



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:42 PM UTC

PDB ID : 6J0W / pdb_00006j0w
Title : Crystal Structure of Yeast Rtt107 and Nse6
Authors : Wan, B.; Wu, J.; Lei, M.
Deposited on : 2018-12-27
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

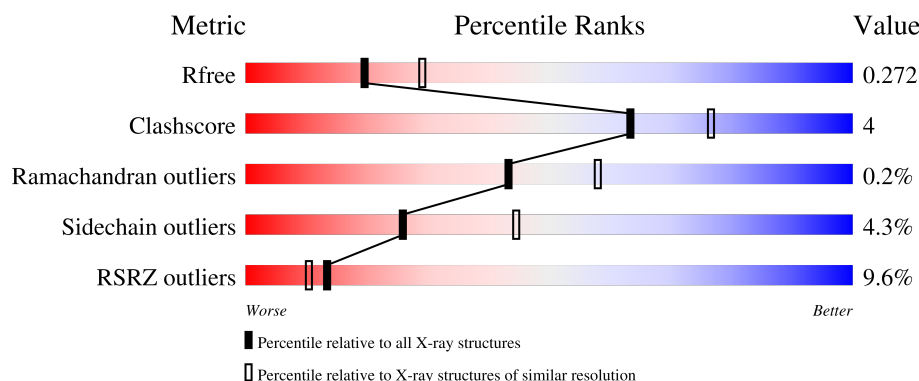
1 Overall quality at a glance

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	513	3% 82% 7% 10%
1	B	513	10% 77% 12% 11%
2	C	29	38% 48% 24% 24%
2	D	29	69% 66% 31%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7963 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 Regulator of Ty1 transposition protein 107.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	S	0	0	0
			3772	2458	617	679	18			
1	B	457	Total	C	N	O	S	0	0	0
			3750	2446	614	672	18			

- Molecule 2 is a protein called Peptide from DNA repair protein KRE29.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	22	Total	C	N	O	0	0	0
			169	105	26	38			
2	D	29	Total	C	N	O	0	0	0
			222	137	33	52			

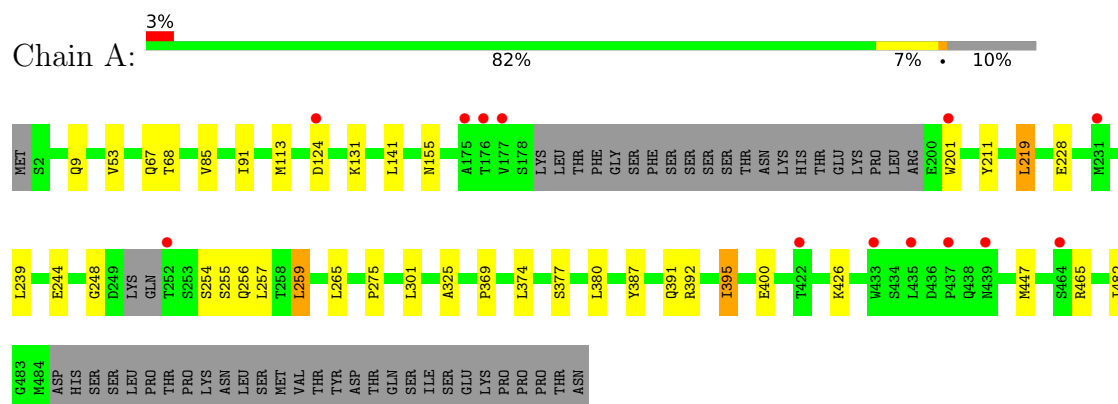
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	26	Total	O	0	0
			26	26		
3	B	18	Total	O	0	0
			18	18		
3	C	2	Total	O	0	0
			2	2		
3	D	4	Total	O	0	0
			4	4		

