAL:CG1	1:B:195:VAL:HG23	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:VAL:HG13	1:A:125:VAL:CG1	2.27	0.64
1:A:104:ARG:NH2	1:E:149:GLU:OE2	2.29	0.64
1:C:130:THR:HG22	1:C:132:SER:N	2.13	0.64
1:D:127:GLY:O	1:D:130:THR:HB	1.98	0.64
1:E:123:CYS:HB2	3:E:401:HOH:O	1.97	0.63
1:B:30:SER:HB2	1:B:57:THR:CG2	2.29	0.63
1:C:165:PHE:CD1	1:C:173:ILE:HD11	2.33	0.63
1:C:47:GLU:OE1	1:C:120:ARG:NH1	2.32	0.62
1:A:168:TYR:CE1	1:E:124:ASP:HB2	2.35	0.62
1:D:124:ASP:HB2	1:E:168:TYR:CE1	2.36	0.61
1:A:94:LYS:HE3	1:B:97:VAL:O	2.01	0.61
1:A:48:VAL:HG12	1:A:50:VAL:HG13	1.84	0.60
1:E:145:HIS:HB3	1:E:150:ILE:HG13	1.84	0.59
1:A:41:VAL:HG22	1:A:48:VAL:HG22	1.86	0.58
1:C:35:PHE:CE1	1:C:199:LEU:HD22	2.39	0.58
1:D:35:PHE:CE2	1:D:199:LEU:HD22	2.38	0.58
1:B:124:ASP:HB2	1:C:168:TYR:CE1	2.39	0.57
1:A:144:THR:HG22	2:A:301:4QW:C16	2.34	0.57
1:C:42:ASN:OD1	1:C:45:THR:N	2.28	0.57
1:D:130:THR:HG22	1:D:133:GLY:H	1.70	0.57
1:A:14:SER:O	1:A:16:PRO:HD3	2.04	0.57
1:E:153:ASP:OD1	1:E:180:LYS:NZ	2.36	0.57
1:A:137:ARG:HH12	1:A:198:SER:CB	2.18	0.56
1:E:38:ILE:HG13	1:E:165:PHE:CZ	2.41	0.56
1:C:33:LEU:HD22	1:C:52:PHE:CD2	2.40	0.56
1:E:38:ILE:HG13	1:E:165:PHE:HZ	1.72	0.55
1:B:130:THR:HG22	1:B:131:GLU:N	2.22	0.54
1:E:105:VAL:HG22	1:E:111:VAL:HG22	1.90	0.54
1:D:160:ASP:HB3	1:D:163:GLU:HB2	1.89	0.54
1:C:39:LEU:HD12	1:C:118:ARG:NE	2.23	0.53
1:A:137:ARG:NH1	1:A:198:SER:CB	2.72	0.53
1:B:152:VAL:HG12	1:B:195:VAL:CG2	2.37	0.52
1:A:124:ASP:HB2	1:B:168:TYR:CE1	2.44	0.52
1:D:128:VAL:HG13	1:D:129:ASP:N	2.24	0.52
1:E:145:HIS:CB	1:E:150:ILE:HG13	2.39	0.51
1:B:19:ILE:HD11	1:B:150:ILE:HG13	1.93	0.51
1:B:65:TRP:NE1	1:B:107:SER:HB3	2.26	0.51
1:C:139:LYS:HE3	1:C:194:ASP:OD1	2.11	0.51
1:D:37:ASN:OD1	1:D:169:SER:HB2	2.11	0.51
1:E:19:ILE:HG22	1:E:82:TRP:CZ2	2.46	0.51
1:E:187:CYS:SG	1:E:188:CYS:N	2.84	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:32:SER:HB3	1:C:155:THR:HG23	1.92	0.51
1:A:54:GLN:O	1:A:114:MET:HA	2.12	0.50
1:A:137:ARG:HH11	1:A:137:ARG:HG2	1.77	0.50
1:B:65:TRP:CE2	1:B:107:SER:HB3	2.46	0.50
1:E:35:PHE:CE1	1:E:199:LEU:HD22	2.47	0.50
1:D:30:SER:HA	1:D:153:ASP:O	2.12	0.49
1:C:33:LEU:HD22	1:C:52:PHE:CE2	2.47	0.49
1:B:19:ILE:HG13	1:B:82:TRP:CZ2	2.47	0.49
1:A:77:PRO:HA	1:A:102:LEU:HD23	1.95	0.48
1:E:106:VAL:HB	1:E:108:ASP:OD1	2.13	0.48
1:A:127:GLY:O	1:A:130:THR:HB	2.12	0.48
1:B:106:VAL:HG12	1:B:107:SER:N	2.29	0.48
1:B:176:VAL:HA	1:B:198:SER:O	2.13	0.48
1:D:152:VAL:CG1	1:D:195:VAL:HG23	2.41	0.48
1:E:188:CYS:HB3	1:E:189:PRO:HD2	1.95	0.48
1:D:36:ILE:HD11	1:D:53:TRP:CE2	2.49	0.48
1:B:16:PRO:HA	1:B:82:TRP:N	2.29	0.48
1:A:33:LEU:HD22	1:A:52:PHE:CD1	2.49	0.47
1:E:48:VAL:HG13	1:E:50:VAL:HG13	1.95	0.47
