



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:02 AM UTC

PDB ID : 6GRK / pdb_00006grk
Title : Structure of the soluble AhlB of the tripartite alpha-pore forming toxin, AHL, from *Aeromonas hydrophila*.
Authors : Churchill-Angus, A.M.; Wilson, J.S.; Baker, P.J.
Deposited on : 2018-06-11
Resolution : 2.33 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

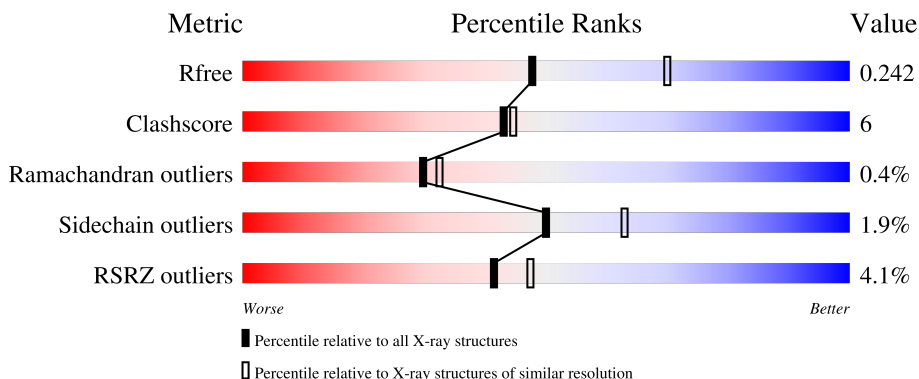
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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}	180053	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	367	3% 87% 9% ..
1	B	367	7% 78% 16% 5%
1	I	367	2% 82% 13% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7875 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 AhlB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	0	0
			2603	1625	452	518	8			
1	I	354	Total	C	N	O	S	0	0	0
			2607	1627	452	520	8			
1	B	347	Total	C	N	O	S	0	0	0
			2564	1600	445	511	8			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	360	LEU	-	expression tag	UNP A0A081US78
A	361	GLU	-	expression tag	UNP A0A081US78
A	362	HIS	-	expression tag	UNP A0A081US78
A	363	HIS	-	expression tag	UNP A0A081US78
A	364	HIS	-	expression tag	UNP A0A081US78
A	365	HIS	-	expression tag	UNP A0A081US78
A	366	HIS	-	expression tag	UNP A0A081US78
A	367	HIS	-	expression tag	UNP A0A081US78
I	360	LEU	-	expression tag	UNP A0A081US78
I	361	GLU	-	expression tag	UNP A0A081US78
I	362	HIS	-	expression tag	UNP A0A081US78
I	363	HIS	-	expression tag	UNP A0A081US78
I	364	HIS	-	expression tag	UNP A0A081US78
I	365	HIS	-	expression tag	UNP A0A081US78
I	366	HIS	-	expression tag	UNP A0A081US78
I	367	HIS	-	expression tag	UNP A0A081US78
B	360	LEU	-	expression tag	UNP A0A081US78
B	361	GLU	-	expression tag	UNP A0A081US78
B	362	HIS	-	expression tag	UNP A0A081US78
B	363	HIS	-	expression tag	UNP A0A081US78
B	364	HIS	-	expression tag	UNP A0A081US78
B	365	HIS	-	expression tag	UNP A0A081US78
B	366	HIS	-	expression tag	UNP A0A081US78

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Chain	Residue	Modelled	Actual	Comment	Reference
B	367	HIS	-	expression tag	UNP A0A081US78

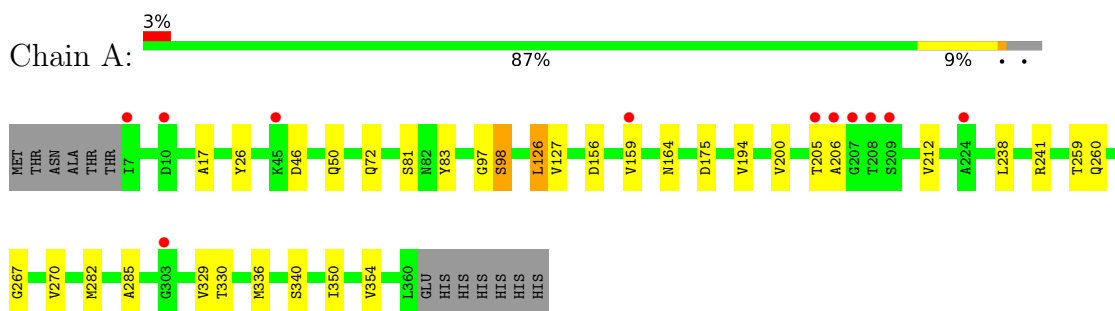
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total 49	O 49	0	0
2	I	39	Total 39	O 39	0	0
2	B	13	Total 13	O 13	0	0

