



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 19, 2024 – 01:24 AM EDT

PDB ID : 4J2J
Title : Crystal structure of AXH domain complex with Capicua
Authors : Song, J.-J.; Kim, E.
Deposited on : 2013-02-04
Resolution : 2.50 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

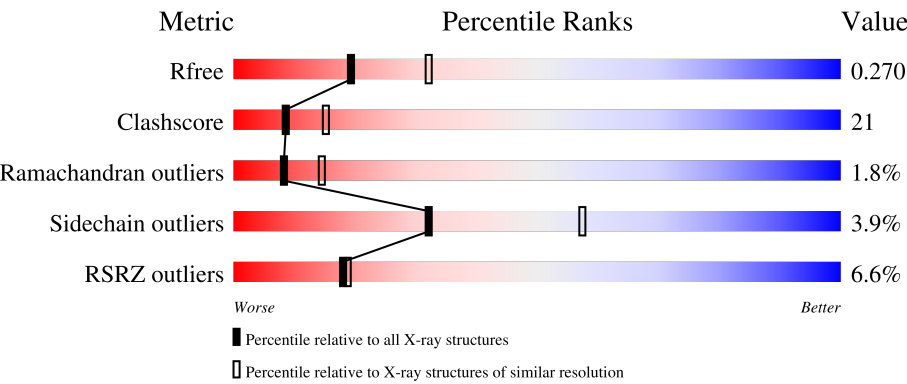
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	129	<div><div>5%</div><div><div></div><div>61%</div><div>31%</div><div>• 6%</div></div></div>
1	B	129	<div><div>13%</div><div><div></div><div>47%</div><div>40%</div><div>5%</div><div>9%</div></div></div>
1	C	129	<div><div>2%</div><div><div></div><div>55%</div><div>31%</div><div>• 11%</div></div></div>
2	D	21	<div><div></div><div><div></div><div>43%</div><div>29%</div><div>29%</div></div></div>
2	E	21	<div><div></div><div><div></div><div>38%</div><div>43%</div><div>19%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	F	21	<div><div></div><div>43%</div><div>24%</div><div>5%</div><div>29%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3161 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 Ataxin-1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	121	Total	C	N	O	S	Se	0	0	0
			937	595	150	185	4	3			
1	B	118	Total	C	N	O	S	Se	0	0	0
			920	585	147	181	4	3			
1	C	115	Total	C	N	O	S	Se	0	0	0
			890	566	143	174	4	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	561	GLU	-	EXPRESSION TAG	UNP P54253
A	562	PHE	-	EXPRESSION TAG	UNP P54253
A	580	MSE	ILE	ENGINEERED MUTATION	UNP P54253
A	605	MSE	ILE	ENGINEERED MUTATION	UNP P54253
B	561	GLU	-	EXPRESSION TAG	UNP P54253
B	562	PHE	-	EXPRESSION TAG	UNP P54253
B	580	MSE	ILE	ENGINEERED MUTATION	UNP P54253
B	605	MSE	ILE	ENGINEERED MUTATION	UNP P54253
C	561	GLU	-	EXPRESSION TAG	UNP P54253
C	562	PHE	-	EXPRESSION TAG	UNP P54253
C	580	MSE	ILE	ENGINEERED MUTATION	UNP P54253
C	605	MSE	ILE	ENGINEERED MUTATION	UNP P54253

- Molecule 2 is a protein called Protein capicua homolog.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	15	Total	C	N	O	0	0	0
			125	87	21	17			
2	E	17	Total	C	N	O	0	0	0
			137	95	23	19			
2	F	15	Total	C	N	O	0	0	0
			124	86	21	17			

