



Full wwPDB X-ray Structure Validation Report

Jul 24, 2014 – 10:19 PM EDT

PDB ID : 4UN1
Title : Sirohaem decarboxylase AhbA/B - an enzyme with structural homology to the Lrp/AsnC transcription factor family that is part of the alternative haem biosynthesis pathway.
Authors : Palmer, D.J.; Brown, D.G.; Warren, M.J.; Pickersgill, R.W.
Deposited on : 2014-05-23
Resolution : 1.97 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>

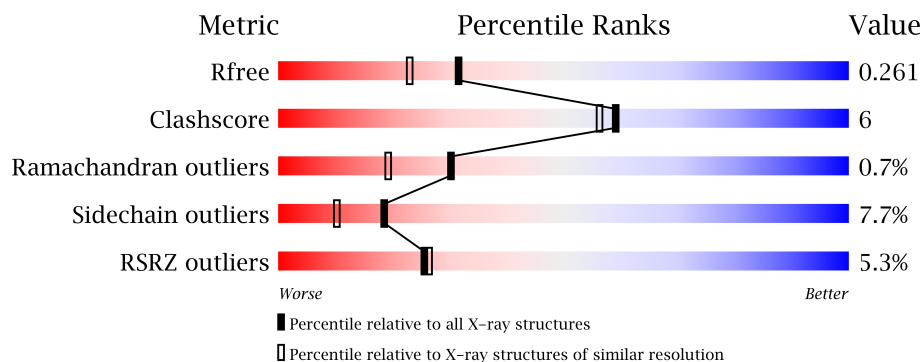
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable23489
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) : stable23489

1 Overall quality at a glance

The reported resolution of this entry is 1.97 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	6577 (2.00-1.96)
Clashscore	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)
RSRZ outliers	66119	6578 (2.00-1.96)

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	173	
1	C	173	
2	B	179	
2	D	179	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	OBV	B	1160	-	X
3	OBV	C	1163	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5119 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 PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	145	Total	C	N	O	S	0	0	0
			1131	719	201	207	4			
1	C	139	Total	C	N	O	S	0	0	0
			1083	687	192	201	3			

- Molecule 2 is a protein called PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	159	Total	C	N	O	S	0	0	0
			1233	772	219	234	8			
2	D	157	Total	C	N	O	S	0	0	0
			1234	775	219	233	7			

There are 40 discrepancies between the modelled and reference sequences:

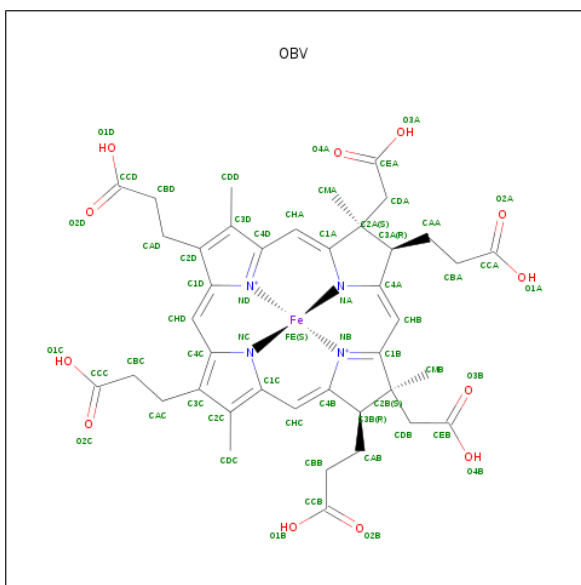
Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MET	-	EXPRESSION TAG	UNP B8J3A4
B	-18	GLY	-	EXPRESSION TAG	UNP B8J3A4
B	-17	SER	-	EXPRESSION TAG	UNP B8J3A4
B	-16	SER	-	EXPRESSION TAG	UNP B8J3A4
B	-15	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-14	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-13	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-12	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-11	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-10	HIS	-	EXPRESSION TAG	UNP B8J3A4
B	-9	SER	-	EXPRESSION TAG	UNP B8J3A4
B	-8	SER	-	EXPRESSION TAG	UNP B8J3A4
B	-7	GLY	-	EXPRESSION TAG	UNP B8J3A4
B	-6	LEU	-	EXPRESSION TAG	UNP B8J3A4
B	-5	VAL	-	EXPRESSION TAG	UNP B8J3A4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	PRO	-	EXPRESSION TAG	UNP B8J3A4
B	-3	ARG	-	EXPRESSION TAG	UNP B8J3A4
B	-2	GLY	-	EXPRESSION TAG	UNP B8J3A4
B	-1	SER	-	EXPRESSION TAG	UNP B8J3A4
B	0	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-19	MET	-	EXPRESSION TAG	UNP B8J3A4
D	-18	GLY	-	EXPRESSION TAG	UNP B8J3A4
D	-17	SER	-	EXPRESSION TAG	UNP B8J3A4
D	-16	SER	-	EXPRESSION TAG	UNP B8J3A4
D	-15	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-14	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-13	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-12	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-11	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-10	HIS	-	EXPRESSION TAG	UNP B8J3A4
D	-9	SER	-	EXPRESSION TAG	UNP B8J3A4
D	-8	SER	-	EXPRESSION TAG	UNP B8J3A4
D	-7	GLY	-	EXPRESSION TAG	UNP B8J3A4
D	-6	LEU	-	EXPRESSION TAG	UNP B8J3A4
D	-5	VAL	-	EXPRESSION TAG	UNP B8J3A4
D	-4	PRO	-	EXPRESSION TAG	UNP B8J3A4
D	-3	ARG	-	EXPRESSION TAG	UNP B8J3A4
D	-2	GLY	-	EXPRESSION TAG	UNP B8J3A4
D	-1	SER	-	EXPRESSION TAG	UNP B8J3A4
D	0	HIS	-	EXPRESSION TAG	UNP B8J3A4

