



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:29 PM GMT

PDB ID : 5ABH
Title : Structure of GH84 with ligand
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Deposited on : 2015-08-05
Resolution : 1.95 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

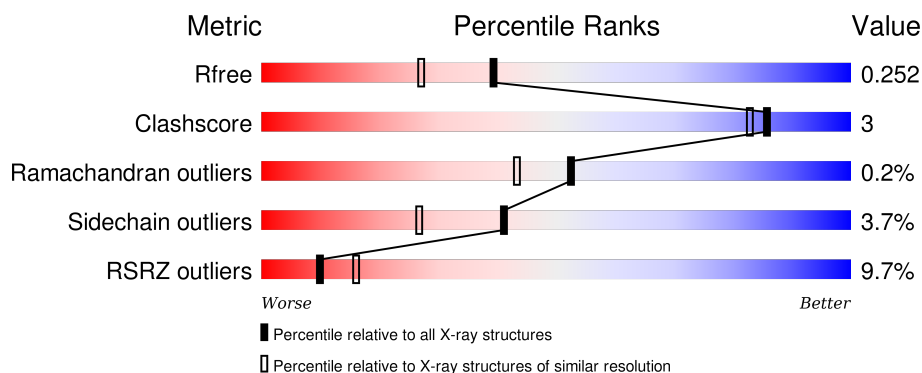
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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. 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	726	<div> <div>8%</div> <div>84%</div> <div>9%</div> <div>7%</div> </div>
1	B	726	<div> <div>10%</div> <div>83%</div> <div>9%</div> <div>8%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	YWN	A	1717	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11667 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 O-GLCNACASE BT_4395.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	677	Total	C	N	O	S	0	4	0
			5515	3529	932	1034	20			
1	B	669	Total	C	N	O	S	0	1	0
			5446	3487	921	1020	18			

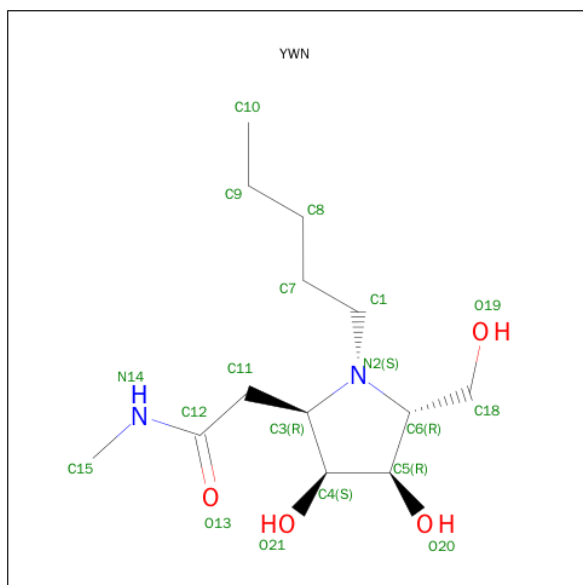
There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	MET	-	EXPRESSION TAG	UNP Q89ZI2
A	-8	GLY	-	EXPRESSION TAG	UNP Q89ZI2
A	-7	SER	-	EXPRESSION TAG	UNP Q89ZI2
A	-6	SER	-	EXPRESSION TAG	UNP Q89ZI2
A	-5	HIS	-	EXPRESSION TAG	UNP Q89ZI2
A	-4	HIS	-	EXPRESSION TAG	UNP Q89ZI2
A	-3	HIS	-	EXPRESSION TAG	UNP Q89ZI2
A	-2	HIS	-	EXPRESSION TAG	UNP Q89ZI2
A	-1	HIS	-	EXPRESSION TAG	UNP Q89ZI2
A	0	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	-9	MET	-	EXPRESSION TAG	UNP Q89ZI2
B	-8	GLY	-	EXPRESSION TAG	UNP Q89ZI2
B	-7	SER	-	EXPRESSION TAG	UNP Q89ZI2
B	-6	SER	-	EXPRESSION TAG	UNP Q89ZI2
B	-5	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	-4	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	-3	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	-2	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	-1	HIS	-	EXPRESSION TAG	UNP Q89ZI2
B	0	HIS	-	EXPRESSION TAG	UNP Q89ZI2

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0

- Molecule 3 is 2-[(2R,3S,4R,5R)-5-(HYDROXYMETHYL)-3,4-BIS(OXIDANYL)-1-PENTYL-PYRROLIDIN-2-YL]-N-METHYL-ETHANAMIDE (three-letter code: YWN) (formula: $C_{13}H_{26}N_2O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 19 13 2 4	0	0
3	B	1	Total C N O 19 13 2 4	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Ca	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	352	Total	O	0	0
			352	352		
6	B	306	Total	O	0	0
			306	306		

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 ($\text{RSRZ} > 2$). 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.