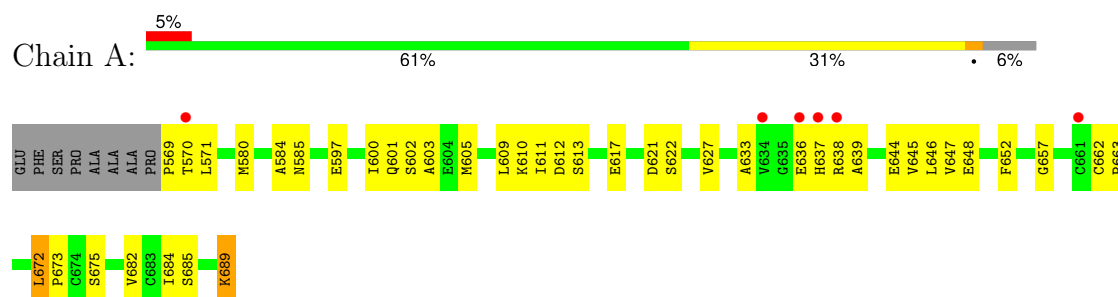
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total 8	O 8	0	0
3	B	5	Total 5	O 5	0	0
3	C	11	Total 11	O 11	0	0
3	E	2	Total 2	O 2	0	0
3	F	2	Total 2	O 2	0	0

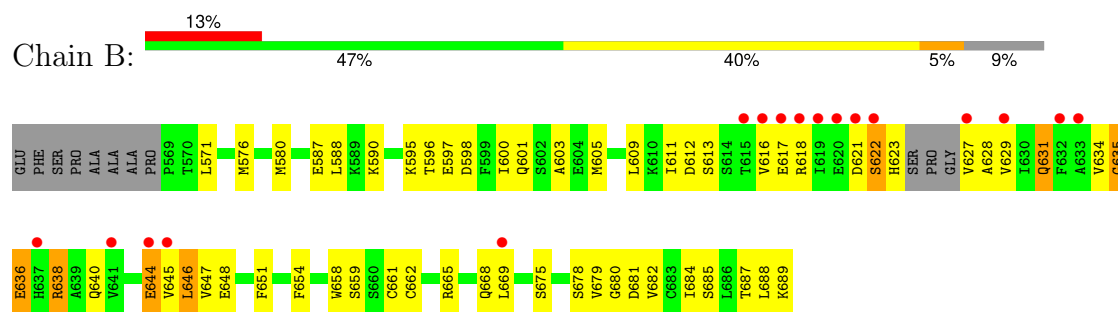
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