- Molecule 3 is 12,18-DIDECARBOXY-SIROHEME (three-letter code: OBV) (formula: $C_{40}H_{44}FeN_4O_{12}$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total 57	C 40	Fe 1	N 4	O 12	0	0
3	C	1	Total 57	C 40	Fe 1	N 4	O 12	0	0

- Molecule 4 is water.

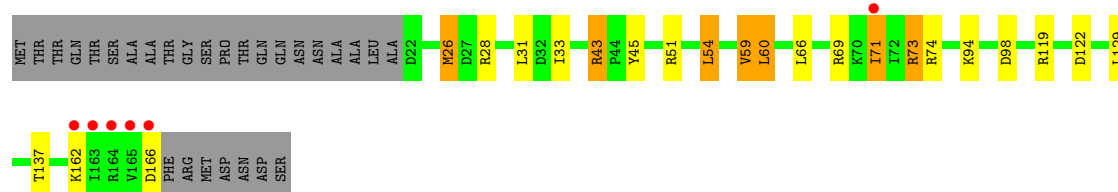
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	85	Total O 85 85	0	0
4	B	67	Total O 67 67	0	0
4	C	97	Total O 97 97	0	0
4	D	75	Total O 75 75	0	0

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.

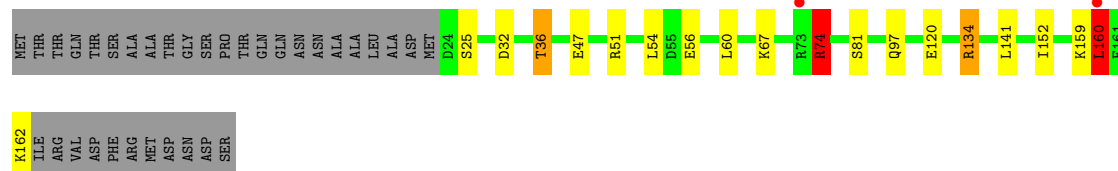
- Molecule 1: PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY

Chain A: 



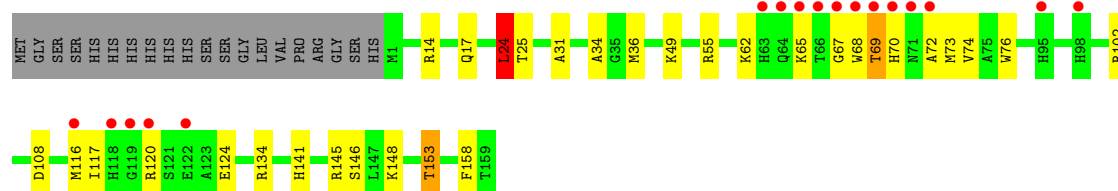
- Molecule 1: PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY

Chain C: 



