



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:51 AM UTC

PDB ID : 8GHJ / pdb_00008ghj
Title : Crystal structure of human AQP2 T125M mutant
Authors : Horsefield, S.; Hagstroemer, C.J.
Deposited on : 2023-03-10
Resolution : 3.90 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

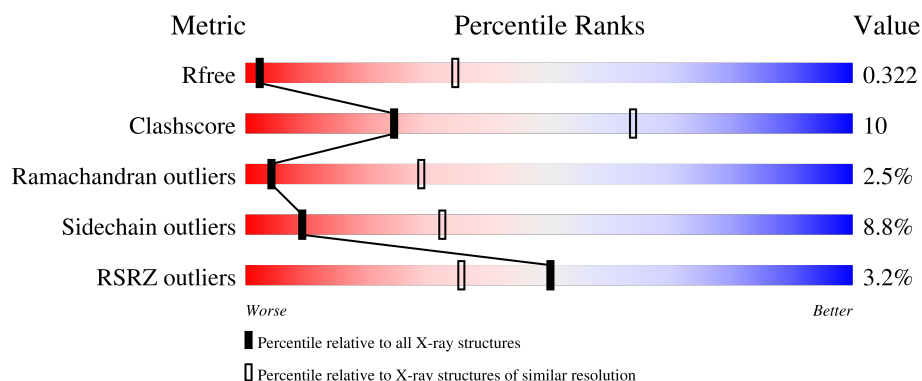
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.90 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	1270 (4.10-3.70)
Clashscore	190562	1034 (4.08-3.72)
Ramachandran outliers	187476	1251 (4.10-3.70)
Sidechain outliers	187428	1243 (4.10-3.70)
RSRZ outliers	180081	1269 (4.10-3.70)

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	241	2% 73% 23% ..
1	B	241	5% 68% 23% 7%
1	C	241	4% 74% 20% ..
1	D	241	% 68% 23% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6844 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 Aquaporin-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1750	1143	294	306	7			
1	B	225	Total	C	N	O	S	0	0	0
			1653	1083	275	288	7			
1	C	236	Total	C	N	O	S	0	0	0
			1737	1136	290	304	7			
1	D	231	Total	C	N	O	S	0	0	0
			1702	1111	284	300	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P41181
A	2	SER	-	expression tag	UNP P41181
A	125	MET	THR	engineered mutation	UNP P41181
B	1	GLY	-	expression tag	UNP P41181
B	2	SER	-	expression tag	UNP P41181
B	125	MET	THR	engineered mutation	UNP P41181
C	1	GLY	-	expression tag	UNP P41181
C	2	SER	-	expression tag	UNP P41181
C	125	MET	THR	engineered mutation	UNP P41181
D	1	GLY	-	expression tag	UNP P41181
D	2	SER	-	expression tag	UNP P41181
D	125	MET	THR	engineered mutation	UNP P41181

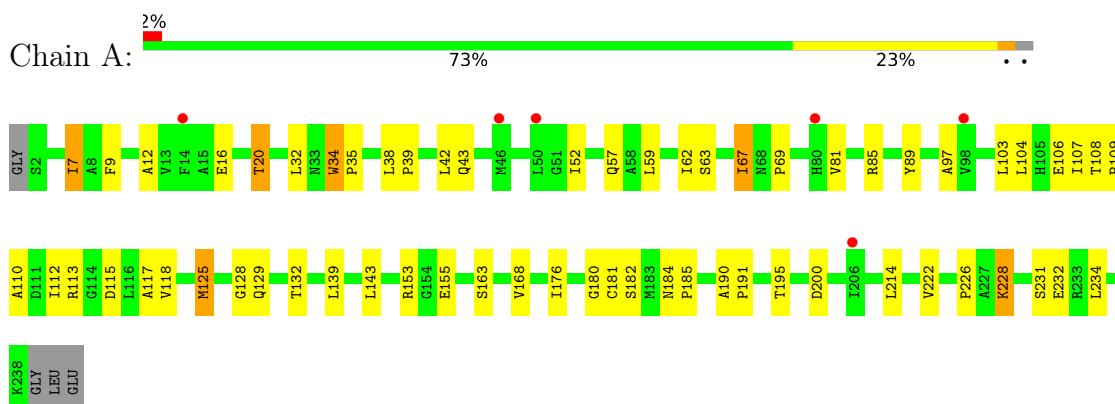
- Molecule 2 is CADMIUM ION (CCD ID: CD) (formula: Cd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cd	1	0
			1	1		
2	D	1	Total	Cd	1	0
			1	1		