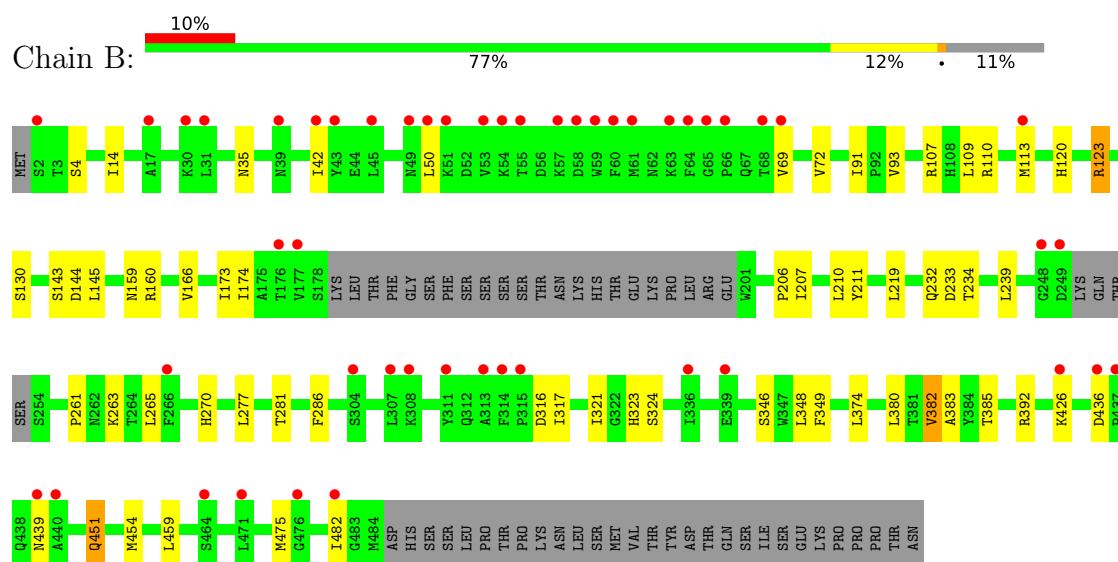
3 Residue-property plots

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.

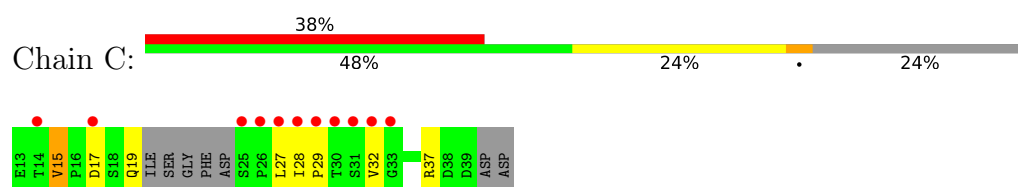
• Molecule 1: Regulator of Ty1 transposition protein 107



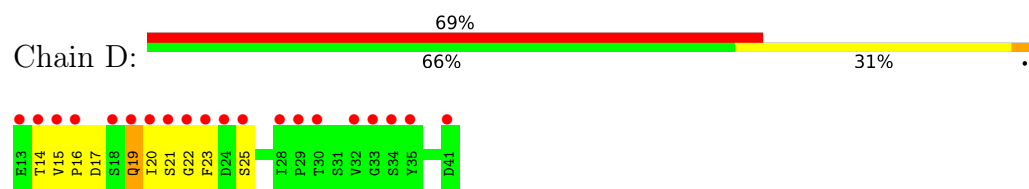
• Molecule 1: Regulator of Ty1 transposition protein 107



• Molecule 2: Peptide from DNA repair protein KRE29



- Molecule 2: Peptide from DNA repair protein KRE29



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.78Å 101.42Å 86.56Å 90.00° 108.57° 90.00°	Depositor
Resolution (Å)	45.71 – 2.40 45.71 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (45.71-2.40) 99.6 (45.71-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.40Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.218 , 0.271 0.222 , 0.272	Depositor DCC
R_{free} test set	2263 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.6	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	7963	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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	0.42	0/3878	0.60	0/5262
1	B	0.39	0/3856	0.60	0/5232
2	C	0.80	1/172 (0.6%)	0.95	1/234 (0.4%)
2	D	0.47	0/227	0.77	0/310
All	All	0.42	1/8133 (0.0%)	0.61	1/11038 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	28	ILE	CA-C	-7.45	1.46	1.53