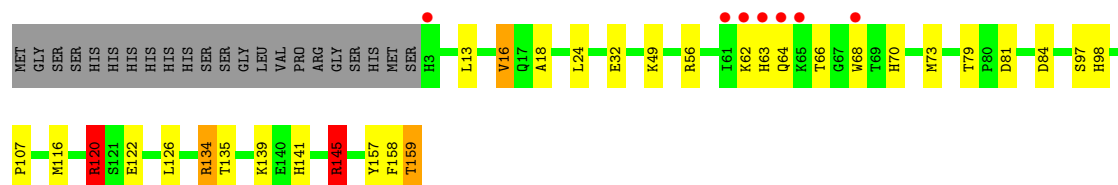
- Molecule 2: PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY

Chain B: 



- Molecule 2: PUTATIVE TRANSCRIPTIONAL REGULATOR, ASNC FAMILY

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.30Å 78.60Å 150.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.23 – 1.97 45.23 – 1.97	Depositor EDS
% Data completeness (in resolution range)	99.2 (45.23-1.97) 99.2 (45.23-1.97)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.07 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.202 , 0.257 0.208 , 0.261	Depositor DCC
R_{free} test set	2333 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 46648 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5119	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: OBV

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	0.87	0/1151	0.99	3/1558 (0.2%)
1	C	0.89	0/1103	1.06	6/1497 (0.4%)
2	B	0.86	0/1259	1.04	9/1704 (0.5%)
2	D	0.96	0/1262	1.07	6/1710 (0.4%)
All	All	0.89	0/4775	1.04	24/6469 (0.4%)

There are no bond length outliers.

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	145	ARG	NE-CZ-NH2	-9.45	115.58	120.30
2	D	145	ARG	NE-CZ-NH1	9.13	124.87	120.30
2	D	120	ARG	NE-CZ-NH1	8.38	124.49	120.30
1	C	134	ARG	NE-CZ-NH1	8.15	124.37	120.30
1	C	134	ARG	NE-CZ-NH2	-7.41	116.59	120.30
1	C	74	ARG	NE-CZ-NH1	7.26	123.93	120.30
2	B	134	ARG	NE-CZ-NH2	-6.79	116.90	120.30
2	D	120	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	C	74	ARG	NE-CZ-NH2	-6.46	117.07	120.30
2	B	102	ARG	NE-CZ-NH2	-6.36	117.12	120.30
2	D	145	ARG	CG-CD-NE	-6.09	99.01	111.80
2	B	55	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	A	28	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	C	51	ARG	NE-CZ-NH2	-5.87	117.37	120.30
2	B	134	ARG	CG-CD-NE	-5.80	99.61	111.80
1	C	160	LEU	CB-CG-CD2	5.76	120.78	111.00
2	B	34	ALA	C-N-CA	-5.75	110.21	122.30
2	D	16	VAL	CG1-CB-CG2	5.60	119.86	110.90
1	A	28	ARG	NE-CZ-NH2	-5.37	117.61	120.30
2	B	134	ARG	NE-CZ-NH1	5.33	122.96	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	ARG	NE-CZ-NH1	5.21	122.90	120.30
2	B	24	LEU	CB-CG-CD1	5.19	119.82	111.00
2	B	145	ARG	NE-CZ-NH1	5.07	122.83	120.30
2	B	73	MET	CG-SD-CE	-5.02	92.16	100.20

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	1131	0	1148	15	0
1	C	1083	0	1087	16	0
2	B	1233	0	1207	14	0
2	D	1234	0	1209	23	0
3	B	57	0	0	2	0
3	C	57	0	0	3	0
4	A	85	0	0	4	0
4	B	67	0	0	2	0
4	C	97	0	0	8	0
4	D	75	0	0	3	0
All	All	5119	0	4651	58	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 6.

