



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:59 AM UTC

PDB ID : 8HUY / pdb_00008huy
Title : N-acetyl-(R)-beta-phenylalanine acylase
Authors : Kato, Y.; Natsume, R.
Deposited on : 2022-12-24
Resolution : 2.45 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

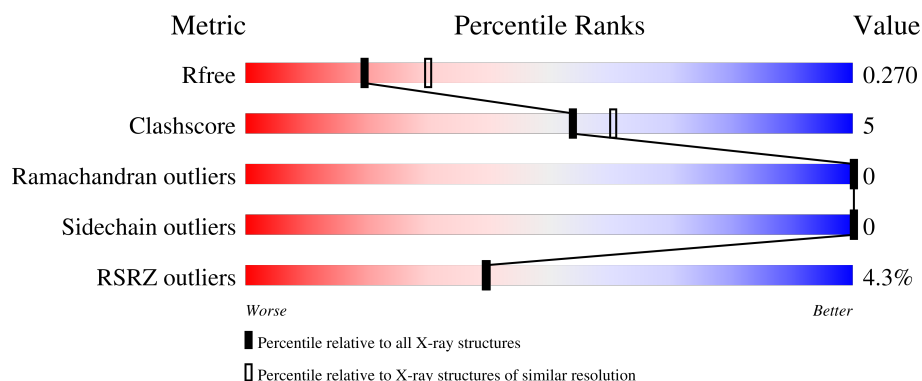
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	772	
1	B	772	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 11099 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.

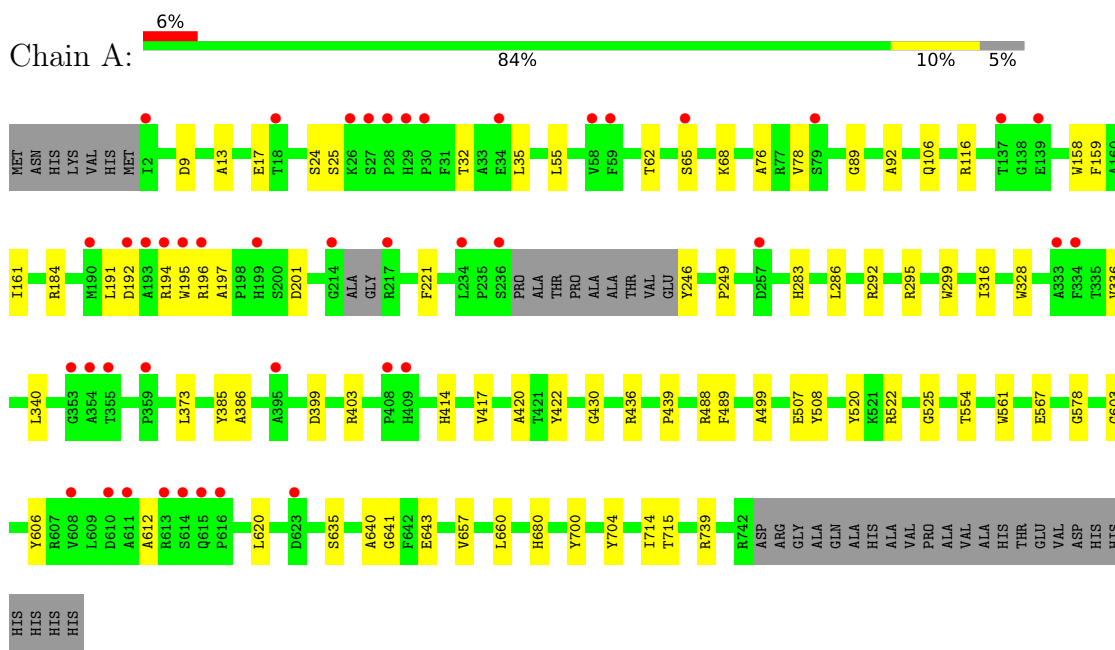
- Molecule 1 is a protein called Chains: A,B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	730	Total	C	N	O	S	0	0	0
			5556	3503	1000	1040	13			
1	B	727	Total	C	N	O	S	0	0	0
			5543	3494	997	1038	14			