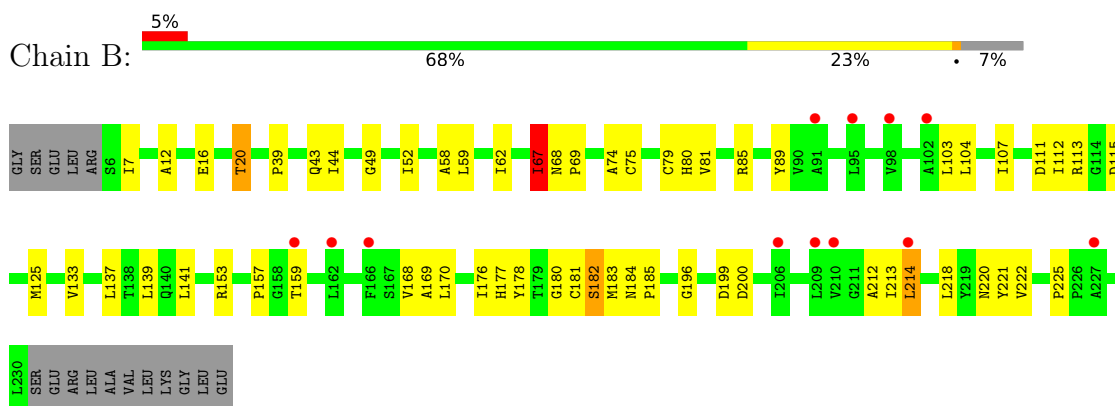
3 Residue-property plots [i](#)

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.

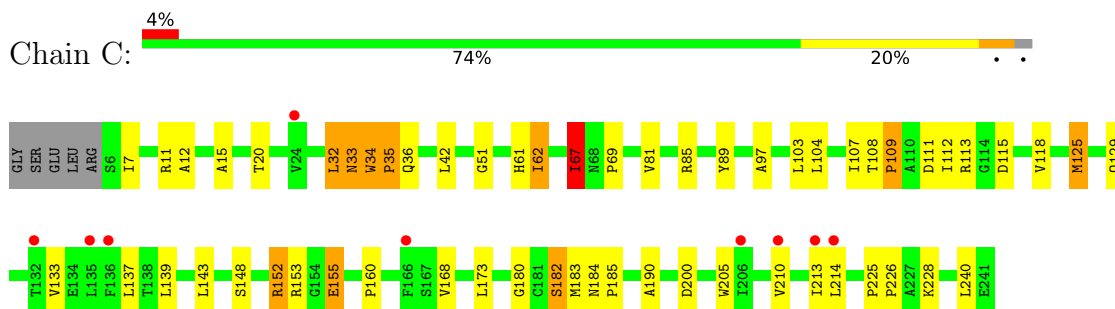
• Molecule 1: Aquaporin-2



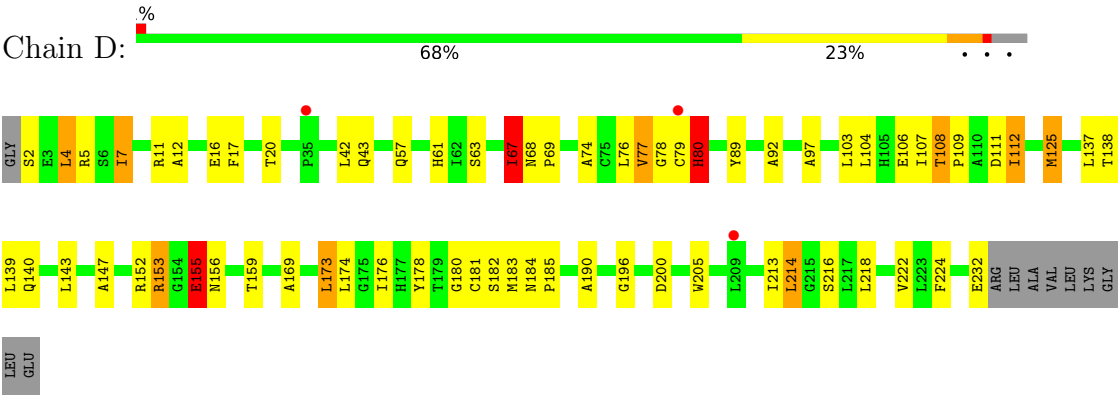
• Molecule 1: Aquaporin-2



• Molecule 1: Aquaporin-2



● Molecule 1: Aquaporin-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	118.31Å 118.31Å 90.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.50 – 3.90 49.50 – 3.90	Depositor EDS
% Data completeness (in resolution range)	99.0 (49.50-3.90) 83.0 (49.50-3.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.28 (at 3.88Å)	Xtriage
Refinement program	PHENIX 1.17_3644	Depositor
R, R_{free}	0.288 , 0.313 0.295 , 0.322	Depositor DCC
R_{free} test set	1157 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	183.4	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 172.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.058 for h,-k,-l	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	6844	wwPDB-VP
Average B, all atoms (Å ²)	214.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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: CD

