



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 02:50 PM GMT

PDB ID : 3D6X
Title : Crystal structure of Campylobacter jejuni FabZ
Authors : Yokoyama, T.; Yeo, H.J.
Deposited on : 2008-05-20
Resolution : 2.59 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

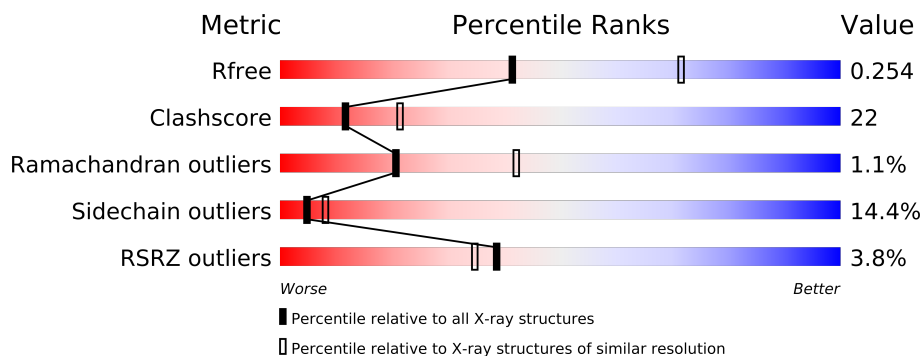
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.59 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}	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	146	
1	B	146	
1	C	146	
1	D	146	
1	E	146	
1	F	146	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6635 atoms, of which 0 are hydrogen and 0 are deuterium.

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 (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	137	Total	C	N	O	Se	0	0	0
			1082	706	180	189	7			
1	B	141	Total	C	N	O	Se	0	0	0
			1114	725	184	197	8			
1	C	140	Total	C	N	O	Se	0	0	0
			1105	720	183	194	8			
1	D	140	Total	C	N	O	Se	0	0	0
			1105	720	183	194	8			
1	E	140	Total	C	N	O	Se	0	0	0
			1107	720	183	196	8			
1	F	139	Total	C	N	O	Se	0	1	0
			1107	721	183	194	9			

- Molecule 2 is water.

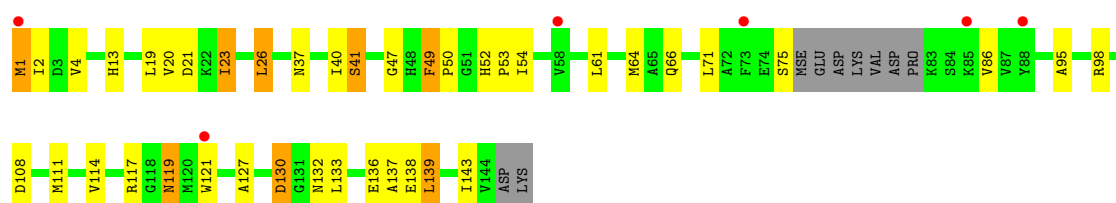
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	O	0	0
			3	3		
2	B	5	Total	O	0	0
			5	5		
2	C	3	Total	O	0	0
			3	3		
2	D	3	Total	O	0	0
			3	3		
2	F	1	Total	O	0	0
			1	1		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

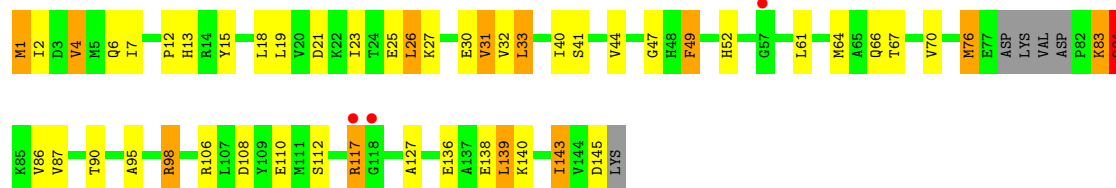
- Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain A: 



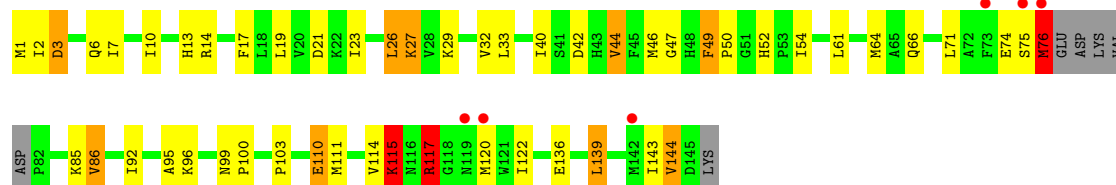
- Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain B: 



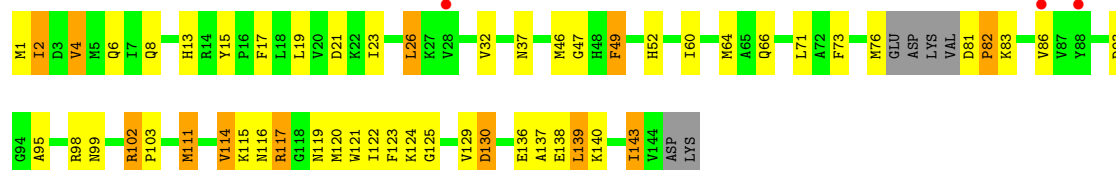
- Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain C: 