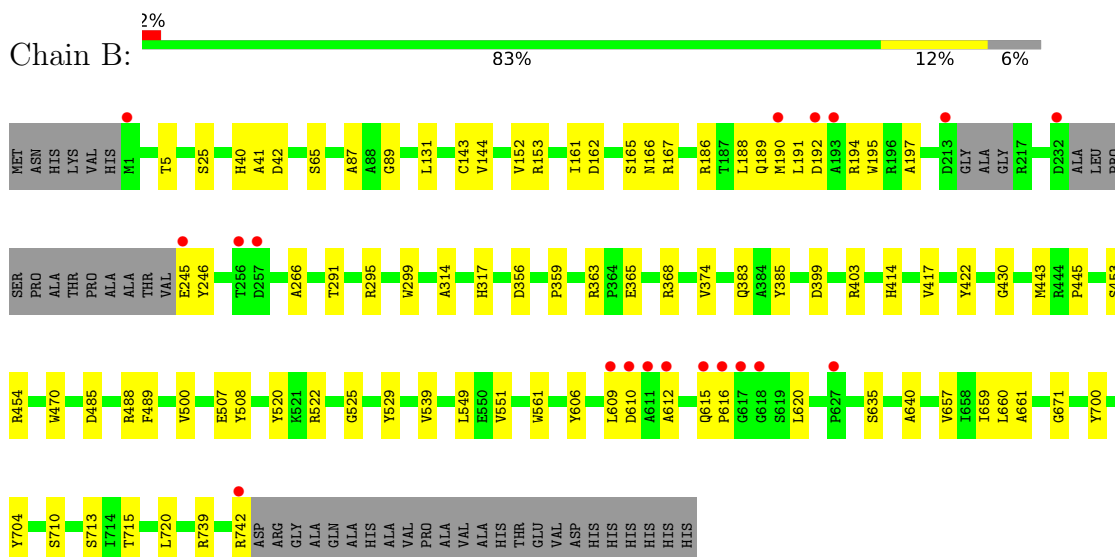
3 Residue-property plots [i](#)

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: Chains: A,B



• Molecule 1: Chains: A,B



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	112.97Å 112.97Å 341.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.12 – 2.45 47.12 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.1 (47.12-2.45) 99.1 (47.12-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.230 , 0.265 0.233 , 0.270	Depositor DCC
R_{free} test set	4066 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.535	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 24.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11099	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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	0.37	0/5708	0.57	0/7773
1	B	0.39	0/5694	0.57	0/7752
All	All	0.38	0/11402	0.57	0/15525

There are no bond length outliers.

There are no bond angle outliers.

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	5556	0	5309	49	0
1	B	5543	0	5296	59	0
All	All	11099	0	10605	107	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 5.

