



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:06 AM GMT

PDB ID : 2VV5
Title : THE OPEN STRUCTURE OF MSCS
Authors : Wang, W.; Dong, C.; Johnson, K.A.; Naismith, J.H.
Deposited on : 2008-06-03
Resolution : 3.45 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

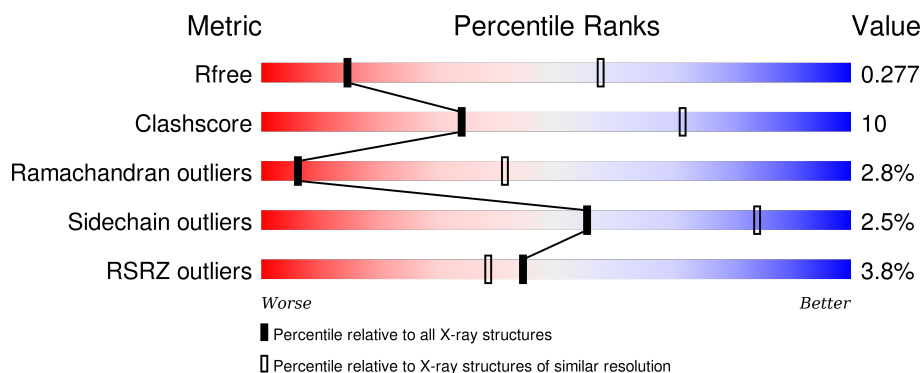
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.45 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}	91344	1000 (3.56-3.36)
Clashscore	102246	1090 (3.56-3.36)
Ramachandran outliers	100387	1057 (3.56-3.36)
Sidechain outliers	100360	1058 (3.56-3.36)
RSRZ outliers	91569	1005 (3.56-3.36)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	286	
1	B	286	
1	C	286	
1	D	286	
1	E	286	

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Mol	Chain	Length	Quality of chain
1	F	286	<div><div><div>4%</div><div>69%</div><div>19%</div><div>•</div><div>10%</div></div></div>
1	G	286	<div><div><div>2%</div><div>71%</div><div>17%</div><div>•</div><div>10%</div></div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 13692 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 SMALL-CONDUCTANCE MECHANOSENSITIVE CHANNEL.

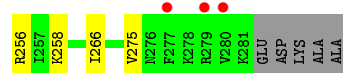
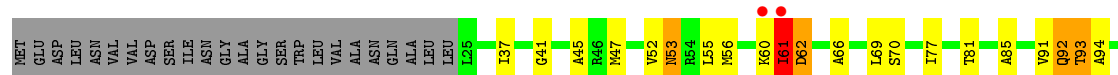
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	B	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	C	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	D	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	E	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	F	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			
1	G	257	Total	C	N	O	S	0	0	1
			1956	1256	342	352	6			

There are 7 discrepancies between the modelled and reference sequences:

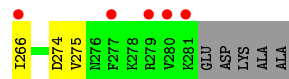
Chain	Residue	Modelled	Actual	Comment	Reference
A	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
B	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
C	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
D	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
E	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
F	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2
G	106	VAL	ALA	ENGINEERED MUTATION	UNP P0C0S2



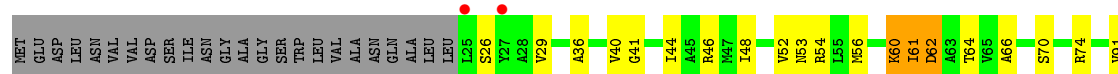
• Molecule 1: SMALL-CONDUCTANCE MECHANOSENSITIVE CHANNEL