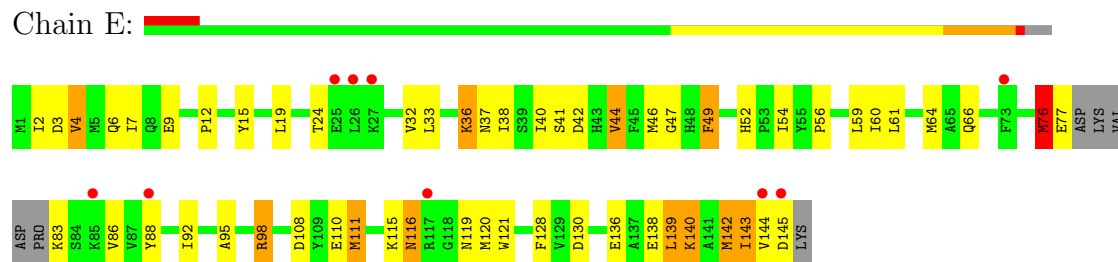
- Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain D: 



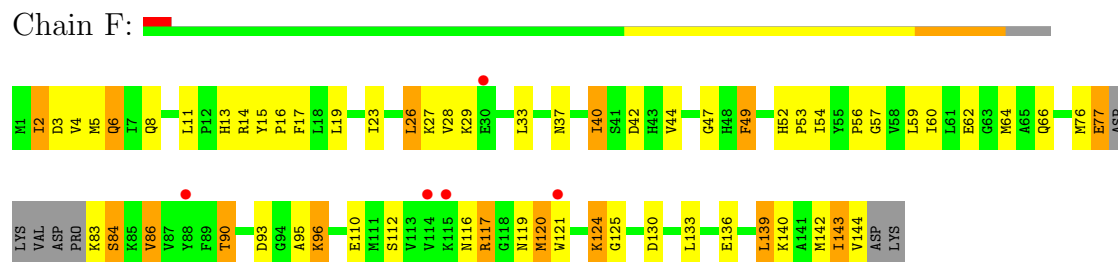
• Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain E:



• Molecule 1: (3R)-hydroxymyristoyl-[acyl-carrier-protein]dehydratase

Chain F:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.59Å 93.40Å 126.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	75.16 – 2.59 33.83 – 2.59	Depositor EDS
% Data completeness (in resolution range)	87.8 (75.16-2.59) 87.9 (33.83-2.59)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.57Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.216 , 0.254 0.216 , 0.254	Depositor DCC
R_{free} test set	1490 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	55.2	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 63.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 29397 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6635	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.63	0/1100	0.75	0/1473
1	B	0.72	0/1132	0.85	1/1514 (0.1%)
1	C	1.22	6/1123 (0.5%)	1.21	6/1502 (0.4%)
1	D	0.60	1/1123 (0.1%)	0.73	0/1503
1	E	0.58	1/1124 (0.1%)	0.70	0/1503
1	F	0.61	0/1124	0.78	0/1502
All	All	0.76	8/6726 (0.1%)	0.85	7/8997 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	117	ARG	NE-CZ	24.36	1.64	1.33
1	C	115	LYS	CE-NZ	19.09	1.96	1.49
1	C	117	ARG	CZ-NH1	8.65	1.44	1.33
1	C	115	LYS	CG-CD	7.97	1.79	1.52
1	C	76	MSE	C-O	7.94	1.38	1.23
1	C	117	ARG	CD-NE	7.73	1.59	1.46
1	E	116	ASN	CG-ND2	6.94	1.50	1.32
1	D	117	ARG	CZ-NH1	5.64	1.40	1.33

