



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:00 PM GMT

PDB ID : 2R1A
Title : Crystal structure of the periplasmic lipopolysaccharide transport protein LptA (YhbN), trigonal form
Authors : Suits, M.D.L.; Polissi, A.; Jia, Z.; Montreal-Kingston Bacterial Structural Genomics Initiative (BSGI)
Deposited on : 2007-08-22
Resolution : 3.26 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>

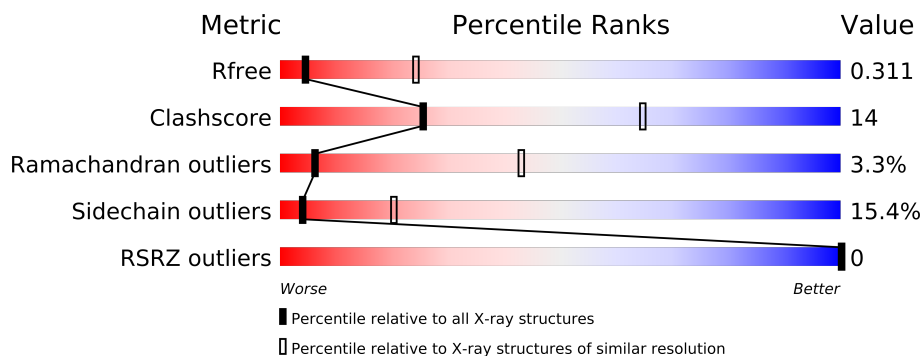
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.26 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	1085 (3.32-3.20)
Clashscore	79885	1374 (3.32-3.20)
Ramachandran outliers	78287	1348 (3.32-3.20)
Sidechain outliers	78261	1346 (3.32-3.20)
RSRZ outliers	66119	1086 (3.32-3.20)

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	159	
1	B	159	
1	C	159	
1	D	159	
1	E	159	
1	F	159	
1	G	159	
1	H	159	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7620 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 Protein yhbN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			927	578	154	191	4			
1	B	134	Total	C	N	O	S	0	0	0
			956	594	160	198	4			
1	C	135	Total	C	N	O	S	0	0	0
			989	617	163	206	3			
1	D	136	Total	C	N	O	S	0	0	0
			923	572	156	193	2			
1	E	132	Total	C	N	O	S	0	0	0
			918	569	155	192	2			
1	F	132	Total	C	N	O	S	0	0	0
			938	584	155	196	3			
1	G	135	Total	C	N	O	S	0	0	0
			980	609	165	203	3			
1	H	131	Total	C	N	O	S	0	0	0
			921	572	152	195	2			

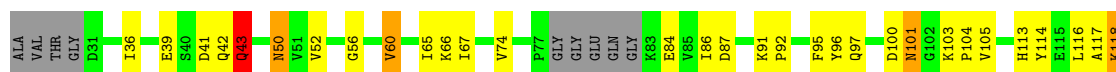
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	9	Total	O	0	0
			9	9		
2	B	8	Total	O	0	0
			8	8		
2	C	2	Total	O	0	0
			2	2		
2	D	8	Total	O	0	0
			8	8		
2	E	14	Total	O	0	0
			14	14		
2	F	9	Total	O	0	0
			9	9		

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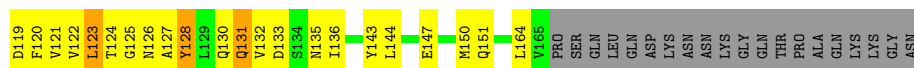
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	14	Total 14	O 14	0	0
2	H	4	Total 4	O 4	0	0



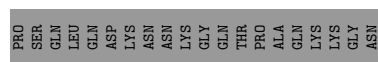
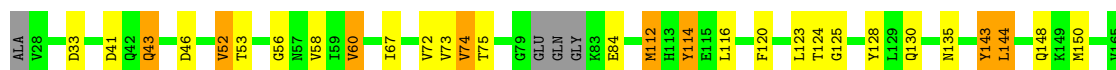
• Molecule 1: Protein yhbN

Chain F:



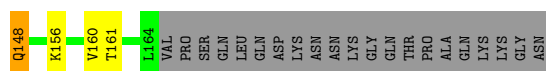
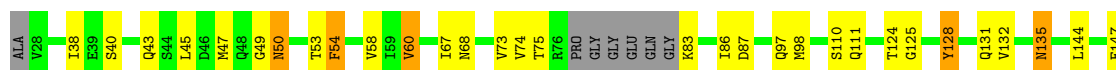
• Molecule 1: Protein yhbN

Chain G:



• Molecule 1: Protein yhbN

Chain H:



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	146.21Å 146.21Å 186.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.98 – 3.26 29.98 – 3.25	Depositor EDS
% Data completeness (in resolution range)	83.6 (29.98-3.26) 97.3 (29.98-3.25)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 3.24Å)	Xtriage
Refinement program	REFMAC 5.4.0065	Depositor
R, R_{free}	0.298 , 0.361 0.308 , 0.311	Depositor DCC
R_{free} test set	1794 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	67.8	Xtriage
Anisotropy	0.935	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 23.6	EDS
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 35752 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7620	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.40 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.4591e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.34	0/940	0.52	0/1283
1	B	0.38	0/971	0.54	0/1327
1	C	0.36	0/1004	0.53	0/1371
1	D	0.36	0/937	0.51	0/1285
1	E	0.36	0/932	0.57	0/1280
1	F	0.39	1/951 (0.1%)	0.58	0/1301
1	G	0.34	0/994	0.52	0/1356
1	H	0.34	0/934	0.50	0/1281
All	All	0.36	1/7663 (0.0%)	0.53	0/10484

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	132	VAL	CB-CG1	-5.18	1.42	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	927	0	798	25	0
1	B	956	0	831	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	989	0	899	28	0
1	D	923	0	762	20	0
1	E	918	0	758	25	0
1	F	938	0	817	36	0
1	G	980	0	882	16	0
1	H	921	0	783	17	0
2	A	9	0	0	1	0
2	B	8	0	0	1	0
2	C	2	0	0	0	0
2	D	8	0	0	0	0
2	E	14	0	0	1	0
2	F	9	0	0	1	0
2	G	14	0	0	1	0
2	H	4	0	0	2	0
All	All	7620	0	6530	191	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 14.