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	0.14	0/1791	0.52	0/2447
1	B	0.18	0/1694	0.56	0/2318
1	C	0.15	0/1778	0.56	0/2430
1	D	0.17	0/1743	0.52	0/2383
All	All	0.16	0/7006	0.54	0/9578

There are no bond length outliers.

There are no bond angle outliers.

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	1750	0	1794	37	0
1	B	1653	0	1686	40	0
1	C	1737	0	1777	40	0
1	D	1702	0	1730	44	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
All	All	6844	0	6987	137	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 (137) 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:B:177:HIS:HE1	1:C:35:PRO:HD3	1.31	0.96
1:B:69:PRO:HD2	1:B:185:PRO:HD2	1.59	0.82
1:B:177:HIS:CE1	1:C:35:PRO:HD3	2.16	0.81
1:A:129:GLN:NE2	1:D:106:GLU:O	2.19	0.76
1:B:68:ASN:ND2	1:B:183:MET:O	2.19	0.75
1:A:125:MET:HE3	1:D:109:PRO:HD3	1.67	0.75
1:B:104:LEU:HD12	1:B:107:ILE:HD11	1.71	0.72
1:D:69:PRO:HD2	1:D:185:PRO:HD2	1.72	0.71
1:C:32:LEU:HD12	1:C:33:ASN:H	1.55	0.70
1:A:104:LEU:HD12	1:A:107:ILE:HD11	1.74	0.70
1:A:109:PRO:HG2	1:A:112:ILE:HD12	1.74	0.69
1:A:42:LEU:HB2	1:D:43:GLN:HG3	1.75	0.69
1:D:2:SER:HB3	1:D:5:ARG:HG3	1.75	0.69
1:A:34:TRP:HB3	1:A:39:PRO:HD3	1.76	0.68
1:C:67:ILE:O	1:C:184:ASN:ND2	2.27	0.67
1:B:178:TYR:HE1	1:C:107:ILE:HD13	1.59	0.67
1:C:104:LEU:HD12	1:C:107:ILE:HD11	1.78	0.64
1:D:67:ILE:O	1:D:184:ASN:ND2	2.32	0.63
1:B:159:THR:HG22	1:D:159:THR:HG21	1.81	0.63
1:D:78:GLY:O	1:D:80:HIS:N	2.28	0.62
1:B:177:HIS:CE1	1:C:33:ASN:HB3	2.35	0.62
1:D:68:ASN:ND2	1:D:183:MET:O	2.22	0.61
1:C:34:TRP:CD1	1:C:34:TRP:H	2.18	0.61
1:A:222:VAL:O	1:D:11:ARG:HD3	2.01	0.60
1:A:69:PRO:HD2	1:A:185:PRO:HD2	1.83	0.60
1:B:67:ILE:O	1:B:184:ASN:ND2	2.32	0.60
1:A:108:THR:O	1:A:113:ARG:NE	2.36	0.59
1:B:133:VAL:HG21	1:B:178:TYR:HD1	1.68	0.59
1:C:34:TRP:CD1	1:C:34:TRP:N	2.71	0.59
1:D:137:LEU:O	1:D:183:MET:HE3	2.02	0.58
1:A:32:LEU:HD21	1:A:104:LEU:HD21	1.86	0.57
1:B:59:LEU:HA	1:B:62:ILE:HD12	1.86	0.57
1:B:141:LEU:HD22	1:B:183:MET:HE2	1.86	0.57
1:C:182:SER:C	1:C:184:ASN:H	2.12	0.57
1:D:176:ILE:HA	1:D:180:GLY:HA2	1.88	0.56
1:C:97:ALA:HB1	1:C:190:ALA:HB2	1.88	0.55
1:B:176:ILE:HA	1:B:180:GLY:HA2	1.90	0.54
1:C:15:ALA:HB2	1:C:62:ILE:HD11	1.91	0.53
1:D:97:ALA:HB1	1:D:190:ALA:HB2	1.90	0.53