• Molecule 1: SMALL-CONDUCTANCE MECHANOSENSITIVE CHANNEL

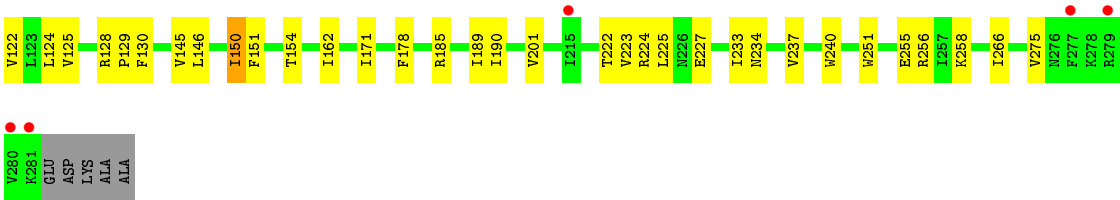


• Molecule 1: SMALL-CONDUCTANCE MECHANOSENSITIVE CHANNEL



• Molecule 1: SMALL-CONDUCTANCE MECHANOSENSITIVE CHANNEL





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	230.88Å 126.62Å 123.22Å 90.00° 90.42° 90.00°	Depositor
Resolution (Å)	123.09 – 3.45 35.48 – 3.45	Depositor EDS
% Data completeness (in resolution range)	99.0 (123.09-3.45) 99.1 (35.48-3.45)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 3.48Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.293 , 0.312 0.263 , 0.277	Depositor DCC
R_{free} test set	2336 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	117.3	Xtriage
Anisotropy	0.130	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 105.9	EDS
Estimated twinning fraction	0.036 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.016 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.030 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.018 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.035 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	3 of 46429 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13692	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.64	0/1981	0.70	1/2687 (0.0%)
1	B	0.64	0/1981	0.70	1/2687 (0.0%)
1	C	0.65	0/1981	0.69	0/2687
1	D	0.58	0/1981	0.68	0/2687
1	E	0.54	0/1981	0.65	0/2687
1	F	0.54	0/1981	0.64	0/2687
1	G	0.59	0/1981	0.65	0/2687
All	All	0.60	0/13867	0.67	2/18809 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	VAL	O-C-N	-5.51	113.89	122.70
1	B	280	VAL	O-C-N	-5.31	114.20	122.70

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	1956	0	2052	47	1
1	B	1956	0	2052	44	0
1	C	1956	0	2052	53	1
1	D	1956	0	2052	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1956	0	2052	38	0
1	F	1956	0	2052	43	0
1	G	1956	0	2052	49	0
All	All	13692	0	14364	282	1

