



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:19 PM UTC

PDB ID : 3C9D / pdb_00003c9d
Title : Crystal structure of Vps75
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Deposited on : 2008-02-15
Resolution : 2.00 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

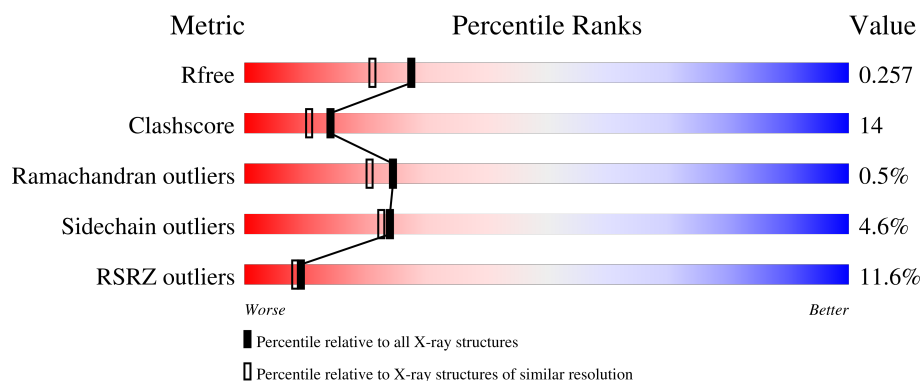
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	259	10% 63% 14% • 20%
1	B	259	8% 64% 13% • 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3761 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 Vacuolar protein sorting-associated protein 75.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	208	Total	C	N	O	S	0	0	0
			1739	1128	283	323	5			
1	B	205	Total	C	N	O	S	0	0	0
			1713	1114	279	315	5			

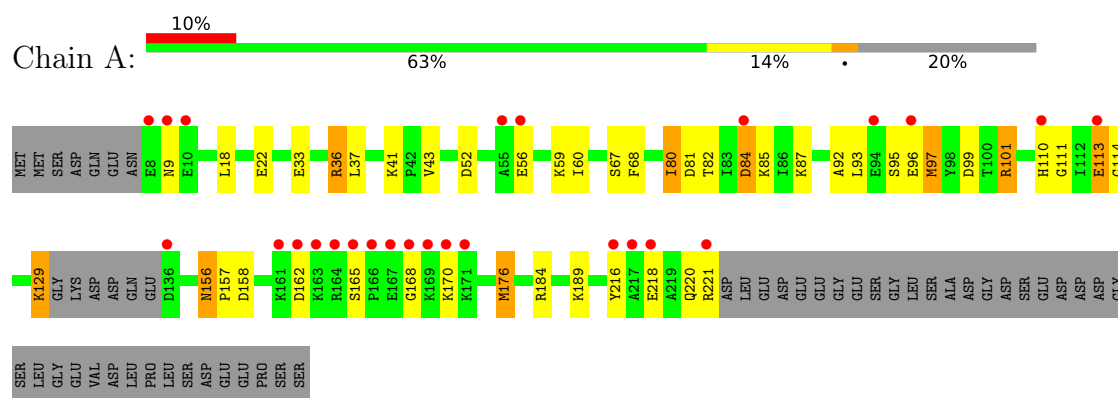
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	135	Total	O	0	0
			135	135		
2	B	174	Total	O	0	0
			174	174		

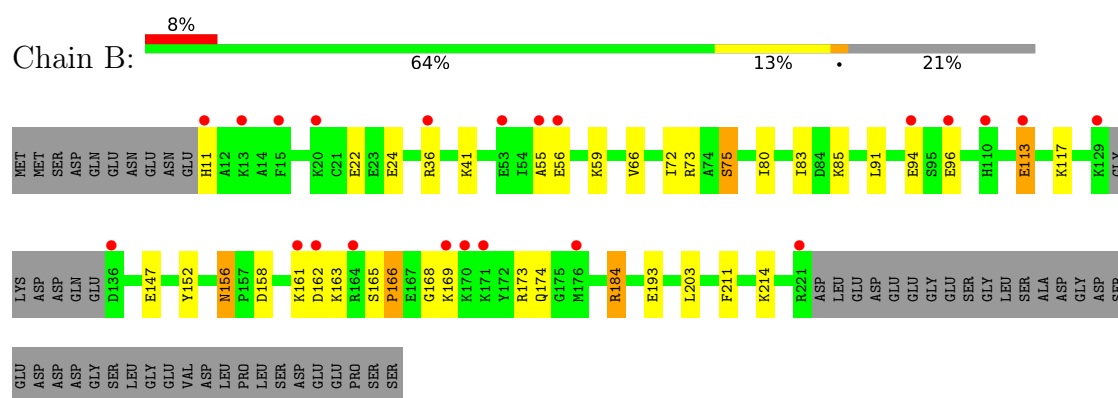
3 Residue-property plots [i](#)