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	28	ILE	O-C-N	-5.14	116.34	121.12

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	3772	0	3745	18	0
1	B	3750	0	3727	41	0
2	C	169	0	153	4	0
2	D	222	0	194	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	26	0	0	0	0
3	B	18	0	0	0	0
3	C	2	0	0	0	0
3	D	4	0	0	0	0
All	All	7963	0	7819	63	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 4.

All (63) 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:B:385:THR:OG1	1:B:426:LYS:HD2	1.41	1.18
1:B:160:ARG:HH12	2:D:22:GLY:HA3	1.38	0.87
1:B:385:THR:OG1	1:B:426:LYS:CD	2.23	0.85
1:A:256:GLN:HB3	1:A:259:LEU:HD22	1.67	0.76
1:B:454:MET:HE1	1:B:482:ILE:HG22	1.70	0.73
1:B:109:LEU:HD23	2:D:20:ILE:HA	1.71	0.71
1:B:385:THR:HG1	1:B:426:LYS:HD2	1.54	0.70
1:B:144:ASP:OD2	1:B:346:SER:OG	2.05	0.68
2:D:16:PRO:HB3	2:D:23:PHE:HD1	1.59	0.68
1:B:107:ARG:NH1	2:D:21:SER:HA	2.14	0.62
1:B:392:ARG:HH12	2:D:19:GLN:HG3	1.63	0.62
1:B:69:VAL:H	1:B:475:MET:HE3	1.64	0.62
1:B:160:ARG:NH1	2:D:22:GLY:HA3	2.14	0.60
1:B:166:VAL:O	1:B:206:PRO:HD3	2.01	0.60
1:B:160:ARG:HH12	2:D:22:GLY:CA	2.12	0.59
1:B:392:ARG:NH1	2:D:19:GLN:HG3	2.17	0.58
2:C:29:PRO:HG2	2:C:32:VAL:HG23	1.86	0.57
1:A:131:LYS:NZ	1:A:400:GLU:OE2	2.35	0.56
1:B:145:LEU:HD21	1:B:349:PHE:CD1	2.40	0.56
1:A:447:MET:HE2	1:A:482:ILE:HD11	1.88	0.55
2:D:17:ASP:OD2	2:D:19:GLN:HG2	2.06	0.55
1:A:447:MET:CE	1:A:482:ILE:HD11	2.38	0.54
1:A:392:ARG:NH1	2:C:19:GLN:OE1	2.40	0.54
1:A:275:PRO:HD3	1:A:301:LEU:HB2	1.90	0.53
1:A:211:TYR:CD2	1:A:219:LEU:HG	2.44	0.53
1:A:387:TYR:O	1:A:392:ARG:HD2	2.07	0.53
1:A:391:GLN:O	1:A:395:ILE:HG12	2.10	0.52
1:B:321:ILE:HD11	1:B:348:LEU:HD21	1.91	0.52
1:B:316:ASP:N	1:B:316:ASP:OD1	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:ILE:HB	1:B:42:ILE:HD13	1.92	0.52
1:B:113:MET:HA	1:B:123:ARG:HG3	1.94	0.50
1:B:270:HIS:HB3	1:B:317:ILE:HA	1.94	0.49
1:B:109:LEU:CD2	2:D:20:ILE:HA	2.40	0.49