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 (282) 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:61:ILE:HD12	1:B:62:ASP:N	1.69	1.06
1:F:61:ILE:HD11	1:F:66:ALA:HB2	1.40	1.03
1:D:56:MET:HE2	1:D:61:ILE:HD13	1.45	0.97
1:E:61:ILE:HD11	1:E:66:ALA:HB2	1.44	0.97
1:B:61:ILE:HD11	1:B:66:ALA:HB2	1.47	0.96
1:B:61:ILE:HD12	1:B:62:ASP:H	1.28	0.95
1:C:61:ILE:HD11	1:C:66:ALA:HB2	1.46	0.94
1:A:61:ILE:HD11	1:A:66:ALA:HB2	1.50	0.91
1:G:56:MET:HE1	1:G:69:LEU:HD12	1.59	0.81
1:D:61:ILE:HD11	1:D:66:ALA:HB2	1.65	0.78
1:G:61:ILE:HD11	1:G:66:ALA:HB2	1.66	0.76
1:G:56:MET:HB3	1:G:61:ILE:HD11	1.64	0.76
1:E:61:ILE:CD1	1:E:66:ALA:HB2	2.16	0.75
1:B:53:ASN:HB3	1:B:70:SER:HB2	1.69	0.75
1:E:62:ASP:OD2	1:E:65:VAL:HG23	1.89	0.72
1:C:53:ASN:HB3	1:C:70:SER:HB2	1.72	0.71
1:C:56:MET:CE	1:C:66:ALA:HA	2.21	0.70
1:D:201:VAL:HG22	1:D:266:ILE:HD13	1.76	0.68
1:B:61:ILE:CD1	1:B:62:ASP:H	2.06	0.67
1:D:53:ASN:HB3	1:D:70:SER:HB2	1.76	0.66
1:E:201:VAL:HG22	1:E:266:ILE:HD13	1.78	0.66
1:B:201:VAL:HG22	1:B:266:ILE:HD13	1.78	0.65
1:C:201:VAL:HG22	1:C:266:ILE:HD13	1.77	0.65
1:G:56:MET:HB3	1:G:61:ILE:CD1	2.29	0.63
1:D:225:LEU:CD1	1:D:233:ILE:HG23	2.28	0.63
1:F:61:ILE:CD1	1:F:66:ALA:HB2	2.24	0.63
1:G:56:MET:HE2	1:G:61:ILE:HD13	1.80	0.62
1:G:56:MET:CE	1:G:61:ILE:HD13	2.29	0.62
1:B:45:ALA:HB2	1:B:77:ILE:HG22	1.81	0.62
1:A:61:ILE:HD12	1:A:62:ASP:N	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:201:VAL:HG22	1:G:266:ILE:HD13	1.82	0.62
1:A:251:TRP:CE3	1:G:224:ARG:HD3	2.35	0.61
1:B:185:ARG:HB2	1:B:240:TRP:CD2	2.35	0.61
1:E:48:ILE:O	1:E:52:VAL:HG23	2.01	0.61
1:C:56:MET:HE2	1:C:66:ALA:HA	1.81	0.61
1:D:61:ILE:CD1	1:D:66:ALA:HB2	2.30	0.60
1:D:56:MET:HB3	1:D:61:ILE:CD1	2.31	0.60
1:C:225:LEU:CD1	1:C:233:ILE:HG23	2.30	0.60
1:E:150:ILE:HG23	1:E:151:PHE:N	2.17	0.60
1:A:150:ILE:HG23	1:A:151:PHE:N	2.17	0.59
1:F:201:VAL:HG22	1:F:266:ILE:HD13	1.83	0.59
1:F:185:ARG:HB2	1:F:240:TRP:CD2	2.38	0.59
1:G:61:ILE:CD1	1:G:66:ALA:HB2	2.32	0.59
1:F:224:ARG:HD3	1:G:251:TRP:CE3	2.38	0.59
1:A:201:VAL:HG22	1:A:266:ILE:HD13	1.86	0.58
1:C:45:ALA:HB2	1:C:77:ILE:HG22	1.85	0.58
1:E:56:MET:HE2	1:E:61:ILE:HD13	1.86	0.58
1:E:53:ASN:HB3	1:E:70:SER:HB2	1.84	0.58
1:B:150:ILE:HG23	1:B:151:PHE:N	2.18	0.58
1:B:124:LEU:HG	1:B:171:ILE:HD13	1.86	0.57
1:D:56:MET:HB3	1:D:61:ILE:HD11	1.86	0.57
1:A:150:ILE:HG23	1:A:151:PHE:CD2	2.39	0.57
1:D:130:PHE:CD2	1:D:145:VAL:HG21	2.40	0.57
1:B:61:ILE:CD1	1:B:62:ASP:N	2.59	0.57
1:C:61:ILE:HD12	1:C:62:ASP:N	2.20	0.57
1:A:185:ARG:HB2	1:A:240:TRP:CD2	2.40	0.57
1:F:56:MET:HB3	1:F:61:ILE:HG12	1.86	0.56
1:C:185:ARG:HB2	1:C:240:TRP:CD2	2.40	0.56
1:C:48:ILE:O	1:C:52:VAL:HG23	2.05	0.56
1:A:91:VAL:O	1:A:93:THR:N	2.38	0.56
1:F:53:ASN:HB3	1:F:70:SER:HB2	1.87	0.56
1:B:130:PHE:CD2	1:B:145:VAL:HG21	2.41	0.56
1:D:91:VAL:O	1:D:93:THR:N	2.40	0.55
1:G:91:VAL:O	1:G:93:THR:N	2.39	0.55
1:D:45:ALA:HB2	1:D:77:ILE:HG22	1.87	0.55
1:C:56:MET:HE3	1:C:69:LEU:HD12	1.88	0.55
1:E:225:LEU:CD1	1:E:233:ILE:HG23	2.37	0.55
1:G:130:PHE:CD2	1:G:145:VAL:HG21	2.41	0.55
1:C:56:MET:HE1	1:C:66:ALA:HA	1.89	0.55