All (58) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:2056:HOH:O	2:B:153:THR:HG21	1.58	1.03
1:A:43:ARG:HH11	1:A:43:ARG:HG2	1.35	0.92
2:D:120:ARG:NE	4:D:2055:HOH:O	2.15	0.80
3:B:1160:OBV:CDB	3:B:1160:OBV:CCB	2.67	0.72
2:B:67:GLY:O	2:B:69:THR:N	2.23	0.72
2:B:31:ALA:HB1	2:B:36:MET:O	1.92	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:162:LYS:HA	4:C:2096:HOH:O	1.92	0.69
1:A:43:ARG:HH11	1:A:43:ARG:CG	2.05	0.68
1:C:162:LYS:C	4:C:2096:HOH:O	2.32	0.66
2:B:17:GLN:HE22	2:B:158:PHE:H	1.45	0.65
4:A:2056:HOH:O	2:B:153:THR:CG2	2.31	0.65
1:C:32:ASP:O	1:C:36:THR:HG23	1.99	0.62
1:A:43:ARG:HG2	1:A:43:ARG:NH1	2.06	0.62
2:D:120:ARG:HH11	2:D:120:ARG:HG2	1.65	0.60
2:D:135:THR:HG21	4:D:2065:HOH:O	2.03	0.59
2:B:49:LYS:NZ	4:B:2034:HOH:O	2.35	0.59
1:A:33:ILE:HD12	1:A:51:ARG:NH2	2.17	0.59
3:C:1163:OBV:O2C	2:D:116:MET:CE	2.51	0.59
2:D:73:MET:HG3	2:D:116:MET:SD	2.42	0.58
2:B:120:ARG:N	2:B:124:GLU:OE1	2.36	0.58
1:C:162:LYS:CA	4:C:2096:HOH:O	2.52	0.58
2:D:139:LYS:O	2:D:141:HIS:HD2	1.88	0.57
2:D:66:THR:HG22	2:D:68:TRP:O	2.06	0.56
2:B:72:ALA:HB3	2:B:117:ILE:HB	1.91	0.53
3:C:1163:OBV:O1C	2:D:98:HIS:HE1	1.93	0.52
1:C:74:ARG:HG2	2:D:18:ALA:O	2.10	0.51
1:A:31:LEU:HD11	1:A:71:ILE:HD11	1.92	0.51
2:D:120:ARG:HH11	2:D:120:ARG:CG	2.24	0.51
1:C:162:LYS:CB	2:D:97:SER:O	2.60	0.50
1:C:160:LEU:CB	4:C:2094:HOH:O	2.60	0.49
1:C:97:GLN:HG3	4:C:2053:HOH:O	2.12	0.49
1:A:45:TYR:HD1	1:A:59:VAL:HG22	1.76	0.49
1:A:73:ARG:HD2	2:B:62:LYS:HG2	1.95	0.49
2:B:69:THR:HG23	2:B:120:ARG:HA	1.95	0.49
1:A:60:LEU:HD12	2:B:24:LEU:HD13	1.94	0.48
1:A:66:LEU:HD22	1:A:71:ILE:HG12	1.96	0.48
1:C:160:LEU:HB2	4:C:2094:HOH:O	2.13	0.48
2:B:14:ARG:NH2	2:D:159:THR:O	2.42	0.47
1:C:81:SER:HB3	2:D:157:TYR:OH	2.15	0.47
1:A:69:ARG:HD3	4:A:2029:HOH:O	2.15	0.46
1:A:166:ASP:N	4:A:2083:HOH:O	2.45	0.46
2:D:134:ARG:HD3	4:D:2063:HOH:O	2.16	0.45
1:C:159:LYS:HE3	2:D:84:ASP:OD1	2.16	0.45
2:D:13:LEU:HD13	2:D:158:PHE:CZ	2.52	0.45
1:C:120:GLU:OE1	2:D:145:ARG:NH1	2.47	0.44
4:B:2009:HOH:O	2:D:159:THR:HG23	2.18	0.44
1:A:119:ARG:NH2	3:B:1160:OBV:O2D	2.51	0.43
2:D:70:HIS:CD2	2:D:122:GLU:OE1	2.70	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:60:LEU:HD22	2:D:24:LEU:CD2	2.49	0.43
1:C:141:LEU:HD22	1:C:152:ILE:HG21	1.99	0.42
1:A:94:LYS:NZ	2:B:108:ASP:OD2	2.42	0.42
2:B:76:TRP:CE3	2:B:141:HIS:HB3	2.55	0.42
1:C:47:GLU:HG3	4:C:2024:HOH:O	2.19	0.41
1:C:56:GLU:HG2	2:D:24:LEU:HD22	2.03	0.41
1:A:26:MET:HG2	1:A:54:LEU:HD21	2.03	0.41
1:A:129:LEU:HD23	1:A:137:THR:HG23	2.02	0.41
3:C:1163:OBV:O2C	2:D:116:MET:HE3	2.19	0.40
4:C:2011:HOH:O	2:D:49:LYS:HE3	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	143/173 (83%)	140 (98%)	3 (2%)	0	100	100
1	C	137/173 (79%)	132 (96%)	5 (4%)	0	100	100
2	B	157/179 (88%)	142 (90%)	12 (8%)	3 (2%)	12	4
2	D	155/179 (87%)	148 (96%)	6 (4%)	1 (1%)	33	23
All	All	592/704 (84%)	562 (95%)	26 (4%)	4 (1%)	30	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	68	TRP
2	B	65	LYS
2	D	62	LYS
2	B	70	HIS

