



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:28 AM UTC

PDB ID : 4YNL / pdb_00004ynl
Title : Crystal structure of the hood domain of Anabaena HetR in complex with the hexapeptide ERGSGR derived from PatS
Authors : Hu, H.X.; Jiang, Y.L.; Zhao, M.X.; Zhang, C.C.; Chen, Y.; Zhou, C.Z.
Deposited on : 2015-03-10
Resolution : 2.10 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

<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

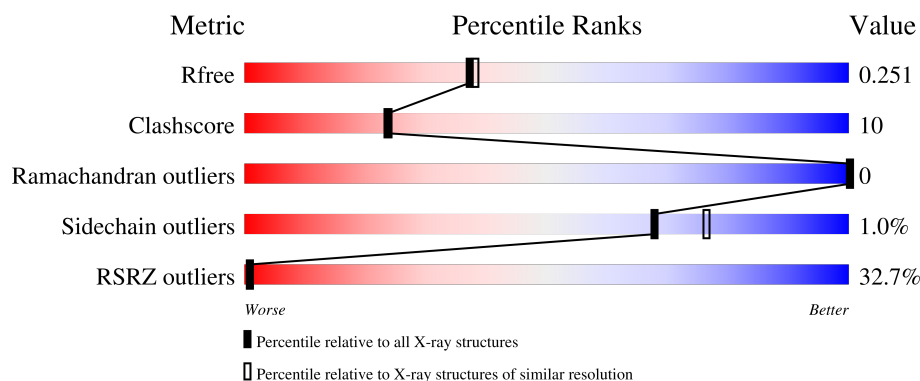
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	90	3% 62% 22% • 14%
1	B	90	7% 60% 20% • 16%
1	M	90	54% 73% 12% 14%
1	N	90	51% 70% 14% 16%
2	C	6	100%

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Mol	Chain	Length	Quality of chain
2	D	6	 33% 83% 17%
2	P	6	 33% 83% 17%
2	R	6	 33% 83% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2958 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 Heterocyst differentiation control protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	76	Total	C	N	O	S	0	1	0
			649	413	113	116	7			
1	A	77	Total	C	N	O	S	0	1	0
			658	418	114	119	7			
1	N	76	Total	C	N	O	S	0	0	0
			645	409	113	116	7			
1	M	77	Total	C	N	O	S	0	0	0
			654	414	114	119	7			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	210	MET	-	expression tag	UNP P27709
B	211	GLY	-	expression tag	UNP P27709
B	212	HIS	-	expression tag	UNP P27709
B	213	HIS	-	expression tag	UNP P27709
B	214	HIS	-	expression tag	UNP P27709
B	215	HIS	-	expression tag	UNP P27709
B	216	HIS	-	expression tag	UNP P27709
B	217	HIS	-	expression tag	UNP P27709
B	218	MET	-	expression tag	UNP P27709
A	210	MET	-	expression tag	UNP P27709
A	211	GLY	-	expression tag	UNP P27709
A	212	HIS	-	expression tag	UNP P27709
A	213	HIS	-	expression tag	UNP P27709
A	214	HIS	-	expression tag	UNP P27709
A	215	HIS	-	expression tag	UNP P27709
A	216	HIS	-	expression tag	UNP P27709
A	217	HIS	-	expression tag	UNP P27709
A	218	MET	-	expression tag	UNP P27709
N	210	MET	-	expression tag	UNP P27709
N	211	GLY	-	expression tag	UNP P27709
N	212	HIS	-	expression tag	UNP P27709

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Chain	Residue	Modelled	Actual	Comment	Reference
N	213	HIS	-	expression tag	UNP P27709
N	214	HIS	-	expression tag	UNP P27709
N	215	HIS	-	expression tag	UNP P27709
N	216	HIS	-	expression tag	UNP P27709
N	217	HIS	-	expression tag	UNP P27709
N	218	MET	-	expression tag	UNP P27709
M	210	MET	-	expression tag	UNP P27709
M	211	GLY	-	expression tag	UNP P27709
M	212	HIS	-	expression tag	UNP P27709
M	213	HIS	-	expression tag	UNP P27709
M	214	HIS	-	expression tag	UNP P27709
M	215	HIS	-	expression tag	UNP P27709
M	216	HIS	-	expression tag	UNP P27709
M	217	HIS	-	expression tag	UNP P27709
M	218	MET	-	expression tag	UNP P27709

