



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:12 AM UTC

PDB ID : 5WM0 / pdb_00005wm0
Title : Crystal structure of apo wild type peptidylglycine alpha-hydroxylating monooxygenase (PHM) soaked with peptide (peptide not observed)
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Deposited on : 2017-07-28
Resolution : 2.40 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

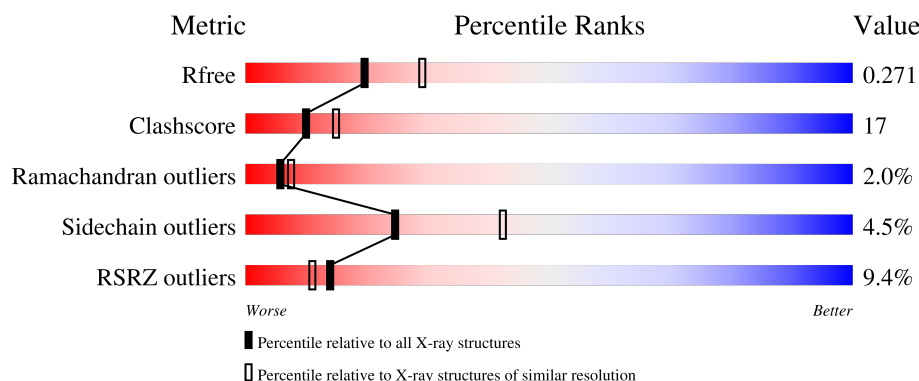
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

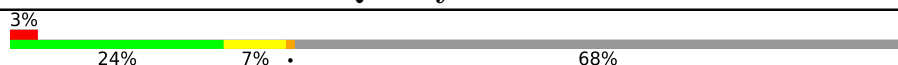
The reported resolution of this entry is 2.40 Å.

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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	976	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2445 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 Peptidyl-glycine alpha-amidating monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	0	1	0
			2418	1541	407	444	26			

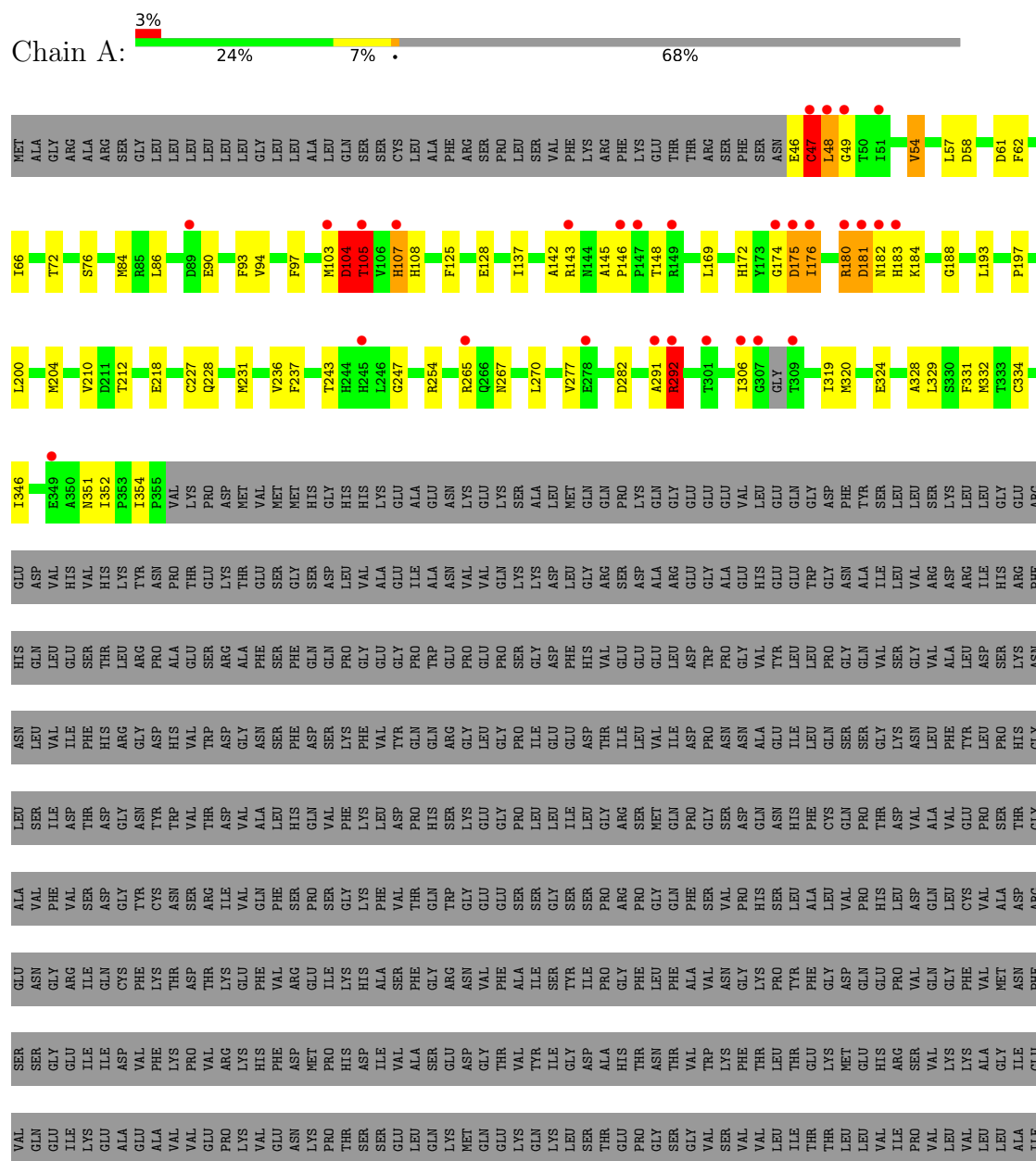
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	27	Total	O	0	0
			27	27		