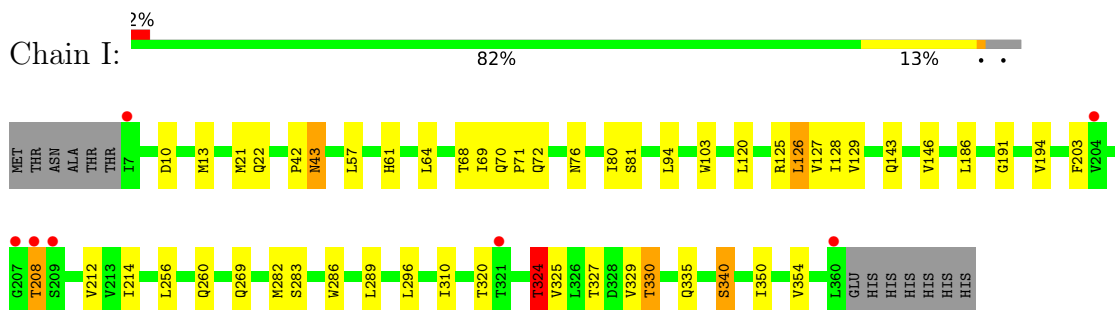
3 Residue-property plots [i](#)

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

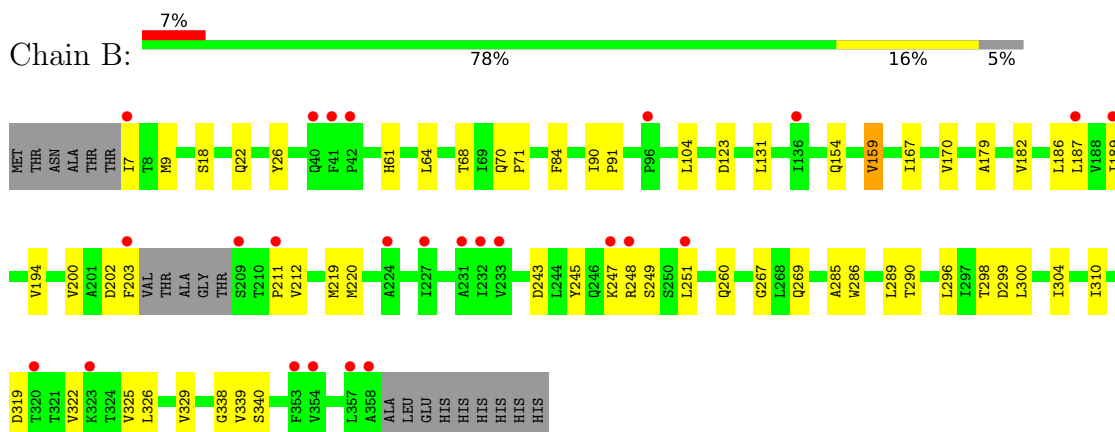
• Molecule 1: AhlB



• Molecule 1: AhlB



• Molecule 1: AhlB



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	133.47Å 79.81Å 111.01Å 90.00° 90.22° 90.00°	Depositor
Resolution (Å)	58.41 – 2.33 58.41 – 2.33	Depositor EDS
% Data completeness (in resolution range)	98.0 (58.41-2.33) 98.0 (58.41-2.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0230	Depositor
R, R_{free}	0.222 , 0.254 (Not available) , 0.242	Depositor DCC
R_{free} test set	2408 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 19.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.024 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.022 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.059 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.076 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.029 for -h,-k,l	Xtriage
Reported twinning fraction	0.895 for H, K, L 0.105 for -1/2H+3/2K, -1/2H-1/2K, L	Depositor
Outliers	0 of 49042 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7875	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	0.86	0/2632	1.23	3/3589 (0.1%)
1	B	0.78	0/2592	1.21	2/3531 (0.1%)
1	I	0.92	2/2636 (0.1%)	1.21	1/3594 (0.0%)
All	All	0.86	2/7860 (0.0%)	1.22	6/10714 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	214	ILE	C-O	-6.96	1.17	1.24
1	I	286	TRP	C-O	-5.09	1.18	1.24