1:B:49:GLY:HA3	1:B:169:ALA:HB2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:ALA:HB1	1:A:190:ALA:HB2	1.91	0.53
1:C:109:PRO:HG2	1:C:112:ILE:HG13	1.90	0.52
1:A:43:GLN:HG3	1:C:42:LEU:HB2	1.89	0.52
1:B:81:VAL:HG23	1:B:85:ARG:HD2	1.91	0.52
1:A:57:GLN:O	1:C:153:ARG:NH1	2.42	0.52
1:D:109:PRO:HG2	1:D:112:ILE:HB	1.91	0.52
1:C:69:PRO:HD2	1:C:185:PRO:HD2	1.91	0.52
1:B:43:GLN:HG3	1:D:42:LEU:HB2	1.91	0.52
1:A:9:PHE:HE2	1:A:85:ARG:HA	1.75	0.51
1:A:34:TRP:HB3	1:A:39:PRO:CD	2.40	0.51
1:A:110:ALA:HA	1:A:113:ARG:HD2	1.94	0.50
1:B:221:TYR:HA	1:B:225:PRO:HB3	1.94	0.50
1:C:148:SER:HB3	1:C:160:PRO:HD2	1.93	0.49
1:B:74:ALA:HB2	1:B:212:ALA:HB1	1.93	0.49
1:D:4:LEU:HA	1:D:7:ILE:HD12	1.94	0.49
1:B:182:SER:C	1:B:184:ASN:H	2.20	0.49
1:C:115:ASP:HB2	1:C:118:VAL:HG22	1.94	0.49
1:D:155:GLU:CD	1:D:156:ASN:H	2.20	0.49
1:B:44:ILE:HB	1:B:176:ILE:HD11	1.95	0.49
1:A:32:LEU:HB3	1:A:117:ALA:HB2	1.95	0.48
1:A:106:GLU:O	1:C:129:GLN:NE2	2.43	0.48
1:A:182:SER:C	1:A:184:ASN:H	2.22	0.48
1:A:12:ALA:HB1	1:A:89:TYR:OH	2.13	0.48
1:C:182:SER:C	1:C:184:ASN:N	2.71	0.48
1:B:16:GLU:O	1:B:20:THR:OG1	2.26	0.48
1:D:125:MET:HE2	1:D:178:TYR:HD1	1.79	0.48
1:A:81:VAL:HG13	1:A:85:ARG:HD2	1.96	0.47
1:A:176:ILE:HA	1:A:180:GLY:HA2	1.97	0.46
1:C:152:ARG:CZ	1:C:226:PRO:HA	2.46	0.46
1:D:153:ARG:O	1:D:153:ARG:HG3	2.15	0.46
1:A:115:ASP:HB2	1:A:118:VAL:HG22	1.97	0.46
1:B:222:VAL:O	1:C:11:ARG:HD3	2.16	0.46
1:D:78:GLY:C	1:D:80:HIS:H	2.18	0.46
1:C:185:PRO:HB3	1:C:205:TRP:CD2	2.51	0.45
1:C:152:ARG:NH2	1:C:225:PRO:O	2.50	0.45
1:A:7:ILE:H	1:A:7:ILE:HG13	1.31	0.45
1:A:34:TRP:HB2	1:A:38:LEU:HD23	1.97	0.45
1:A:155:GLU:OE1	1:D:57:GLN:NE2	2.45	0.45
1:B:220:ASN:O	1:B:225:PRO:HA	2.17	0.45
1:A:231:SER:HA	1:A:234:LEU:HG	1.99	0.45
1:B:177:HIS:ND1	1:C:33:ASN:HB3	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:176:ILE:HA	1:D:180:GLY:CA	2.48	0.44
1:B:58:ALA:HA	1:D:147:ALA:CB	2.47	0.44
1:B:52:ILE:HG13	1:B:168:VAL:HG21	1.98	0.44
1:C:125:MET:HB3	1:C:125:MET:HE3	1.54	0.44
1:D:4:LEU:HD22	1:D:4:LEU:H	1.83	0.44
1:D:137:LEU:HD21	1:D:174:LEU:HD23	2.00	0.44
1:D:152:ARG:HD2	1:D:224:PHE:CE2	2.53	0.44
1:D:169:ALA:O	1:D:173:LEU:HD22	2.18	0.44
1:C:81:VAL:HG13	1:C:85:ARG:HD2	1.99	0.44
1:A:81:VAL:HG13	1:A:85:ARG:HB3	2.00	0.43