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	117	ARG	NE-CZ-NH2	-32.60	104.00	120.30
1	C	117	ARG	NH1-CZ-NH2	9.55	129.91	119.40
1	C	117	ARG	CD-NE-CZ	-7.45	113.17	123.60
1	C	117	ARG	NE-CZ-NH1	5.64	123.12	120.30
1	B	26	LEU	CA-CB-CG	5.42	127.75	115.30
1	C	115	LYS	CD-CE-NZ	-5.36	99.36	111.70
1	C	40	ILE	CB-CA-C	-5.05	101.50	111.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	84	SER	Peptide
1	C	117	ARG	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1082	0	1107	39	0
1	B	1114	0	1134	63	0
1	C	1105	0	1128	60	0
1	D	1105	0	1127	50	0
1	E	1107	0	1126	45	0
1	F	1107	0	1130	51	0
2	A	3	0	0	0	0
2	B	5	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
2	F	1	0	0	0	0
All	All	6635	0	6752	293	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (293) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:115:LYS:CD	1:C:115:LYS:CG	1.79	1.57
1:C:115:LYS:CE	1:C:115:LYS:NZ	1.96	1.27
1:C:117:ARG:HD2	1:C:120:MSE:HE3	1.26	1.18
1:D:111:MSE:HE2	1:D:123:PHE:HB3	1.27	1.10
1:C:61:LEU:HD23	1:C:64:MSE:CE	1.83	1.08
1:B:83:LYS:CG	1:B:84:SER:H	1.69	1.03
1:B:83:LYS:CD	1:B:84:SER:H	1.72	1.03
1:B:61:LEU:HD23	1:B:64:MSE:CE	1.88	1.03
1:E:98:ARG:CG	1:E:98:ARG:HH11	1.74	0.99
1:B:61:LEU:HD23	1:B:64:MSE:HE2	1.43	0.98
1:A:53:PRO:HD3	1:F:5[B]:MSE:HE3	1.42	0.98
1:B:98:ARG:HH11	1:B:98:ARG:HG2	1.29	0.95
1:C:61:LEU:HD23	1:C:64:MSE:HE1	1.50	0.94
1:D:111:MSE:CE	1:D:123:PHE:HB3	1.96	0.94
1:B:98:ARG:HH11	1:B:98:ARG:CG	1.79	0.94
1:C:117:ARG:HD2	1:C:120:MSE:CE	1.97	0.93
1:D:102:ARG:HH11	1:D:102:ARG:CG	1.82	0.93
1:B:15:TYR:CE2	1:C:46:MSE:HE2	2.05	0.92
1:C:19:LEU:H	1:C:66:GLN:HE22	1.15	0.92
1:D:102:ARG:HH11	1:D:102:ARG:HG2	1.33	0.91
1:E:76:MSE:HA	1:E:76:MSE:HE3	1.53	0.91
1:B:83:LYS:HG3	1:B:84:SER:H	1.36	0.88
1:F:19:LEU:H	1:F:66:GLN:HE22	1.22	0.88
1:F:64:MSE:HE3	1:F:125:GLY:HA3	1.56	0.87
1:E:98:ARG:HG2	1:E:98:ARG:HH11	1.36	0.87
1:D:2:ILE:HG23	1:D:23:ILE:HB	1.57	0.87
1:B:83:LYS:HG3	1:B:84:SER:N	1.90	0.86
1:F:2:ILE:HG23	1:F:23:ILE:HB	1.58	0.85
1:B:83:LYS:CG	1:B:84:SER:N	2.39	0.84
1:C:61:LEU:HA	1:C:64:MSE:HE3	1.59	0.83
1:B:117:ARG:HH11	1:B:117:ARG:HG3	1.42	0.83
1:E:42:ASP:OD1	1:E:44:VAL:HG23	1.77	0.83
1:A:19:LEU:H	1:A:66:GLN:HE22	1.26	0.83
1:D:138:GLU:C	1:D:139:LEU:HD23	1.99	0.82
1:B:61:LEU:HA	1:B:64:MSE:HE2	1.61	0.81
1:A:111:MSE:HE1	1:A:139:LEU:HD12	1.63	0.80
1:A:119:ASN:ND2	1:A:143:ILE:O	2.15	0.79
1:F:117:ARG:HB2	1:F:120:MSE:SE	2.31	0.79
1:A:52:HIS:CE1	1:A:54:ILE:HD11	2.18	0.79
1:A:119:ASN:ND2	1:A:143:ILE:HB	1.97	0.79
1:D:19:LEU:H	1:D:66:GLN:HE22	1.30	0.79
1:B:12:PRO:HD2	1:B:13:HIS:HD2	1.48	0.78
1:D:103:PRO:HG2	1:E:4:VAL:HG21	1.66	0.78