Chain A:

Position	Amino Acid	Bits
1	T670	0.08
2	V671	0.04
3	D672	0.04
4	L673	0.04
5	LYS	0.04
6	GLN	0.04
7	LYS	0.04
8	GLU	0.04
9	SER	0.04
10	ARG	0.04
11	LEU	0.04
12	S681	0.04
13	A682	0.04
14	V689	0.04
15	ILE	0.04
16	K603	0.04
17	H604	0.04
18	L605	0.04
19	ASN	0.04
20	VAL	0.04
21	SER	0.04
22	ASP	0.04
23	GLU	0.04
24	N620	0.04
25	E621	0.04
26	VAL	0.04
27	TYR	0.04
28	LEU	0.04
29	ARG	0.04
30	GLN	0.04
31	F709	0.04
32	K745	0.04
33	LYS	0.04
34	R519	0.04
35	K562	0.04
36	M591	0.04
37	P592	0.04
38	H593	0.04
39	K594	0.04
40	M595	0.04
41	I596	0.04
42	S597	0.04
43	M598	0.04
44	V599	0.04
45	GLU	0.04
46	G164	0.04
47	P165	0.04
48	P181	0.04
49	D182	0.04
50	K183	0.04
51	E184	0.04
52	K213	0.04
53	A238	0.04
54	P270	0.04
55	V276	0.04
56	M308	0.04
57	I323	0.04
58	V335	0.04
59	I336	0.04
60	H337	0.04
61	V370	0.04
62	T371	0.04
63	Y385	0.04
64	A406	0.04
65	M423	0.04
66	M441	0.04
67	D442	0.04
68	I443	0.04
69	K444	0.04
70	E448	0.04
71	G457	0.04
72	E510	0.04
73	K514	0.04
74	E519	0.04
75	K520	0.04
76	E521	0.04
77	K522	0.04
78	K53	0.04
79	L56	0.04
80	I57	0.04
81	S58	0.04
82	I59	0.04
83	G60	0.04
84	K65	0.04
85	R68	0.04
86	S71	0.04
87	R72	0.04
88	Y81	0.04
89	E87	0.04
90	K88	0.04
91	E89	0.04
92	I90	0.04
93	V91	0.04
94	I121	0.04
95	K122	0.04
96	D123	0.04
97	G131	0.04
98	Q149	0.04
99	Y153	0.04
100	S4	0.04

[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	51.41Å 161.31Å 223.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	130.74 – 1.95 65.37 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.9 (130.74-1.95) 100.0 (65.37-1.95)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0124	Depositor
R, R_{free}	0.212 , 0.252 0.217 , 0.252	Depositor DCC
R_{free} test set	6816 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	0.801	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	5 of 136405 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11667	wwPDB-VP
Average B, all atoms (Å ²)	42.0	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 20.81 % 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 8.0175e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: YWN, CA, EDO, CL

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	0.57	1/5649 (0.0%)	0.67	2/7650 (0.0%)
1	B	0.60	1/5579 (0.0%)	0.70	4/7555 (0.1%)
All	All	0.59	2/11228 (0.0%)	0.68	6/15205 (0.0%)

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	A	0	1
1	B	0	5
All	All	0	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	442	ASP	CB-CG	6.22	1.64	1.51
1	B	442	ASP	CB-CG	5.03	1.62	1.51

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	258	LEU	CA-CB-CG	-7.11	98.96	115.30
1	B	442	ASP	CB-CG-OD1	7.08	124.68	118.30
1	B	168	ASP	CB-CG-OD1	5.94	123.64	118.30
1	A	442	ASP	CB-CG-OD1	5.79	123.51	118.30
1	B	168	ASP	CB-CG-OD2	-5.35	113.49	118.30
1	A	689	VAL	CB-CA-C	-5.07	101.76	111.40