1:D:9:ASN:O	1:D:13:THR:HB	2.14	0.47
1:E:19:ILE:O	1:E:21:THR:N	2.48	0.47
1:A:76:VAL:HG11	1:A:81:LEU:HD11	1.96	0.47
1:D:16:PRO:HB3	1:D:81:LEU:O	2.14	0.47
1:E:183:VAL:HG23	1:E:185:TYR:CE2	2.50	0.47
1:C:89:TYR:CD2	1:C:139:LYS:HG2	2.50	0.47
1:C:139:LYS:HE3	1:C:181:ASN:HD22	1.79	0.47
1:C:83:VAL:HG13	1:C:84:PRO:HD2	1.96	0.47
1:C:48:VAL:O	1:C:120:ARG:HG3	2.14	0.47
1:D:2:ASP:HA	1:D:5:ASP:OD2	2.16	0.46
1:E:48:VAL:O	1:E:120:ARG:HA	2.16	0.46
1:E:180:LYS:HG2	1:E:181:ASN:N	2.30	0.46
1:C:130:THR:CG2	1:C:131:GLU:N	2.78	0.46
1:E:83:VAL:HG13	1:E:84:PRO:HD2	1.99	0.45
1:B:143:TRP:CE2	1:C:99:THR:HG21	2.52	0.45
2:B:301:4QW:C15	1:C:104:ARG:HB2	2.47	0.45
1:E:30:SER:HB2	1:E:57:THR:OG1	2.17	0.45
1:C:2:ASP:OD1	1:C:5:ASP:N	2.40	0.45
1:C:90:ASN:OD1	1:C:90:ASN:N	2.50	0.44
1:A:99:THR:HG23	1:A:116:SER:HB3	1.98	0.44
1:E:134:ALA:O	1:E:200:ASN:ND2	2.46	0.44
1:D:60:ASP:OD1	1:D:62:THR:OG1	2.35	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:36:ILE:HD11	1:E:53:TRP:CD2	2.53	0.44
1:C:85:ASP:HB2	1:D:100:PRO:HG2	2.00	0.44
1:C:125:VAL:O	1:C:125:VAL:HG12	2.17	0.44
1:E:125:VAL:HG12	1:E:125:VAL:O	2.17	0.44
1:B:130:THR:CG2	1:B:131:GLU:N	2.81	0.44
1:A:125:VAL:HG12	1:A:125:VAL:O	2.18	0.43
1:B:127:GLY:O	1:B:130:THR:HB	2.18	0.43
1:D:32:SER:HA	1:D:178:GLN:HE22	1.83	0.43
1:C:188:CYS:HA	1:C:189:PRO:HD3	1.67	0.42
1:D:149:GLU:OE1	1:E:104:ARG:NH2	2.50	0.42
1:B:151:SER:HB2	1:B:193:GLU:OE1	2.19	0.42
1:A:137:ARG:NH1	1:A:137:ARG:HG2	2.34	0.42
1:D:125:VAL:O	1:D:125:VAL:HG12	2.20	0.42
1:A:32:SER:HA	1:A:178:GLN:HE22	1.83	0.42
1:A:57:THR:HA	1:A:111:VAL:O	2.20	0.42
1:A:130:THR:CG2	1:A:132:SER:H	2.20	0.42
1:C:77:PRO:HA	1:C:102:LEU:HD23	2.01	0.42
1:D:144:THR:HG22	2:D:301:4QW:C16	2.50	0.42
1:C:45:THR:O	1:C:46:ASN:HB2	2.20	0.41
1:D:128:VAL:CG1	1:D:129:ASP:N	2.83	0.41
1:D:108:ASP:OD1	1:D:108:ASP:N	2.54	0.41
1:A:143:TRP:CE2	1:B:99:THR:HG21	2.56	0.41
1:C:39:LEU:HD12	1:C:118:ARG:HE	1.86	0.41
1:D:3:ARG:HG3	1:D:71:PRO:HG2	2.01	0.41
1:E:49:ASP:OD2	1:E:118:ARG:NH2	2.31	0.41
1:A:48:VAL:HG12	1:A:50:VAL:CG1	2.49	0.41
1:A:137:ARG:NH1	1:A:198:SER:HB3	2.35	0.41
1:B:137:ARG:HA	1:B:197:VAL:O	2.21	0.41
1:A:55:GLN:HA	1:A:114:MET:HG3	2.03	0.41
1:C:72:ASP:C	1:C:73:GLN:HG3	2.46	0.40
1:D:36:ILE:HD11	1:D:53:TRP:CD2	2.57	0.40
1:E:84:PRO:HB2	1:E:86:LEU:HG	2.03	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	189/218 (87%)	183 (97%)	5 (3%)	1 (0%)	24	46
1	B	184/218 (84%)	181 (98%)	3 (2%)	0	100	100
1	C	194/218 (89%)	188 (97%)	6 (3%)	0	100	100
1	D	195/218 (89%)	193 (99%)	2 (1%)	0	100	100
1	E	195/218 (89%)	189 (97%)	6 (3%)	0	100	100
All	All	957/1090 (88%)	934 (98%)	22 (2%)	1 (0%)	48	70