3 Residue-property plots

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: Peptidyl-glycine alpha-amidating monooxygenase



VAL	GLY
MET	GLY
PHE	SER
ILE	ASP
ARG	GLN
TRP	GLU
LYS	LYS
LYS	ASP
SER	GLU
ARG	ASP
ALA	ASP
PHE	GLY
GLY	THR
ASP	GLU
HIS	SER
ASP	GLU
ASP	GLU
LYS	GLU
LEU	TYR
GLU	SER
SER	ALA
SER	PRO
SER	LEU
GLY	PRO
ARG	LYS
VAL	PRO
LEU	ALA
GLY	PRO
ARG	SER
PHE	SER
ARG	SER
GLY	
LYS	
GLY	
SER	
GLY	
GLY	
LEU	
ASN	
LEU	
GLY	
ASN	
PHE	
PHE	
ALA	
SER	
ARG	
LYS	
GLY	
TYR	
SER	
ARG	
LYS	
GLY	
PHE	
ASP	
ARG	
VAL	
SER	
THR	

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.10Å 65.93Å 69.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.93 – 2.40 47.93 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.93-2.40) 99.1 (47.93-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.200 , 0.281 0.233 , 0.271	Depositor DCC
R_{free} test set	562 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2445	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.16% 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 [i](#)

5.1 Standard geometry [i](#)

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	1.21	3/2489 (0.1%)	1.19	9/3385 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	146	PRO	N-CD	6.44	1.56	1.47
1	A	97	PHE	CA-C	6.08	1.59	1.52
1	A	86	LEU	CA-C	-5.55	1.47	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	104	ASP	N-CA-C	9.28	124.23	113.19
1	A	328	ALA	N-CA-C	9.05	122.77	110.55
1	A	176	ILE	N-CA-C	7.67	119.89	108.46
1	A	146	PRO	O-C-N	6.87	124.47	121.31
1	A	277	VAL	N-CA-C	-6.15	99.56	108.36
1	A	267	ASN	N-CA-C	-6.11	101.95	109.65
1	A	54	VAL	CB-CA-C	-5.64	102.42	110.83
1	A	107	HIS	CB-CA-C	-5.19	101.00	109.56
1	A	47	CYS	CA-CB-SG	-5.00	102.89	114.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2418	0	2342	82	0
2	A	27	0	0	2	0
All	All	2445	0	2342	82	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 17.