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	120/146 (82%)	110 (92%)	10 (8%)	16	8
1	C	115/146 (79%)	108 (94%)	7 (6%)	26	17
2	B	130/152 (86%)	122 (94%)	8 (6%)	26	17
2	D	131/152 (86%)	118 (90%)	13 (10%)	11	5
All	All	496/596 (83%)	458 (92%)	38 (8%)	18	10

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	MET
1	A	43	ARG
1	A	54	LEU
1	A	59	VAL
1	A	60	LEU
1	A	71	ILE
1	A	73	ARG
1	A	98	ASP
1	A	122	ASP
1	A	162	LYS
2	B	24	LEU
2	B	25	THR
2	B	69	THR
2	B	74	VAL
2	B	116	MET
2	B	146	SER
2	B	148	LYS
2	B	153	THR
1	C	25	SER
1	C	36	THR
1	C	54	LEU
1	C	67	LYS
1	C	74	ARG
1	C	134	ARG
1	C	160	LEU
2	D	16	VAL

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Mol	Chain	Res	Type
2	D	32	GLU
2	D	56	ARG
2	D	63	HIS
2	D	64	GLN
2	D	79	THR
2	D	81	ASP
2	D	107	PRO
2	D	120	ARG
2	D	126	LEU
2	D	134	ARG
2	D	145	ARG
2	D	159	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
2	B	10	GLN
2	B	17	GLN
2	B	95	HIS
2	B	141	HIS
2	D	98	HIS
2	D	141	HIS

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 ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	OBV	B	1160	1	64,64,64	2.19	18 (28%)	90,104,104	2.63	32 (35%)
3	OBV	C	1163	-	64,64,64	1.96	15 (23%)	90,104,104	2.59	40 (44%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	OBV	B	1160	1	-	0/30/138/138	0/0/8/8
3	OBV	C	1163	-	-	0/30/138/138	0/0/8/8

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1160	OBV	C4A-NA	-8.24	1.25	1.36
3	C	1163	OBV	C4A-NA	-7.52	1.26	1.36
3	B	1160	OBV	C2D-C3D	5.32	1.48	1.36
3	C	1163	OBV	C2D-C3D	4.92	1.47	1.36
3	B	1160	OBV	C1C-C2C	4.02	1.50	1.43
3	B	1160	OBV	C3C-C2C	3.90	1.49	1.37
3	C	1163	OBV	FE-ND	3.83	2.11	1.96
3	B	1160	OBV	C4C-C3C	3.76	1.49	1.43
3	C	1163	OBV	CHD-C1D	3.67	1.47	1.35
3	B	1160	OBV	CHD-C1D	3.67	1.47	1.35
3	C	1163	OBV	C1D-C2D	3.54	1.51	1.45
3	B	1160	OBV	C1C-CHC	3.51	1.49	1.39
3	B	1160	OBV	CHB-C4A	3.41	1.44	1.36
3	B	1160	OBV	C1D-ND	-3.39	1.32	1.39
3	C	1163	OBV	C3C-C2C	3.38	1.47	1.37
3	B	1160	OBV	C4C-CHD	3.27	1.48	1.39
3	C	1163	OBV	C1C-C2C	3.24	1.48	1.43
3	B	1160	OBV	FE-NB	-3.19	1.83	1.96
3	B	1160	OBV	C4B-NB	-3.12	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1160	OBV	CHA-C4D	3.10	1.46	1.39
3	B	1160	OBV	FE-ND	3.04	2.08	1.96
3	C	1163	OBV	CHB-C1B	-3.01	1.32	1.41
3	C	1163	OBV	C4C-CHD	3.00	1.48	1.39
3	C	1163	OBV	CHB-C4A	2.90	1.43	1.36
3	C	1163	OBV	CHA-C4D	2.85	1.46	1.39
3	C	1163	OBV	C4C-C3C	2.83	1.48	1.43
3	B	1160	OBV	C1D-C2D	2.60	1.49	1.45
3	C	1163	OBV	C4D-C3D	2.53	1.49	1.44
3	C	1163	OBV	C1C-CHC	2.43	1.46	1.39
3	B	1160	OBV	CHB-C1B	-2.40	1.34	1.41
3	B	1160	OBV	CHC-C4B	2.32	1.36	1.33
3	B	1160	OBV	C4D-ND	-2.16	1.34	1.38
3	C	1163	OBV	C4D-ND	-2.11	1.34	1.38