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	PRO

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	143/204 (70%)	137 (96%)	6 (4%)	26	52
1	B	140/204 (69%)	136 (97%)	4 (3%)	37	65
1	C	160/204 (78%)	156 (98%)	4 (2%)	42	69
1	D	154/204 (76%)	150 (97%)	4 (3%)	40	68
1	E	153/204 (75%)	144 (94%)	9 (6%)	18	38
All	All	750/1020 (74%)	723 (96%)	27 (4%)	31	58

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

Mol	Chain	Res	Type
1	A	18	VAL
1	A	40	GLU
1	A	41	VAL

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Mol	Chain	Res	Type
1	A	56	THR
1	A	106	VAL
1	A	188	CYS
1	B	31	VAL
1	B	41	VAL
1	B	188	CYS
1	B	198	SER
1	C	99	THR
1	C	128	VAL
1	C	152	VAL
1	C	188	CYS
1	D	13	THR
1	D	14	SER
1	D	129	ASP
1	D	130	THR
1	E	19	ILE
1	E	41	VAL
1	E	43	GLU
1	E	48	VAL
1	E	132	SER
1	E	147	SER
1	E	177	THR
1	E	183	VAL
1	E	187	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	C	69	HIS
1	C	181	ASN
1	D	55	GLN
1	D	200	ASN
1	E	101	GLN
1	E	145	HIS

