Full wwPDB X-ray Structure Validation Report ⓘ

Jun 24, 2024 – 08:54 AM EDT

PDB ID	:	6N64
Title	:	Crystal structure of mouse SMCHD1 hinge domain
Authors	:	Birkinshaw, R.W.; Chen, K.; Czabotar, P.E.; Blewitt, M.E.; Murphy, J.M.
Deposited on	:	2018-11-25
Resolution	:	3.30 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:

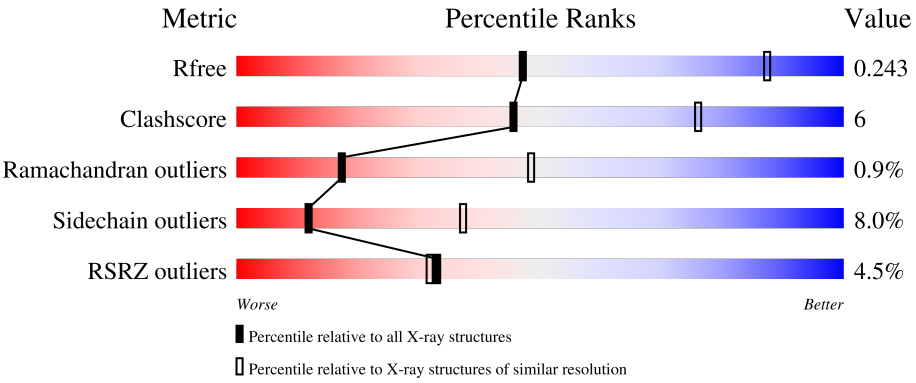
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 <=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	222	6% 64% 12% • 23%
1	B	222	4% 60% 16% • 22%
1	C	222	62% 15% • 21%
1	D	222	% 66% 12% 21%
1	E	222	5% 64% 12% • 22%

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Mol	Chain	Length	Quality of chain
1	F	222	4% 58% 17% 23%
2	G	21	57% 43%
2	H	21	100%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 8412 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 Structural maintenance of chromosomes flexible hinge domain-containing protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	Se	0	0	0
			1353	855	245	246	3	4			
1	B	173	Total	C	N	O	S	Se	0	0	0
			1375	870	248	250	3	4			
1	C	175	Total	C	N	O	S	Se	0	0	0
			1389	879	250	252	4	4			
1	D	175	Total	C	N	O	S	Se	0	0	0
			1389	879	250	252	4	4			
1	E	173	Total	C	N	O	S	Se	0	0	0
			1373	867	248	251	3	4			
1	F	172	Total	C	N	O	S	Se	0	0	0
			1368	866	247	248	3	4			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1678	GLY	-	expression tag	UNP Q6P5D8
A	1679	ALA	-	expression tag	UNP Q6P5D8
A	1680	MSE	-	expression tag	UNP Q6P5D8
A	1681	GLY	-	expression tag	UNP Q6P5D8
A	1682	SER	-	expression tag	UNP Q6P5D8
B	1678	GLY	-	expression tag	UNP Q6P5D8
B	1679	ALA	-	expression tag	UNP Q6P5D8
B	1680	MSE	-	expression tag	UNP Q6P5D8
B	1681	GLY	-	expression tag	UNP Q6P5D8
B	1682	SER	-	expression tag	UNP Q6P5D8
C	1678	GLY	-	expression tag	UNP Q6P5D8
C	1679	ALA	-	expression tag	UNP Q6P5D8
C	1680	MSE	-	expression tag	UNP Q6P5D8
C	1681	GLY	-	expression tag	UNP Q6P5D8
C	1682	SER	-	expression tag	UNP Q6P5D8
D	1678	GLY	-	expression tag	UNP Q6P5D8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1679	ALA	-	expression tag	UNP Q6P5D8
D	1680	MSE	-	expression tag	UNP Q6P5D8
D	1681	GLY	-	expression tag	UNP Q6P5D8
D	1682	SER	-	expression tag	UNP Q6P5D8
E	1678	GLY	-	expression tag	UNP Q6P5D8
E	1679	ALA	-	expression tag	UNP Q6P5D8
E	1680	MSE	-	expression tag	UNP Q6P5D8
E	1681	GLY	-	expression tag	UNP Q6P5D8
E	1682	SER	-	expression tag	UNP Q6P5D8
F	1678	GLY	-	expression tag	UNP Q6P5D8
F	1679	ALA	-	expression tag	UNP Q6P5D8
F	1680	MSE	-	expression tag	UNP Q6P5D8
F	1681	GLY	-	expression tag	UNP Q6P5D8
F	1682	SER	-	expression tag	UNP Q6P5D8