All (107) 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:32:THR:HG22	1:A:62:THR:HG22	1.63	0.81
1:B:615:GLN:HB3	1:B:616:PRO:HD2	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:LEU:HD11	1:B:195:TRP:HB3	1.73	0.69
1:A:507:GLU:HB3	1:A:508:TYR:CD2	2.32	0.65
1:A:13:ALA:HB1	1:A:17:GLU:HB2	1.80	0.64
1:A:292:ARG:HG2	1:A:316:ILE:HG12	1.81	0.62
1:B:609:LEU:HD13	1:B:659:ILE:HG22	1.80	0.61
1:A:191:LEU:HD11	1:A:195:TRP:HB3	1.81	0.61
1:B:507:GLU:HB3	1:B:508:TYR:CD2	2.36	0.60
1:A:417:VAL:HG12	1:A:439:PRO:HG2	1.84	0.60
1:A:196:ARG:HG2	1:A:197:ALA:H	1.67	0.60
1:B:40:HIS:CE1	1:B:42:ASP:HB3	2.38	0.59
1:A:116:ARG:HH12	1:A:201:ASP:HB3	1.68	0.58
1:A:422:TYR:OH	1:A:507:GLU:OE1	2.21	0.58
1:B:610:ASP:OD1	1:B:610:ASP:N	2.38	0.57
1:A:9:ASP:O	1:A:436:ARG:NH1	2.37	0.57
1:A:76:ALA:HB2	1:A:286:LEU:HD23	1.86	0.56
1:A:184:ARG:NH2	1:A:246:TYR:O	2.38	0.55
1:B:186:ARG:HG2	1:B:246:TYR:CZ	2.42	0.55
1:A:620:LEU:HD21	1:A:739:ARG:HD3	1.88	0.55
1:B:167:ARG:NH1	1:B:190:MET:HG3	2.22	0.55
1:B:739:ARG:HA	1:B:742:ARG:HE	1.72	0.54
1:B:422:TYR:OH	1:B:507:GLU:OE1	2.25	0.54
1:A:196:ARG:HG2	1:A:197:ALA:N	2.22	0.54
1:A:24:SER:HB2	1:A:68:LYS:HB2	1.89	0.54
1:B:189:GLN:HG2	1:B:194:ARG:HH22	1.72	0.54
1:A:116:ARG:HH12	1:A:201:ASP:CB	2.21	0.53
1:B:609:LEU:CD1	1:B:659:ILE:HG22	2.38	0.53
1:A:606:TYR:CZ	1:A:640:ALA:HB2	2.45	0.52
1:B:385:TYR:CZ	1:B:422:TYR:CZ	2.98	0.52
1:B:609:LEU:HD12	1:B:661:ALA:HA	1.90	0.52
1:B:383:GLN:HB3	1:B:549:LEU:HD13	1.92	0.52
1:A:385:TYR:CZ	1:A:422:TYR:CZ	2.98	0.51
1:A:25:SER:OG	1:A:65:SER:HB3	2.11	0.51
1:A:643:GLU:HB2	1:A:714:ILE:HD12	1.93	0.51
1:B:190:MET:HE2	1:B:192:ASP:HB2	1.92	0.51
1:B:414:HIS:O	1:B:417:VAL:HG22	2.11	0.51
1:A:522:ARG:HG2	1:A:704:TYR:CZ	2.45	0.51
1:B:609:LEU:HD12	1:B:661:ALA:CA	2.41	0.51
1:A:520:TYR:CZ	1:A:525:GLY:HA3	2.46	0.50
1:B:295:ARG:HD2	1:B:299:TRP:CD2	2.47	0.50
1:B:529:TYR:HD2	1:B:710:SER:HB3	1.76	0.50
1:B:635:SER:HB2	1:B:715:THR:HB	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:HIS:O	1:A:417:VAL:HG22	2.13	0.49
1:B:470:TRP:CD1	1:B:720:LEU:HB3	2.48	0.49
1:A:295:ARG:HD2	1:A:299:TRP:CD2	2.48	0.49
1:B:245:GLU:HG3	1:B:246:TYR:N	2.26	0.49
1:B:190:MET:O	1:B:194:ARG:NH2	2.44	0.49
1:B:144:VAL:HG22	1:B:189:GLN:OE1	2.14	0.48
1:A:680:HIS:CD2	1:B:671:GLY:HA3	2.49	0.47
1:B:620:LEU:HD21	1:B:739:ARG:HD3	1.96	0.47
1:B:453:SER:OG	1:B:454:ARG:HG3	2.15	0.47
1:B:186:ARG:CZ	1:B:188:LEU:HD11	2.45	0.47
1:A:192:ASP:C	1:A:194:ARG:H	2.23	0.46
1:A:89:GLY:HA3	1:A:161:ILE:O	2.15	0.46
1:A:635:SER:HB2	1:A:715:THR:HB	1.98	0.46
1:B:430:GLY:HA3	1:B:561:TRP:CD2	2.50	0.46
1:A:430:GLY:HA3	1:A:561:TRP:CD2	2.51	0.45
1:B:539:VAL:HG12	1:B:551:VAL:HG22	1.98	0.45
1:A:603:GLY:HA2	1:A:641:GLY:HA3	1.98	0.45
1:B:522:ARG:HG2	1:B:704:TYR:CZ	2.51	0.45
1:B:41:ALA:HB1	1:B:445:PRO:HA	1.99	0.45
1:A:158:TRP:CG	1:A:249:PRO:HD2	2.52	0.44
1:B:192:ASP:C	1:B:194:ARG:H	2.25	0.44
1:A:24:SER:HB3	1:A:328:TRP:CH2	2.53	0.44
1:B:166:ASN:O	1:B:167:ARG:HD3	2.18	0.44
1:A:78:VAL:HG13	1:A:283:HIS:O	2.18	0.43
1:A:158:TRP:CD2	1:A:249:PRO:HD2	2.53	0.43
1:A:35:LEU:HD23	1:A:55:LEU:HB2	2.00	0.43
1:B:606:TYR:CZ	1:B:640:ALA:HB2	2.53	0.43
1:B:266:ALA:HB2	1:B:314:ALA:HB2	2.00	0.43
1:B:657:VAL:HA	1:B:700:TYR:HB3	2.01	0.43
1:A:488:ARG:HD3	1:A:489:PHE:CZ	2.53	0.43
1:B:131:LEU:O	1:B:143:CYS:HA	2.19	0.43
1:B:520:TYR:CZ	1:B:525:GLY:HA3	2.54	0.43
1:A:68:LYS:HB3	1:A:328:TRP:CD1	2.54	0.43
1:B:25:SER:HB3	1:B:65:SER:HB3	2.01	0.43
1:B:5:THR:HA	1:B:356:ASP:OD2	2.19	0.42
1:B:365:GLU:HB3	1:B:368:ARG:HB2	2.00	0.42
1:B:399:ASP:O	1:B:403:ARG:HG3	2.19	0.42
1:A:336:VAL:HG13	1:A:340:LEU:HD23	2.00	0.42
1:A:192:ASP:O	1:A:194:ARG:N	2.52	0.42
1:B:363:ARG:HD3	1:B:485:ASP:CG	2.44	0.42
1:A:373:LEU:HD22	1:A:499:ALA:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:ASP:OD2	1:B:165:SER:OG	2.23	0.42
1:B:291:THR:HB	1:B:317:HIS:HB3	2.01	0.42
1:B:89:GLY:HA3	1:B:161:ILE:O	2.20	0.41
1:A:106:GLN:HB3	1:A:221:PHE:HA	2.02	0.41
1:B:190:MET:CE	1:B:192:ASP:HB2	2.49	0.41
1:A:612:ALA:HB1	1:A:660:LEU:HD23	2.02	0.41
1:B:507:GLU:HB3	1:B:508:TYR:HD2	1.84	0.41
1:A:657:VAL:HA	1:A:700:TYR:HB3	2.02	0.41
1:B:87:ALA:HA	1:B:197:ALA:HB2	2.02	0.41
1:B:152:VAL:O	1:B:153:ARG:HG2	2.20	0.41
1:A:68:LYS:HB3	1:A:328:TRP:NE1	2.36	0.41
1:B:359:PRO:HG2	1:B:443:MET:O	2.21	0.41
1:A:295:ARG:HG3	1:A:567:GLU:OE1	2.21	0.41
1:A:399:ASP:O	1:A:403:ARG:HG3	2.21	0.41
1:B:606:TYR:CE2	1:B:640:ALA:HB2	2.56	0.41
1:A:554:THR:OG1	1:A:578:GLY:HA3	2.20	0.41
1:B:374:VAL:CG1	1:B:500:VAL:HG22	2.51	0.40
1:B:488:ARG:HD3	1:B:489:PHE:CE2	2.56	0.40
1:B:191:LEU:O	1:B:194:ARG:HB2	2.21	0.40
1:B:612:ALA:HB1	1:B:660:LEU:HD23	2.03	0.40
1:B:640:ALA:HA	1:B:713:SER:OG	2.22	0.40
1:A:92:ALA:HB3	1:A:159:PHE:CE1	2.57	0.40
1:A:386:ALA:O	1:A:420:ALA:HB2	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	724/772 (94%)	702 (97%)	22 (3%)	0	100	100
1	B	721/772 (93%)	698 (97%)	23 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1445/1544 (94%)	1400 (97%)	45 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	554/585 (95%)	554 (100%)	0	100	100
1	B	553/585 (94%)	553 (100%)	0	100	100
All	All	1107/1170 (95%)	1107 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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	409	HIS
1	A	466	HIS
1	B	280	HIS
1	B	722	HIS