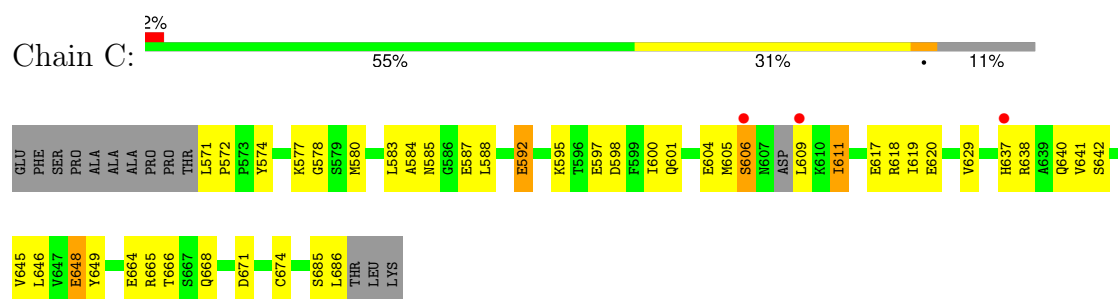
• Molecule 1: Ataxin-1



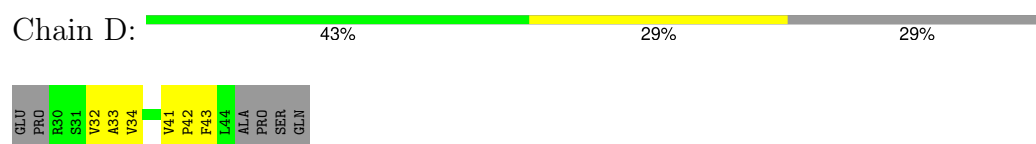
• Molecule 1: Ataxin-1



• Molecule 1: Ataxin-1



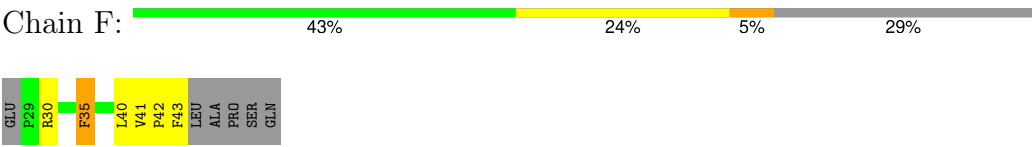
• Molecule 2: Protein capicua homolog



● Molecule 2: Protein capicua homolog



● Molecule 2: Protein capicua homolog



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	77.72Å 89.10Å 132.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.90 – 2.50 43.90 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.1 (43.90-2.50) 98.7 (43.90-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.22 (at 2.51Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.223 , 0.270 0.231 , 0.270	Depositor DCC
R_{free} test set	1525 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3161	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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](#)

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	0/953	0.69	0/1283
1	B	0.36	0/934	0.64	0/1255
1	C	0.40	0/904	0.61	0/1216
2	D	0.47	0/131	0.67	0/180
2	E	0.50	0/144	0.71	0/198
2	F	0.38	0/131	0.62	0/180
All	All	0.40	0/3197	0.65	0/4312

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	937	0	925	28	0
1	B	920	0	909	61	0
1	C	890	0	874	39	0
2	D	125	0	125	7	0
2	E	137	0	138	9	0
2	F	124	0	122	8	0
3	A	8	0	0	0	0
3	B	5	0	0	0	0
3	C	11	0	0	1	0
3	E	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	2	0	0	0	0
All	All	3161	0	3093	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 21.