1:E:130:PHE:CD2	1:E:145:VAL:HG21	2.42	0.55
1:E:91:VAL:O	1:E:93:THR:N	2.40	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:TRP:CD2	1:G:224:ARG:HD3	2.42	0.54
1:G:56:MET:HE2	1:G:66:ALA:HA	1.88	0.54
1:G:185:ARG:HB2	1:G:240:TRP:CD2	2.42	0.54
1:G:56:MET:CE	1:G:69:LEU:HD12	2.33	0.54
1:A:130:PHE:CD2	1:A:145:VAL:HG21	2.43	0.54
1:F:150:ILE:HG23	1:F:151:PHE:N	2.22	0.54
1:D:150:ILE:HG23	1:D:151:PHE:N	2.23	0.54
1:E:124:LEU:HG	1:E:171:ILE:HD13	1.90	0.54
1:G:150:ILE:HG23	1:G:151:PHE:CD2	2.43	0.54
1:G:190:ILE:N	1:G:190:ILE:HD12	2.23	0.53
1:E:224:ARG:HD3	1:F:251:TRP:CE3	2.43	0.53
1:A:225:LEU:CD1	1:A:233:ILE:HG23	2.37	0.53
1:A:224:ARG:HD3	1:B:251:TRP:CE3	2.44	0.53
1:D:150:ILE:HG23	1:D:151:PHE:CD2	2.43	0.53
1:C:150:ILE:HG23	1:C:151:PHE:CD2	2.44	0.53
1:E:150:ILE:HG23	1:E:151:PHE:CD2	2.44	0.53
1:B:225:LEU:CD1	1:B:233:ILE:HG23	2.39	0.53
1:F:91:VAL:O	1:F:93:THR:N	2.41	0.53
1:B:150:ILE:HG23	1:B:151:PHE:CD2	2.45	0.52
1:G:225:LEU:CD1	1:G:233:ILE:HG23	2.40	0.52
1:F:48:ILE:O	1:F:52:VAL:HG23	2.09	0.52
1:B:115:LEU:HD21	1:C:110:ALA:HB2	1.90	0.52
1:G:124:LEU:HG	1:G:171:ILE:HD13	1.92	0.52
1:C:150:ILE:HG23	1:C:151:PHE:N	2.24	0.52
1:A:53:ASN:HB3	1:A:70:SER:HB2	1.91	0.52
1:B:91:VAL:O	1:B:93:THR:N	2.42	0.52
1:A:92:GLN:O	1:A:94:ALA:N	2.43	0.52
1:C:91:VAL:O	1:C:93:THR:N	2.43	0.52
1:F:150:ILE:HG23	1:F:151:PHE:CD2	2.45	0.51
1:A:275:VAL:HG23	1:G:275:VAL:HG13	1.92	0.51
1:C:61:ILE:O	1:C:62:ASP:HB2	2.11	0.51
1:F:130:PHE:CD2	1:F:145:VAL:HG21	2.45	0.51
1:C:56:MET:HB3	1:C:61:ILE:HG12	1.93	0.51
1:E:185:ARG:HB2	1:E:240:TRP:CD2	2.46	0.51
1:B:92:GLN:O	1:B:94:ALA:N	2.44	0.51
1:C:130:PHE:CD2	1:C:145:VAL:HG21	2.46	0.51
1:F:225:LEU:CD1	1:F:233:ILE:HG23	2.40	0.51
1:B:190:ILE:N	1:B:190:ILE:HD12	2.26	0.51
1:E:92:GLN:O	1:E:94:ALA:N	2.44	0.50
1:A:124:LEU:HG	1:A:171:ILE:HD13	1.93	0.50
1:A:146:LEU:HD12	1:A:154:THR:HG22	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:THR:HG21	1:C:92:GLN:HE21	1.77	0.50
1:A:190:ILE:N	1:A:190:ILE:HD12	2.27	0.50
1:C:224:ARG:HD3	1:D:251:TRP:CE3	2.47	0.50
1:F:124:LEU:HG	1:F:171:ILE:HD13	1.94	0.50
1:E:41:GLY:HA2	1:E:44:ILE:HG22	1.94	0.50
1:G:150:ILE:HG23	1:G:151:PHE:N	2.28	0.49
1:G:189:ILE:C	1:G:190:ILE:HD12	2.33	0.49
1:D:185:ARG:HB2	1:D:240:TRP:CD2	2.46	0.49
1:F:53:ASN:HB2	1:F:66:ALA:HB1	1.95	0.49
1:A:275:VAL:CG2	1:G:275:VAL:HG13	2.42	0.49
1:F:189:ILE:C	1:F:190:ILE:HD12	2.33	0.49
1:B:56:MET:HB3	1:B:61:ILE:HG12	1.95	0.49
1:A:233:ILE:HD11	1:B:258:LYS:HE2	1.94	0.49
1:G:45:ALA:HB2	1:G:77:ILE:HG22	1.93	0.49
1:C:92:GLN:O	1:C:94:ALA:N	2.45	0.49
1:C:124:LEU:HG	1:C:171:ILE:HD13	1.94	0.49
1:A:251:TRP:HB3	1:G:224:ARG:HH11	1.77	0.48
1:C:61:ILE:HD11	1:C:66:ALA:CB	2.32	0.48
1:F:190:ILE:N	1:F:190:ILE:HD12	2.28	0.48
1:B:185:ARG:HB2	1:B:240:TRP:CE2	2.49	0.48
1:C:115:LEU:HD21	1:D:110:ALA:HB2	1.96	0.47
1:A:37:ILE:HG22	1:A:85:ALA:HB2	1.96	0.47
1:D:37:ILE:CG2	1:D:85:ALA:HB2	2.43	0.47
1:C:190:ILE:HD12	1:C:190:ILE:N	2.29	0.47
1:G:56:MET:HE1	1:G:69:LEU:CD1	2.36	0.47
1:B:227:GLU:HB3	1:B:234:ASN:HB2	1.96	0.47
1:D:56:MET:HE2	1:D:61:ILE:CD1	2.32	0.47
1:B:41:GLY:HA3	1:B:81:THR:HG21	1.97	0.47