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: Vacuolar protein sorting-associated protein 75



- Molecule 1: Vacuolar protein sorting-associated protein 75



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.81Å 86.08Å 85.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 20.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.2 (20.00-2.00) 99.0 (20.00-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.229 , 0.251 0.235 , 0.257	Depositor DCC
R_{free} test set	1868 reflections (3.63%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.759	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -h,l,k	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3761	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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	0.53	0/1786	0.71	0/2405
1	B	0.58	0/1760	0.71	0/2370
All	All	0.55	0/3546	0.71	0/4775

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

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	1739	0	1693	58	0
1	B	1713	0	1674	40	5
2	A	135	0	0	11	1
2	B	174	0	0	14	4
All	All	3761	0	3367	97	7

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 14.

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:B:117:LYS:NZ	1:B:147:GLU:HG3	1.59	1.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:117:LYS:HZ3	1:B:147:GLU:HG3	1.07	1.15
1:A:101:ARG:HG3	1:A:101:ARG:HH11	1.11	1.14
1:B:41:LYS:HB2	2:B:467:HOH:O	1.61	0.98
1:B:91:LEU:O	1:B:94:GLU:HG2	1.65	0.96
1:A:176:MET:HA	1:A:176:MET:HE2	1.47	0.95
1:B:173:ARG:NH2	2:B:491:HOH:O	2.02	0.91
1:A:110:HIS:CE1	2:A:341:HOH:O	2.26	0.87
1:A:82:THR:HA	1:A:110:HIS:CE1	2.11	0.85
1:A:101:ARG:HG3	1:A:101:ARG:NH1	1.91	0.83
1:B:73:ARG:NH1	2:B:356:HOH:O	2.13	0.82
1:A:113:GLU:OE1	2:A:394:HOH:O	2.05	0.74
1:A:99:ASP:OD1	1:A:101:ARG:NH1	2.24	0.70
1:B:117:LYS:HZ1	1:B:147:GLU:HG3	1.58	0.68
1:A:101:ARG:HH11	1:A:101:ARG:CG	1.96	0.66
1:A:113:GLU:OE1	1:A:113:GLU:HA	1.95	0.66
1:A:189:LYS:HD3	2:A:348:HOH:O	1.96	0.65
1:A:82:THR:CB	1:A:110:HIS:CE1	2.81	0.63
1:A:84:ASP:OD1	1:A:110:HIS:HB3	2.00	0.62
1:A:82:THR:CA	1:A:110:HIS:CE1	2.82	0.62
1:A:56:GLU:O	1:A:60:ILE:HG12	2.00	0.62
1:B:72:ILE:CD1	1:B:80:ILE:CD1	2.77	0.62
1:B:41:LYS:CB	2:B:467:HOH:O	2.34	0.61
1:A:82:THR:HG22	1:A:110:HIS:CE1	2.35	0.61
1:A:176:MET:HA	1:A:176:MET:CE	2.25	0.61
1:B:184:ARG:NH1	2:B:331:HOH:O	2.32	0.61
1:B:66:VAL:HG22	2:B:401:HOH:O	2.01	0.60
1:A:113:GLU:OE1	1:A:114:GLY:N	2.35	0.60
1:B:156:ASN:C	1:B:156:ASN:HD22	2.09	0.60
1:B:161:LYS:HG3	1:B:162:ASP:H	1.67	0.59
1:A:82:THR:HB	1:A:110:HIS:ND1	2.18	0.58
1:A:129:LYS:HA	1:A:129:LYS:HE3	1.85	0.58
1:A:110:HIS:CG	2:A:341:HOH:O	2.56	0.58
1:A:36:ARG:NH2	2:A:323:HOH:O	2.37	0.57