- Molecule 2 is a protein called Heterocyst inhibition-signaling peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	6	Total	C	N	O	0	0	0
			45	24	12	9			
2	C	6	Total	C	N	O	0	0	0
			45	24	12	9			
2	P	6	Total	C	N	O	0	0	0
			45	24	12	9			
2	R	6	Total	C	N	O	0	0	0
			45	24	12	9			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	68	Total	O	0	0
			68	68		
3	A	82	Total	O	0	0
			82	82		
3	D	10	Total	O	0	0
			10	10		
3	C	11	Total	O	0	0
			11	11		
3	M	1	Total	O	0	0
			1	1		

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.

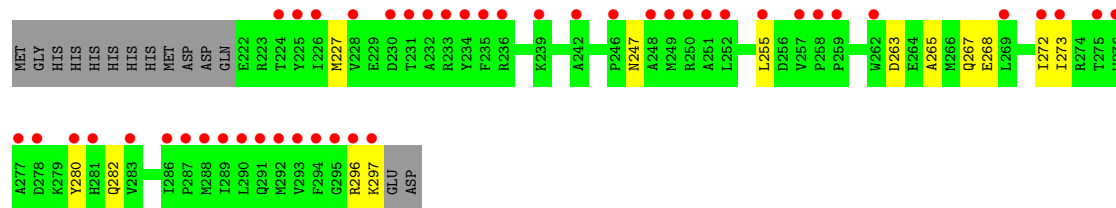
- Molecule 1: Heterocyst differentiation control protein



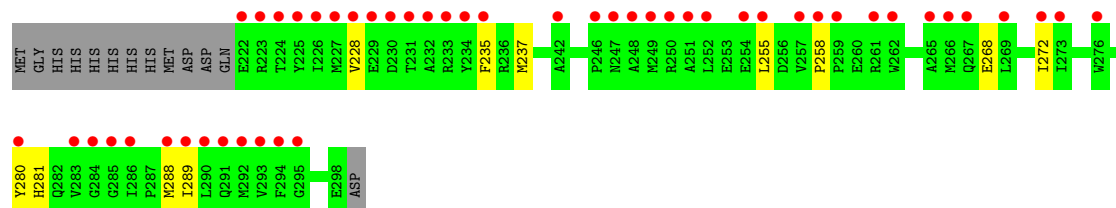
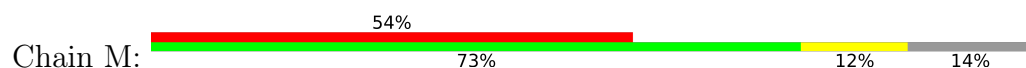
- Molecule 1: Heterocyst differentiation control protein



- Molecule 1: Heterocyst differentiation control protein



- Molecule 1: Heterocyst differentiation control protein



- Molecule 2: Heterocyst inhibition-signaling peptide

Chain D:  83% 17%



- Molecule 2: Heterocyst inhibition-signaling peptide

Chain C:  100%

There are no outlier residues recorded for this chain.

- Molecule 2: Heterocyst inhibition-signaling peptide

Chain P:  33% 83% 17%



- Molecule 2: Heterocyst inhibition-signaling peptide

Chain R:  33% 83% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	218.22Å 43.46Å 55.11Å 90.00° 97.54° 90.00°	Depositor
Resolution (Å)	42.65 – 2.10 42.65 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.3 (42.65-2.10) 97.1 (42.65-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.212 , 0.256 0.213 , 0.251	Depositor DCC
R_{free} test set	1511 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtriage
Anisotropy	0.773	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2958	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.40% 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	1.07	2/676 (0.3%)	1.05	3/911 (0.3%)
1	B	1.02	2/667 (0.3%)	1.14	2/899 (0.2%)
1	M	0.46	0/669	0.78	0/900
1	N	0.44	0/660	0.80	0/888
2	C	1.46	0/44	1.12	0/55
2	D	1.39	0/44	1.01	0/55
2	P	0.44	0/44	0.44	0/55
2	R	0.51	0/44	0.49	0/55
All	All	0.83	4/2848 (0.1%)	0.95	5/3818 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	289	ILE	C-O	5.57	1.29	1.24
1	A	254	GLU	CA-C	-5.39	1.46	1.52
1	B	254	GLU	CA-C	-5.18	1.46	1.52
1	B	289	ILE	C-O	5.11	1.29	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	286	ILE	CA-C-N	-6.18	113.58	119.76
1	B	286	ILE	C-N-CA	-6.18	113.58	119.76
1	A	286	ILE	CA-C-N	-5.90	113.71	119.90
1	A	286	ILE	C-N-CA	-5.90	113.71	119.90
1	A	285	GLY	N-CA-C	-5.34	104.00	112.66