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1160	OBV	C3A-C2A-C1A	-8.49	90.53	100.57
3	C	1163	OBV	C2C-C1C-NC	7.82	114.72	109.50
3	B	1160	OBV	C4A-CHB-C1B	-7.23	111.75	125.84
3	C	1163	OBV	C1C-C2C-C3C	-6.72	102.32	107.00
3	B	1160	OBV	CDB-C2B-C3B	6.53	126.32	108.18
3	C	1163	OBV	C3D-C4D-ND	5.98	117.51	109.99
3	B	1160	OBV	C3B-C2B-C1B	-5.71	91.00	101.07
3	B	1160	OBV	C4A-NA-C1A	-5.58	104.16	107.94
3	B	1160	OBV	CDA-C2A-C3A	5.52	123.52	108.18
3	C	1163	OBV	C4A-NA-C1A	-5.45	104.25	107.94
3	C	1163	OBV	C2D-C1D-ND	5.34	116.37	110.14
3	B	1160	OBV	C3D-C4D-ND	5.05	116.34	109.99
3	B	1160	OBV	CDA-C2A-C1A	4.90	115.06	104.94
3	B	1160	OBV	C4D-C3D-C2D	-4.81	100.90	106.96
3	B	1160	OBV	C2C-C1C-NC	4.67	112.62	109.50
3	C	1163	OBV	O2C-CCC-CBC	-4.67	107.28	123.06
3	B	1160	OBV	C4B-NB-C1B	4.65	109.30	106.43
3	B	1160	OBV	C4C-C3C-C2C	-4.33	102.43	106.92
3	C	1163	OBV	CAB-C3B-C4B	4.17	118.20	111.19
3	C	1163	OBV	CHC-C4B-NB	4.15	128.50	123.98
3	B	1160	OBV	C1C-C2C-C3C	-4.13	104.12	107.00
3	C	1163	OBV	C2A-C1A-CHA	-4.05	117.71	123.40
3	C	1163	OBV	C3A-C4A-NA	3.96	117.69	110.78
3	C	1163	OBV	C4D-C3D-C2D	-3.80	102.18	106.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1163	OBV	C2B-CDB-CEB	-3.73	106.08	115.19
3	C	1163	OBV	C2B-C1B-CHB	-3.72	116.59	123.46
3	C	1163	OBV	CBC-CAC-C3C	-3.69	106.48	112.63
3	B	1160	OBV	CAB-C3B-C4B	-3.68	105.00	111.19
3	C	1163	OBV	CDB-C2B-C3B	3.68	118.39	108.18
3	C	1163	OBV	CDA-C2A-C1A	3.64	112.46	104.94
3	C	1163	OBV	CDD-C3D-C4D	3.61	130.10	124.95
3	B	1160	OBV	C1C-CHC-C4B	-3.60	122.99	130.12
3	C	1163	OBV	CHA-C4D-ND	-3.52	120.38	124.45
3	B	1160	OBV	CHB-C4A-NA	-3.43	121.27	125.48
3	C	1163	OBV	C4C-C3C-C2C	-3.30	103.51	106.92
3	B	1160	OBV	CHC-C4B-NB	3.20	127.47	123.98
3	C	1163	OBV	CAD-C2D-C1D	3.12	130.12	124.51
3	C	1163	OBV	C1D-C2D-C3D	-3.11	101.83	106.93
3	C	1163	OBV	C1D-ND-C4D	-3.00	101.99	105.11
3	B	1160	OBV	CAD-C2D-C1D	3.00	129.91	124.51
3	C	1163	OBV	O1C-CCC-CBC	2.94	124.36	114.19
3	C	1163	OBV	CDA-C2A-C3A	2.94	116.34	108.18
3	B	1160	OBV	CBC-CAC-C3C	2.91	117.47	112.63
3	B	1160	OBV	CDD-C3D-C4D	2.89	129.07	124.95
3	C	1163	OBV	CHD-C1D-ND	-2.89	120.21	124.28
3	B	1160	OBV	CHC-C1C-NC	-2.82	119.92	124.70
3	C	1163	OBV	C1C-CHC-C4B	-2.81	124.54	130.12
3	C	1163	OBV	C3C-C4C-NC	2.77	113.59	109.73
3	B	1160	OBV	CMA-C2A-CDA	-2.73	106.46	110.78
3	B	1160	OBV	CMB-C2B-CDB	-2.69	106.51	110.78
3	B	1160	OBV	CHD-C1D-ND	-2.67	120.53	124.28
3	C	1163	OBV	C3C-C4C-CHD	-2.66	120.95	126.00
3	B	1160	OBV	C3B-C4B-NB	-2.65	104.83	111.22
3	B	1160	OBV	C2D-C1D-ND	2.59	113.17	110.14
3	B	1160	OBV	C3C-C4C-NC	2.57	113.31	109.73
3	B	1160	OBV	CMB-C2B-C3B	-2.54	107.31	112.15
3	C	1163	OBV	O1D-CCD-CBD	2.49	122.79	114.19
3	C	1163	OBV	CHA-C1A-NA	2.48	128.16	124.09
3	B	1160	OBV	C2B-C1B-NB	-2.47	106.27	111.81
3	B	1160	OBV	CHA-C4D-C3D	-2.43	120.80	124.86
3	C	1163	OBV	CAA-C3A-C4A	-2.38	106.04	111.05
3	B	1160	OBV	C2A-C3A-C4A	-2.30	96.99	101.20
3	C	1163	OBV	C4A-CHB-C1B	-2.28	121.40	125.84
3	C	1163	OBV	CAC-CBC-CCC	-2.27	106.27	113.66
3	C	1163	OBV	C3A-C4A-CHB	-2.21	117.29	123.13
3	C	1163	OBV	O3B-CEB-CDB	-2.18	115.85	122.75