1:F:92:GLN:O	1:F:94:ALA:N	2.47	0.47
1:D:122:VAL:O	1:D:125:VAL:HG22	2.15	0.47
1:A:37:ILE:CG2	1:A:85:ALA:HB2	2.46	0.47
1:A:56:MET:HB3	1:A:61:ILE:HG12	1.97	0.46
1:D:52:VAL:O	1:D:55:LEU:HB3	2.14	0.46
1:A:178:PHE:CE2	1:B:161:LYS:HD3	2.50	0.46
1:B:150:ILE:HG23	1:B:151:PHE:H	1.81	0.46
1:A:46:ARG:HE	1:A:74:ARG:HH21	1.63	0.46
1:C:201:VAL:HG11	1:C:235:PHE:CD2	2.51	0.46
1:A:189:ILE:C	1:A:190:ILE:HD12	2.36	0.46
1:E:150:ILE:HG23	1:E:151:PHE:H	1.81	0.46
1:A:251:TRP:HB3	1:G:224:ARG:NH1	2.30	0.46
1:D:92:GLN:O	1:D:94:ALA:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:128:ARG:N	1:G:129:PRO:CD	2.79	0.46
1:E:56:MET:HB3	1:E:61:ILE:CD1	2.46	0.45
1:G:223:VAL:O	1:G:224:ARG:HG3	2.16	0.45
1:A:227:GLU:HB3	1:A:234:ASN:HB2	1.97	0.45
1:B:201:VAL:HG11	1:B:235:PHE:CD2	2.52	0.45
1:E:56:MET:HE3	1:E:69:LEU:HD12	1.98	0.45
1:D:227:GLU:HB3	1:D:234:ASN:HB2	1.98	0.45
1:D:190:ILE:N	1:D:190:ILE:HD12	2.32	0.45
1:G:92:GLN:O	1:G:94:ALA:N	2.49	0.45
1:D:124:LEU:HG	1:D:171:ILE:HD13	1.98	0.45
1:F:61:ILE:O	1:F:62:ASP:HB2	2.17	0.45
1:B:224:ARG:HD3	1:C:251:TRP:CE3	2.51	0.45
1:F:185:ARG:HB2	1:F:240:TRP:CE2	2.52	0.45
1:D:61:ILE:O	1:D:62:ASP:HB2	2.17	0.45
1:E:274:ASP:HB2	1:F:272:GLN:NE2	2.32	0.45
1:C:53:ASN:HB2	1:C:66:ALA:HB1	1.99	0.45
1:C:122:VAL:O	1:C:125:VAL:HG22	2.17	0.45
1:E:45:ALA:HB2	1:E:77:ILE:HG22	1.98	0.44
1:F:114:SER:O	1:F:115:LEU:C	2.55	0.44
1:G:61:ILE:O	1:G:62:ASP:HB2	2.17	0.44
1:C:233:ILE:HD11	1:D:258:LYS:HE2	1.99	0.44
1:F:233:ILE:HD11	1:G:258:LYS:HE2	2.00	0.44
1:D:224:ARG:HD3	1:E:251:TRP:CE3	2.52	0.44
1:B:93:THR:HG21	1:C:92:GLN:NE2	2.32	0.44
1:E:227:GLU:HB3	1:E:234:ASN:HB2	2.00	0.44
1:C:45:ALA:HB2	1:C:77:ILE:CG2	2.47	0.44
1:D:45:ALA:HB2	1:D:77:ILE:CG2	2.47	0.44
1:D:146:LEU:HD12	1:D:154:THR:HG22	2.00	0.44
1:B:189:ILE:C	1:B:190:ILE:HD12	2.37	0.44
1:D:223:VAL:O	1:D:224:ARG:HG3	2.18	0.44
1:F:224:ARG:HD3	1:G:251:TRP:CD2	2.52	0.44
1:C:157:THR:O	1:C:159:ASP:N	2.50	0.44
1:A:161:LYS:HD3	1:G:178:PHE:CE2	2.53	0.44
1:D:37:ILE:HG22	1:D:85:ALA:HB2	2.01	0.43
1:A:122:VAL:O	1:A:123:LEU:C	2.57	0.43
1:D:53:ASN:HB2	1:D:66:ALA:HB1	2.00	0.43
1:B:45:ALA:HB2	1:B:77:ILE:CG2	2.46	0.43
1:C:189:ILE:C	1:C:190:ILE:HD12	2.39	0.43
1:F:227:GLU:HB3	1:F:234:ASN:HB2	2.00	0.43
1:G:53:ASN:HB3	1:G:70:SER:HB2	2.01	0.43
1:F:41:GLY:HA2	1:F:44:ILE:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ILE:HG23	1:A:151:PHE:H	1.82	0.43
1:C:227:GLU:HB3	1:C:234:ASN:HB2	1.99	0.43
1:D:56:MET:HE3	1:D:69:LEU:HD12	2.01	0.43
1:F:94:ALA:HA	1:F:97:ILE:HG12	2.01	0.43
1:F:137:ASP:HB3	1:F:176:ILE:HB	2.01	0.43
1:D:128:ARG:N	1:D:129:PRO:CD	2.81	0.43
1:A:224:ARG:HD3	1:B:251:TRP:CD2	2.53	0.43
1:A:45:ALA:HB2	1:A:77:ILE:HG22	2.01	0.43
1:B:178:PHE:CE2	1:C:161:LYS:HD3	2.54	0.42
1:C:37:ILE:CG2	1:C:85:ALA:HB2	2.49	0.42
1:E:190:ILE:HD12	1:E:190:ILE:N	2.34	0.42
1:E:275:VAL:HG13	1:F:275:VAL:HG23	2.00	0.42
1:E:61:ILE:O	1:E:62:ASP:HB2	2.20	0.42
1:G:41:GLY:HA2	1:G:44:ILE:HG22	2.00	0.42
1:C:128:ARG:N	1:C:129:PRO:CD	2.81	0.42
1:C:185:ARG:HB2	1:C:240:TRP:CE2	2.54	0.42
1:G:122:VAL:O	1:G:125:VAL:HG22	2.18	0.42
1:F:150:ILE:HG23	1:F:151:PHE:H	1.85	0.42