All (191) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:148:GLN:HG3	1:B:47:MET:HB2	1.48	0.95
1:F:98:MET:HA	1:F:99:GLN:CB	1.98	0.93
1:F:97:GLN:HG2	1:F:98:MET:H	1.43	0.83
1:B:148:GLN:HB3	1:C:47:MET:HB2	1.60	0.81
1:C:143:TYR:CE1	1:C:145:VAL:HA	2.15	0.81
1:A:115:GLU:HG3	2:A:205:HOH:O	1.82	0.78
1:C:148:GLN:HB3	1:D:47:MET:HB2	1.66	0.78
1:B:65:ILE:HA	1:B:96:TYR:O	1.84	0.77
1:B:42:GLN:HB2	1:B:55:THR:HG23	1.67	0.77
1:D:160:VAL:HG12	1:D:161:THR:H	1.54	0.73
1:G:84:GLU:HB2	1:G:116:LEU:HD12	1.70	0.72
1:C:143:TYR:HE1	1:C:145:VAL:HA	1.53	0.72
1:B:150:MET:HG2	1:B:151:GLN:H	1.54	0.72
1:B:106:GLU:HG3	1:B:107:GLY:H	1.54	0.71
1:F:130:GLN:HG2	1:F:131:GLN:H	1.55	0.71
1:E:118:LYS:O	1:E:119:ASP:HB2	1.92	0.70
1:G:46:ASP:HB2	2:G:201:HOH:O	1.91	0.70
1:C:41:ASP:HB2	1:C:56:GLY:HA3	1.72	0.70
1:C:135:ASN:ND2	1:C:135:ASN:O	2.24	0.69
1:B:125:GLY:HA3	1:B:126:ASN:C	2.12	0.69
1:B:42:GLN:O	1:B:43:GLN:HB3	1.91	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:28:VAL:N	1:F:98:MET:O	2.27	0.68
1:C:60:VAL:HG13	1:C:67:ILE:HB	1.75	0.68
1:F:136:ILE:HG12	1:G:43:GLN:HE21	1.59	0.68
1:B:29:THR:HB	1:B:100:ASP:HB3	1.77	0.67
1:D:160:VAL:HG12	1:D:161:THR:N	2.10	0.67
1:B:76:ARG:HB2	2:B:203:HOH:O	1.94	0.66
1:E:42:GLN:O	1:E:43:GLN:HB3	1.95	0.65
1:A:143:TYR:CE1	1:A:145:VAL:HA	2.32	0.64
1:A:143:TYR:HE1	1:A:145:VAL:HA	1.61	0.64
1:B:150:MET:HG2	1:B:151:GLN:N	2.11	0.63
1:E:131:GLN:HB2	2:E:209:HOH:O	1.96	0.63
1:E:86:ILE:HB	1:E:114:TYR:HB3	1.80	0.63
1:H:111:GLN:HB3	1:H:124:THR:HB	1.82	0.62
1:G:60:VAL:HG13	1:G:67:ILE:HB	1.82	0.62
1:C:95:PHE:HB2	1:C:112:MET:CE	2.31	0.61
1:F:98:MET:CA	1:F:99:GLN:CB	2.77	0.60
1:H:110:SER:HB2	1:H:125:GLY:HA3	1.83	0.60
1:G:112:MET:SD	1:G:123:LEU:HD22	2.42	0.60
1:G:58:VAL:HG11	1:G:72:VAL:HG21	1.84	0.59
1:F:52:VAL:HG13	1:F:74:VAL:HG13	1.83	0.59
1:A:92:PRO:HB2	1:A:108:HIS:HB2	1.85	0.58
1:E:154:SER:HB3	1:E:160:VAL:HG23	1.84	0.58
1:E:117:ALA:H	1:E:118:LYS:CB	2.17	0.57
1:F:136:ILE:HG12	1:G:43:GLN:NE2	2.18	0.57
1:A:98:MET:CE	1:A:98:MET:HA	2.35	0.56
1:A:132:VAL:HA	1:A:133:ASP:C	2.24	0.56
1:F:76:ARG:HG2	2:F:201:HOH:O	2.05	0.56
1:F:65:ILE:HG12	1:F:97:GLN:HG3	1.86	0.56
1:D:97:GLN:HE21	1:D:98:MET:H	1.53	0.56
1:F:111:GLN:HB3	1:F:124:THR:HB	1.87	0.56
1:F:59:ILE:HG22	1:F:59:ILE:O	2.06	0.56
1:B:106:GLU:HB3	1:B:130:GLN:HB3	1.87	0.56
1:A:60:VAL:HG13	1:A:67:ILE:HB	1.88	0.56
1:A:113:HIS:HB3	1:A:122:VAL:HB	1.89	0.55
1:G:128:TYR:HE2	1:G:130:GLN:HE21	1.54	0.54
1:H:40:SER:HB3	1:H:58:VAL:HG13	1.90	0.54
1:D:160:VAL:CG1	1:D:161:THR:H	2.19	0.54
1:E:101:ASN:N	1:E:101:ASN:HD22	2.06	0.54
1:E:152:ALA:HB3	1:F:43:GLN:HG3	1.89	0.54
1:C:29:THR:N	1:C:100:ASP:OD2	2.40	0.53
1:B:62:GLN:HE21	1:B:65:ILE:HD12	1.74	0.53
1:B:135:ASN:HD22	1:B:135:ASN:C	2.12	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:67:ILE:C	1:B:68:ASN:HD22	2.12	0.53
1:E:41:ASP:HB2	1:E:56:GLY:HA3	1.90	0.53
1:E:65:ILE:HG22	1:E:66:LYS:N	2.22	0.53