5.3.3 RNA ⓘ

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](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	4QW	A	301	-	17,20,20	1.53	4 (23%)	17,27,27	2.08	5 (29%)
2	4QW	E	301	-	17,20,20	1.52	3 (17%)	17,27,27	2.19	6 (35%)
2	4QW	D	301	-	17,20,20	1.52	1 (5%)	17,27,27	2.23	6 (35%)
2	4QW	B	301	-	17,20,20	1.44	3 (17%)	17,27,27	2.55	5 (29%)
2	4QW	C	301	-	17,20,20	1.49	4 (23%)	17,27,27	2.61	6 (35%)

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	4QW	A	301	-	-	6/8/18/18	0/3/3/3
2	4QW	E	301	-	-	6/8/18/18	0/3/3/3
2	4QW	D	301	-	-	4/8/18/18	0/3/3/3
2	4QW	B	301	-	-	4/8/18/18	0/3/3/3
2	4QW	C	301	-	-	4/8/18/18	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	301	4QW	C07-C06	4.22	1.53	1.44
2	A	301	4QW	C07-C06	3.97	1.52	1.44
2	E	301	4QW	C07-C06	3.78	1.52	1.44
2	C	301	4QW	C07-C06	3.58	1.51	1.44
2	B	301	4QW	C07-C06	3.50	1.51	1.44
2	B	301	4QW	C01-N02	-2.61	1.41	1.46
2	E	301	4QW	C06-N02	-2.60	1.33	1.38
2	B	301	4QW	C06-N02	-2.57	1.33	1.38
2	C	301	4QW	C06-N02	-2.43	1.34	1.38
2	A	301	4QW	C06-N02	-2.31	1.34	1.38
2	E	301	4QW	C05-N04	-2.28	1.33	1.37
2	C	301	4QW	C13-C12	2.18	1.51	1.48
2	A	301	4QW	C05-N04	-2.17	1.33	1.37
2	C	301	4QW	C09-C08	-2.14	1.46	1.50
2	A	301	4QW	C13-C12	2.03	1.51	1.48

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	4QW	N02-C03-N04	-7.34	108.57	112.94
2	C	301	4QW	N02-C03-N04	-5.96	109.39	112.94
2	D	301	4QW	N02-C03-N04	-5.84	109.47	112.94
2	A	301	4QW	N02-C03-N04	-5.71	109.54	112.94
2	C	301	4QW	C09-C08-C07	-5.59	121.27	131.17
2	C	301	4QW	C09-C10-N11	-4.50	102.74	106.50
2	E	301	4QW	N02-C03-N04	-4.36	110.34	112.94
2	E	301	4QW	C09-C10-N11	-4.23	102.97	106.50
2	E	301	4QW	C09-C08-C07	-4.08	123.94	131.17
2	B	301	4QW	C03-N02-C06	3.93	108.24	106.39
2	B	301	4QW	C09-C08-C07	-3.88	124.30	131.17
2	D	301	4QW	C09-C10-N11	-3.60	103.49	106.50
2	D	301	4QW	C09-C08-C07	-3.46	125.05	131.17
2	A	301	4QW	C09-C08-C07	-3.33	125.27	131.17
2	D	301	4QW	C03-N02-C06	3.23	107.91	106.39
2	C	301	4QW	C03-N02-C06	2.71	107.66	106.39
2	A	301	4QW	C03-N02-C06	2.64	107.63	106.39
2	B	301	4QW	C05-N04-C03	2.50	108.51	105.24
2	E	301	4QW	C16-N17-C18	2.47	121.19	116.85
2	E	301	4QW	C13-C18-N17	-2.37	119.97	123.50
2	C	301	4QW	C05-N04-C03	2.32	108.27	105.24
2	B	301	4QW	C16-N17-C18	2.28	120.84	116.85
2	C	301	4QW	C16-N17-C18	2.19	120.68	116.85
2	E	301	4QW	C14-C13-C18	2.16	120.02	117.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	301	4QW	C16-N17-C18	2.16	120.63	116.85
2	A	301	4QW	C16-N17-C18	2.09	120.50	116.85
2	A	301	4QW	C14-C13-C18	2.03	119.88	117.61
2	D	301	4QW	C05-N04-C03	2.02	107.88	105.24

There are no chirality outliers.

