



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:47 PM UTC

PDB ID : 2CKH / pdb_00002ckh
Title : SENP1-SUMO2 complex
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Deposited on : 2006-04-18
Resolution : 3.20 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

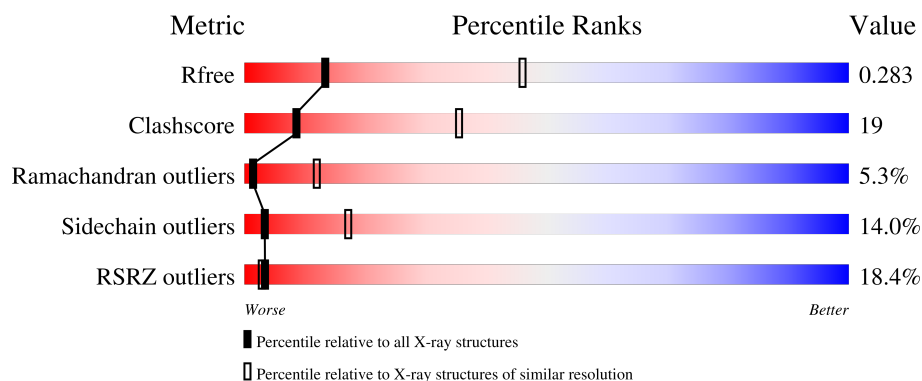
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	225	17% 49% 41% 8%
2	B	79	23% 70% 20% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2515 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 SENTRIN-SPECIFIC PROTEASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	225	Total	C	N	O	S	0	0	0
			1882	1208	326	334	14			

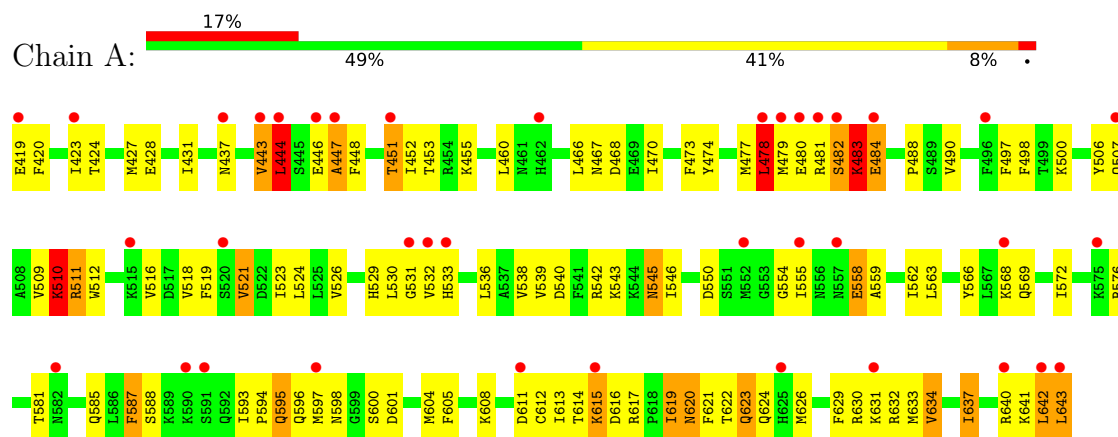
- Molecule 2 is a protein called SMALL UBIQUITIN-RELATED MODIFIER 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	79	Total	C	N	O	S	0	0	0
			633	390	115	124	4			

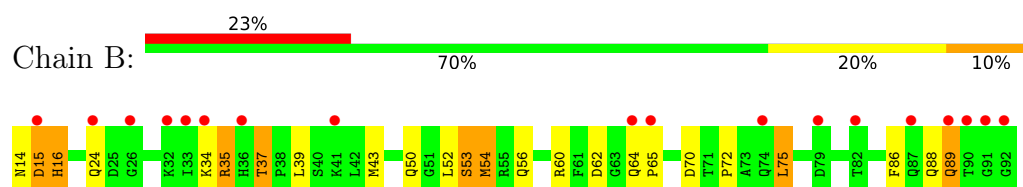
3 Residue-property plots [i](#)

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: SENTRIN-SPECIFIC PROTEASE 1



• Molecule 2: SMALL UBIQUITIN-RELATED MODIFIER 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	143.38Å 143.38Å 71.89Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	124.03 – 3.20 124.03 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (124.03-3.20) 99.0 (124.03-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.245 , 0.294 0.238 , 0.283	Depositor DCC
R_{free} test set	716 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	92.9	Xtriage
Anisotropy	0.893	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 140.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	2515	wwPDB-VP
Average B, all atoms (Å ²)	129.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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	0.60	1/1925 (0.1%)	0.94	9/2587 (0.3%)
2	B	0.44	0/642	0.83	0/861
All	All	0.56	1/2567 (0.0%)	0.91	9/3448 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	615	LYS	CE-NZ	7.51	1.71	1.49