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	270	PRO	Peptide
1	B	46	SER	Peptide
1	B	646	ILE	Peptide
1	B	663	THR	Peptide
1	B	673	LEU	Peptide
1	B	674	LYS	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5515	0	5431	36	0
1	B	5446	0	5363	25	0
2	A	1	0	0	0	0
3	A	19	0	26	0	0
3	B	19	0	26	0	0
4	A	4	0	6	0	0
4	B	4	0	6	0	0
5	B	1	0	0	1	0
6	A	352	0	0	2	0
6	B	306	0	0	3	0
All	All	11667	0	10858	62	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (62) 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:A:423[A]:MET:CE	1:A:441[A]:MET:HG3	2.18	0.74
5:B:1716:CA:CA	6:B:2034:HOH:O	1.67	0.71
1:A:423[A]:MET:HE1	1:A:441[A]:MET:HG3	1.74	0.70
1:B:223:LEU:CD1	1:B:260:TYR:HE2	2.08	0.67
1:A:11:GLN:HG2	1:A:122:LYS:HB2	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:VAL:O	1:B:399:THR:HG23	1.95	0.66
1:B:454:PHE:C	1:B:456:GLU:O	2.36	0.63
1:A:597:SER:HB3	1:A:631:VAL:HA	1.83	0.61
1:A:598:ASN:HB3	1:A:599:VAL:HG23	1.85	0.58
1:B:70:TYR:OH	1:B:89:GLU:OE2	2.19	0.57
1:B:109:GLN:O	1:B:112:LYS:HE2	2.05	0.56
1:B:223:LEU:HD11	1:B:260:TYR:CE2	2.44	0.54
1:A:457:GLY:HA2	6:A:2250:HOH:O	2.07	0.53
1:A:597:SER:CB	1:A:631:VAL:HA	2.39	0.53
1:A:444[B]:GLN:NE2	1:A:448:GLU:OE1	2.41	0.53
1:B:661:ILE:HD12	1:B:692:VAL:HG13	1.91	0.52
1:B:661:ILE:HG23	1:B:662:SER:N	2.24	0.52
1:B:662:SER:HB3	1:B:691:PHE:HB2	1.93	0.51
1:B:81:TYR:CE2	1:B:123:ASP:HB3	2.45	0.51
1:A:181:PRO:HG2	1:A:184:GLU:OE1	2.11	0.50
1:A:661:ILE:HD12	1:A:671:VAL:HG21	1.93	0.50
1:B:128:ARG:HD2	6:B:2050:HOH:O	2.12	0.50
1:A:18:THR:O	1:A:19:ILE:HD12	2.14	0.48
1:A:597:SER:OG	1:A:632:GLU:N	2.47	0.47
1:A:385:TYR:CD1	1:A:406:ALA:HB2	2.50	0.47
1:A:423[A]:MET:HE2	1:A:441[A]:MET:H	1.80	0.47
1:A:238:ALA:HA	1:A:276:VAL:O	2.13	0.47
1:B:223:LEU:CD1	1:B:260:TYR:CE2	2.91	0.47
1:A:131:GLY:O	1:A:370:VAL:HA	2.15	0.46
1:B:308:MET:HA	1:B:335:TYR:O	2.15	0.46
1:B:81:TYR:CZ	1:B:123:ASP:HB3	2.50	0.46
1:A:28:ASN:HB2	1:A:56:LEU:HD11	1.98	0.46
1:A:308:MET:HA	1:A:335:TYR:O	2.16	0.46
1:B:131:GLY:HA3	1:B:160:THR:O	2.17	0.45
1:A:595:MET:O	1:A:595:MET:SD	2.75	0.45
1:B:354:PRO:HB2	1:B:399:THR:HG22	1.98	0.45
1:A:605:LEU:HD13	1:A:617:SER:O	2.17	0.45
1:A:11:GLN:HG3	1:A:11:GLN:O	2.16	0.44
1:B:661:ILE:CD1	1:B:692:VAL:HG13	2.48	0.44
1:B:238:ALA:HA	1:B:276:VAL:O	2.17	0.44
1:B:557:ALA:HB1	1:B:561:ILE:HB	1.99	0.44
1:A:12:LEU:HD13	1:A:121:ILE:HG12	1.99	0.44
1:A:562[A]:LYS:HD3	6:A:2281:HOH:O	2.16	0.44
1:A:164:GLY:N	1:A:165:PRO:CD	2.81	0.43
1:A:81:TYR:CZ	1:A:123:ASP:HB3	2.54	0.43
1:B:356:TYR:HB3	1:B:399:THR:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:LYS:HD2	1:A:89:GLU:HB3	2.01	0.43
1:A:423[A]:MET:CE	1:A:441[A]:MET:H	2.32	0.43
1:B:385:TYR:CD1	1:B:406:ALA:HB2	2.54	0.43
1:A:423[A]:MET:O	1:A:423[A]:MET:HG3	2.19	0.42
1:B:225:LYS:O	1:B:229:MET:HG2	2.19	0.42
1:B:349:HIS:HE1	6:B:2078:HOH:O	2.02	0.42
1:A:12:LEU:HD13	1:A:121:ILE:CG1	2.49	0.42
1:A:510:GLU:O	1:A:514:LYS:HG3	2.19	0.42
1:A:33:ALA:HB2	1:A:60:GLY:HA2	2.02	0.42
1:A:591:MET:HG3	1:A:593:HIS:O	2.20	0.41
1:B:223:LEU:HD12	1:B:260:TYR:HE2	1.85	0.41
1:A:28:ASN:HB3	1:A:58:SER:HA	2.01	0.41
1:A:81:TYR:CE2	1:A:123:ASP:HB3	2.56	0.41
1:A:149:GLN:HB3	1:A:153:TYR:CZ	2.56	0.41
1:A:12:LEU:HD12	1:A:13:ILE:N	2.37	0.40
1:B:591:MET:SD	1:B:592:PRO:HD2	2.61	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	669/726 (92%)	640 (96%)	29 (4%)	0	100	100
1	B	658/726 (91%)	628 (95%)	28 (4%)	2 (0%)	46	35
All	All	1327/1452 (91%)	1268 (96%)	57 (4%)	2 (0%)	52	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	518	GLY
1	B	664	ASP