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:117:ARG:CD	1:C:120:MSE:HE3	2.10	0.76
1:B:117:ARG:HH11	1:B:117:ARG:CG	1.98	0.76
1:B:4:VAL:HG21	1:C:103:PRO:HB2	1.69	0.75
1:B:83:LYS:HD2	1:B:84:SER:H	1.51	0.74
1:C:115:LYS:CD	1:C:115:LYS:CB	2.65	0.74
1:B:61:LEU:CD2	1:B:64:MSE:CE	2.65	0.74
1:F:42:ASP:OD1	1:F:44:VAL:HG22	1.87	0.74
1:D:102:ARG:NH1	1:D:102:ARG:HG2	1.96	0.74
1:B:76:MSE:SE	1:B:143:ILE:HD11	2.39	0.73
1:F:52:HIS:CD2	1:F:54:ILE:HD13	2.24	0.72
1:B:1:MSE:HB3	1:B:2:ILE:HD12	1.71	0.72
1:E:33:LEU:HD22	1:E:110:GLU:HG2	1.72	0.72
1:D:111:MSE:CE	1:D:123:PHE:CB	2.68	0.72
1:C:6:GLN:O	1:C:10:ILE:HD12	1.90	0.71
1:C:61:LEU:HD23	1:C:64:MSE:HE3	1.71	0.71
1:B:61:LEU:HD23	1:B:64:MSE:HE1	1.73	0.71
1:B:19:LEU:H	1:B:66:GLN:HE22	1.35	0.71
1:B:83:LYS:CD	1:B:84:SER:N	2.52	0.71
1:C:114:VAL:HG22	1:C:122:ILE:HG22	1.73	0.70
1:C:19:LEU:H	1:C:66:GLN:NE2	1.89	0.70
1:E:98:ARG:HG3	1:E:98:ARG:HH11	1.56	0.70
1:B:15:TYR:HE2	1:C:46:MSE:HE2	1.53	0.70
1:B:12:PRO:HD2	1:B:13:HIS:CD2	2.27	0.69
1:E:98:ARG:HG2	1:E:98:ARG:NH1	2.06	0.69
1:B:98:ARG:NH1	1:B:98:ARG:HG2	2.08	0.68
1:D:98:ARG:HD3	1:D:99:ASN:HD21	1.57	0.68
1:B:83:LYS:HD2	1:B:84:SER:N	2.09	0.67
1:F:19:LEU:H	1:F:66:GLN:NE2	1.91	0.67
1:F:2:ILE:CG2	1:F:23:ILE:HB	2.25	0.67
1:B:32:VAL:CG1	1:B:67:THR:HG22	2.24	0.67
1:F:2:ILE:HG23	1:F:23:ILE:O	1.95	0.67
1:F:142:MSE:O	1:F:143:ILE:HD12	1.94	0.67
1:A:117:ARG:HH11	1:A:117:ARG:HG3	1.60	0.66
1:D:139:LEU:HD23	1:D:139:LEU:N	2.09	0.66
1:B:98:ARG:NH1	1:B:98:ARG:CG	2.49	0.66
1:F:57:GLY:HA2	1:F:60:ILE:HD12	1.79	0.65
1:C:13:HIS:O	1:C:14:ARG:HD3	1.97	0.65
1:E:76:MSE:HA	1:E:76:MSE:CE	2.24	0.65
1:E:111:MSE:HE1	1:E:139:LEU:HG	1.78	0.65
1:B:2:ILE:HG23	1:B:6:GLN:HB2	1.80	0.64
1:C:26:LEU:CD1	1:C:71:LEU:HD13	2.29	0.63
1:D:111:MSE:HE3	1:D:125:GLY:CA	2.28	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:61:LEU:CD2	1:B:64:MSE:HE1	2.29	0.63
1:C:114:VAL:CG2	1:C:122:ILE:HG22	2.28	0.63
1:C:26:LEU:HD13	1:C:71:LEU:HD13	1.81	0.63
1:B:25:GLU:HB2	1:B:33:LEU:HB2	1.79	0.63
1:A:111:MSE:HE1	1:A:139:LEU:CD1	2.29	0.63
1:C:13:HIS:CD2	1:C:13:HIS:H	2.15	0.63
1:B:49:PHE:HB2	1:B:52:HIS:O	1.98	0.63
1:D:19:LEU:H	1:D:66:GLN:NE2	1.95	0.62
1:F:49:PHE:HB2	1:F:52:HIS:O	2.00	0.61
1:E:98:ARG:NH1	1:E:98:ARG:CG	2.47	0.61
1:D:81:ASP:N	1:D:82:PRO:HD3	2.15	0.61
1:F:64:MSE:CE	1:F:125:GLY:HA3	2.29	0.61
1:E:49:PHE:HB2	1:E:52:HIS:O	2.01	0.61
1:A:49:PHE:HB2	1:A:52:HIS:O	2.01	0.60
1:C:49:PHE:HB2	1:C:52:HIS:O	2.02	0.60
1:F:86:VAL:HG13	1:F:144:VAL:HB	1.84	0.60
1:D:26:LEU:C	1:D:26:LEU:HD12	2.22	0.59
1:D:116:ASN:ND2	1:D:121:TRP:CZ2	2.70	0.59
1:B:32:VAL:HG13	1:B:67:THR:HG22	1.85	0.59
1:B:13:HIS:CD2	1:B:13:HIS:H	2.20	0.59
1:E:139:LEU:N	1:E:139:LEU:HD22	2.18	0.59
1:E:19:LEU:H	1:E:66:GLN:HE22	1.49	0.59