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	THR	CB-CA-C	8.54	124.29	110.88
1	A	206	ALA	N-CA-C	-5.95	102.85	110.53
1	B	123	ASP	CA-CB-CG	5.93	118.53	112.60
1	A	175	ASP	CA-CB-CG	5.38	117.98	112.60
1	I	324	THR	CB-CA-C	5.09	118.86	110.88
1	B	319	ASP	CA-CB-CG	5.09	117.69	112.60

There are no chirality outliers.

There are no planarity outliers.

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	2603	0	2637	25	0
1	B	2564	0	2593	36	0
1	I	2607	0	2641	36	0
2	A	49	0	0	1	0
2	B	13	0	0	0	0
2	I	39	0	0	1	0
All	All	7875	0	7871	97	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 6.

All (97) 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:I:13:MET:HA	1:I:282:MET:HE1	1.49	0.92
1:A:81:SER:HB2	1:A:329:VAL:HG11	1.54	0.88
1:B:9:MET:HE1	1:B:286:TRP:HA	1.59	0.85
1:I:81:SER:HB3	1:I:329:VAL:HG11	1.62	0.81
1:A:46:ASP:O	1:A:50:GLN:HG2	1.86	0.75
1:I:76:ASN:ND2	1:I:282:MET:HE3	2.03	0.73
1:A:126:LEU:C	1:A:126:LEU:HD23	2.18	0.69
1:I:42:PRO:O	1:I:43:ASN:CB	2.40	0.68
1:I:191:GLY:C	1:I:335:GLN:HE21	2.02	0.68
1:I:13:MET:HA	1:I:282:MET:CE	2.22	0.66
1:I:126:LEU:C	1:I:126:LEU:HD23	2.21	0.66
1:I:13:MET:CA	1:I:282:MET:HE1	2.26	0.65
1:B:202:ASP:O	1:B:203:PHE:CD1	2.51	0.64
1:A:238:LEU:HD12	1:A:354:VAL:HG21	1.79	0.64
1:I:194:VAL:HA	1:I:340:SER:O	1.97	0.64
1:I:76:ASN:HD21	1:I:282:MET:HE3	1.62	0.63
1:I:42:PRO:O	1:I:43:ASN:HB3	1.96	0.63
1:I:327:THR:O	1:I:330:THR:HG22	1.99	0.62
1:I:320:THR:O	1:I:324:THR:HG22	1.99	0.61
1:B:200:VAL:HG23	1:B:212:VAL:O	2.01	0.60
1:A:17:ALA:HA	1:A:336:MET:CE	2.31	0.59
1:B:18:SER:HB2	1:B:182:VAL:HG23	1.83	0.59
1:A:17:ALA:HB1	1:A:336:MET:CE	2.35	0.57
1:B:159:VAL:HG13	1:B:251:LEU:HD11	1.87	0.57
1:A:270:VAL:CG2	2:A:434:HOH:O	2.52	0.56
1:B:104:LEU:HD23	1:B:300:LEU:HD23	1.87	0.56
1:I:289:LEU:C	1:I:289:LEU:HD23	2.31	0.56
1:A:194:VAL:HA	1:A:340:SER:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:MET:HE3	1:A:285:ALA:HB3	1.87	0.56
1:A:17:ALA:CB	1:A:336:MET:CE	2.83	0.55
1:B:296:LEU:HD11	1:B:310:ILE:HG12	1.89	0.54
1:I:13:MET:SD	1:I:329:VAL:HG22	2.47	0.54
1:A:17:ALA:HA	1:A:336:MET:HE1	1.88	0.54
1:I:212:VAL:HG11	1:I:260:GLN:HB3	1.90	0.53
1:I:22:GLN:NE2	2:I:403:HOH:O	2.41	0.53
1:B:9:MET:HE3	1:B:289:LEU:CD2	2.39	0.52
1:I:203:PHE:O	1:I:208:THR:O	2.27	0.52
1:B:9:MET:HE3	1:B:289:LEU:HD22	1.92	0.52
1:B:219:MET:O	1:B:220:MET:HG3	2.10	0.51
1:B:212:VAL:HG11	1:B:260:GLN:HB3	1.93	0.50
1:B:64:LEU:HD23	1:B:131:LEU:HD13	1.92	0.49
1:A:350:ILE:O	1:A:354:VAL:HG13	2.13	0.49
1:I:186:LEU:HA	1:I:335:GLN:OE1	2.12	0.49
1:B:325:VAL:O	1:B:329:VAL:HG12	2.13	0.48
1:B:219:MET:C	1:B:220:MET:HG3	2.38	0.48