1:E:178:PHE:CE2	1:F:161:LYS:HD3	2.55	0.42
1:D:215:ILE:HD11	1:D:239:VAL:HG21	2.02	0.42
1:C:61:ILE:HD12	1:C:62:ASP:O	2.18	0.42
1:B:150:ILE:CG2	1:B:151:PHE:N	2.82	0.42
1:F:275:VAL:HG13	1:G:275:VAL:HG23	2.02	0.42
1:A:52:VAL:O	1:A:55:LEU:HB3	2.20	0.42
1:F:46:ARG:HE	1:F:74:ARG:HH21	1.68	0.42
1:B:61:ILE:O	1:B:62:ASP:HB2	2.20	0.42
1:C:112:GLN:O	1:C:115:LEU:N	2.53	0.42
1:G:94:ALA:HA	1:G:97:ILE:HG12	2.02	0.42
1:G:227:GLU:HB3	1:G:234:ASN:HB2	2.01	0.42
1:B:222:THR:O	1:B:237:VAL:HA	2.20	0.42
1:B:201:VAL:HG11	1:B:235:PHE:CE2	2.55	0.42
1:D:233:ILE:HD11	1:E:258:LYS:HE2	2.02	0.42
1:F:225:LEU:HD23	1:G:255:GLU:HA	2.02	0.42
1:D:94:ALA:HA	1:D:97:ILE:HG12	2.02	0.42
1:C:146:LEU:HD12	1:C:154:THR:HG22	2.02	0.42
1:G:146:LEU:HD12	1:G:154:THR:HG22	2.00	0.42
1:C:76:GLY:O	1:C:79:ALA:HB3	2.19	0.42
1:E:56:MET:HB3	1:E:61:ILE:HG12	2.02	0.41
1:E:150:ILE:CG2	1:E:151:PHE:N	2.82	0.41
1:C:223:VAL:O	1:C:224:ARG:HG3	2.19	0.41
1:C:275:VAL:HG13	1:D:275:VAL:HG23	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:146:LEU:HD12	1:F:154:THR:HG22	2.02	0.41
1:A:122:VAL:O	1:A:125:VAL:HG22	2.21	0.41
1:D:115:LEU:HD21	1:E:110:ALA:HB2	2.02	0.41
1:A:61:ILE:HD12	1:A:62:ASP:H	1.83	0.41
1:F:128:ARG:N	1:F:129:PRO:CD	2.82	0.41
1:C:41:GLY:HA3	1:C:81:THR:HG21	2.02	0.41
1:E:94:ALA:HA	1:E:97:ILE:HG12	2.02	0.41
1:D:112:GLN:O	1:D:113:GLY:C	2.59	0.41
1:D:41:GLY:HA3	1:D:81:THR:HG21	2.03	0.41
1:E:37:ILE:HG22	1:E:81:THR:HG23	2.02	0.41
1:C:120:ALA:HB3	1:C:167:ASN:HB3	2.02	0.41
1:F:61:ILE:HD12	1:F:62:ASP:N	2.36	0.41
1:A:225:LEU:HD23	1:B:255:GLU:HA	2.02	0.41
1:B:163:ILE:HG22	1:B:165:ILE:HG13	2.03	0.41
1:C:52:VAL:O	1:C:55:LEU:HB3	2.20	0.41
1:F:36:ALA:O	1:F:40:VAL:HG23	2.21	0.41
1:C:112:GLN:O	1:C:113:GLY:C	2.57	0.41
1:F:222:THR:O	1:F:237:VAL:HA	2.21	0.41
1:A:110:ALA:HB2	1:G:115:LEU:HD21	2.03	0.41
1:C:222:THR:O	1:C:237:VAL:HA	2.21	0.41
1:E:233:ILE:HD11	1:F:258:LYS:HE2	2.03	0.41
1:G:185:ARG:HB2	1:G:240:TRP:CE2	2.56	0.41
1:G:112:GLN:O	1:G:115:LEU:N	2.54	0.41
1:A:112:GLN:O	1:A:115:LEU:N	2.54	0.41
1:B:128:ARG:N	1:B:129:PRO:CD	2.84	0.41
1:E:137:ASP:HB3	1:E:176:ILE:HB	2.01	0.41
1:G:222:THR:O	1:G:237:VAL:HA	2.21	0.41
1:F:120:ALA:HB3	1:F:167:ASN:HB3	2.03	0.41
1:A:150:ILE:CG2	1:A:151:PHE:N	2.81	0.40
1:E:223:VAL:O	1:E:224:ARG:HG3	2.21	0.40
1:A:128:ARG:N	1:A:129:PRO:CD	2.84	0.40
1:C:201:VAL:HG11	1:C:235:PHE:CE2	2.56	0.40
1:A:185:ARG:HB2	1:A:240:TRP:CE2	2.56	0.40
1:F:188:PHE:HB3	1:F:190:ILE:CD1	2.52	0.40
1:D:112:GLN:O	1:D:115:LEU:N	2.54	0.40
1:A:114:SER:O	1:A:115:LEU:C	2.59	0.40
1:C:163:ILE:HG22	1:C:165:ILE:HG13	2.04	0.40
1:B:61:ILE:HD12	1:B:62:ASP:O	2.21	0.40
1:D:201:VAL:HG11	1:D:235:PHE:CD2	2.56	0.40
1:A:223:VAL:O	1:A:224:ARG:HG3	2.22	0.40
1:A:94:ALA:HA	1:A:97:ILE:HG12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:GLY:O	1:B:79:ALA:HB3	2.21	0.40
1:B:149:GLN:O	1:B:150:ILE:C	2.59	0.40
1:E:122:VAL:O	1:E:125:VAL:HG22	2.21	0.40
1:G:162:ILE:HD13	1:G:162:ILE:HA	1.94	0.40
1:A:76:GLY:O	1:A:79:ALA:HB3	2.21	0.40
1:G:112:GLN:O	1:G:113:GLY:C	2.59	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ARG:NH1	1:C:213:ASP:OD2[4_546]	2.10	0.10