1:F:53:PRO:O	1:F:54:ILE:HD12	2.03	0.58
1:F:96:LYS:NZ	1:F:96:LYS:HB3	2.17	0.58
1:B:19:LEU:H	1:B:66:GLN:NE2	2.00	0.58
1:D:49:PHE:HB2	1:D:52:HIS:O	2.04	0.57
1:C:26:LEU:HD11	1:C:71:LEU:CD1	2.33	0.57
1:D:119:ASN:HA	1:D:143:ILE:HD11	1.84	0.57
1:B:90:THR:OG1	1:B:140:LYS:HE2	2.04	0.57
1:C:61:LEU:CD2	1:C:64:MSE:HE1	2.30	0.57
1:A:4:VAL:HG21	1:F:40:ILE:HG12	1.85	0.57
1:B:4:VAL:HG22	1:B:21:ASP:O	2.04	0.57
1:C:115:LYS:CD	1:C:115:LYS:NZ	2.67	0.57
1:F:5[B]:MSE:CE	1:F:8:GLN:HE22	2.17	0.56
1:B:32:VAL:HG11	1:B:67:THR:HG22	1.87	0.56
1:F:144:VAL:O	1:F:144:VAL:HG12	2.05	0.56
1:D:120:MSE:HE2	1:D:122:ILE:HD11	1.86	0.56
1:F:33:LEU:CD2	1:F:110:GLU:HG2	2.36	0.56
1:C:2:ILE:HB	1:C:23:ILE:HB	1.87	0.56
1:C:26:LEU:CD1	1:C:71:LEU:CD1	2.84	0.56
1:F:2:ILE:HG13	1:F:2:ILE:O	2.03	0.56
1:B:87:VAL:HG22	1:B:143:ILE:HB	1.89	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:53:PRO:HD3	1:F:5[B]:MSE:CE	2.28	0.55
1:B:117:ARG:NH1	1:B:117:ARG:CG	2.64	0.55
1:C:86:VAL:HG12	1:C:144:VAL:O	2.06	0.55
1:D:111:MSE:HE1	1:D:123:PHE:CB	2.36	0.55
1:B:95:ALA:HA	1:B:136:GLU:O	2.07	0.55
1:B:76:MSE:SE	1:B:143:ILE:CD1	3.04	0.55
1:B:23:ILE:CD1	1:B:70:VAL:HG11	2.37	0.55
1:A:2:ILE:HB	1:A:23:ILE:HB	1.89	0.54
1:C:115:LYS:CE	1:C:115:LYS:CG	2.79	0.54
1:B:139:LEU:CD2	1:B:139:LEU:N	2.70	0.54
1:F:64:MSE:HE2	1:F:139:LEU:HD21	1.90	0.54
1:F:144:VAL:CG1	1:F:144:VAL:O	2.55	0.54
1:C:3:ASP:O	1:C:7:ILE:HD12	2.08	0.54
1:E:3:ASP:O	1:E:7:ILE:HG13	2.08	0.54
1:B:33:LEU:HD22	1:B:110:GLU:HG2	1.90	0.53
1:E:33:LEU:CD2	1:E:110:GLU:HG2	2.36	0.53
1:F:52:HIS:HD2	1:F:54:ILE:HD13	1.70	0.52
1:D:26:LEU:HD13	1:D:71:LEU:HD13	1.91	0.52
1:D:119:ASN:HA	1:D:143:ILE:CD1	2.40	0.52
1:C:42:ASP:OD1	1:C:44:VAL:HG22	2.09	0.52
1:D:47:GLY:C	1:D:49:PHE:H	2.11	0.52
1:A:19:LEU:H	1:A:66:GLN:NE2	2.00	0.52
1:C:13:HIS:CD2	1:C:13:HIS:N	2.77	0.52
1:A:121:TRP:HE1	1:A:143:ILE:HD11	1.76	0.51
1:A:40:ILE:HD12	1:A:41:SER:N	2.25	0.51
1:E:12:PRO:HG2	1:F:49:PHE:CE2	2.45	0.51
1:F:33:LEU:HD22	1:F:110:GLU:HG2	1.91	0.51
1:D:60:ILE:O	1:D:64:MSE:HG3	2.11	0.51
1:C:75:SER:C	1:C:76:MSE:HG3	2.32	0.51
1:B:61:LEU:HA	1:B:64:MSE:CE	2.37	0.50
1:D:2:ILE:HG13	1:D:6:GLN:HB2	1.94	0.50
1:C:29:LYS:HG3	1:C:75:SER:HB3	1.93	0.50
1:C:26:LEU:HD12	1:C:26:LEU:C	2.32	0.50
1:D:13:HIS:H	1:D:13:HIS:CD2	2.28	0.50
1:F:117:ARG:CB	1:F:120:MSE:SE	3.06	0.50
1:E:144:VAL:HG12	1:E:145:ASP:H	1.77	0.50
1:E:56:PRO:HD2	1:E:59:LEU:HD12	1.94	0.50
1:F:76:MSE:O	1:F:77:GLU:C	2.50	0.50
1:E:144:VAL:HG12	1:E:145:ASP:N	2.27	0.50
1:D:111:MSE:HE2	1:D:123:PHE:CB	2.17	0.50
1:D:111:MSE:HE3	1:D:124:LYS:C	2.32	0.50
1:D:26:LEU:CD1	1:D:71:LEU:HD13	2.42	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:13:HIS:O	1:F:14:ARG:HD3	2.12	0.50
1:D:116:ASN:ND2	1:D:121:TRP:CE2	2.80	0.50