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	658	0	646	30	0
1	B	649	0	640	31	0
1	M	654	0	637	12	0
1	N	645	0	631	15	0
2	C	45	0	45	0	0
2	D	45	0	45	1	0
2	P	45	0	45	1	0
2	R	45	0	45	1	0
3	A	82	0	0	2	0
3	B	68	0	0	1	2
3	C	11	0	0	0	2
3	D	10	0	0	0	0
3	M	1	0	0	0	0
All	All	2958	0	2734	56	2

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 10.

All (56) 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:N:268:GLU:O	1:N:272:ILE:HG13	1.77	0.84
1:B:288:MET:HG3	1:A:257[A]:VAL:HG12	1.71	0.73
1:B:272:ILE:HG21	1:A:272:ILE:CG2	2.21	0.71
1:B:292:MET:HE2	1:A:292:MET:CE	2.22	0.69
1:B:272:ILE:CG2	1:A:272:ILE:CG2	2.71	0.69
1:B:257[A]:VAL:HG12	1:A:288:MET:HG3	1.75	0.68
1:B:288:MET:HG3	1:A:257[A]:VAL:CG1	2.23	0.68
1:N:272:ILE:HG21	1:M:272:ILE:HG21	1.74	0.68
1:B:281:HIS:O	3:B:301:HOH:O	2.12	0.68
1:B:272:ILE:CG2	1:A:272:ILE:HG21	2.24	0.67
1:B:292:MET:HE2	1:A:292:MET:HE3	1.77	0.67
1:B:292:MET:CE	1:A:292:MET:HE2	2.27	0.64
1:A:271:GLU:HG2	3:A:369:HOH:O	1.97	0.64
1:N:263:ASP:O	1:N:267:GLN:OE1	2.15	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:291:GLN:HG2	3:A:362:HOH:O	1.97	0.63
1:B:292:MET:CE	1:A:292:MET:CE	2.75	0.63
1:B:257[A]:VAL:CG1	1:A:288:MET:HG3	2.29	0.62
1:B:257[A]:VAL:HG12	1:A:288:MET:CG	2.30	0.61
1:B:288:MET:CG	1:A:257[A]:VAL:HG12	2.32	0.60
1:B:241:TRP:O	1:A:223:ARG:NH1	2.28	0.58
1:N:282:GLN:HE21	1:M:258:PRO:HG3	1.69	0.58
1:N:272:ILE:CG2	1:M:272:ILE:HG21	2.35	0.57
1:B:295:GLY:HA2	1:A:247:ASN:O	2.07	0.54
1:A:224:THR:HG22	1:N:227:MET:CE	2.38	0.54
1:B:232:ALA:HB1	1:B:236:ARG:HH12	1.72	0.54
1:A:261:ARG:NH1	1:A:264:GLU:OE2	2.42	0.53
1:A:257[A]:VAL:HG13	1:A:262:TRP:HE1	1.74	0.53
1:N:282:GLN:NE2	1:M:258:PRO:HG3	2.24	0.53
1:B:257[A]:VAL:HG13	1:B:262:TRP:CD1	2.45	0.51
1:B:257[A]:VAL:HG13	1:B:262:TRP:NE1	2.27	0.50
1:A:257[A]:VAL:HG13	1:A:262:TRP:NE1	2.27	0.49
1:M:237:MET:HE1	1:M:289:ILE:HB	1.95	0.49
1:A:224:THR:HG22	1:N:227:MET:HE3	1.96	0.48
1:B:261:ARG:NH1	1:B:264:GLU:OE1	2.48	0.46
1:B:230:ASP:OD1	1:B:233:ARG:NH2	2.48	0.46
1:N:265:ALA:HA	1:M:280:TYR:CE1	2.51	0.46
1:B:254:GLU:HG2	2:D:1:GLU:O	2.17	0.45
1:B:257[A]:VAL:HG13	1:B:262:TRP:HE1	1.81	0.45
1:N:255:LEU:O	1:M:288:MET:N	2.47	0.45
1:B:251:ALA:HB1	1:A:273:ILE:HD11	1.98	0.45
1:N:272:ILE:HG21	1:M:272:ILE:CG2	2.45	0.44
1:B:242:ALA:HA	1:A:223:ARG:HB3	2.00	0.44
1:N:273:ILE:HD11	2:R:6:ARG:NH2	2.33	0.44
1:A:235:PHE:CE2	1:M:228:VAL:HA	2.53	0.43
1:B:257[B]:VAL:HG22	1:A:280:TYR:O	2.19	0.43
1:M:281:HIS:CE1	2:P:2:ARG:HB2	2.54	0.42
1:B:250:ARG:HG2	1:B:251:ALA:N	2.33	0.42
1:N:280:TYR:OH	1:M:268:GLU:OE2	2.34	0.42
1:B:237:MET:HE1	1:B:289:ILE:O	2.19	0.41
1:A:228:VAL:HG13	1:M:235:PHE:HD1	1.86	0.41
1:N:296:ARG:HG2	1:N:297:LYS:N	2.35	0.41
1:B:286:ILE:O	1:B:288:MET:HG2	2.21	0.41
1:N:247:ASN:O	1:N:247:ASN:OD1	2.39	0.41
1:B:290:LEU:HD11	1:A:292:MET:HG3	2.03	0.40
1:A:257[A]:VAL:HG13	1:A:262:TRP:CD1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:ILE:HD12	1:A:259:PRO:HD3	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:322:HOH:O	3:C:107:HOH:O[4_545]	2.06	0.14
3:B:322:HOH:O	3:C:105:HOH:O[4_545]	2.14	0.06