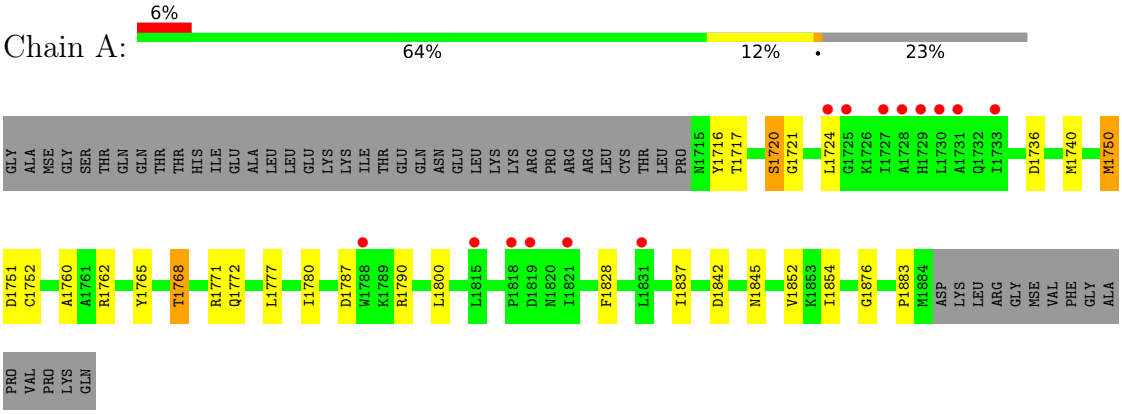
- Molecule 2 is a protein called Uncharacterized peptide from Structural maintenance of chromosomes flexible hinge domain-containing protein 1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	12	Total	C	N	O	0	0	0
			60	36	12	12			
2	H	21	Total	C	N	O	0	0	0
			105	63	21	21			

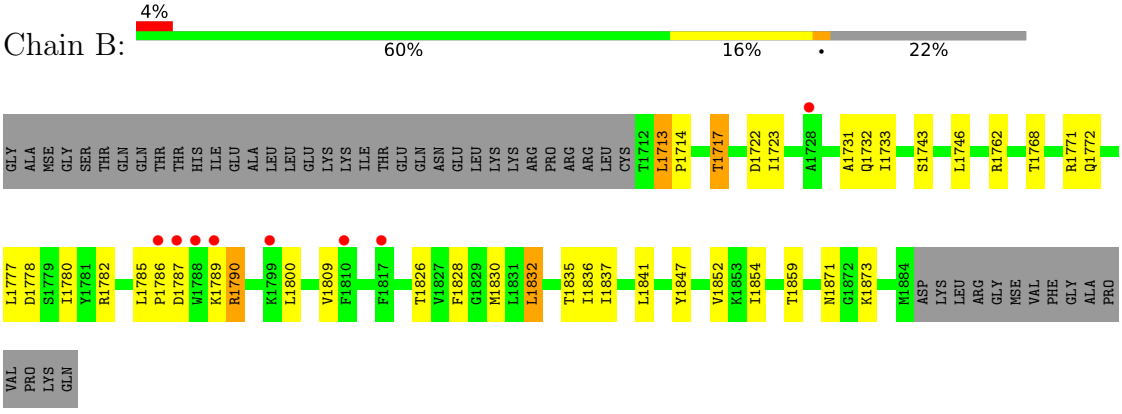
3 Residue-property plots (i)

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.