1:A:53:PRO:CD	1:F:5[B]:MSE:HE3	2.30	0.50
1:D:4:VAL:HG13	1:D:21:ASP:O	2.12	0.50
1:B:23:ILE:HD12	1:B:70:VAL:HG11	1.93	0.49
1:E:120:MSE:HG3	1:E:142:MSE:SE	2.62	0.49
1:E:36:LYS:HD2	1:E:38:ILE:HD13	1.95	0.49
1:B:83:LYS:HE2	1:B:84:SER:OG	2.12	0.49
1:C:95:ALA:HA	1:C:136:GLU:O	2.12	0.49
1:B:108:ASP:O	1:B:127:ALA:HA	2.13	0.49
1:F:56:PRO:HD2	1:F:59:LEU:HD12	1.95	0.49
1:A:40:ILE:C	1:A:40:ILE:HD12	2.33	0.49
1:D:37:ASN:HB3	1:E:37:ASN:HB3	1.94	0.49
1:C:13:HIS:HD2	1:C:13:HIS:H	1.59	0.49
1:E:138:GLU:C	1:E:139:LEU:HD22	2.34	0.49
1:B:2:ILE:HG23	1:B:6:GLN:CB	2.42	0.48
1:A:13:HIS:CD2	1:A:13:HIS:H	2.31	0.48
1:C:26:LEU:HD11	1:C:71:LEU:HD12	1.94	0.48
1:A:121:TRP:HE1	1:A:143:ILE:CD1	2.26	0.48
1:D:111:MSE:HE1	1:D:123:PHE:HB2	1.95	0.48
1:A:26:LEU:HD13	1:A:71:LEU:CD1	2.43	0.48
1:B:18:LEU:C	1:B:19:LEU:HD23	2.34	0.48
1:B:30:GLU:HB3	1:B:31:VAL:HG12	1.95	0.48
1:B:40:ILE:C	1:B:40:ILE:HD12	2.34	0.48
1:A:121:TRP:NE1	1:A:143:ILE:HD11	2.28	0.48
1:C:3:ASP:C	1:C:3:ASP:OD1	2.52	0.48
1:E:92:ILE:HG12	1:E:139:LEU:HD13	1.96	0.48
1:C:92:ILE:HG12	1:C:139:LEU:HD13	1.95	0.48
1:D:73:PHE:HD2	1:D:76:MSE:HE2	1.79	0.47
1:F:62:GLU:O	1:F:66:GLN:HG3	2.13	0.47
1:D:111:MSE:HE3	1:D:125:GLY:N	2.30	0.47
1:F:47:GLY:C	1:F:49:PHE:H	2.18	0.47
1:C:86:VAL:O	1:C:143:ILE:HA	2.13	0.47
1:F:2:ILE:CG2	1:F:23:ILE:O	2.62	0.47
1:C:96:LYS:NZ	1:C:96:LYS:HB3	2.28	0.47
1:E:139:LEU:N	1:E:139:LEU:CD2	2.77	0.47
1:D:102:ARG:NH1	1:D:102:ARG:CG	2.52	0.47
1:B:47:GLY:C	1:B:49:PHE:H	2.18	0.47
1:E:120:MSE:SE	1:E:140:LYS:HE3	2.65	0.47
1:B:13:HIS:HD2	1:B:13:HIS:H	1.62	0.47
1:E:120:MSE:C	1:E:121:TRP:HD1	2.18	0.47
1:F:2:ILE:HD12	1:F:6:GLN:HB3	1.97	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:26:LEU:HD13	1:A:71:LEU:HD13	1.97	0.46
1:E:60:ILE:O	1:E:64:MSE:HG3	2.15	0.46
1:B:1:MSE:HE3	1:B:1:MSE:HA	1.97	0.46
1:C:96:LYS:NZ	1:C:96:LYS:CB	2.77	0.46
1:D:114:VAL:HG23	1:D:122:ILE:O	2.15	0.46
1:D:4:VAL:O	1:D:8:GLN:HG3	2.16	0.46
1:F:5[B]:MSE:HE1	1:F:8:GLN:HE22	1.81	0.46
1:B:139:LEU:N	1:B:139:LEU:HD23	2.30	0.46
1:B:40:ILE:HD12	1:B:41:SER:N	2.30	0.46
1:B:13:HIS:CD2	1:B:13:HIS:N	2.82	0.46
1:E:95:ALA:HA	1:E:136:GLU:O	2.16	0.45
1:C:96:LYS:HZ3	1:C:96:LYS:HB3	1.81	0.45
1:D:95:ALA:HA	1:D:136:GLU:O	2.15	0.45
1:A:49:PHE:HA	1:A:50:PRO:HD2	1.82	0.45
1:A:20:VAL:HG11	1:A:23:ILE:HD11	1.97	0.45
1:E:32:VAL:CG1	1:E:111:MSE:HB3	2.46	0.45
1:C:33:LEU:CD2	1:C:110:GLU:HG3	2.47	0.45
1:A:23:ILE:N	1:A:23:ILE:HD13	2.31	0.45
1:F:95:ALA:HA	1:F:136:GLU:O	2.16	0.45
1:C:13:HIS:C	1:C:14:ARG:HD3	2.38	0.45
1:D:115:LYS:HB2	1:D:115:LYS:HE2	1.79	0.45
1:C:120:MSE:HG3	1:C:120:MSE:O	2.16	0.44
1:D:15:TYR:CE2	1:E:46:MSE:HE2	2.51	0.44
1:F:116:ASN:OD1	1:F:121:TRP:CE2	2.70	0.44
1:B:2:ILE:HG22	1:B:7:ILE:HG13	1.99	0.44
1:D:98:ARG:HD3	1:D:99:ASN:ND2	2.26	0.44