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	598/639 (94%)	580 (97%)	18 (3%)	48	36
1	B	591/639 (92%)	565 (96%)	26 (4%)	35	19
All	All	1189/1278 (93%)	1145 (96%)	44 (4%)	41	27

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	17	LYS
1	A	32	GLU
1	A	49	GLN
1	A	68	ARG
1	A	71	SER
1	A	88	LYS
1	A	183	LYS
1	A	213	LYS
1	A	323	ILE
1	A	337	TRP
1	A	371	THR
1	A	519	ARG
1	A	620	ASN
1	A	648	PHE
1	A	650	LYS
1	A	661	ILE
1	A	666	LYS
1	B	17	LYS
1	B	18	THR
1	B	62	LYS
1	B	112	LYS
1	B	258	LEU
1	B	290	ASN
1	B	292	ASN
1	B	300	LYS
1	B	324	SER

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Mol	Chain	Res	Type
1	B	337	TRP
1	B	371	THR
1	B	408	ARG
1	B	438	GLU
1	B	448	GLU
1	B	496	THR
1	B	564	LEU
1	B	574	LYS
1	B	642	GLU
1	B	660	GLU
1	B	661	ILE
1	B	664	ASP
1	B	666	LYS
1	B	672	ASP
1	B	676	LYS
1	B	694	PHE
1	B	715	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	620	ASN
1	B	349	HIS
1	B	433	HIS
1	B	459	ASN
1	B	608	GLN
1	B	685	GLN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 for Mogul analysis.

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	YWN	A	1717	-	19,19,19	0.80	1 (5%)	16,25,25	0.51	0
4	EDO	A	1718	-	3,3,3	0.58	0	2,2,2	0.48	0
3	YWN	B	1717	-	19,19,19	0.75	0	16,25,25	0.81	0
4	EDO	B	1718	-	3,3,3	0.52	0	2,2,2	0.54	0

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	YWN	A	1717	-	-	0/13/33/33	0/1/1/1
4	EDO	A	1718	-	-	0/1/1/1	0/0/0/0
3	YWN	B	1717	-	-	0/13/33/33	0/1/1/1
4	EDO	B	1718	-	-	0/1/1/1	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1717	YWN	C4-C3	-2.17	1.50	1.53

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	677/726 (93%)	0.47	58 (8%)	13 21	24, 38, 78, 102	0
1	B	669/726 (92%)	0.49	72 (10%)	8 12	22, 35, 76, 97	0
All	All	1346/1452 (92%)	0.48	130 (9%)	10 16	22, 37, 77, 102	0

All (130) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	668	TRP	9.7
1	B	673	LEU	9.5
1	B	706	LEU	8.6
1	A	599	VAL	7.6
1	A	653	PRO	7.1
1	A	673	LEU	6.9
1	A	646	ILE	6.4
1	B	605	LEU	6.1
1	B	53	LYS	5.9
1	A	671	VAL	5.7
1	B	618	PRO	5.6
1	B	680	LEU	5.6
1	A	668	TRP	5.3
1	B	679	ARG	5.3
1	B	631	VAL	5.3
1	B	661	ILE	5.3
1	B	290	ASN	5.2
1	A	672	ASP	5.2
1	B	685	GLN	4.7
1	B	676	LYS	4.7
1	B	671	VAL	4.6
1	B	688	PRO	4.6
1	A	709	PHE	4.6
1	A	651	ASP	4.6