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	510	LYS	CA-C-N	6.76	133.88	121.70
1	A	510	LYS	C-N-CA	6.76	133.88	121.70
1	A	478	LEU	CA-C-N	6.14	128.79	120.44
1	A	478	LEU	C-N-CA	6.14	128.79	120.44
1	A	420	PHE	N-CA-C	5.86	118.10	109.42
1	A	587	PHE	N-CA-C	5.80	116.21	108.38
1	A	545	ASN	N-CA-C	5.73	117.58	107.61
1	A	443	VAL	N-CA-C	5.06	115.83	110.72
1	A	444	LEU	N-CA-C	5.04	121.55	110.80

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	1882	0	1883	83	0
2	B	633	0	618	16	0
All	All	2515	0	2501	95	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 19.

All (95) 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:615:LYS:CE	1:A:615:LYS:NZ	1.71	1.48
1:A:446:GLU:HG3	1:A:447:ALA:H	1.16	1.08
1:A:593:ILE:HB	1:A:594:PRO:HD2	1.42	1.00
1:A:446:GLU:CG	1:A:447:ALA:H	1.74	0.99
1:A:478:LEU:HD22	1:A:481:ARG:HH12	1.25	0.99
2:B:70:ASP:HB3	2:B:75:LEU:HD21	1.45	0.97
1:A:633:MET:O	1:A:637:ILE:HG12	1.77	0.85
1:A:446:GLU:HG3	1:A:447:ALA:N	1.95	0.81
1:A:593:ILE:HB	1:A:594:PRO:CD	2.09	0.81
1:A:510:LYS:HB3	1:A:511:ARG:HB2	1.62	0.81
1:A:615:LYS:C	1:A:617:ARG:H	1.93	0.78
1:A:510:LYS:CB	1:A:511:ARG:HB2	2.13	0.77
1:A:524:LEU:HB2	1:A:539:VAL:HB	1.66	0.77
1:A:526:VAL:O	1:A:536:LEU:HD12	1.89	0.72
1:A:545:ASN:HB3	1:A:585:GLN:HB3	1.72	0.71
1:A:477:MET:HB3	1:A:642:LEU:HD21	1.72	0.69
1:A:451:THR:C	1:A:452:ILE:HD12	2.17	0.69
1:A:478:LEU:HD22	1:A:481:ARG:NH1	2.03	0.68
1:A:595:GLN:HE21	1:A:595:GLN:HA	1.61	0.66
2:B:70:ASP:HB3	2:B:75:LEU:CD2	2.23	0.66
1:A:633:MET:HG2	1:A:637:ILE:HD11	1.78	0.66
1:A:447:ALA:O	1:A:448:PHE:HB2	1.97	0.64
1:A:478:LEU:HD11	1:A:611:ASP:HA	1.79	0.64
1:A:539:VAL:HG22	1:A:546:ILE:HG12	1.79	0.63
1:A:598:ASN:HD21	1:A:601:ASP:HB2	1.62	0.62
1:A:500:LYS:HE2	2:B:86:PHE:HE2	1.64	0.62
1:A:615:LYS:O	1:A:617:ARG:N	2.35	0.60
1:A:478:LEU:CD2	1:A:481:ARG:HH12	2.08	0.59
1:A:558:GLU:O	1:A:562:ILE:HG13	2.03	0.59
1:A:477:MET:CB	1:A:642:LEU:HD21	2.33	0.58
1:A:490:VAL:HG22	1:A:523:ILE:HB	1.84	0.58
1:A:424:THR:H	1:A:427:MET:HE2	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:615:LYS:C	1:A:617:ARG:N	2.62	0.57
1:A:615:LYS:NZ	1:A:615:LYS:CD	2.62	0.57
1:A:444:LEU:HB2	1:A:452:ILE:O	2.05	0.57
1:A:460:LEU:HD21	1:A:470:ILE:HG21	1.87	0.56
1:A:643:LEU:HD23	1:A:643:LEU:N	2.20	0.56
1:A:480:GLU:O	1:A:483:LYS:HB3	2.06	0.56
1:A:446:GLU:CG	1:A:447:ALA:N	2.51	0.55
2:B:39:LEU:HD12	2:B:70:ASP:O	2.06	0.55
2:B:60:ARG:NH2	2:B:65:PRO:HG3	2.22	0.55
1:A:558:GLU:CG	1:A:562:ILE:HD11	2.39	0.53