1:A:95:ALA:HA	1:A:136:GLU:O	2.17	0.44
1:A:37:ASN:HB3	1:F:37:ASN:HB3	1.98	0.44
1:A:47:GLY:C	1:A:49:PHE:H	2.20	0.44
1:C:32:VAL:HG13	1:C:111:MSE:HB2	2.00	0.44
1:E:3:ASP:C	1:E:3:ASP:OD2	2.55	0.44
1:F:90:THR:HG23	1:F:140:LYS:HD2	1.99	0.44
1:F:5[B]:MSE:CE	1:F:8:GLN:NE2	2.80	0.44
1:C:99:ASN:HB3	1:C:100:PRO:HD2	2.00	0.44
1:C:26:LEU:HD12	1:C:27:LYS:N	2.33	0.44
1:D:120:MSE:CE	1:D:122:ILE:HD11	2.48	0.44
1:F:11:LEU:HB3	1:F:13:HIS:HD2	1.83	0.44
1:A:21:ASP:OD1	1:F:40:ILE:CD1	2.66	0.43
1:E:119:ASN:O	1:E:143:ILE:HG23	2.17	0.43
1:C:61:LEU:HA	1:C:64:MSE:CE	2.37	0.43
1:E:76:MSE:HB3	1:E:77:GLU:H	1.57	0.43
1:E:47:GLY:C	1:E:49:PHE:H	2.22	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:120:MSE:CE	1:D:140:LYS:HE2	2.48	0.43
1:A:117:ARG:HG3	1:A:117:ARG:NH1	2.32	0.43
1:B:98:ARG:HH11	1:B:98:ARG:HG3	1.71	0.43
1:C:47:GLY:C	1:C:49:PHE:H	2.22	0.43
1:C:86:VAL:CG1	1:C:144:VAL:O	2.65	0.43
1:C:32:VAL:HG12	1:C:71:LEU:HD22	2.00	0.43
1:C:61:LEU:CD2	1:C:64:MSE:CE	2.76	0.43
1:A:1:MSE:C	1:A:2:ILE:HG13	2.39	0.42
1:E:61:LEU:HA	1:E:64:MSE:HE2	2.00	0.42
1:E:83:LYS:HB3	1:E:83:LYS:HE3	1.70	0.42
1:A:108:ASP:O	1:A:127:ALA:HA	2.19	0.42
1:F:119:ASN:HB2	1:F:143:ILE:HB	2.00	0.42
1:A:40:ILE:CD1	1:A:40:ILE:C	2.87	0.42
1:B:40:ILE:HG13	1:C:21:ASP:OD1	2.19	0.42
1:D:102:ARG:HH11	1:D:102:ARG:HG3	1.75	0.42
1:C:49:PHE:HA	1:C:50:PRO:HD2	1.84	0.42
1:C:52:HIS:NE2	1:C:54:ILE:HD11	2.35	0.42
1:F:26:LEU:CD1	1:F:26:LEU:C	2.89	0.42
1:A:64:MSE:HE1	1:A:137:ALA:HB3	2.02	0.42
1:F:83:LYS:HB2	1:F:84:SER:H	1.63	0.42
1:A:139:LEU:HD23	1:A:139:LEU:N	2.35	0.42
1:D:32:VAL:CG1	1:D:111:MSE:HB2	2.50	0.41
1:B:98:ARG:NH1	1:B:98:ARG:HG3	2.31	0.41
1:F:124:LYS:HG3	1:F:124:LYS:O	2.16	0.41
1:A:13:HIS:CD2	1:A:13:HIS:N	2.88	0.41
1:A:138:GLU:C	1:A:139:LEU:HD23	2.40	0.41
1:E:120:MSE:C	1:E:121:TRP:CD1	2.93	0.41
1:A:61:LEU:HA	1:A:64:MSE:HE2	2.02	0.41
1:F:19:LEU:N	1:F:66:GLN:HE22	2.03	0.41
1:F:15:TYR:CD1	1:F:16:PRO:HA	2.55	0.41
1:C:6:GLN:O	1:C:7:ILE:C	2.59	0.41
1:E:19:LEU:H	1:E:66:GLN:NE2	2.18	0.41
1:E:6:GLN:O	1:E:9:GLU:HB3	2.21	0.41
1:A:130:ASP:N	1:A:130:ASP:OD1	2.53	0.41
1:E:52:HIS:NE2	1:E:54:ILE:HD11	2.37	0.40
1:E:59:LEU:HA	1:E:59:LEU:HD23	1.87	0.40
1:E:108:ASP:HB2	1:E:128:PHE:HB2	2.02	0.40
1:D:137:ALA:HB3	1:D:139:LEU:HD21	2.03	0.40
1:D:46:MSE:HE2	1:E:15:TYR:CE2	2.56	0.40
1:D:129:VAL:HG12	1:D:130:ASP:OD2	2.22	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	133/146 (91%)	127 (96%)	5 (4%)	1 (1%)	27	53
1	B	137/146 (94%)	129 (94%)	7 (5%)	1 (1%)	30	58
1	C	136/146 (93%)	127 (93%)	8 (6%)	1 (1%)	30	58
1	D	136/146 (93%)	129 (95%)	4 (3%)	3 (2%)	10	18
1	E	136/146 (93%)	127 (93%)	7 (5%)	2 (2%)	15	30
1	F	136/146 (93%)	130 (96%)	5 (4%)	1 (1%)	30	58
All	All	814/876 (93%)	769 (94%)	36 (4%)	9 (1%)	21	42