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	76/90 (84%)	76 (100%)	0	0	100	100
1	B	75/90 (83%)	74 (99%)	1 (1%)	0	100	100
1	M	75/90 (83%)	74 (99%)	1 (1%)	0	100	100
1	N	74/90 (82%)	74 (100%)	0	0	100	100
2	C	4/6 (67%)	4 (100%)	0	0	100	100
2	D	4/6 (67%)	4 (100%)	0	0	100	100
2	P	4/6 (67%)	4 (100%)	0	0	100	100
2	R	4/6 (67%)	4 (100%)	0	0	100	100
All	All	316/384 (82%)	314 (99%)	2 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	69/80 (86%)	68 (99%)	1 (1%)	59	67
1	B	68/80 (85%)	67 (98%)	1 (2%)	57	65
1	M	68/80 (85%)	67 (98%)	1 (2%)	57	65
1	N	67/80 (84%)	67 (100%)	0	100	100
2	C	4/4 (100%)	4 (100%)	0	100	100
2	D	4/4 (100%)	4 (100%)	0	100	100
2	P	4/4 (100%)	4 (100%)	0	100	100
2	R	4/4 (100%)	4 (100%)	0	100	100
All	All	288/336 (86%)	285 (99%)	3 (1%)	68	76

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	272	ILE
1	A	228	VAL
1	M	255	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	N	282	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 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	77/90 (85%)	0.34	3 (3%) 43 45	12, 27, 52, 69	1 (1%)
1	B	76/90 (84%)	0.40	6 (7%) 18 20	12, 28, 58, 72	1 (1%)
1	M	77/90 (85%)	2.69	49 (63%) 0 0	76, 119, 154, 162	0
1	N	76/90 (84%)	2.53	46 (60%) 0 0	71, 113, 157, 168	0
2	C	6/6 (100%)	-0.00	0 100 100	20, 21, 25, 35	0
2	D	6/6 (100%)	-0.06	0 100 100	21, 22, 25, 38	0
2	P	6/6 (100%)	1.17	2 (33%) 1 1	95, 105, 121, 126	0
2	R	6/6 (100%)	1.61	2 (33%) 1 1	73, 87, 108, 126	0
All	All	330/384 (85%)	1.43	108 (32%) 1 1	12, 71, 147, 168	2 (0%)