1:A:17:ALA:HB1	1:A:336:MET:HE3	1.95	0.48
1:B:299:ASP:OD1	1:B:304:ILE:HD12	2.13	0.48
1:I:70:GLN:HB3	1:I:71:PRO:HD3	1.95	0.47
1:B:167:ILE:O	1:B:170:VAL:N	2.47	0.47
1:B:22:GLN:OE1	1:B:179:ALA:HA	2.15	0.47
1:I:72:GLN:OE1	1:I:120:LEU:HD11	2.14	0.47
1:B:194:VAL:HA	1:B:340:SER:O	2.15	0.47
1:I:296:LEU:HD11	1:I:310:ILE:HG12	1.98	0.46
1:I:350:ILE:O	1:I:354:VAL:HG23	2.14	0.46
1:B:64:LEU:HD12	1:B:68:THR:HB	1.97	0.46
1:I:76:ASN:HD21	1:I:282:MET:HG2	1.81	0.46
1:B:9:MET:CE	1:B:286:TRP:HA	2.37	0.46
1:B:248:ARG:NH2	1:B:249:SER:O	2.49	0.45
1:A:83:TYR:CD2	1:A:83:TYR:C	2.93	0.45
1:A:238:LEU:CD1	1:A:354:VAL:HG21	2.44	0.45
1:B:203:PHE:CE1	1:B:211:PRO:HB3	2.52	0.45
1:B:212:VAL:HG21	1:B:260:GLN:HB2	1.99	0.45
1:I:94:LEU:HD23	1:I:103:TRP:CE2	2.52	0.45
1:I:64:LEU:O	1:I:68:THR:HB	2.17	0.45
1:A:26:TYR:OH	1:A:267:GLY:HA3	2.16	0.45
1:B:186:LEU:O	1:B:187:LEU:HB2	2.15	0.45
1:I:21:MET:HE1	1:I:186:LEU:HD13	1.98	0.44
1:I:94:LEU:HD23	1:I:103:TRP:CD2	2.52	0.44
1:B:26:TYR:OH	1:B:267:GLY:HA3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:126:LEU:HD23	1:I:127:VAL:N	2.32	0.44
1:B:338:GLY:O	1:B:339:VAL:C	2.59	0.44
1:A:17:ALA:CA	1:A:336:MET:CE	2.94	0.44
1:B:7:ILE:HG23	1:B:7:ILE:O	2.18	0.44
1:B:243:ASP:O	1:B:247:LYS:HG2	2.17	0.44
1:A:156:ASP:HB3	1:A:159:VAL:HG12	1.99	0.44
1:I:57:LEU:O	1:I:61:HIS:CD2	2.71	0.43
1:B:90:ILE:O	1:B:91:PRO:C	2.60	0.43
1:B:9:MET:HE2	1:B:285:ALA:O	2.19	0.43
1:A:97:GLY:O	1:A:98:SER:C	2.62	0.43
1:A:97:GLY:O	1:A:98:SER:O	2.36	0.43
1:I:68:THR:O	1:I:72:GLN:HG2	2.18	0.43
1:B:84:PHE:HB3	1:B:322:VAL:HG13	2.01	0.43
1:A:212:VAL:HG11	1:A:260:GLN:HB3	2.00	0.43
1:B:61:HIS:HB3	1:B:131:LEU:HD11	2.00	0.43
1:B:326:LEU:HA	1:B:329:VAL:HG12	2.01	0.42
1:B:70:GLN:HB3	1:B:71:PRO:HD3	2.00	0.42
1:A:164:ASN:OD1	1:A:259:THR:CG2	2.68	0.42
1:I:125:ARG:NH1	1:I:129:VAL:HG21	2.35	0.42
1:I:143:GLN:O	1:I:146:VAL:HG12	2.20	0.42
1:B:245:TYR:CE1	1:B:260:GLN:HG2	2.55	0.42
1:A:126:LEU:HD23	1:A:127:VAL:N	2.35	0.41
1:I:80:ILE:HG22	1:I:329:VAL:HG21	2.02	0.41
1:I:13:MET:SD	1:I:329:VAL:CG2	3.08	0.41
1:A:72:GLN:OE1	1:A:72:GLN:HA	2.20	0.41
1:I:69:ILE:HD13	1:I:128:ILE:HG12	2.01	0.41
1:B:64:LEU:O	1:B:68:THR:HB	2.21	0.41
1:A:200:VAL:HG13	1:A:241:ARG:CD	2.51	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	352/367 (96%)	336 (96%)	15 (4%)	1 (0%)	36	41
1	B	343/367 (94%)	329 (96%)	13 (4%)	1 (0%)	36	41
1	I	352/367 (96%)	338 (96%)	12 (3%)	2 (1%)	21	22
All	All	1047/1101 (95%)	1003 (96%)	40 (4%)	4 (0%)	30	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	98	SER
1	I	43	ASN
1	B	154	GLN
1	I	208	THR