1:A:85:VAL:HG13	1:A:91:ILE:HB	1.95	0.49
1:B:385:THR:CB	1:B:426:LYS:CD	2.92	0.48
1:B:436:ASP:HB3	1:B:439:ASN:HB2	1.97	0.47
1:A:244:GLU:O	1:A:248:GLY:HA3	2.15	0.47
1:B:211:TYR:CD2	1:B:219:LEU:HG	2.50	0.47
1:B:277:LEU:HD23	1:B:323:HIS:HB2	1.96	0.46
1:B:120:HIS:O	1:B:123:ARG:HB3	2.17	0.45
1:B:385:THR:CB	1:B:426:LYS:HD3	2.47	0.45
1:B:277:LEU:CD2	1:B:323:HIS:HB2	2.47	0.44
1:A:426:LYS:HG3	2:C:15:VAL:HG11	1.97	0.44
1:A:392:ARG:NH2	2:C:17:ASP:OD2	2.49	0.44
1:A:254:SER:OG	1:A:255:SER:N	2.49	0.44
1:A:325:ALA:HB2	1:A:369:PRO:O	2.18	0.44
1:B:210:LEU:HG	1:B:286:PHE:CZ	2.53	0.44
2:D:16:PRO:HB3	2:D:23:PHE:CD1	2.46	0.43
1:B:166:VAL:CG1	1:B:173:ILE:HD11	2.48	0.43
1:A:395:ILE:HG21	1:A:395:ILE:HD12	1.81	0.43
1:B:385:THR:HB	1:B:426:LYS:HD3	2.00	0.43
1:B:382:VAL:HG23	1:B:383:ALA:O	2.19	0.43
1:B:72:VAL:O	1:B:93:VAL:HA	2.19	0.43
1:B:91:ILE:HD11	1:B:475:MET:HE2	2.00	0.43
1:B:110:ARG:HB3	2:D:20:ILE:HD11	2.01	0.42
1:B:166:VAL:HG11	1:B:174:ILE:HD11	2.01	0.42
1:B:232:GLN:O	1:B:234:THR:N	2.52	0.42
1:A:141:LEU:C	1:A:141:LEU:HD23	2.44	0.41
1:B:261:PRO:O	1:B:263:LYS:HG3	2.20	0.41
1:A:447:MET:H	1:A:447:MET:HG3	1.71	0.41
1:B:4:SER:HB3	1:B:35:ASN:OD1	2.21	0.40
1:B:145:LEU:HD21	1:B:349:PHE:CE1	2.56	0.40
1:B:451:GLN:HB3	1:B:459:LEU:HD12	2.02	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	454/513 (88%)	442 (97%)	12 (3%)	0	100	100
1	B	451/513 (88%)	417 (92%)	33 (7%)	1 (0%)	43	58
2	C	18/29 (62%)	16 (89%)	2 (11%)	0	100	100
2	D	27/29 (93%)	23 (85%)	3 (11%)	1 (4%)	2	2
All	All	950/1084 (88%)	898 (94%)	50 (5%)	2 (0%)	43	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	19	GLN
1	B	233	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	428/480 (89%)	409 (96%)	19 (4%)	25	43
1	B	425/480 (88%)	411 (97%)	14 (3%)	33	55
2	C	21/27 (78%)	18 (86%)	3 (14%)	3	4
2	D	27/27 (100%)	24 (89%)	3 (11%)	6	9
All	All	901/1014 (89%)	862 (96%)	39 (4%)	26	44