All (24) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	4QW	N02-C06-C07-C08
2	A	301	4QW	C05-C06-C07-C08
2	A	301	4QW	C06-C07-C08-C09
2	A	301	4QW	C06-C07-C08-C12
2	B	301	4QW	N02-C06-C07-C08
2	B	301	4QW	C05-C06-C07-C08
2	B	301	4QW	C06-C07-C08-C09
2	B	301	4QW	C06-C07-C08-C12
2	C	301	4QW	N02-C06-C07-C08
2	C	301	4QW	C05-C06-C07-C08
2	C	301	4QW	C06-C07-C08-C09
2	C	301	4QW	C06-C07-C08-C12
2	D	301	4QW	N02-C06-C07-C08
2	D	301	4QW	C05-C06-C07-C08
2	D	301	4QW	C06-C07-C08-C09
2	D	301	4QW	C06-C07-C08-C12
2	E	301	4QW	N02-C06-C07-C08
2	E	301	4QW	C05-C06-C07-C08
2	E	301	4QW	C06-C07-C08-C09
2	E	301	4QW	C06-C07-C08-C12
2	E	301	4QW	C08-C12-C13-C18
2	E	301	4QW	C08-C12-C13-C14
2	A	301	4QW	N11-C12-C13-C18
2	A	301	4QW	N11-C12-C13-C14

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	4QW	1	0
2	D	301	4QW	1	0
2	B	301	4QW	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	195/218 (89%)	0.55	14 (7%) 21 17	32, 50, 79, 99	0
1	B	192/218 (88%)	0.38	11 (5%) 29 24	25, 45, 73, 88	0
1	C	198/218 (90%)	-0.09	3 (1%) 72 68	26, 36, 56, 76	0
1	D	199/218 (91%)	0.16	7 (3%) 47 41	27, 45, 67, 84	0
1	E	199/218 (91%)	0.31	10 (5%) 34 28	31, 46, 79, 90	0
All	All	983/1090 (90%)	0.26	45 (4%) 37 32	25, 45, 74, 99	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	4	ALA	4.7
1	E	155	THR	3.6
1	B	63	LEU	3.6
1	B	3	ARG	3.5
1	E	22	GLN	3.4
1	B	132	SER	3.4
1	B	176	VAL	3.4
1	E	2	ASP	3.2
1	B	24	ASP	3.0
1	C	168	TYR	3.0
1	A	129	ASP	3.0
1	E	190	GLU	2.9
1	C	23	ARG	2.9
1	D	187	CYS	2.9
1	B	200	ASN	2.8
1	D	24	ASP	2.8
1	A	190	GLU	2.8
1	D	186	SER	2.8
1	A	3	ARG	2.8
1	A	59	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	5	ASP	2.7
1	D	22	GLN	2.7
1	B	70	SER	2.7
1	D	155	THR	2.7
1	C	188	CYS	2.6
1	A	31	VAL	2.6
1	D	2	ASP	2.5
1	A	30	SER	2.5
1	A	184	THR	2.5
1	A	86	LEU	2.4
1	E	187	CYS	2.4
1	D	188	CYS	2.4
1	A	8	TYR	2.3
1	A	69	HIS	2.3
1	A	66	ASN	2.3
1	B	5	ASP	2.3
1	A	188	CYS	2.3
1	B	130	THR	2.3
1	E	14	SER	2.3
1	A	67	SER	2.3
1	B	129	ASP	2.2
1	B	131	GLU	2.2
1	A	152	VAL	2.1
1	E	3	ARG	2.1
1	E	8	TYR	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	4QW	A	301	18/18	0.94	0.09	46,55,67,73	0
2	4QW	D	301	18/18	0.94	0.09	42,52,62,68	0
2	4QW	E	301	18/18	0.94	0.09	42,51,64,68	0
2	4QW	B	301	18/18	0.95	0.08	26,35,43,44	0
2	4QW	C	301	18/18	0.95	0.07	28,34,40,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.