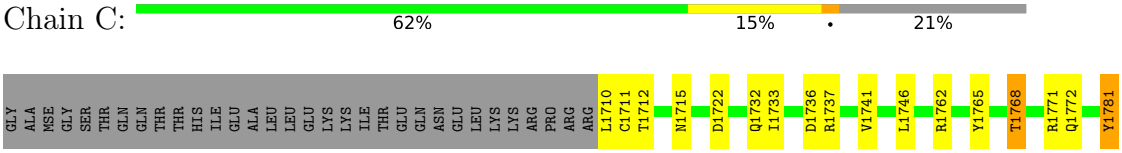
• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1

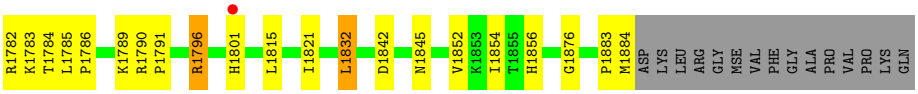


• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1



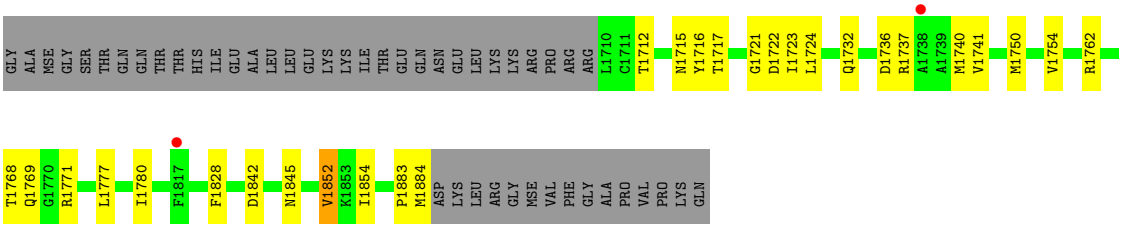
• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1





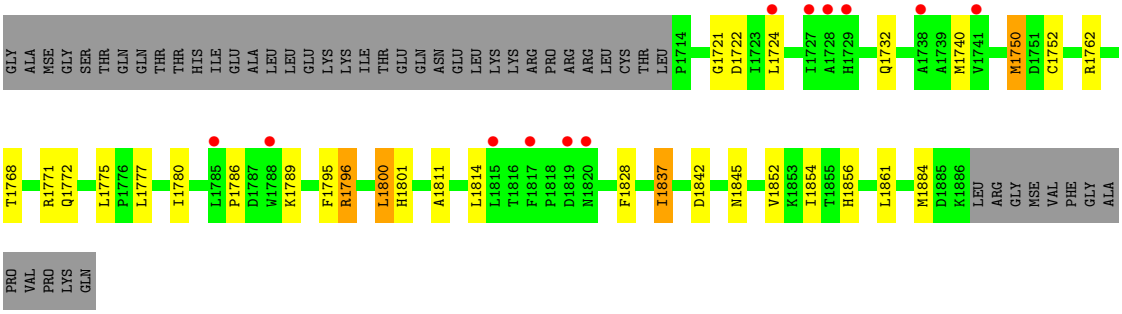
• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1

Chain D:



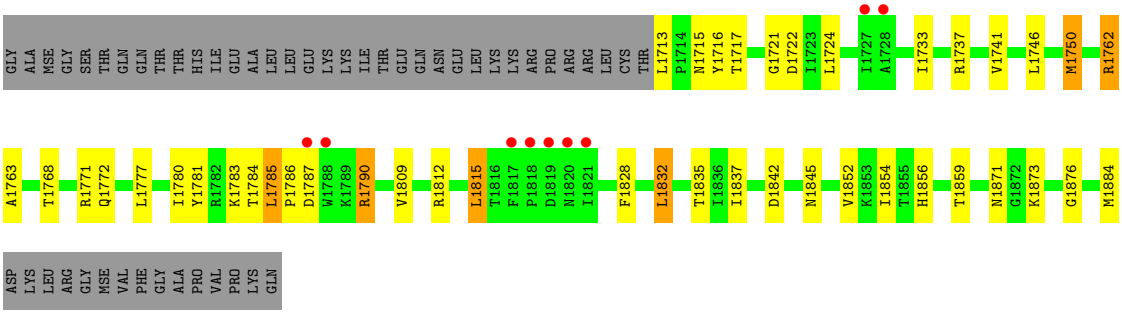
• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1

Chain E:



• Molecule 1: Structural maintenance of chromosomes flexible hinge domain-containing protein 1

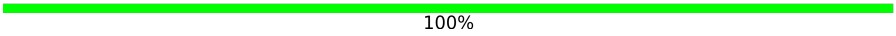
Chain F:



• Molecule 2: Uncharacterized peptide from Structural maintenance of chromosomes flexible hinge domain-containing protein 1

Chain G:

• Molecule 2: Uncharacterized peptide from Structural maintenance of chromosomes flexible hinge domain-containing protein 1

Chain H: 

There are no outlier residues recorded for this chain.