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	282/295 (96%)	280 (99%)	2 (1%)	76	84
1	B	279/295 (95%)	274 (98%)	5 (2%)	51	64
1	I	283/295 (96%)	274 (97%)	9 (3%)	34	44
All	All	844/885 (95%)	828 (98%)	16 (2%)	50	63

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	LEU
1	A	330	THR
1	I	10	ASP
1	I	126	LEU
1	I	256	LEU
1	I	269	GLN
1	I	283	SER
1	I	324	THR
1	I	325	VAL

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Mol	Chain	Res	Type
1	I	330	THR
1	I	340	SER
1	B	159	VAL
1	B	189	ILE
1	B	269	GLN
1	B	290	THR
1	B	298	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	47	ASN
1	A	132	ASN
1	A	273	GLN
1	A	346	GLN
1	I	11	GLN
1	I	22	GLN
1	I	106	GLN
1	I	132	ASN
1	I	143	GLN
1	I	273	GLN
1	I	335	GLN
1	B	15	ASN
1	B	24	GLN
1	B	40	GLN
1	B	50	GLN
1	B	53	GLN
1	B	132	ASN
1	B	143	GLN
1	B	236	ASN
1	B	246	GLN
1	B	269	GLN
1	B	273	GLN
1	B	274	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	354/367 (96%)	0.32	11 (3%) 51 58	29, 45, 69, 109	0
1	B	347/367 (94%)	0.61	25 (7%) 21 25	29, 55, 81, 104	0
1	I	354/367 (96%)	0.26	7 (1%) 65 70	26, 43, 73, 107	0
All	All	1055/1101 (95%)	0.40	43 (4%) 41 47	26, 47, 79, 109	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	206	ALA	5.6
1	B	203	PHE	4.9
1	B	357	LEU	4.5
1	A	208	THR	4.4
1	I	7	ILE	3.8
1	A	7	ILE	3.4
1	I	208	THR	3.3
1	B	227	ILE	3.2
1	I	209	SER	3.1
1	B	232	ILE	3.1
1	B	248	ARG	3.0
1	I	207	GLY	3.0
1	A	207	GLY	2.9
1	A	209	SER	2.9
1	B	40	GLN	2.8
1	B	136	ILE	2.8
1	I	360	LEU	2.7
1	B	358	ALA	2.7
1	B	189	ILE	2.6
1	B	209	SER	2.6
1	B	251	LEU	2.6
1	A	205	THR	2.6
1	B	224	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	I	321	THR	2.5
1	A	10	ASP	2.5
1	B	320	THR	2.4
1	A	303	GLY	2.4
1	A	45	LYS	2.3
1	B	187	LEU	2.3
1	B	233	VAL	2.3
1	B	354	VAL	2.2
1	A	224	ALA	2.2
1	B	353	PHE	2.2
1	B	7	ILE	2.2
1	B	41	PHE	2.2
1	B	231	ALA	2.1
1	B	211	PRO	2.1
1	I	204	VAL	2.1
1	B	96	PRO	2.1
1	B	42	PRO	2.1
1	A	159	VAL	2.0
1	B	247	LYS	2.0
1	B	323	LYS	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.