1:E:60:VAL:HG13	1:E:67:ILE:HB	1.89	0.53
1:A:105:VAL:HG22	1:A:131:GLN:HG3	1.92	0.52
1:B:128:TYR:C	1:B:128:TYR:CD2	2.82	0.52
1:B:143:TYR:CZ	1:B:145:VAL:HA	2.45	0.52
1:C:135:ASN:HD22	1:C:135:ASN:C	2.14	0.51
1:D:36:ILE:HG23	1:D:62:GLN:HB2	1.90	0.51
1:C:143:TYR:C	1:C:143:TYR:CD1	2.84	0.51
1:F:143:TYR:HD1	1:F:150:MET:HG3	1.75	0.51
1:A:40:SER:HB3	1:A:58:VAL:HG13	1.93	0.51
1:C:45:LEU:HD23	1:C:46:ASP:H	1.76	0.51
1:F:123:LEU:HB3	1:F:127:ALA:HB1	1.92	0.51
1:A:121:VAL:HB	1:A:143:TYR:HB3	1.92	0.50
1:E:164:LEU:HD21	1:F:76:ARG:HH22	1.76	0.50
1:G:150:MET:HE3	1:H:47:MET:SD	2.51	0.50
1:D:130:GLN:HG2	1:D:131:GLN:H	1.76	0.50
1:H:131:GLN:HG3	1:H:132:VAL:H	1.77	0.50
1:D:60:VAL:HG13	1:D:67:ILE:HB	1.93	0.50
1:D:52:VAL:HB	1:D:74:VAL:HG13	1.94	0.50
1:B:106:GLU:HG3	1:B:107:GLY:N	2.26	0.49
1:B:159:ARG:HH21	1:B:159:ARG:HB2	1.77	0.49
1:B:159:ARG:HD3	1:C:39:GLU:HG3	1.93	0.49
1:F:41:ASP:HB2	1:F:56:GLY:HA3	1.94	0.49
1:F:113:HIS:HB3	1:F:122:VAL:HB	1.93	0.49
1:D:107:GLY:O	1:D:108:HIS:HB3	2.13	0.49
1:H:83:LYS:N	2:H:201:HOH:O	2.46	0.49
1:C:99:GLN:HB2	1:C:103:LYS:O	2.13	0.49
1:C:83:LYS:N	1:C:84:GLU:OE2	2.45	0.49
1:F:60:VAL:HG13	1:F:67:ILE:HB	1.93	0.49
1:H:147:GLU:O	1:H:148:GLN:HB2	2.13	0.48
1:F:120:PHE:HE1	1:F:122:VAL:HG23	1.79	0.48
1:F:97:GLN:HG2	1:F:98:MET:N	2.22	0.48
1:B:83:LYS:HA	1:B:117:ALA:HB2	1.96	0.48
1:C:163:VAL:HA	1:D:36:ILE:O	2.14	0.48
1:G:143:TYR:HD1	1:G:144:LEU:N	2.11	0.48
1:A:135:ASN:HB3	1:A:163:VAL:HG22	1.95	0.48
1:B:101:ASN:O	1:B:103:LYS:N	2.47	0.47
1:C:147:GLU:O	1:C:148:GLN:HB2	2.13	0.47
1:A:128:TYR:CZ	1:A:130:GLN:HB2	2.49	0.47
1:F:55:THR:OG1	1:F:56:GLY:N	2.48	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:125:GLY:O	1:D:127:ALA:N	2.47	0.47
1:D:62:GLN:HE21	1:D:65:ILE:HD12	1.80	0.47
1:H:135:ASN:C	1:H:135:ASN:ND2	2.67	0.47
1:B:111:GLN:HB3	1:B:124:THR:HB	1.95	0.47
1:E:101:ASN:H	1:E:101:ASN:HD22	1.62	0.47
1:B:40:SER:HB2	1:B:58:VAL:HG12	1.97	0.47
1:A:114:TYR:CD2	1:A:114:TYR:C	2.89	0.47
1:C:143:TYR:HD1	1:C:143:TYR:C	2.17	0.47
1:D:76:ARG:N	1:D:77:PRO:HD3	2.30	0.47
1:G:53:THR:HG23	1:G:73:VAL:HG22	1.97	0.46
1:C:77:PRO:HD2	1:C:83:LYS:O	2.15	0.46
1:C:122:VAL:HG12	1:C:122:VAL:O	2.15	0.46
1:H:53:THR:HG23	1:H:73:VAL:HG12	1.97	0.46
1:G:41:ASP:HB2	1:G:56:GLY:HA3	1.96	0.46
1:B:52:VAL:HB	1:B:74:VAL:HG13	1.97	0.46
1:C:113:HIS:HB3	1:C:122:VAL:HB	1.97	0.46
1:B:50:ASN:N	1:B:50:ASN:OD1	2.49	0.46
1:D:128:TYR:C	1:D:128:TYR:CD2	2.89	0.46
1:F:72:VAL:HG12	1:F:73:VAL:N	2.31	0.46
1:E:117:ALA:N	1:E:118:LYS:CB	2.78	0.45
1:E:65:ILE:CG2	1:E:66:LYS:N	2.79	0.45
1:F:72:VAL:CG1	1:F:73:VAL:N	2.79	0.45
1:E:128:TYR:HA	1:E:137:LYS:HA	1.97	0.45
1:F:128:TYR:HE2	1:F:135:ASN:HB2	1.82	0.45
1:F:42:GLN:HB2	1:F:55:THR:HG23	1.99	0.45
1:G:143:TYR:HE1	1:G:148:GLN:HA	1.82	0.45
1:B:91:LYS:HA	1:B:92:PRO:HA	1.71	0.45
1:F:135:ASN:O	1:F:136:ILE:HG13	2.17	0.44
1:E:114:TYR:OH	1:E:116:LEU:HD13	2.17	0.44
1:B:113:HIS:HB3	1:B:122:VAL:HB	2.00	0.44