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	123.55Å 123.55Å 232.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.45 – 3.30 48.45 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.45-3.30) 99.9 (48.45-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.06 (at 3.33Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.204 , 0.239 0.212 , 0.243	Depositor DCC
R_{free} test set	1399 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	107.5	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 82.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8412	wwPDB-VP
Average B, all atoms (Å^2)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.79% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.47	0/1380	0.70	0/1864
1	B	0.49	0/1403	0.69	0/1897
1	C	0.48	0/1417	0.71	0/1916
1	D	0.49	0/1417	0.72	0/1916
1	E	0.49	0/1401	0.71	0/1893
1	F	0.50	0/1396	0.70	0/1887
All	All	0.49	0/8414	0.71	0/11373

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

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	1353	0	1349	17	0
1	B	1375	0	1374	25	0
1	C	1389	0	1390	25	0
1	D	1389	0	1388	12	0
1	E	1373	0	1363	18	0
1	F	1368	0	1367	25	0
2	G	60	0	14	0	0
2	H	105	0	23	0	0
All	All	8412	0	8268	106	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 6.

All (106) 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:1835:THR:HG22	1:B:1859:THR:HB	1.58	0.85
1:C:1783:LYS:HB2	1:C:1790:ARG:CZ	2.09	0.81
1:C:1783:LYS:HB2	1:C:1790:ARG:NH1	1.99	0.77
1:F:1785:LEU:HG	1:F:1786:PRO:HD3	1.71	0.72
1:A:1768:THR:HG21	1:A:1772:GLN:HB3	1.74	0.68
1:C:1768:THR:HG21	1:C:1772:GLN:HB3	1.75	0.66
1:A:1768:THR:CG2	1:A:1772:GLN:H	2.09	0.65
1:C:1768:THR:CG2	1:C:1772:GLN:H	2.11	0.64
1:C:1783:LYS:N	1:C:1783:LYS:HD3	2.15	0.61
1:E:1796:ARG:HB3	1:E:1801:HIS:CD2	2.35	0.61
1:F:1746:LEU:HD21	1:F:1832:LEU:HD13	1.82	0.61
1:A:1765:TYR:O	1:A:1768:THR:HG22	2.00	0.61
1:B:1826:THR:HG22	1:B:1830:MSE:HE3	1.82	0.60
1:C:1765:TYR:O	1:C:1768:THR:HG22	2.01	0.60
1:B:1787:ASP:HA	1:B:1790:ARG:CG	2.32	0.59
1:B:1871:ASN:HD21	1:B:1873:LYS:HE2	1.66	0.59
1:C:1783:LYS:CB	1:C:1790:ARG:CZ	2.80	0.59
1:A:1787:ASP:HA	1:A:1790:ARG:HD2	1.85	0.59
1:A:1768:THR:HG21	1:A:1772:GLN:H	1.68	0.58