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	53	VAL
1	A	67	GLN
1	A	68	THR
1	A	113	MET
1	A	124	ASP
1	A	155	ASN
1	A	201	TRP
1	A	219	LEU
1	A	228	GLU
1	A	239	LEU
1	A	257	LEU
1	A	259	LEU
1	A	265	LEU
1	A	374	LEU
1	A	377	SER
1	A	380	LEU
1	A	395	ILE
1	A	465	ARG
1	B	50	LEU
1	B	123	ARG
1	B	130	SER
1	B	143	SER
1	B	159	ASN
1	B	207	ILE
1	B	239	LEU
1	B	265	LEU
1	B	281	THR
1	B	324	SER
1	B	374	LEU
1	B	380	LEU
1	B	382	VAL
1	B	451	GLN
2	C	15	VAL
2	C	27	LEU
2	C	37	ARG
2	D	14	THR
2	D	15	VAL
2	D	25	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	GLN
1	A	78	ASN
1	A	112	ASN
1	A	362	GLN
1	A	386	ASN
1	B	67	GLN
1	B	78	ASN
1	B	159	ASN
1	B	268	ASN
1	B	355	GLN
1	B	371	HIS
1	B	438	GLN

5.3.3 RNA [i](#)

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	460/513 (89%)	-0.05	13 (2%) 55 51	16, 37, 72, 114	0
1	B	457/513 (89%)	0.79	49 (10%) 11 8	23, 51, 90, 111	0
2	C	22/29 (75%)	1.99	11 (50%) 0 0	49, 78, 103, 107	0
2	D	29/29 (100%)	2.49	20 (68%) 0 0	53, 83, 95, 99	0
All	All	968/1084 (89%)	0.47	93 (9%) 13 10	16, 45, 90, 114	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	19	GLN	5.2
2	D	20	ILE	4.9
2	D	21	SER	4.8
2	D	22	GLY	4.3
2	C	32	VAL	4.3
2	D	23	PHE	4.3
1	A	176	THR	4.2
1	B	307	LEU	4.1
2	C	30	THR	3.9
1	B	314	PHE	3.8
2	C	28	ILE	3.7
2	C	33	GLY	3.7
2	C	27	LEU	3.4
2	D	15	VAL	3.4
2	C	29	PRO	3.2
1	B	53	VAL	3.2
1	A	201	TRP	3.2
1	B	64	PHE	3.1
2	D	41	ASP	3.1
1	B	476	GLY	3.1
1	A	175	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
2	C	26	PRO	3.0
1	B	339	GLU	3.0
1	B	176	THR	3.0
1	B	61	MET	3.0
1	B	43	TYR	2.9
2	D	30	THR	2.9
1	B	65	GLY	2.8
1	B	311	TYR	2.8
2	D	33	GLY	2.7
1	B	482	ILE	2.7
1	B	60	PHE	2.7
1	B	42	ILE	2.7
1	B	248	GLY	2.6
2	D	32	VAL	2.6
2	D	24	ASP	2.6
1	B	68	THR	2.6
1	B	57	LYS	2.6
1	B	440	ALA	2.6
1	B	66	PRO	2.5
1	B	249	ASP	2.5
1	B	336	ILE	2.5
1	B	45	LEU	2.5
1	B	59	TRP	2.5
2	D	16	PRO	2.4
2	D	34	SER	2.4
1	B	69	VAL	2.4
1	A	437	PRO	2.4
2	D	18	SER	2.4
1	B	54	LYS	2.4
1	B	308	LYS	2.4
2	C	31	SER	2.3
1	A	422	THR	2.3
1	B	55	THR	2.3
1	A	433	TRP	2.3
1	B	464	SER	2.3
2	D	28	ILE	2.3
1	B	58	ASP	2.3
1	B	31	LEU	2.3
1	B	113	MET	2.3
1	B	315	PRO	2.3
2	C	25	SER	2.3
1	B	51	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	177	VAL	2.3
1	B	437	PRO	2.3
1	B	313	ALA	2.2
1	A	435	LEU	2.2
1	A	464	SER	2.2
1	B	471	LEU	2.2
2	D	29	PRO	2.2
1	B	39	ASN	2.2
1	A	124	ASP	2.2
1	B	436	ASP	2.2
1	A	439	ASN	2.2
1	B	50	LEU	2.2
2	D	25	SER	2.2
1	B	17	ALA	2.2
1	B	49	ASN	2.2
2	C	14	THR	2.2
1	B	2	SER	2.2
2	D	13	GLU	2.2
1	A	252	THR	2.1
1	B	30	LYS	2.1
2	D	35	TYR	2.1
2	C	17	ASP	2.1
1	A	231	MET	2.1
2	D	14	THR	2.1
1	B	426	LYS	2.1
1	B	304	SER	2.1
1	B	439	ASN	2.0
1	B	63	LYS	2.0
1	B	177	VAL	2.0
1	B	266	PHE	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.