1:A:446:GLU:CD	1:A:447:ALA:H	2.17	0.53
1:A:555:ILE:HG23	1:A:588:SER:HB2	1.91	0.53
1:A:594:PRO:HD3	1:A:619:ILE:HG21	1.93	0.51
1:A:473:PHE:CD2	1:A:637:ILE:HG23	2.45	0.50
1:A:608:LYS:HG3	1:A:629:PHE:CD2	2.46	0.49
1:A:468:ASP:HB3	2:B:89:GLN:HG2	1.95	0.49
1:A:512:TRP:CD1	2:B:62:ASP:HA	2.48	0.49
2:B:14:ASN:HD22	2:B:35:ARG:HD3	1.78	0.48
1:A:563:LEU:O	1:A:566:TYR:HB3	2.14	0.48
1:A:604:MET:HE2	1:A:630:ARG:HA	1.96	0.48
1:A:481:ARG:NH1	1:A:642:LEU:HD22	2.29	0.48
2:B:37:THR:O	2:B:72:PRO:HD2	2.13	0.47
1:A:519:PHE:C	1:A:521:VAL:H	2.23	0.47
1:A:624:GLN:H	1:A:624:GLN:CD	2.22	0.47
2:B:53:SER:O	2:B:54:MET:CB	2.61	0.47
1:A:478:LEU:HD13	1:A:614:THR:OG1	2.15	0.47
1:A:538:VAL:HG21	1:A:613:ILE:HD13	1.95	0.47
1:A:540:ASP:OD2	1:A:543:LYS:HD3	2.15	0.47
1:A:453:THR:OG1	1:A:455:LYS:HG2	2.16	0.46
2:B:39:LEU:O	2:B:43:MET:HG3	2.15	0.46
1:A:509:VAL:O	1:A:510:LYS:C	2.59	0.46
1:A:478:LEU:CD1	1:A:611:ASP:HA	2.45	0.45
1:A:478:LEU:HD11	1:A:611:ASP:CA	2.46	0.45
1:A:608:LYS:HE2	1:A:629:PHE:CZ	2.52	0.45
2:B:16:HIS:HA	2:B:34:LYS:HA	1.98	0.45
1:A:424:THR:OG1	1:A:427:MET:HG3	2.16	0.45
1:A:510:LYS:CA	1:A:511:ARG:HB2	2.46	0.45
1:A:482:SER:HB2	1:A:488:PRO:O	2.17	0.44
1:A:506:TYR:O	1:A:507:GLN:C	2.60	0.44
1:A:490:VAL:HG21	1:A:614:THR:HG22	2.00	0.44
1:A:452:ILE:HD12	1:A:452:ILE:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:623:GLN:O	1:A:626:MET:HB2	2.18	0.43
1:A:529:HIS:NE2	2:B:89:GLN:HB2	2.32	0.43
1:A:623:GLN:O	1:A:626:MET:N	2.34	0.43
1:A:474:TYR:CE1	1:A:633:MET:HE2	2.54	0.43
1:A:594:PRO:HB2	1:A:605:PHE:HE2	1.84	0.43
1:A:447:ALA:O	1:A:448:PHE:CB	2.67	0.42
2:B:15:ASP:O	2:B:35:ARG:HG2	2.19	0.42
2:B:70:ASP:CB	2:B:75:LEU:HD21	2.33	0.42
1:A:550:ASP:OD2	1:A:554:GLY:N	2.52	0.42
1:A:559:ALA:HA	1:A:562:ILE:HD12	2.00	0.42
2:B:56:GLN:O	2:B:88:GLN:HB2	2.20	0.42
1:A:428:GLU:HA	1:A:431:ILE:HD12	2.02	0.42
1:A:488:PRO:HB3	1:A:542:ARG:HH21	1.84	0.42
1:A:530:LEU:O	1:A:533:HIS:HB3	2.20	0.41
1:A:497:PHE:O	1:A:498:PHE:C	2.63	0.41
1:A:481:ARG:O	1:A:484:GLU:CG	2.69	0.41
1:A:569:GLN:HA	1:A:572:ILE:HD12	2.02	0.41
1:A:631:LYS:O	1:A:634:VAL:HG23	2.21	0.41
1:A:558:GLU:HG3	1:A:562:ILE:HD11	2.02	0.41
1:A:605:PHE:HA	1:A:626:MET:HE1	2.02	0.41
1:A:620:ASN:HD22	1:A:621:PHE:N	2.19	0.41
1:A:612:CYS:SG	1:A:621:PHE:HE2	2.44	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	223/225 (99%)	185 (83%)	25 (11%)	13 (6%)	1	10
2	B	77/79 (98%)	69 (90%)	5 (6%)	3 (4%)	2	18
All	All	300/304 (99%)	254 (85%)	30 (10%)	16 (5%)	1	12