All (82) 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:62:PHE:CE1	1:A:193:LEU:HD12	1.76	1.20
1:A:265:ARG:NH1	1:A:354:ILE:HD11	1.52	1.20
1:A:174:GLY:O	1:A:175:ASP:HB2	1.44	1.16
1:A:227:CYS:SG	1:A:334:CYS:SG	1.36	1.13
1:A:231:MET:HE3	2:A:1012:HOH:O	1.45	1.13
1:A:265:ARG:NH1	1:A:354:ILE:CD1	2.11	1.12
1:A:227:CYS:SG	1:A:334:CYS:CB	2.44	1.05
1:A:329:LEU:HD11	1:A:332:MET:HG3	1.44	0.99
1:A:103:MET:O	1:A:143:ARG:NH2	1.97	0.97
1:A:72:THR:HG23	1:A:182:ASN:O	1.65	0.97
1:A:265:ARG:HH11	1:A:354:ILE:HD11	1.20	0.95
1:A:46:GLU:O	1:A:48:LEU:N	2.01	0.93
1:A:105:THR:HG21	1:A:180:ARG:NH1	1.82	0.93
1:A:62:PHE:CZ	1:A:193:LEU:HD12	2.03	0.93
1:A:174:GLY:O	1:A:175:ASP:CB	2.22	0.87
1:A:105:THR:HG23	1:A:180:ARG:HH22	1.37	0.87
1:A:72:THR:CG2	1:A:182:ASN:O	2.27	0.82
1:A:181:ASP:O	2:A:1001:HOH:O	2.01	0.79
1:A:105:THR:HG21	1:A:180:ARG:HH12	1.45	0.78
1:A:105:THR:HG23	1:A:180:ARG:NH2	2.00	0.75
1:A:47:CYS:C	1:A:48:LEU:HG	2.12	0.74
1:A:265:ARG:NH1	1:A:354:ILE:CG1	2.51	0.74
1:A:62:PHE:CZ	1:A:193:LEU:CD1	2.72	0.72
1:A:292:ARG:HD2	1:A:346:ILE:HD13	1.70	0.72
1:A:105:THR:CG2	1:A:180:ARG:HH12	2.05	0.70
1:A:104:ASP:O	1:A:105:THR:HG23	1.92	0.69
1:A:329:LEU:CD1	1:A:332:MET:HG3	2.18	0.69
1:A:231:MET:HE2	1:A:329:LEU:HD23	1.74	0.68
1:A:108:HIS:HB3	1:A:172:HIS:HB3	1.76	0.67
1:A:292:ARG:HD2	1:A:346:ILE:CD1	2.25	0.66
1:A:94:VAL:HG22	1:A:193:LEU:HD23	1.79	0.65
1:A:105:THR:CG2	1:A:180:ARG:NH1	2.57	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:265:ARG:HH12	1:A:354:ILE:HG12	1.62	0.65
1:A:329:LEU:HD13	1:A:331:PHE:O	1.97	0.65
1:A:265:ARG:HH11	1:A:354:ILE:CD1	1.94	0.64
1:A:66:ILE:HG23	1:A:84:MET:HG3	1.79	0.63
1:A:291:ALA:O	1:A:292:ARG:HB2	1.98	0.63
1:A:265:ARG:NH1	1:A:354:ILE:HG12	2.13	0.63
1:A:265:ARG:HH12	1:A:354:ILE:CD1	2.06	0.63
1:A:46:GLU:O	1:A:46:GLU:HG3	2.00	0.62
1:A:183:HIS:O	1:A:184:LYS:HB2	2.00	0.61
1:A:200:LEU:HD23	1:A:324:GLU:HA	1.82	0.60
1:A:94:VAL:HG22	1:A:193:LEU:CD2	2.32	0.60
1:A:180:ARG:HD3	1:A:183:HIS:ND1	2.17	0.58
1:A:62:PHE:HE1	1:A:193:LEU:HD12	1.56	0.58
1:A:254:ARG:NH1	1:A:282:ASP:O	2.30	0.56
1:A:76:SER:OG	1:A:176:ILE:N	2.37	0.55
1:A:107:HIS:ND1	1:A:108:HIS:HB2	2.22	0.54
1:A:62:PHE:CE1	1:A:193:LEU:CD1	2.71	0.54
1:A:105:THR:HG22	1:A:175:ASP:O	2.07	0.54
1:A:103:MET:C	1:A:143:ARG:HH22	2.13	0.54
1:A:57:LEU:HG	1:A:62:PHE:HA	1.90	0.53
1:A:329:LEU:C	1:A:329:LEU:HD12	2.34	0.53
1:A:243:THR:HG21	1:A:247:GLY:HA3	1.92	0.52
1:A:103:MET:O	1:A:143:ARG:CZ	2.58	0.51
1:A:105:THR:CG2	1:A:180:ARG:NH2	2.72	0.51
1:A:47:CYS:C	1:A:49:GLY:H	2.18	0.50
1:A:292:ARG:CD	1:A:346:ILE:HD13	2.39	0.50
1:A:105:THR:HG21	1:A:180:ARG:CZ	2.40	0.50
1:A:329:LEU:HD11	1:A:332:MET:CG	2.29	0.50
1:A:265:ARG:HH12	1:A:354:ILE:CG1	2.17	0.49
1:A:125:PHE:O	1:A:128:GLU:HB2	2.13	0.48
1:A:265:ARG:HG2	1:A:354:ILE:HD11	1.97	0.47
1:A:58:ASP:OD1	1:A:61:ASP:HB2	2.15	0.47
1:A:137:ILE:HG23	1:A:204:MET:HE3	1.97	0.47
1:A:105:THR:CG2	1:A:180:ARG:CZ	2.93	0.46
1:A:212:THR:O	1:A:306:ILE:HG12	2.16	0.46
1:A:46:GLU:C	1:A:48:LEU:N	2.73	0.46
1:A:227:CYS:CB	1:A:334:CYS:SG	2.82	0.46
1:A:180:ARG:HD3	1:A:183:HIS:HD1	1.80	0.45
1:A:236:VAL:HG13	1:A:319:ILE:HG23	1.99	0.44
1:A:93:PHE:CE2	1:A:197:PRO:HA	2.51	0.44
1:A:72:THR:HG21	1:A:182:ASN:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:LEU:HD21	1:A:188:GLY:HA2	2.02	0.42
1:A:175:ASP:O	1:A:176:ILE:HD13	2.19	0.42
1:A:292:ARG:NH1	1:A:351:ASN:OD1	2.53	0.41
1:A:265:ARG:HH11	1:A:354:ILE:CG1	2.27	0.41
1:A:237:PHE:CZ	1:A:320:MET:SD	3.13	0.41
1:A:47:CYS:C	1:A:49:GLY:N	2.79	0.40
1:A:47:CYS:O	1:A:48:LEU:CB	2.69	0.40
1:A:142:ALA:HB3	1:A:145:ALA:HB3	2.04	0.40
1:A:329:LEU:HD12	1:A:329:LEU:O	2.21	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	306/976 (31%)	278 (91%)	22 (7%)	6 (2%)	6 7