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1163	OBV	CHC-C1C-NC	-2.17	121.02	124.70
3	C	1163	OBV	CAD-CBD-CCD	-2.16	109.45	113.53
3	C	1163	OBV	CMA-C2A-C3A	-2.15	108.05	112.15
3	C	1163	OBV	CDC-C2C-C3C	2.11	128.92	124.94
3	C	1163	OBV	CMB-C2B-C3B	-2.07	108.20	112.15
3	B	1160	OBV	CAC-C3C-C4C	2.04	131.65	125.50

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	145/173 (83%)	0.19	6 (4%) 35 36	23, 35, 55, 107	0
1	C	139/173 (80%)	0.09	2 (1%) 72 74	22, 35, 53, 63	0
2	B	159/179 (88%)	0.72	17 (10%) 6 6	23, 39, 93, 150	0
2	D	157/179 (87%)	0.19	7 (4%) 32 33	23, 34, 69, 115	0
All	All	600/704 (85%)	0.31	32 (5%) 25 26	22, 36, 67, 150	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	68	TRP	14.0
1	A	163	ILE	12.8
2	B	67	GLY	10.2
2	B	66	THR	7.2
1	A	164	ARG	6.4
2	B	63	HIS	6.3
2	B	65	LYS	6.0
2	B	72	ALA	4.6
1	A	165	VAL	4.3
2	B	71	ASN	4.3
2	B	64	GLN	4.1
2	D	3	HIS	3.9
2	D	64	GLN	3.2
2	D	63	HIS	3.2
2	D	65	LYS	3.1
2	B	119	GLY	3.1
2	B	69	THR	3.0
2	B	70	HIS	3.0
2	B	118	HIS	3.0
2	B	98	HIS	2.8
2	B	120	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
2	B	95	HIS	2.7
1	A	162	LYS	2.5
2	B	122	GLU	2.5
1	C	160	LEU	2.4
2	D	62	LYS	2.4
2	D	68	TRP	2.4
2	D	61	ILE	2.3
2	B	116	MET	2.2
1	A	71	ILE	2.1
1	A	166	ASP	2.1
1	C	73	ARG	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 ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	OBV	C	1163	57/57	0.31	3.33	30,50,74,88	57
3	OBV	B	1160	57/57	0.42	2.45	41,56,71,77	57

6.5 Other polymers ⓘ

There are no such residues in this entry.