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	225	TYR	6.6
1	M	265	ALA	6.4
1	N	272	ILE	6.1
1	M	294	PHE	6.0
1	M	292	MET	5.9
1	N	252	LEU	5.8
1	N	276	TRP	5.7
1	M	272	ILE	5.7
1	M	290	LEU	5.5
1	N	251	ALA	5.1
1	M	252	LEU	5.0
1	M	251	ALA	4.9
1	M	235	PHE	4.8
1	M	276	TRP	4.7
1	M	285	GLY	4.5
1	N	289	ILE	4.5

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Mol	Chain	Res	Type	RSRZ
1	N	290	LEU	4.4
1	N	294	PHE	4.4
1	M	286	ILE	4.4
1	M	269	LEU	4.3
1	N	297	LYS	4.3
1	N	292	MET	4.3
1	N	293	VAL	4.1
1	M	224	THR	4.1
1	N	249	MET	4.0
1	M	273	ILE	4.0
1	M	289	ILE	3.9
1	N	287	PRO	3.9
1	M	262	TRP	3.9
1	N	269	LEU	3.6
1	M	293	VAL	3.6
1	N	283	VAL	3.5
1	N	226	ILE	3.5
1	N	286	ILE	3.5
1	N	235	PHE	3.5
1	M	232	ALA	3.4
1	N	250	ARG	3.4
1	N	225	TYR	3.4
1	N	262	TRP	3.3
1	N	242	ALA	3.3
1	M	230	ASP	3.3
1	B	225	TYR	3.2
1	N	281	HIS	3.2
1	M	229	GLU	3.2
1	M	231	THR	3.2
1	N	291	GLN	3.2
1	M	283	VAL	3.1
1	M	257	VAL	3.1
1	N	273	ILE	3.1
1	M	249	MET	3.1
1	N	259	PRO	3.1
1	M	255	LEU	3.1
1	M	258	PRO	3.0
1	N	224	THR	3.0
1	M	227	MET	3.0
1	N	248	ALA	3.0
1	N	295	GLY	3.0
1	N	280	TYR	2.9

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Mol	Chain	Res	Type	RSRZ
1	N	257	VAL	2.8
1	M	228	VAL	2.8
1	N	233	ARG	2.7
1	M	295	GLY	2.7
1	M	226	ILE	2.7
1	N	255	LEU	2.7
1	B	228	VAL	2.7
1	N	234	TYR	2.7
1	N	288	MET	2.6
1	N	246	PRO	2.6
1	M	222	GLU	2.5
1	M	284	GLY	2.5
1	N	239	LYS	2.5
1	N	232	ALA	2.5
1	M	242	ALA	2.5
1	M	259	PRO	2.5
1	M	266	MET	2.5
2	P	5	GLY	2.5
1	M	233	ARG	2.5
1	M	288	MET	2.5
1	M	234	TYR	2.4
1	N	231	THR	2.4
1	A	225	TYR	2.4
1	N	275	THR	2.4
1	N	228	VAL	2.4
1	B	229	GLU	2.3
1	M	248	ALA	2.3
1	B	231	THR	2.3
1	M	247	ASN	2.3
1	N	230	ASP	2.3
1	B	235	PHE	2.3
1	M	254	GLU	2.3
1	M	261	ARG	2.3
1	N	258	PRO	2.3
1	A	273	ILE	2.2
1	N	236	ARG	2.2
1	M	267	GLN	2.2
1	M	246	PRO	2.2
1	N	296	ARG	2.2
1	M	223	ARG	2.2
1	M	280	TYR	2.2
1	B	226	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
2	R	3	GLY	2.1
1	N	277	ALA	2.1
1	M	291	GLN	2.1
1	A	239	LYS	2.1
1	M	250	ARG	2.1
2	R	5	GLY	2.1
1	N	278	ASP	2.1
2	P	6	ARG	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.