1:B:1713:LEU:HB3	1:B:1714:PRO:CD	2.33	0.58
1:B:1786:PRO:HB2	1:B:1789:LYS:HB2	1.86	0.57
1:D:1740:MSE:HE3	1:D:1884:MSE:HG2	1.85	0.57
1:B:1786:PRO:HG2	1:B:1790:ARG:HH21	1.70	0.57
1:F:1835:THR:HG22	1:F:1859:THR:HB	1.86	0.57
1:C:1768:THR:HG21	1:C:1772:GLN:H	1.70	0.56
1:E:1777:LEU:HA	1:E:1780:ILE:HD12	1.88	0.56
1:E:1811:ALA:O	1:E:1814:LEU:HB2	2.05	0.55
1:F:1777:LEU:HA	1:F:1780:ILE:HD12	1.89	0.55
1:F:1768:THR:HG21	1:F:1772:GLN:HB2	1.88	0.55
1:C:1746:LEU:HD21	1:C:1832:LEU:HD13	1.89	0.55
1:D:1712:THR:HA	1:D:1715:ASN:OD1	2.07	0.54
1:F:1871:ASN:HD21	1:F:1873:LYS:HE2	1.73	0.53
1:B:1768:THR:HG21	1:B:1772:GLN:HB2	1.89	0.53
1:C:1762:ARG:HD2	1:D:1845:ASN:OD1	2.09	0.53
1:E:1768:THR:HG21	1:E:1772:GLN:HB2	1.90	0.53
1:A:1777:LEU:HA	1:A:1780:ILE:HD12	1.90	0.53
1:F:1832:LEU:O	1:F:1835:THR:HG23	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1787:ASP:HA	1:B:1790:ARG:HG2	1.91	0.52
1:D:1715:ASN:HA	1:D:1754:VAL:O	2.10	0.52
1:C:1783:LYS:HB2	1:C:1790:ARG:NE	2.25	0.51
1:E:1724:LEU:HD12	1:E:1750:MSE:HG3	1.92	0.51
1:B:1777:LEU:HA	1:B:1780:ILE:HD12	1.92	0.51
1:B:1778:ASP:HA	1:B:1782:ARG:HH22	1.75	0.51
1:C:1783:LYS:CB	1:C:1790:ARG:NE	2.74	0.50
1:A:1724:LEU:HD12	1:A:1750:MSE:HG3	1.94	0.50
1:E:1768:THR:HG21	1:E:1772:GLN:CB	2.42	0.49
1:A:1768:THR:HG23	1:A:1772:GLN:H	1.77	0.49
1:D:1777:LEU:HA	1:D:1780:ILE:HD12	1.95	0.49
1:F:1724:LEU:HD12	1:F:1750:MSE:HG3	1.94	0.49
1:F:1733:ILE:HD11	1:F:1815:LEU:HD23	1.95	0.49
1:B:1768:THR:HG21	1:B:1772:GLN:CB	2.43	0.48
1:A:1876:GLY:HA2	1:B:1771:ARG:O	2.13	0.48
1:B:1787:ASP:HA	1:B:1790:ARG:HG3	1.92	0.48
1:F:1768:THR:HG22	1:F:1771:ARG:HB2	1.95	0.48
1:F:1783:LYS:HG2	1:F:1784:THR:H	1.78	0.48
1:C:1876:GLY:HA2	1:D:1771:ARG:O	2.14	0.48
1:B:1768:THR:HG22	1:B:1771:ARG:HB2	1.95	0.47
1:C:1768:THR:HG23	1:C:1772:GLN:H	1.79	0.47
1:C:1784:THR:HG22	1:C:1785:LEU:HD12	1.95	0.47
1:C:1712:THR:HA	1:C:1715:ASN:OD1	2.15	0.47
1:C:1796:ARG:HB2	1:C:1801:HIS:HB2	1.96	0.47
1:A:1768:THR:HG23	1:A:1771:ARG:N	2.30	0.46
1:B:1782:ARG:HG2	1:B:1830:MSE:HE2	1.97	0.46
1:C:1789:LYS:O	1:C:1791:PRO:HD3	2.15	0.46
1:E:1837:ILE:HA	1:E:1861:LEU:O	2.16	0.46
1:E:1845:ASN:OD1	1:F:1762:ARG:HD2	2.16	0.46
1:D:1737:ARG:O	1:D:1741:VAL:HG23	2.16	0.46