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	255/286 (89%)	231 (91%)	17 (7%)	7 (3%)	6	42
1	B	255/286 (89%)	229 (90%)	19 (8%)	7 (3%)	6	42
1	C	255/286 (89%)	230 (90%)	17 (7%)	8 (3%)	5	39
1	D	255/286 (89%)	231 (91%)	17 (7%)	7 (3%)	6	42
1	E	255/286 (89%)	230 (90%)	18 (7%)	7 (3%)	6	42
1	F	255/286 (89%)	229 (90%)	19 (8%)	7 (3%)	6	42
1	G	255/286 (89%)	231 (91%)	17 (7%)	7 (3%)	6	42
All	All	1785/2002 (89%)	1611 (90%)	124 (7%)	50 (3%)	6	41

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	LYS
1	A	61	ILE
1	A	62	ASP
1	A	92	GLN
1	A	93	THR
1	A	112	GLN
1	B	60	LYS
1	B	61	ILE
1	B	62	ASP
1	B	92	GLN
1	B	93	THR
1	B	112	GLN
1	C	60	LYS
1	C	61	ILE
1	C	62	ASP
1	C	92	GLN
1	C	93	THR
1	C	112	GLN
1	C	150	ILE
1	D	60	LYS
1	D	61	ILE
1	D	62	ASP
1	D	92	GLN
1	D	112	GLN
1	E	60	LYS
1	E	61	ILE
1	E	62	ASP
1	E	92	GLN
1	E	93	THR
1	E	112	GLN
1	F	60	LYS
1	F	61	ILE
1	F	62	ASP
1	F	92	GLN
1	F	93	THR
1	F	112	GLN
1	G	60	LYS
1	G	61	ILE
1	G	62	ASP
1	G	92	GLN
1	G	93	THR
1	G	112	GLN
1	G	150	ILE