1:B:11:HIS:CB	2:B:484:HOH:O	2.54	0.56
1:B:73:ARG:NH2	1:B:174:GLN:OE1	2.29	0.56
1:A:82:THR:HA	1:A:110:HIS:HE1	1.67	0.56
1:A:110:HIS:ND1	2:A:341:HOH:O	2.33	0.56
1:A:113:GLU:OE1	1:A:113:GLU:CA	2.53	0.55
1:A:82:THR:HB	1:A:110:HIS:CE1	2.41	0.55
1:A:110:HIS:CD2	2:A:341:HOH:O	2.60	0.55
1:A:97:MET:HA	1:A:97:MET:HE2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:THR:HA	1:A:110:HIS:ND1	2.22	0.54
1:B:117:LYS:NZ	1:B:147:GLU:CG	2.52	0.54
1:B:161:LYS:HG3	1:B:162:ASP:N	2.24	0.53
1:A:41:LYS:HE3	1:A:93:LEU:CD1	2.40	0.51
1:A:113:GLU:CD	2:A:394:HOH:O	2.51	0.51
1:B:56:GLU:HB3	1:B:59:LYS:HE2	1.92	0.51
1:A:157:PRO:HG3	1:A:176:MET:HE3	1.92	0.50
1:A:59:LYS:HE2	1:A:81:ASP:HB3	1.94	0.50
1:B:163:LYS:O	1:B:169:LYS:HG3	2.12	0.49
1:A:218:GLU:HA	1:A:221:ARG:HG2	1.95	0.49
1:A:220:GLN:HA	1:A:220:GLN:OE1	2.13	0.48
1:A:110:HIS:CD2	1:A:111:GLY:O	2.66	0.48
1:B:72:ILE:HD12	1:B:80:ILE:HD12	1.96	0.48
1:B:72:ILE:HD13	1:B:80:ILE:CD1	2.44	0.47
1:B:165:SER:O	1:B:168:GLY:N	2.47	0.47
1:A:156:ASN:HD22	1:A:158:ASP:H	1.61	0.47
1:B:36:ARG:NH2	2:B:414:HOH:O	2.47	0.47
1:B:72:ILE:CD1	1:B:80:ILE:HD12	2.43	0.47
1:A:96:GLU:H	1:A:96:GLU:CD	2.23	0.47
1:A:158:ASP:CG	1:A:184:ARG:HH22	2.23	0.47
1:A:41:LYS:HE3	1:A:93:LEU:HD13	1.96	0.47
1:A:165:SER:O	1:A:168:GLY:N	2.49	0.46
1:B:211:PHE:HB3	1:B:214:LYS:HD3	1.97	0.46
1:A:82:THR:CA	1:A:110:HIS:ND1	2.77	0.46
1:A:82:THR:CG2	1:A:110:HIS:CE1	2.98	0.46
1:A:162:ASP:CG	1:A:165:SER:HB2	2.40	0.46
1:B:162:ASP:O	1:B:168:GLY:HA3	2.15	0.46
1:B:193:GLU:OE1	2:B:375:HOH:O	2.21	0.46
1:A:96:GLU:CD	1:A:96:GLU:N	2.74	0.46
1:B:165:SER:O	1:B:166:PRO:C	2.58	0.46
1:B:85:LYS:HE3	2:B:460:HOH:O	2.15	0.46
1:B:158:ASP:OD1	1:B:184:ARG:NH2	2.49	0.46
1:B:80:ILE:HG23	1:B:83:ILE:HD12	1.98	0.46
1:A:101:ARG:NH1	1:A:101:ARG:CG	2.63	0.45
1:A:36:ARG:HH12	1:A:37:LEU:HD23	1.82	0.45
1:A:92:ALA:HA	1:A:95:SER:O	2.16	0.45
1:A:18:LEU:O	1:A:22:GLU:HG2	2.17	0.45
1:A:156:ASN:HD22	1:A:156:ASN:C	2.25	0.45
1:A:170:LYS:HE2	2:A:387:HOH:O	2.17	0.44
1:B:72:ILE:HD12	1:B:80:ILE:CD1	2.46	0.44
1:B:113:GLU:HG3	2:B:387:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:HIS:HB3	2:B:484:HOH:O	2.15	0.44
1:B:11:HIS:HB2	2:B:484:HOH:O	2.16	0.44
1:B:75:SER:HB2	1:B:152:TYR:OH	2.18	0.44
1:B:117:LYS:HE3	2:B:473:HOH:O	2.18	0.43
1:A:216:TYR:O	1:A:220:GLN:HG2	2.17	0.43
1:A:43:VAL:HG22	1:B:24:GLU:HG2	2.00	0.43
1:B:156:ASN:HD22	1:B:158:ASP:H	1.66	0.43
1:A:36:ARG:HH12	1:A:37:LEU:CD2	2.32	0.42
1:A:52:ASP:OD2	1:A:85:LYS:NZ	2.47	0.42
1:A:33:GLU:OE2	1:A:36:ARG:NH1	2.53	0.41
1:A:82:THR:CB	1:A:110:HIS:ND1	2.83	0.41
2:A:384:HOH:O	1:B:22:GLU:HB2	2.20	0.41
1:A:68:PHE:CZ	1:A:80:ILE:HD12	2.55	0.41
1:A:110:HIS:NE2	2:A:341:HOH:O	2.42	0.41