1:F:1768:THR:HG21	1:F:1772:GLN:CB	2.46	0.46
1:C:1768:THR:HG23	1:C:1771:ARG:N	2.31	0.45
1:D:1724:LEU:HD12	1:D:1750:MSE:HG3	1.98	0.45
1:B:1746:LEU:HD21	1:B:1832:LEU:HD13	1.99	0.45
1:C:1845:ASN:OD1	1:D:1762:ARG:HD2	2.16	0.45
1:E:1768:THR:HG22	1:E:1771:ARG:HB2	1.98	0.45
1:F:1737:ARG:O	1:F:1741:VAL:HG23	2.18	0.44
1:E:1721:GLY:N	1:E:1750:MSE:HB2	2.33	0.44
1:E:1786:PRO:HB2	1:E:1789:LYS:HB2	2.00	0.43
1:F:1809:VAL:HG13	1:F:1837:ILE:HG23	1.99	0.43
1:F:1790:ARG:HD2	1:F:1812:ARG:NH2	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1771:ARG:O	1:F:1876:GLY:HA2	2.19	0.43
1:E:1795:PHE:CE1	1:E:1800:LEU:HD12	2.53	0.43
1:C:1781:TYR:CB	1:D:1852:VAL:HG22	2.48	0.43
1:E:1762:ARG:HD2	1:F:1845:ASN:OD1	2.18	0.43
1:A:1717:THR:HB	1:F:1717:THR:HG23	2.00	0.43
1:C:1733:ILE:HD11	1:C:1815:LEU:HD23	2.00	0.43
1:B:1836:ILE:HD13	1:B:1847:TYR:HE2	1.83	0.42
1:A:1720:SER:HB2	1:A:1752:CYS:O	2.20	0.42
1:C:1737:ARG:O	1:C:1741:VAL:HG23	2.19	0.42
1:B:1809:VAL:HG13	1:B:1837:ILE:HG23	2.01	0.42
1:D:1717:THR:HG21	1:D:1723:ILE:CD1	2.50	0.42
1:B:1731:ALA:O	1:B:1733:ILE:HG13	2.19	0.42
1:F:1721:GLY:N	1:F:1750:MSE:HB2	2.35	0.42
1:C:1786:PRO:HG3	1:E:1800:LEU:HD23	2.02	0.42
1:B:1713:LEU:HB3	1:B:1714:PRO:HD3	2.02	0.41
1:B:1832:LEU:HD12	1:B:1832:LEU:HA	1.93	0.41
1:A:1845:ASN:OD1	1:B:1762:ARG:HD2	2.19	0.41
1:E:1752:CYS:SG	1:F:1873:LYS:HD3	2.61	0.41
1:F:1871:ASN:ND2	1:F:1873:LYS:HE2	2.36	0.41
1:C:1783:LYS:CG	1:C:1790:ARG:CZ	2.98	0.41
1:D:1721:GLY:N	1:D:1750:MSE:HB2	2.36	0.41
1:A:1716:TYR:HB2	1:F:1716:TYR:O	2.21	0.41
1:B:1717:THR:OG1	1:B:1723:ILE:HD11	2.21	0.41
1:E:1856:HIS:ND1	1:F:1781:TYR:CD2	2.89	0.41
1:A:1760:ALA:HA	1:F:1763:ALA:HB1	2.02	0.40
1:A:1721:GLY:N	1:A:1750:MSE:HB2	2.37	0.40
1:A:1762:ARG:NH1	1:B:1841:LEU:HB3	2.36	0.40
1:E:1775:LEU:HD11	1:E:1780:ILE:HD11	2.03	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	168/222 (76%)	155 (92%)	12 (7%)	1 (1%)	25	57
1	B	171/222 (77%)	160 (94%)	9 (5%)	2 (1%)	13	42
1	C	173/222 (78%)	160 (92%)	11 (6%)	2 (1%)	13	42
1	D	173/222 (78%)	161 (93%)	10 (6%)	2 (1%)	13	42
1	E	171/222 (77%)	164 (96%)	5 (3%)	2 (1%)	13	42
1	F	170/222 (77%)	158 (93%)	12 (7%)	0	100	100
All	All	1026/1332 (77%)	958 (93%)	59 (6%)	9 (1%)	17	48