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	447	ALA
1	A	511	ARG
1	A	616	ASP
1	A	640	ARG
1	A	482	SER
1	A	510	LYS
1	A	531	GLY
1	A	576	ARG
1	A	623	GLN
1	A	641	LYS
2	B	54	MET
1	A	444	LEU
1	A	483	LYS
2	B	53	SER
2	B	89	GLN
1	A	437	ASN

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	208/208 (100%)	178 (86%)	30 (14%)	3	16
2	B	70/70 (100%)	61 (87%)	9 (13%)	4	20
All	All	278/278 (100%)	239 (86%)	39 (14%)	3	17

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

Mol	Chain	Res	Type
1	A	419	GLU
1	A	423	ILE
1	A	443	VAL
1	A	451	THR
1	A	466	LEU
1	A	467	ASN
1	A	478	LEU

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Mol	Chain	Res	Type
1	A	479	MET
1	A	483	LYS
1	A	484	GLU
1	A	516	VAL
1	A	518	VAL
1	A	521	VAL
1	A	532	VAL
1	A	558	GLU
1	A	568	LYS
1	A	581	THR
1	A	587	PHE
1	A	595	GLN
1	A	596	GLN
1	A	597	MET
1	A	600	SER
1	A	619	ILE
1	A	620	ASN
1	A	622	THR
1	A	632	ARG
1	A	634	VAL
1	A	637	ILE
1	A	642	LEU
1	A	643	LEU
2	B	15	ASP
2	B	16	HIS
2	B	24	GLN
2	B	35	ARG
2	B	37	THR
2	B	50	GLN
2	B	52	LEU
2	B	64	GLN
2	B	75	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	437	ASN
1	A	491	HIS
1	A	545	ASN
1	A	595	GLN
1	A	596	GLN
1	A	620	ASN

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Mol	Chain	Res	Type
1	A	624	GLN
1	A	625	HIS
2	B	14	ASN
2	B	18	ASN
2	B	36	HIS
2	B	50	GLN
2	B	64	GLN
2	B	67	ASN
2	B	88	GLN
2	B	89	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	225/225 (100%)	1.28	38 (16%) 4 3	102, 124, 134, 146	0
2	B	79/79 (100%)	1.54	18 (22%) 2 2	127, 148, 153, 154	0
All	All	304/304 (100%)	1.35	56 (18%) 3 3	102, 126, 150, 154	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	91	GLY	8.9
2	B	92	GLY	5.2
2	B	90	THR	4.6
1	A	423	ILE	3.8
1	A	482	SER	3.7
2	B	89	GLN	3.6
1	A	520	SER	3.6
1	A	478	LEU	3.4
1	A	443	VAL	3.2
1	A	591	SER	3.2
2	B	33	ILE	3.1
1	A	590	LYS	2.9
1	A	447	ALA	2.9
1	A	642	LEU	2.8
2	B	24	GLN	2.8
1	A	479	MET	2.7
1	A	582	ASN	2.7
1	A	446	GLU	2.6
2	B	34	LYS	2.6
1	A	531	GLY	2.6
1	A	496	PHE	2.5
1	A	515	LYS	2.4
1	A	444	LEU	2.4
2	B	82	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	532	VAL	2.4
1	A	555	ILE	2.4
2	B	41	LYS	2.4
1	A	419	GLU	2.4
2	B	15	ASP	2.3
2	B	79	ASP	2.3
2	B	32	LYS	2.3
2	B	65	PRO	2.3
1	A	568	LYS	2.3
1	A	643	LEU	2.3
1	A	557	ASN	2.3
1	A	484	GLU	2.2
1	A	552	MET	2.2
2	B	74	GLN	2.2
1	A	480	GLU	2.2
1	A	597	MET	2.2
1	A	611	ASP	2.2
1	A	462	HIS	2.2
1	A	533	HIS	2.2
2	B	87	GLN	2.1
1	A	615	LYS	2.1
1	A	640	ARG	2.1
1	A	625	HIS	2.1
2	B	36	HIS	2.1
1	A	481	ARG	2.1
1	A	575	LYS	2.1
1	A	507	GLN	2.1
2	B	64	GLN	2.1
2	B	26	GLY	2.1
1	A	631	LYS	2.1
1	A	437	ASN	2.0
1	A	451	THR	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.