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:B:580:MSE:HE1	1:B:590:LYS:HG2	1.14	1.11
1:B:646:LEU:HB3	1:B:648:GLU:OE2	1.77	0.84
1:B:685:SER:HA	2:E:41:VAL:HG13	1.65	0.78
1:A:647:VAL:HB	1:A:662:CYS:SG	2.25	0.77
1:B:617:GLU:HA	1:B:679:VAL:HG13	1.68	0.76
1:B:662:CYS:SG	1:B:665:ARG:HB2	2.29	0.73
1:B:627:VAL:HG22	1:B:646:LEU:HD23	1.73	0.70
1:C:571:LEU:N	1:C:572:PRO:HD3	2.08	0.69
1:C:685:SER:HA	2:F:41:VAL:HG13	1.74	0.68
1:C:685:SER:HA	2:F:41:VAL:CG1	2.23	0.68
1:B:636:GLU:H	1:B:636:GLU:CD	1.97	0.68
1:B:576:MSE:CE	1:C:629:VAL:HB	2.25	0.67
1:C:666:THR:HG21	1:C:674:CYS:SG	2.36	0.65
1:A:571:LEU:HD11	1:A:648:GLU:HG2	1.78	0.64
1:B:629:VAL:HG22	1:B:644:GLU:HG3	1.80	0.63
1:B:634:VAL:HG11	2:E:44:LEU:HD13	1.80	0.63
1:B:627:VAL:HG12	1:B:628:ALA:N	2.14	0.63
1:B:636:GLU:C	1:B:638:ARG:H	2.02	0.62
1:B:616:VAL:HG13	1:B:680:GLY:H	1.64	0.62
1:B:685:SER:HA	2:E:41:VAL:CG1	2.30	0.61
1:A:571:LEU:HG	1:A:646:LEU:HD12	1.82	0.61
1:B:648:GLU:H	1:B:648:GLU:CD	2.03	0.61
1:B:631:GLN:HE22	1:B:640:GLN:HG3	1.64	0.61
1:B:687:THR:HG22	2:E:44:LEU:HB2	1.84	0.59
1:C:645:VAL:HG21	1:C:649:TYR:CG	2.37	0.59
1:B:609:LEU:O	1:B:689:LYS:HE3	2.02	0.59
1:B:601:GLN:O	1:B:605:MSE:HG2	2.02	0.59
1:C:578:GLY:H	1:C:592:GLU:CD	2.05	0.58
1:B:576:MSE:HE1	1:C:629:VAL:HB	1.85	0.58
1:C:585:ASN:OD1	1:C:587:GLU:HG2	2.04	0.57
1:A:684:ILE:O	2:D:41:VAL:HG13	2.05	0.57
1:B:580:MSE:CE	1:B:590:LYS:HG2	2.10	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:570:THR:HA	1:A:646:LEU:HD11	1.88	0.56
1:B:628:ALA:O	1:B:644:GLU:HG2	2.06	0.56
1:B:597:GLU:H	1:B:597:GLU:CD	2.09	0.56
1:C:664:GLU:O	1:C:668:GLN:HG3	2.07	0.55
1:C:600:ILE:HA	1:C:611:ILE:HD12	1.88	0.55
1:B:616:VAL:HG13	1:B:680:GLY:N	2.21	0.54
1:B:622:SER:CB	1:B:629:VAL:HG23	2.37	0.54
1:B:654:PHE:CZ	1:B:684:ILE:HG12	2.43	0.54
1:A:610:LYS:HD2	1:A:689:LYS:HB2	1.89	0.54
1:A:663:PRO:HG3	1:A:675:SER:O	2.08	0.54
1:C:618:ARG:HG2	3:C:704:HOH:O	2.09	0.53
1:B:622:SER:HB3	1:B:629:VAL:HG23	1.89	0.52
1:B:645:VAL:HG22	1:B:646:LEU:N	2.25	0.52
1:C:617:GLU:OE1	1:C:640:GLN:OE1	2.26	0.52
1:A:569:PRO:HG3	1:A:644:GLU:O	2.10	0.52
1:A:603:ALA:HB2	1:A:611:ILE:HG13	1.93	0.51
1:B:618:ARG:HB3	1:B:631:GLN:HB2	1.92	0.51
2:F:35:PHE:CD1	2:F:35:PHE:N	2.79	0.51
1:B:616:VAL:HG11	1:B:678:SER:O	2.11	0.50
1:A:617:GLU:CD	1:A:633:ALA:HB2	2.32	0.50
1:B:576:MSE:HE3	1:C:620:GLU:CB	2.41	0.50
1:A:585:ASN:HB3	1:A:605:MSE:HE3	1.92	0.50
2:F:30:ARG:HH11	2:F:30:ARG:HB3	1.76	0.50
1:C:574:TYR:CD1	2:F:40:LEU:HD23	2.47	0.50
2:D:41:VAL:HG13	2:D:42:PRO:HD2	1.94	0.49
1:B:576:MSE:HE2	1:C:629:VAL:CG1	2.43	0.49
1:B:576:MSE:HE2	1:C:629:VAL:HB	1.94	0.49
1:A:600:ILE:HA	1:A:611:ILE:HD12	1.94	0.49
1:A:580:MSE:HE3	1:B:588:LEU:HD12	1.94	0.49
1:A:637:HIS:O	1:A:638:ARG:HB2	2.13	0.49
1:C:571:LEU:N	1:C:572:PRO:CD	2.75	0.49
1:C:597:GLU:O	1:C:601:GLN:HG3	2.14	0.48
1:C:648:GLU:OE1	1:C:665:ARG:HD3	2.14	0.48
1:A:637:HIS:CD2	1:A:639:ALA:HB2	2.50	0.47
1:C:641:VAL:HG22	1:C:642:SER:N	2.29	0.47
1:C:646:LEU:HB3	1:C:648:GLU:OE2	2.15	0.47
1:B:616:VAL:HG12	1:B:681:ASP:O	2.15	0.47