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	47	CYS
1	A	175	ASP
1	A	292	ARG
1	A	105	THR
1	A	148	THR
1	A	181	ASP

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	266/841 (32%)	254 (96%)	12 (4%)	24	42

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LEU
1	A	54	VAL
1	A	90	GLU
1	A	104	ASP
1	A	105	THR
1	A	180	ARG
1	A	210	VAL
1	A	218	GLU
1	A	228	GLN
1	A	270	LEU
1	A	292	ARG
1	A	352	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	ASN
1	A	182	ASN
1	A	245	HIS
1	A	316	ASN

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 [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	309/976 (31%)	0.47	29 (9%) 14 11	27, 46, 90, 125	1 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	291	ALA	4.3
1	A	307	GLY	3.9
1	A	180	ARG	3.6
1	A	48	LEU	3.4
1	A	183	HIS	3.3
1	A	265	ARG	3.3
1	A	47	CYS	3.1
1	A	175	ASP	3.0
1	A	105	THR	2.8
1	A	278	GLU	2.7
1	A	181	ASP	2.6
1	A	107	HIS	2.6
1	A	103	MET	2.5
1	A	146	PRO	2.5
1	A	301	THR	2.5
1	A	147	PRO	2.5
1	A	349	GLU	2.4
1	A	51	ILE	2.4
1	A	182	ASN	2.3
1	A	309	THR	2.3
1	A	174	GLY	2.3
1	A	143	ARG	2.3
1	A	292	ARG	2.2
1	A	245	HIS	2.2
1	A	149	ARG	2.1
1	A	176	ILE	2.1
1	A	89[A]	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	49	GLY	2.0
1	A	306	ILE	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.