1:D:213:ILE:HG13	1:D:214:LEU:N	2.33	0.43
1:C:108:THR:HG21	1:C:113:ARG:HG3	2.00	0.43
1:C:213:ILE:HG13	1:C:214:LEU:N	2.33	0.43
1:D:12:ALA:HB1	1:D:89:TYR:OH	2.18	0.43
1:D:182:SER:C	1:D:184:ASN:H	2.26	0.43
1:A:153:ARG:HG2	1:D:61:HIS:CD2	2.54	0.43
1:D:109:PRO:HB2	1:D:111:ASP:OD1	2.18	0.43
1:C:133:VAL:O	1:C:137:LEU:HG	2.19	0.43
1:D:182:SER:C	1:D:184:ASN:N	2.77	0.43
1:D:104:LEU:O	1:D:108:THR:OG1	2.36	0.43
1:D:185:PRO:HB3	1:D:205:TRP:CD2	2.54	0.43
1:B:170:LEU:HD22	1:C:51:GLY:HA2	2.01	0.42
1:B:177:HIS:CE1	1:C:33:ASN:C	2.97	0.42
1:B:213:ILE:HG13	1:B:214:LEU:N	2.32	0.42
1:C:108:THR:HG21	1:C:113:ARG:CG	2.49	0.42
1:D:77:VAL:HG22	1:D:216:SER:OG	2.19	0.42
1:D:218:LEU:HD12	1:D:222:VAL:HG21	2.01	0.42
1:B:133:VAL:HG21	1:B:178:TYR:CD1	2.51	0.42
1:B:181:CYS:SG	1:B:183:MET:SD	3.17	0.42
1:D:137:LEU:O	1:D:140:GLN:HG3	2.19	0.42
1:A:163:SER:HB3	1:D:57:GLN:HG2	2.00	0.42
1:A:16:GLU:O	1:A:20:THR:OG1	2.35	0.42
1:D:17:PHE:HB2	1:D:92:ALA:HB1	2.02	0.42
1:A:12:ALA:HA	1:A:63:SER:HB3	2.01	0.42
1:B:58:ALA:HA	1:D:147:ALA:HB3	2.02	0.42
1:A:128:GLY:O	1:A:132:THR:OG1	2.30	0.41
1:B:39:PRO:HB2	1:B:44:ILE:HD11	2.02	0.41
1:A:228:LYS:HD3	1:A:228:LYS:HA	1.60	0.41
1:C:12:ALA:HB1	1:C:89:TYR:OH	2.21	0.41
1:D:16:GLU:HG3	1:D:89:TYR:CG	2.55	0.41
1:D:74:ALA:HB1	1:D:216:SER:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:ARG:HD2	1:C:61:HIS:CD2	2.55	0.41
1:B:177:HIS:CE1	1:C:35:PRO:CD	2.96	0.41
1:A:52:ILE:HG13	1:A:168:VAL:HG21	2.02	0.41
1:C:35:PRO:HB2	1:C:36:GLN:H	1.63	0.41
1:B:79:CYS:C	1:B:81:VAL:H	2.29	0.41
1:B:112:ILE:O	1:B:113:ARG:C	2.63	0.41
1:B:153:ARG:NE	1:B:157:PRO:HG3	2.36	0.41
1:C:168:VAL:HA	1:C:183:MET:HE1	2.02	0.41
1:D:12:ALA:HA	1:D:63:SER:HB3	2.02	0.41
1:B:12:ALA:HB1	1:B:89:TYR:OH	2.21	0.41
1:A:190:ALA:HB3	1:A:191:PRO:HD3	2.03	0.40
1:B:218:LEU:O	1:B:222:VAL:HB	2.21	0.40
1:A:59:LEU:HA	1:A:62:ILE:HD12	2.04	0.40
1:C:32:LEU:CD1	1:C:33:ASN:H	2.29	0.40
1:C:210:VAL:O	1:C:214:LEU:HD12	2.20	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	235/241 (98%)	221 (94%)	10 (4%)	4 (2%)	7	35
1	B	223/241 (92%)	196 (88%)	22 (10%)	5 (2%)	5	31
1	C	234/241 (97%)	214 (92%)	12 (5%)	8 (3%)	3	24
1	D	229/241 (95%)	215 (94%)	8 (4%)	6 (3%)	4	28
All	All	921/964 (96%)	846 (92%)	52 (6%)	23 (2%)	4	29