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Mol	Chain	Res	Type	RSRZ
1	B	715	LYS	4.6
1	A	619	ALA	4.5
1	B	709	PHE	4.5
1	B	646	ILE	4.4
1	A	620	ASN	4.4
1	A	659	LEU	4.3
1	B	674	LYS	4.2
1	B	48	LYS	4.2
1	B	687	ALA	4.2
1	B	678	SER	4.2
1	A	616	ILE	4.2
1	B	707	ARG	4.1
1	A	667	GLU	4.1
1	A	694	PHE	4.1
1	A	631	VAL	4.1
1	B	666	LYS	4.0
1	B	644	ILE	3.9
1	A	658	ARG	3.9
1	B	672	ASP	3.9
1	A	666	LYS	3.9
1	A	88	LYS	3.9
1	B	669	LYS	3.9
1	A	652	ALA	3.7
1	B	684	LEU	3.7
1	A	53	LYS	3.7
1	A	604	ASN	3.7
1	B	616	ILE	3.7
1	B	286	TRP	3.6
1	B	693	ARG	3.5
1	B	24	VAL	3.5
1	B	689	VAL	3.4
1	A	598	ASN	3.4
1	A	16	ASN	3.4
1	B	694	PHE	3.4
1	A	665	GLY	3.4
1	B	663	THR	3.4
1	B	647	ASN	3.3
1	A	27	LEU	3.3
1	B	617	SER	3.3
1	B	606	PRO	3.3
1	A	49	GLN	3.2
1	B	660	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	597	SER	3.2
1	A	65	LYS	3.2
1	A	50	SER	3.1
1	A	648	PHE	3.1
1	B	692	VAL	3.1
1	B	633	ILE	3.1
1	B	46	SER	3.1
1	A	660	GLU	3.1
1	B	677	GLU	3.1
1	A	52	LYS	3.1
1	B	708	GLN	3.1
1	B	51	SER	3.0
1	A	661	ILE	3.0
1	B	683	GLY	3.0
1	B	670	THR	3.0
1	B	682	ALA	2.9
1	A	24	VAL	2.9
1	B	681	SER	2.9
1	A	670	THR	2.9
1	A	72	ARG	2.8
1	B	2	ASN	2.8
1	B	645	GLN	2.8
1	A	68	ARG	2.8
1	A	693	ARG	2.8
1	A	47	GLY	2.8
1	A	596	ILE	2.7
1	A	695	THR	2.7
1	B	291	GLY	2.7
1	B	638	ILE	2.7
1	B	675	GLN	2.7
1	A	650	LYS	2.6
1	B	567	ARG	2.6
1	A	649	GLY	2.6
1	A	17	LYS	2.6
1	B	52	LYS	2.6
1	B	49	GLN	2.5
1	A	682	ALA	2.5
1	A	657	GLY	2.4
1	A	423[A]	MET	2.4
1	B	632	GLU	2.4
1	A	647	ASN	2.4
1	B	16	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	87	GLU	2.4
1	B	643	ASN	2.4
1	B	667	GLU	2.3
1	A	21	LEU	2.3
1	B	451	LEU	2.3
1	B	25	TYR	2.3
1	A	91	VAL	2.3
1	B	441	MET	2.3
1	B	54	GLY	2.3
1	B	613	ARG	2.3
1	A	14	VAL	2.2
1	A	605	LEU	2.2
1	A	669	LYS	2.2
1	B	21	LEU	2.1
1	B	454	PHE	2.1
1	A	595	MET	2.1
1	B	595	MET	2.1
1	A	621	GLU	2.0
1	B	640	PRO	2.0
1	B	642	GLU	2.0
1	A	51	SER	2.0
1	B	583	HIS	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](#)

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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	YWN	A	1717	19/19	0.95	0.17	3.21	26,36,51,59	0
4	EDO	B	1718	4/4	0.96	0.13	1.32	35,40,41,42	0
4	EDO	A	1718	4/4	0.96	0.16	1.30	33,34,36,39	0
3	YWN	B	1717	19/19	0.95	0.16	1.27	29,36,50,54	0
5	CA	B	1716	1/1	0.99	0.09	-0.09	36,36,36,36	0
2	CL	A	1716	1/1	0.94	0.10	-1.82	69,69,69,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.