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	PHE
1	B	49	PHE
1	C	49	PHE
1	D	49	PHE
1	D	82	PRO
1	E	49	PHE
1	D	83	LYS
1	F	49	PHE
1	E	76	MSE

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	117/118 (99%)	104 (89%)	13 (11%)	9	16
1	B	121/118 (102%)	102 (84%)	19 (16%)	4	6

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	120/118 (102%)	106 (88%)	14 (12%)	8	14
1	D	120/118 (102%)	106 (88%)	14 (12%)	8	14
1	E	120/118 (102%)	101 (84%)	19 (16%)	4	6
1	F	120/118 (102%)	96 (80%)	24 (20%)	2	3
All	All	718/708 (101%)	615 (86%)	103 (14%)	5	8

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	23	ILE
1	A	26	LEU
1	A	41	SER
1	A	75	SER
1	A	86	VAL
1	A	98	ARG
1	A	114	VAL
1	A	119	ASN
1	A	130	ASP
1	A	132	ASN
1	A	133	LEU
1	A	139	LEU
1	B	1	MSE
1	B	4	VAL
1	B	26	LEU
1	B	27	LYS
1	B	31	VAL
1	B	33	LEU
1	B	44	VAL
1	B	76	MSE
1	B	83	LYS
1	B	84	SER
1	B	86	VAL
1	B	98	ARG
1	B	106	ARG
1	B	112	SER
1	B	117	ARG
1	B	138	GLU
1	B	139	LEU
1	B	143	ILE
1	B	145	ASP

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Mol	Chain	Res	Type
1	C	1	MSE
1	C	3	ASP
1	C	17	PHE
1	C	26	LEU
1	C	27	LYS
1	C	44	VAL
1	C	74	GLU
1	C	76	MSE
1	C	85	LYS
1	C	86	VAL
1	C	110	GLU
1	C	115	LYS
1	C	139	LEU
1	C	144	VAL
1	D	1	MSE
1	D	2	ILE
1	D	4	VAL
1	D	17	PHE
1	D	26	LEU
1	D	86	VAL
1	D	93	ASP
1	D	102	ARG
1	D	111	MSE
1	D	114	VAL
1	D	117	ARG
1	D	130	ASP
1	D	139	LEU
1	D	143	ILE
1	E	2	ILE
1	E	4	VAL
1	E	24	THR
1	E	36	LYS
1	E	40	ILE
1	E	41	SER
1	E	44	VAL
1	E	76	MSE
1	E	86	VAL
1	E	88	TYR
1	E	98	ARG
1	E	111	MSE
1	E	115	LYS
1	E	116	ASN

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Mol	Chain	Res	Type
1	E	130	ASP
1	E	139	LEU
1	E	140	LYS
1	E	142	MSE
1	E	143	ILE
1	F	2	ILE
1	F	3	ASP
1	F	4	VAL
1	F	6	GLN
1	F	17	PHE
1	F	26	LEU
1	F	27	LYS
1	F	28	VAL
1	F	29	LYS
1	F	40	ILE
1	F	77	GLU
1	F	84	SER
1	F	86	VAL
1	F	90	THR
1	F	93	ASP
1	F	96	LYS
1	F	112	SER
1	F	117	ARG
1	F	120	MSE
1	F	124	LYS
1	F	130	ASP
1	F	133	LEU
1	F	139	LEU
1	F	143	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	HIS
1	A	66	GLN
1	A	119	ASN
1	A	132	ASN
1	B	13	HIS
1	B	66	GLN
1	B	116	ASN
1	B	132	ASN
1	C	8	GLN

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Mol	Chain	Res	Type
1	C	13	HIS
1	C	66	GLN
1	D	13	HIS
1	D	52	HIS
1	D	66	GLN
1	D	99	ASN
1	E	66	GLN
1	F	8	GLN
1	F	13	HIS
1	F	52	HIS
1	F	66	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	137/146 (93%)	0.24	6 (4%)	33 29	51, 56, 63, 83	0
1	B	141/146 (96%)	0.01	3 (2%)	60 58	52, 56, 63, 84	0
1	C	140/146 (95%)	0.15	6 (4%)	34 30	52, 56, 63, 78	0
1	D	140/146 (95%)	0.20	3 (2%)	60 58	51, 56, 64, 79	0
1	E	140/146 (95%)	0.32	9 (6%)	19 16	51, 56, 63, 77	0
1	F	139/146 (95%)	0.17	5 (3%)	41 37	49, 56, 63, 80	0
All	All	837/876 (95%)	0.18	32 (3%)	38 35	49, 56, 64, 84	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	26	LEU	4.6
1	E	144	VAL	4.3
1	C	119	ASN	4.2
1	A	73	PHE	4.0
1	F	121	TRP	3.6
1	E	27	LYS	3.3
1	C	120	MSE	3.3
1	B	117	ARG	3.1
1	F	114	VAL	3.1
1	A	121	TRP	3.0
1	A	85	LYS	3.0
1	A	1	MSE	2.9
1	D	28	VAL	2.9
1	E	117	ARG	2.8
1	F	88	TYR	2.6
1	C	142	MSE	2.5
1	C	73	PHE	2.5
1	B	57	GLY	2.4
1	F	30	GLU	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	86	VAL	2.4
1	C	76	MSE	2.4
1	E	88	TYR	2.4
1	B	118	GLY	2.3
1	E	85	LYS	2.2
1	A	58	VAL	2.2
1	C	75	SER	2.2
1	F	115	LYS	2.1
1	E	145	ASP	2.1
1	A	88	TYR	2.1
1	E	25	GLU	2.1
1	E	73	PHE	2.1
1	D	88	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.