1:F:67:ILE:C	1:F:68:ASN:HD22	2.21	0.44
1:A:95:PHE:HB2	1:A:112:MET:HE1	2.00	0.44
1:C:40:SER:HB3	1:C:58:VAL:HG13	2.00	0.44
1:E:52:VAL:HB	1:E:74:VAL:HG23	2.00	0.44
1:A:98:MET:HE3	1:A:98:MET:HA	2.00	0.44
1:B:135:ASN:ND2	1:B:135:ASN:C	2.70	0.44
1:A:98:MET:HE2	1:A:98:MET:HA	2.00	0.44
1:F:62:GLN:HE22	1:F:116:LEU:HD21	1.83	0.44
1:C:99:GLN:HG2	1:C:105:VAL:HG23	1.99	0.43
1:A:135:ASN:O	1:A:135:ASN:ND2	2.44	0.43
1:A:114:TYR:C	1:A:114:TYR:HD2	2.22	0.43
1:G:52:VAL:HB	1:G:74:VAL:HG13	2.00	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:163:VAL:HG12	1:F:37:HIS:HA	2.01	0.43
1:F:120:PHE:CE1	1:F:122:VAL:HG23	2.54	0.43
1:C:52:VAL:HB	1:C:74:VAL:HG13	2.00	0.43
1:F:96:TYR:C	1:F:96:TYR:CD2	2.92	0.42
1:H:97:GLN:HG3	1:H:98:MET:N	2.34	0.42
1:G:114:TYR:C	1:G:114:TYR:CD2	2.93	0.42
1:H:128:TYR:HE2	1:H:135:ASN:HB2	1.85	0.42
1:A:95:PHE:HB2	1:A:112:MET:CE	2.49	0.42
1:D:97:GLN:HG3	1:D:98:MET:N	2.33	0.42
1:D:38:ILE:HD12	1:D:38:ILE:N	2.34	0.42
1:B:107:GLY:HA2	1:B:129:LEU:HA	2.00	0.42
1:C:119:ASP:O	1:C:145:VAL:HG23	2.20	0.42
1:B:106:GLU:CG	1:B:107:GLY:H	2.20	0.42
1:A:135:ASN:CB	1:A:163:VAL:HG22	2.49	0.42
1:C:84:GLU:HB2	1:C:116:LEU:HD12	2.02	0.42
1:A:68:ASN:HB2	1:A:94:THR:OG1	2.20	0.42
1:D:46:ASP:HB3	1:D:51:VAL:HB	2.02	0.42
1:C:29:THR:HG23	1:C:29:THR:O	2.19	0.42
1:B:64:THR:HG21	1:B:98:MET:HB3	2.02	0.42
1:E:36:ILE:HD12	1:E:36:ILE:H	1.85	0.42
1:F:150:MET:HG2	1:F:151:GLN:N	2.34	0.42
1:B:103:LYS:HA	1:B:104:PRO:HD3	1.84	0.41
1:D:91:LYS:HA	1:D:92:PRO:HA	1.91	0.41
1:H:54:PHE:CD1	1:H:54:PHE:N	2.87	0.41
1:E:103:LYS:HA	1:E:104:PRO:HD3	1.90	0.41
1:E:113:HIS:HB3	1:E:122:VAL:HB	2.02	0.41
1:A:143:TYR:HE1	1:A:145:VAL:CA	2.28	0.41
1:F:57:ASN:HD22	1:F:57:ASN:C	2.24	0.41
1:E:36:ILE:HD12	1:E:36:ILE:N	2.35	0.41
1:H:75:THR:HB	2:H:202:HOH:O	2.20	0.41
1:H:60:VAL:HG13	1:H:67:ILE:HB	2.02	0.41
1:C:61:THR:HG22	1:C:66:LYS:HG3	2.03	0.41
1:E:95:PHE:CD2	1:E:96:TYR:N	2.89	0.41
1:H:49:GLY:O	1:H:50:ASN:HB2	2.20	0.41
1:F:130:GLN:HG2	1:F:131:GLN:N	2.29	0.41
1:A:108:HIS:CE1	1:A:128:TYR:HD2	2.39	0.41
1:B:68:ASN:N	1:B:68:ASN:HD22	2.17	0.41
1:E:91:LYS:HA	1:E:92:PRO:HA	1.79	0.41
1:H:86:ILE:HG22	1:H:87:ASP:N	2.36	0.41
1:F:83:LYS:HA	1:F:117:ALA:HB2	2.02	0.40
1:H:160:VAL:HG12	1:H:161:THR:N	2.36	0.40
1:G:114:TYR:C	1:G:114:TYR:HD2	2.25	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:52:VAL:HB	1:D:74:VAL:CG1	2.52	0.40
1:B:112:MET:SD	1:B:113:HIS:N	2.94	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	125/159 (79%)	98 (78%)	23 (18%)	4 (3%)	6	44
1	B	130/159 (82%)	95 (73%)	31 (24%)	4 (3%)	7	45
1	C	131/159 (82%)	117 (89%)	12 (9%)	2 (2%)	15	66
1	D	132/159 (83%)	105 (80%)	19 (14%)	8 (6%)	2	22
1	E	128/159 (80%)	101 (79%)	19 (15%)	8 (6%)	2	20
1	F	128/159 (80%)	100 (78%)	24 (19%)	4 (3%)	7	45
1	G	131/159 (82%)	116 (88%)	14 (11%)	1 (1%)	27	79
1	H	127/159 (80%)	109 (86%)	15 (12%)	3 (2%)	9	53
All	All	1032/1272 (81%)	841 (82%)	157 (15%)	34 (3%)	6	43