All (7) 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)
2:A:376:HOH:O	2:B:393:HOH:O[4_555]	1.59	0.61
1:B:96:GLU:OE1	1:B:113:GLU:OE2[3_444]	1.76	0.44
1:B:96:GLU:OE2	2:B:489:HOH:O[3_444]	1.98	0.22
2:B:492:HOH:O	2:B:495:HOH:O[3_454]	2.02	0.18
1:B:96:GLU:OE2	1:B:113:GLU:OE2[3_444]	2.09	0.11
1:B:96:GLU:CD	1:B:113:GLU:OE2[3_444]	2.12	0.08
1:B:96:GLU:CD	2:B:489:HOH:O[3_444]	2.17	0.03

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	204/259 (79%)	199 (98%)	5 (2%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	201/259 (78%)	196 (98%)	3 (2%)	2 (1%)	12	8
All	All	405/518 (78%)	395 (98%)	8 (2%)	2 (0%)	24	21

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	55	ALA
1	B	166	PRO

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	186/230 (81%)	174 (94%)	12 (6%)	15	12
1	B	183/230 (80%)	178 (97%)	5 (3%)	39	42
All	All	369/460 (80%)	352 (95%)	17 (5%)	24	22

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	36	ARG
1	A	67	SER
1	A	80	ILE
1	A	84	ASP
1	A	87	LYS
1	A	97	MET
1	A	101	ARG
1	A	113	GLU
1	A	129	LYS
1	A	156	ASN
1	A	176	MET
1	B	75	SER
1	B	113	GLU

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Mol	Chain	Res	Type
1	B	156	ASN
1	B	184	ARG
1	B	203	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	108	HIS
1	A	156	ASN
1	A	174	GLN
1	B	70	ASN
1	B	150	GLN
1	B	156	ASN

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	208/259 (80%)	0.78	26 (12%) 8 7	22, 35, 57, 73	0
1	B	205/259 (79%)	0.48	22 (10%) 11 10	19, 29, 46, 63	0
All	All	413/518 (79%)	0.63	48 (11%) 9 8	19, 32, 54, 73	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	162	ASP	6.1
1	B	56	GLU	5.2
1	A	167	GLU	4.9
1	A	165	SER	4.7
1	A	110	HIS	4.7
1	B	164	ARG	4.5
1	A	166	PRO	3.9
1	A	8	GLU	3.8
1	A	136	ASP	3.7
1	B	176	MET	3.7
1	B	136	ASP	3.5
1	B	161	LYS	3.5
1	A	84	ASP	3.4
1	B	53	GLU	3.4
1	A	55	ALA	3.4
1	B	55	ALA	3.3
1	A	163	LYS	3.3
1	B	20	LYS	3.2
1	B	96	GLU	3.2
1	B	113	GLU	3.1
1	A	168	GLY	3.1
1	B	110	HIS	3.0
1	A	164	ARG	3.0
1	B	94	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	170	LYS	2.9
1	A	96	GLU	2.9
1	B	11	HIS	2.9
1	A	10	GLU	2.8
1	A	9	ASN	2.8
1	A	221	ARG	2.8
1	A	217	ALA	2.8
1	B	169	LYS	2.7
1	B	162	ASP	2.7
1	A	113	GLU	2.7
1	A	161	LYS	2.7
1	B	170	LYS	2.5
1	B	36	ARG	2.5
1	B	221	ARG	2.5
1	B	15	PHE	2.5
1	B	171	LYS	2.5
1	A	94	GLU	2.4
1	A	169	LYS	2.4
1	B	13	LYS	2.2
1	A	171	LYS	2.1
1	A	56	GLU	2.1
1	A	218	GLU	2.1
1	B	129	LYS	2.0
1	A	216	TYR	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.