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	1883	PRO
1	E	1884	MSE
1	B	1732	GLN
1	D	1732	GLN
1	E	1732	GLN
1	A	1883	PRO
1	C	1732	GLN
1	C	1883	PRO
1	B	1713	LEU

5.3.2 Protein sidechains (i)

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	147/186 (79%)	135 (92%)	12 (8%)	11	36
1	B	150/186 (81%)	140 (93%)	10 (7%)	16	45
1	C	152/186 (82%)	137 (90%)	15 (10%)	8	28
1	D	152/186 (82%)	143 (94%)	9 (6%)	19	49
1	E	149/186 (80%)	139 (93%)	10 (7%)	16	45
1	F	149/186 (80%)	133 (89%)	16 (11%)	6	25
All	All	899/1116 (81%)	827 (92%)	72 (8%)	12	37

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1720	SER
1	A	1736	ASP
1	A	1740	MSE
1	A	1750	MSE
1	A	1751	ASP
1	A	1768	THR
1	A	1800	LEU
1	A	1828	PHE
1	A	1837	ILE
1	A	1842	ASP
1	A	1852	VAL
1	A	1854	ILE
1	B	1717	THR
1	B	1722	ASP
1	B	1743	SER
1	B	1785	LEU
1	B	1790	ARG
1	B	1800	LEU
1	B	1828	PHE
1	B	1832	LEU
1	B	1852	VAL
1	B	1854	ILE
1	C	1710	LEU
1	C	1711	CYS
1	C	1722	ASP
1	C	1736	ASP
1	C	1768	THR
1	C	1781	TYR
1	C	1782	ARG
1	C	1796	ARG
1	C	1821	ILE
1	C	1832	LEU
1	C	1842	ASP
1	C	1852	VAL
1	C	1854	ILE
1	C	1856	HIS
1	C	1884	MSE
1	D	1716	TYR
1	D	1722	ASP
1	D	1736	ASP
1	D	1768	THR
1	D	1769	GLN

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Mol	Chain	Res	Type
1	D	1828	PHE
1	D	1842	ASP
1	D	1852	VAL
1	D	1854	ILE
1	E	1722	ASP
1	E	1740	MSE
1	E	1750	MSE
1	E	1796	ARG
1	E	1800	LEU
1	E	1828	PHE
1	E	1837	ILE
1	E	1842	ASP
1	E	1852	VAL
1	E	1854	ILE
1	F	1713	LEU
1	F	1715	ASN
1	F	1722	ASP
1	F	1750	MSE
1	F	1762	ARG
1	F	1785	LEU
1	F	1787	ASP
1	F	1790	ARG
1	F	1815	LEU
1	F	1828	PHE
1	F	1832	LEU
1	F	1842	ASP
1	F	1852	VAL
1	F	1854	ILE
1	F	1856	HIS
1	F	1884	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1797	ASN
1	F	1715	ASN
1	F	1823	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 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, 95^{th} 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	166/222 (74%)	0.32	14 (8%)	11	10	80, 105, 175, 188	0
1	B	169/222 (76%)	0.36	8 (4%)	31	29	76, 122, 181, 198	0
1	C	171/222 (77%)	0.04	1 (0%)	89	90	69, 99, 139, 157	0
1	D	171/222 (77%)	0.07	2 (1%)	79	78	72, 98, 159, 175	0
1	E	169/222 (76%)	0.38	12 (7%)	16	16	71, 104, 184, 198	0
1	F	168/222 (75%)	0.19	9 (5%)	25	24	76, 106, 172, 194	0
2	G	0/21	-	-			-	-
2	H	0/21	-	-			-	-
All	All	1014/1374 (73%)	0.23	46 (4%)	33	32	69, 106, 174, 198	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1727	ILE	4.7
1	A	1728	ALA	4.4
1	F	1788	TRP	4.2
1	B	1788	TRP	4.0
1	E	1728	ALA	3.8
1	B	1817	PHE	3.8
1	E	1817	PHE	3.7
1	E	1815	LEU	3.2
1	B	1786	PRO	3.1
1	A	1724	LEU	3.1
1	B	1728	ALA	3.0
1	F	1727	ILE	2.9
1	E	1788	TRP	2.9
1	E	1819	ASP	2.9
1	A	1731	ALA	2.9
1	A	1819	ASP	2.8
1	E	1820	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	1733	ILE	2.8
1	E	1727	ILE	2.6
1	A	1818	PRO	2.6
1	A	1725	GLY	2.6
1	C	1801	HIS	2.6
1	E	1785	LEU	2.6
1	D	1817	PHE	2.6
1	D	1738	ALA	2.5
1	B	1789	LYS	2.5
1	A	1788	TRP	2.5
1	E	1741	VAL	2.5
1	A	1730	LEU	2.5
1	B	1787	ASP	2.4
1	E	1738	ALA	2.4
1	F	1817	PHE	2.4
1	E	1724	LEU	2.3
1	F	1819	ASP	2.3
1	A	1831	LEU	2.3
1	E	1729	HIS	2.3
1	B	1799	LYS	2.3
1	A	1815	LEU	2.2
1	A	1729	HIS	2.2
1	F	1787	ASP	2.2
1	B	1810	PHE	2.2
1	F	1728	ALA	2.1
1	F	1818	PRO	2.1
1	F	1820	ASN	2.0
1	F	1821	ILE	2.0
1	A	1821	ILE	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 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.