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	156	LYS
1	D	77	PRO
1	D	119	ASP
1	E	118	LYS
1	F	99	GLN
1	A	156	LYS
1	B	43	GLN
1	B	102	GLY
1	D	118	LYS
1	D	126	ASN

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Mol	Chain	Res	Type
1	E	50	ASN
1	E	119	ASP
1	E	166	PRO
1	H	50	ASN
1	A	110	SER
1	A	135	ASN
1	B	63	GLY
1	D	117	ALA
1	E	43	GLN
1	H	156	LYS
1	D	76	ARG
1	E	84	GLU
1	F	102	GLY
1	H	148	GLN
1	D	63	GLY
1	F	56	GLY
1	F	125	GLY
1	B	49	GLY
1	D	51	VAL
1	A	56	GLY
1	E	105	VAL
1	E	125	GLY
1	G	125	GLY
1	C	125	GLY

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	88/135 (65%)	77 (88%)	11 (12%)	7	32
1	B	91/135 (67%)	73 (80%)	18 (20%)	2	10
1	C	101/135 (75%)	89 (88%)	12 (12%)	8	35
1	D	82/135 (61%)	72 (88%)	10 (12%)	7	33
1	E	82/135 (61%)	71 (87%)	11 (13%)	6	28
1	F	91/135 (67%)	65 (71%)	26 (29%)	0	2
1	G	98/135 (73%)	85 (87%)	13 (13%)	6	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	87/135 (64%)	77 (88%)	10 (12%)	8	36
All	All	720/1080 (67%)	609 (85%)	111 (15%)	4	20