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Mol	Chain	Res	Type
1	A	150	ILE
1	B	150	ILE
1	D	93	THR
1	D	150	ILE
1	E	150	ILE
1	F	150	ILE
1	C	158	ALA

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	209/232 (90%)	202 (97%)	7 (3%)	45	79
1	B	209/232 (90%)	204 (98%)	5 (2%)	57	85
1	C	209/232 (90%)	204 (98%)	5 (2%)	57	85
1	D	209/232 (90%)	205 (98%)	4 (2%)	65	87
1	E	209/232 (90%)	204 (98%)	5 (2%)	57	85
1	F	209/232 (90%)	203 (97%)	6 (3%)	50	82
1	G	209/232 (90%)	205 (98%)	4 (2%)	65	87
All	All	1463/1624 (90%)	1427 (98%)	36 (2%)	55	84

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	VAL
1	A	37	ILE
1	A	42	LEU
1	A	47	MET
1	A	53	ASN
1	A	54	ARG
1	A	256	ARG
1	B	47	MET
1	B	53	ASN
1	B	54	ARG

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Mol	Chain	Res	Type
1	B	219	ARG
1	B	256	ARG
1	C	48	ILE
1	C	49	SER
1	C	54	ARG
1	C	219	ARG
1	C	256	ARG
1	D	47	MET
1	D	53	ASN
1	D	61	ILE
1	D	256	ARG
1	E	26	SER
1	E	53	ASN
1	E	54	ARG
1	E	61	ILE
1	E	256	ARG
1	F	26	SER
1	F	29	VAL
1	F	54	ARG
1	F	60	LYS
1	F	64	THR
1	F	256	ARG
1	G	26	SER
1	G	53	ASN
1	G	54	ARG
1	G	256	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	53	ASN

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 carbohydrates 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	257/286 (89%)	-0.08	7 (2%) 58 51	110, 113, 118, 134	0
1	B	257/286 (89%)	-0.12	5 (1%) 70 63	110, 113, 118, 135	0
1	C	257/286 (89%)	-0.05	8 (3%) 52 46	110, 113, 118, 135	0
1	D	257/286 (89%)	-0.03	12 (4%) 35 30	110, 113, 118, 134	0
1	E	257/286 (89%)	0.12	20 (7%) 16 15	110, 113, 118, 135	0
1	F	257/286 (89%)	0.05	12 (4%) 35 30	110, 113, 118, 134	0
1	G	257/286 (89%)	-0.04	5 (1%) 70 63	110, 113, 118, 135	0
All	All	1799/2002 (89%)	-0.02	69 (3%) 44 38	110, 113, 118, 135	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	281	LYS	8.2
1	G	281	LYS	8.1
1	G	280	VAL	6.8
1	C	280	VAL	5.6
1	A	277	PHE	5.3
1	C	279	ARG	4.5
1	G	277	PHE	4.1
1	F	279	ARG	4.1
1	F	278	LYS	4.0
1	F	277	PHE	3.9
1	E	281	LYS	3.9
1	E	61	ILE	3.9
1	E	216	LEU	3.7
1	D	277	PHE	3.6
1	D	280	VAL	3.4
1	A	278	LYS	3.4
1	D	61	ILE	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	280	VAL	3.4
1	A	279	ARG	3.2
1	F	216	LEU	3.1
1	E	242	ASN	3.0
1	E	239	VAL	3.0
1	G	279	ARG	3.0
1	B	281	LYS	2.9
1	F	280	VAL	2.9
1	F	242	ASN	2.9
1	B	277	PHE	2.9
1	E	177	ASN	2.9
1	E	215	ILE	2.9
1	D	279	ARG	2.7
1	C	61	ILE	2.7
1	E	60	LYS	2.7
1	D	60	LYS	2.6
1	D	221	MET	2.6
1	E	277	PHE	2.6
1	A	253	VAL	2.5
1	E	178	PHE	2.5
1	E	130	PHE	2.5
1	D	219	ARG	2.5
1	F	281	LYS	2.5
1	F	231	SER	2.4
1	F	27	TYR	2.4
1	D	183	VAL	2.4
1	B	25	LEU	2.4
1	C	221	MET	2.3
1	F	177	ASN	2.3
1	A	280	VAL	2.3
1	D	242	ASN	2.3
1	E	221	MET	2.3
1	E	279	ARG	2.2
1	D	240	TRP	2.2
1	F	25	LEU	2.2
1	F	130	PHE	2.2
1	E	240	TRP	2.2
1	E	176	ILE	2.2
1	A	25	LEU	2.1
1	E	266	ILE	2.1
1	D	241	SER	2.1
1	D	216	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	E	243	SER	2.1
1	B	240	TRP	2.1
1	C	220	GLU	2.1
1	C	239	VAL	2.1
1	E	241	SER	2.1
1	C	58	SER	2.1
1	E	57	ILE	2.1
1	A	27	TYR	2.0
1	B	213	ASP	2.0
1	G	215	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 carbohydrates 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.