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

There are no ligands in this entry.

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	730/772 (94%)	0.56	43 (5%) 28 25	22, 36, 56, 74	0
1	B	727/772 (94%)	0.30	19 (2%) 57 58	21, 32, 50, 78	0
All	All	1457/1544 (94%)	0.43	62 (4%) 40 39	21, 34, 53, 78	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	214	GLY	6.8
1	B	616	PRO	4.8
1	A	236	SER	4.4
1	B	256	THR	4.3
1	A	610	ASP	4.2
1	B	615	GLN	4.1
1	A	234	LEU	4.1
1	B	742	ARG	3.8
1	A	616	PRO	3.3
1	A	2	ILE	3.3
1	A	27	SER	3.2
1	A	29	HIS	3.1
1	A	611	ALA	3.1
1	A	192	ASP	3.0
1	B	192	ASP	3.0
1	B	617	GLY	3.0
1	B	213	ASP	3.0
1	B	1	MET	2.9
1	B	610	ASP	2.9
1	A	65	SER	2.9
1	A	28	PRO	2.8
1	A	137	THR	2.8
1	A	199	HIS	2.8
1	A	409	HIS	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	618	GLY	2.7
1	A	615	GLN	2.7
1	A	193	ALA	2.7
1	B	611	ALA	2.7
1	A	139	GLU	2.7
1	B	257	ASP	2.6
1	A	194	ARG	2.6
1	A	217	ARG	2.5
1	A	195	TRP	2.5
1	A	623	ASP	2.5
1	A	354	ALA	2.4
1	A	334	PHE	2.4
1	B	627	PRO	2.3
1	A	257	ASP	2.3
1	B	609	LEU	2.3
1	A	395	ALA	2.3
1	A	196	ARG	2.3
1	A	614	SER	2.3
1	A	58	VAL	2.3
1	A	333	ALA	2.3
1	B	193	ALA	2.2
1	B	612	ALA	2.2
1	A	355	THR	2.2
1	B	232	ASP	2.2
1	B	245	GLU	2.2
1	A	190	MET	2.2
1	B	190	MET	2.2
1	A	359	PRO	2.2
1	A	34	GLU	2.2
1	A	59	PHE	2.2
1	A	18	THR	2.1
1	A	353	GLY	2.1
1	A	613	ARG	2.1
1	A	26	LYS	2.1
1	A	408	PRO	2.0
1	A	608	VAL	2.0
1	A	30	PRO	2.0
1	A	79	SER	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 [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.