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	PRO
1	C	33	ASN
1	C	35	PRO
1	D	155	GLU
1	B	196	GLY
1	C	155	GLU
1	D	79	CYS
1	D	80	HIS
1	D	196	GLY
1	A	181	CYS
1	A	226	PRO
1	B	125	MET
1	C	180	GLY
1	D	181	CYS
1	C	182	SER
1	B	80	HIS
1	C	109	PRO
1	C	111	ASP
1	B	182	SER
1	B	67	ILE
1	C	67	ILE
1	A	67	ILE
1	D	67	ILE

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	181/183 (99%)	168 (93%)	13 (7%)	13	39
1	B	170/183 (93%)	158 (93%)	12 (7%)	13	40
1	C	179/183 (98%)	163 (91%)	16 (9%)	9	32
1	D	176/183 (96%)	155 (88%)	21 (12%)	5	21
All	All	706/732 (96%)	644 (91%)	62 (9%)	9	32

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ILE
1	A	20	THR
1	A	34	TRP
1	A	67	ILE
1	A	103	LEU
1	A	125	MET
1	A	139	LEU
1	A	143	LEU
1	A	195	THR
1	A	200	ASP
1	A	214	LEU
1	A	228	LYS
1	A	232	GLU
1	B	7	ILE
1	B	20	THR
1	B	67	ILE
1	B	75	CYS
1	B	103	LEU
1	B	111	ASP
1	B	115	ASP
1	B	137	LEU
1	B	139	LEU
1	B	199	ASP
1	B	200	ASP
1	B	214	LEU
1	C	7	ILE
1	C	20	THR
1	C	32	LEU
1	C	34	TRP
1	C	62	ILE
1	C	67	ILE
1	C	103	LEU
1	C	125	MET
1	C	139	LEU
1	C	143	LEU
1	C	152	ARG
1	C	155	GLU
1	C	173	LEU
1	C	200	ASP
1	C	228	LYS
1	C	240	LEU
1	D	4	LEU
1	D	7	ILE

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Mol	Chain	Res	Type
1	D	20	THR
1	D	67	ILE
1	D	76	LEU
1	D	77	VAL
1	D	80	HIS
1	D	103	LEU
1	D	107	ILE
1	D	108	THR
1	D	112	ILE
1	D	125	MET
1	D	138	THR
1	D	139	LEU
1	D	143	LEU
1	D	153	ARG
1	D	155	GLU
1	D	173	LEU
1	D	200	ASP
1	D	214	LEU
1	D	232	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	123	ASN
1	A	177	HIS
1	A	220	ASN
1	B	119	ASN
1	B	177	HIS
1	B	220	ASN
1	C	66	HIS
1	C	105	HIS
1	D	129	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 [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	237/241 (98%)	-0.16	6 (2%) 58 41	166, 208, 257, 268	0
1	B	225/241 (93%)	-0.10	12 (5%) 32 25	167, 214, 246, 278	0
1	C	236/241 (97%)	0.01	9 (3%) 44 32	169, 210, 249, 264	0
1	D	231/241 (95%)	-0.23	3 (1%) 75 54	164, 217, 250, 283	1 (0%)
All	All	929/964 (96%)	-0.12	30 (3%) 50 35	164, 213, 251, 283	1 (0%)

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	210	VAL	8.3
1	C	206	ILE	5.7
1	D	209	LEU	5.5
1	B	210	VAL	5.0
1	C	213	ILE	4.9
1	A	14	PHE	4.4
1	B	102	ALA	4.1
1	C	136	PHE	3.6
1	B	209	LEU	3.4
1	C	214	LEU	3.3
1	B	95	LEU	3.3
1	B	214	LEU	3.1
1	C	24	VAL	3.0
1	B	162	LEU	3.0
1	B	159	THR	3.0
1	D	79	CYS	3.0
1	D	35	PRO	2.8
1	A	98	VAL	2.8
1	C	166	PHE	2.7
1	A	46	MET	2.6
1	B	91	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	132	THR	2.4
1	B	227	ALA	2.4
1	A	206	ILE	2.3
1	C	135	LEU	2.3
1	A	80	HIS	2.3
1	B	206	ILE	2.3
1	B	98	VAL	2.2
1	B	166	PHE	2.2
1	A	50	LEU	2.2

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
2	CD	A	301	1/1	-	-	181,181,181,181	1
2	CD	D	301	1/1	-	-	212,212,212,212	1

6.5 Other polymers [i](#)

There are no such residues in this entry.