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	GLU
1	A	43	GLN
1	A	55	THR
1	A	57	ASN
1	A	76	ARG
1	A	97	GLN
1	A	98	MET
1	A	114	TYR
1	A	116	LEU
1	A	135	ASN
1	A	164	LEU
1	B	43	GLN
1	B	45	LEU
1	B	50	ASN
1	B	60	VAL
1	B	61	THR
1	B	73	VAL
1	B	101	ASN
1	B	112	MET
1	B	119	ASP
1	B	121	VAL
1	B	128	TYR
1	B	131	GLN
1	B	133	ASP
1	B	134	SER
1	B	135	ASN
1	B	147	GLU
1	B	161	THR
1	B	164	LEU
1	C	33	ASP
1	C	39	GLU
1	C	43	GLN
1	C	45	LEU
1	C	74	VAL
1	C	84	GLU
1	C	111	GLN

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Mol	Chain	Res	Type
1	C	112	MET
1	C	119	ASP
1	C	135	ASN
1	C	143	TYR
1	C	144	LEU
1	D	43	GLN
1	D	45	LEU
1	D	60	VAL
1	D	87	ASP
1	D	97	GLN
1	D	100	ASP
1	D	120	PHE
1	D	124	THR
1	D	135	ASN
1	D	144	LEU
1	E	39	GLU
1	E	43	GLN
1	E	50	ASN
1	E	60	VAL
1	E	87	ASP
1	E	97	GLN
1	E	100	ASP
1	E	101	ASN
1	E	131	GLN
1	E	135	ASN
1	E	144	LEU
1	F	31	ASP
1	F	43	GLN
1	F	44	SER
1	F	45	LEU
1	F	55	THR
1	F	57	ASN
1	F	58	VAL
1	F	60	VAL
1	F	61	THR
1	F	73	VAL
1	F	74	VAL
1	F	98	MET
1	F	100	ASP
1	F	106	GLU
1	F	112	MET
1	F	114	TYR

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Mol	Chain	Res	Type
1	F	119	ASP
1	F	121	VAL
1	F	123	LEU
1	F	126	ASN
1	F	128	TYR
1	F	131	GLN
1	F	133	ASP
1	F	144	LEU
1	F	147	GLU
1	F	164	LEU
1	G	33	ASP
1	G	43	GLN
1	G	52	VAL
1	G	60	VAL
1	G	74	VAL
1	G	75	THR
1	G	112	MET
1	G	114	TYR
1	G	120	PHE
1	G	124	THR
1	G	135	ASN
1	G	143	TYR
1	G	144	LEU
1	H	38	ILE
1	H	43	GLN
1	H	45	LEU
1	H	54	PHE
1	H	60	VAL
1	H	68	ASN
1	H	74	VAL
1	H	128	TYR
1	H	135	ASN
1	H	144	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	68	ASN
1	A	101	ASN
1	A	108	HIS
1	A	131	GLN

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Mol	Chain	Res	Type
1	B	42	GLN
1	B	62	GLN
1	B	68	ASN
1	B	130	GLN
1	B	135	ASN
1	B	148	GLN
1	C	42	GLN
1	C	50	ASN
1	C	57	ASN
1	C	62	GLN
1	C	108	HIS
1	C	113	HIS
1	C	148	GLN
1	D	62	GLN
1	D	97	GLN
1	D	101	ASN
1	D	130	GLN
1	E	62	GLN
1	E	68	ASN
1	E	101	ASN
1	F	57	ASN
1	F	62	GLN
1	F	68	ASN
1	F	108	HIS
1	G	42	GLN
1	G	62	GLN
1	G	130	GLN
1	G	135	ASN
1	H	50	ASN
1	H	62	GLN
1	H	68	ASN
1	H	130	GLN
1	H	131	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	131/159 (82%)	-0.09	0 100 100	44, 64, 75, 77	0
1	B	134/159 (84%)	-0.01	0 100 100	54, 63, 74, 80	0
1	C	135/159 (84%)	-0.18	0 100 100	32, 41, 61, 65	0
1	D	136/159 (85%)	-0.08	0 100 100	47, 69, 76, 78	0
1	E	132/159 (83%)	-0.02	0 100 100	63, 71, 77, 80	0
1	F	132/159 (83%)	-0.04	0 100 100	46, 58, 68, 70	0
1	G	135/159 (84%)	-0.16	0 100 100	34, 46, 58, 60	0
1	H	131/159 (82%)	-0.12	0 100 100	35, 54, 68, 71	0
All	All	1066/1272 (83%)	-0.09	0 100 100	32, 60, 75, 80	0

There are no RSRZ outliers to report.

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.