1:B:627:VAL:CG1	1:B:628:ALA:N	2.77	0.46
1:B:638:ARG:HG3	1:B:638:ARG:HH11	1.80	0.46
1:C:572:PRO:HB3	1:C:574:TYR:CZ	2.51	0.46
1:B:571:LEU:CD1	1:B:669:LEU:HD12	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:659:SER:HA	1:B:675:SER:O	2.16	0.46
2:E:36:PRO:HG2	2:E:39:SER:OG	2.15	0.46
1:A:609:LEU:HD22	2:D:43:PHE:HZ	1.81	0.46
1:A:627:VAL:HG11	1:A:644:GLU:CG	2.47	0.45
1:B:571:LEU:HD11	1:B:669:LEU:HD12	1.97	0.45
2:F:42:PRO:O	2:F:43:PHE:HB2	2.15	0.45
1:B:596:THR:HG22	1:B:600:ILE:HD12	1.99	0.44
1:B:621:ASP:O	1:B:622:SER:HB2	2.18	0.44
1:B:654:PHE:HD1	1:B:682:VAL:O	2.00	0.44
1:A:684:ILE:HG22	2:D:41:VAL:HG11	1.99	0.44
1:B:665:ARG:O	1:B:668:GLN:HB3	2.17	0.44
1:B:612:ASP:O	1:B:613:SER:C	2.56	0.44
1:A:613:SER:HB3	1:A:682:VAL:CG1	2.47	0.44
1:A:672:LEU:O	1:A:673:PRO:C	2.53	0.44
1:C:618:ARG:HG3	1:C:619:ILE:N	2.32	0.44
2:D:32:VAL:HG12	2:D:33:ALA:N	2.33	0.44
1:A:621:ASP:OD1	1:A:622:SER:N	2.48	0.43
1:B:609:LEU:HD23	1:B:688:LEU:HD23	2.00	0.43
1:B:576:MSE:HE3	1:C:620:GLU:HB3	2.00	0.43
2:D:34:VAL:HG12	2:E:30:ARG:HG3	2.00	0.43
1:C:645:VAL:HG22	1:C:649:TYR:HB2	2.01	0.43
1:B:636:GLU:C	1:B:638:ARG:N	2.69	0.43
1:B:645:VAL:HG22	1:B:646:LEU:H	1.84	0.43
1:C:592:GLU:CD	1:C:592:GLU:H	2.20	0.43
2:F:30:ARG:CB	2:F:30:ARG:NH1	2.81	0.43
1:C:637:HIS:O	1:C:638:ARG:HB2	2.19	0.43
1:B:603:ALA:HB2	1:B:611:ILE:CD1	2.48	0.43
1:B:627:VAL:HG12	1:B:628:ALA:H	1.82	0.43
1:C:580:MSE:HE3	1:C:588:LEU:HD13	2.00	0.43
1:C:584:ALA:HB1	1:C:605:MSE:HG2	2.01	0.43
1:B:595:LYS:O	1:B:598:ASP:HB2	2.19	0.42
1:B:576:MSE:CE	1:C:620:GLU:HB3	2.49	0.42
1:C:577:LYS:HG3	1:C:592:GLU:HG2	2.00	0.42
1:B:647:VAL:HG13	1:B:661:CYS:HB2	2.02	0.42
1:C:595:LYS:HB2	1:C:598:ASP:OD2	2.20	0.42
1:C:606:SER:OG	1:C:609:LEU:HB2	2.20	0.42
1:B:634:VAL:CG1	2:E:44:LEU:HD13	2.49	0.41
1:C:645:VAL:HG22	1:C:646:LEU:N	2.35	0.41
1:A:685:SER:HA	2:D:41:VAL:CG1	2.50	0.41
1:B:596:THR:HG22	1:B:600:ILE:CD1	2.50	0.41
1:C:601:GLN:HA	1:C:604:GLU:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:32:VAL:CG1	2:E:33:ALA:N	2.82	0.41
1:B:634:VAL:HG12	1:B:635:GLY:N	2.35	0.41
1:B:587:GLU:HG2	1:B:588:LEU:N	2.35	0.41
1:C:645:VAL:HG21	1:C:649:TYR:CB	2.50	0.41
2:F:41:VAL:HA	2:F:42:PRO:HD3	1.78	0.41
1:A:569:PRO:CG	1:A:645:VAL:HA	2.50	0.41
1:B:576:MSE:HE2	1:C:629:VAL:HG11	2.02	0.41
1:C:645:VAL:CG2	1:C:649:TYR:CB	2.99	0.41
1:A:603:ALA:CB	1:A:611:ILE:HG13	2.51	0.41
1:A:652:PHE:CZ	1:A:657:GLY:HA2	2.56	0.41
1:A:597:GLU:HB2	1:A:601:GLN:NE2	2.36	0.41
1:A:584:ALA:HB3	1:A:602:SER:HA	2.03	0.40
1:A:584:ALA:CB	1:A:605:MSE:HB3	2.51	0.40
1:B:651:PHE:O	1:B:658:TRP:HA	2.20	0.40
1:B:688:LEU:HG	2:E:43:PHE:HZ	1.85	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	119/129 (92%)	109 (92%)	9 (8%)	1 (1%)	19	35
1	B	114/129 (88%)	97 (85%)	13 (11%)	4 (4%)	3	4
1	C	111/129 (86%)	104 (94%)	5 (4%)	2 (2%)	8	14
2	D	13/21 (62%)	12 (92%)	1 (8%)	0	100	100
2	E	15/21 (71%)	13 (87%)	2 (13%)	0	100	100
2	F	13/21 (62%)	12 (92%)	1 (8%)	0	100	100
All	All	385/450 (86%)	347 (90%)	31 (8%)	7 (2%)	8	14

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	583	LEU
1	B	622	SER
1	B	638	ARG
1	A	636	GLU
1	C	606	SER
1	B	635	GLY
1	B	646	LEU

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	109/111 (98%)	106 (97%)	3 (3%)	43	70
1	B	107/111 (96%)	103 (96%)	4 (4%)	34	60
1	C	103/111 (93%)	98 (95%)	5 (5%)	25	47
2	D	14/19 (74%)	14 (100%)	0	100	100
2	E	15/19 (79%)	14 (93%)	1 (7%)	16	31
2	F	14/19 (74%)	13 (93%)	1 (7%)	14	28
All	All	362/390 (93%)	348 (96%)	14 (4%)	32	57

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	612	ASP
1	A	672	LEU
1	A	689	LYS
1	B	623	HIS
1	B	631	GLN
1	B	636	GLU
1	B	644	GLU
1	C	592	GLU
1	C	611	ILE
1	C	648	GLU
1	C	671	ASP
1	C	686	LEU

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Mol	Chain	Res	Type
2	E	34	VAL
2	F	35	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	601	GLN
1	A	640	GLN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	118/129 (91%)	0.17	6 (5%) 28 29	29, 44, 88, 163	0
1	B	115/129 (89%)	0.52	17 (14%) 2 2	32, 55, 120, 150	0
1	C	112/129 (86%)	-0.03	3 (2%) 54 58	30, 52, 93, 124	1 (0%)
2	D	15/21 (71%)	0.05	0 100 100	33, 40, 68, 74	0
2	E	17/21 (80%)	0.03	0 100 100	29, 45, 99, 107	0
2	F	15/21 (71%)	-0.09	0 100 100	48, 53, 78, 80	0
All	All	392/450 (87%)	0.19	26 (6%) 18 19	29, 51, 101, 163	1 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	619	ILE	8.6
1	A	638	ARG	5.5
1	B	618	ARG	3.7
1	A	637	HIS	3.7
1	C	637	HIS	3.7
1	B	621	ASP	3.7
1	B	632	PHE	3.5
1	B	641	VAL	3.5
1	B	622	SER	3.4
1	B	620	GLU	3.1
1	B	616	VAL	2.8
1	A	636	GLU	2.7
1	B	645	VAL	2.6
1	B	633	ALA	2.6
1	A	661	CYS	2.6
1	B	644	GLU	2.6
1	C	609	LEU	2.6
1	B	627	VAL	2.5
1	B	617	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	606	SER	2.4
1	B	615	THR	2.4
1	B	637	HIS	2.4
1	A	634	VAL	2.3
1	B	629	VAL	2.3
1	B	669	LEU	2.3
1	A	570	THR	2.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.