



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:29 AM UTC

PDB ID : 3G06 / pdb_00003g06
Title : The Salmonella Virulence Effector SspH2 Functions As A Novel E3 Ligase
Authors : Quezada, C.M.; Stebbins, C.E.
Deposited on : 2009-01-27
Resolution : 1.90 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

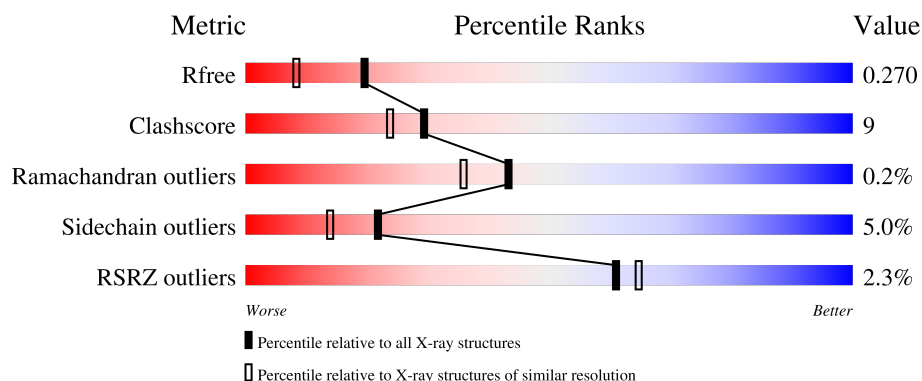
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

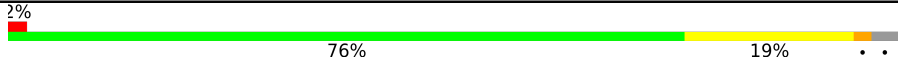
The reported resolution of this entry is 1.90 Å.

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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	622	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5069 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 SspH2 (Leucine-rich repeat protein).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	601	Total	C	N	O	S	0	0	0
			4663	2951	803	891	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	162	GLY	-	expression tag	UNP Q9RPH0
A	163	PRO	-	expression tag	UNP Q9RPH0
A	164	VAL	-	expression tag	UNP Q9RPH0
A	165	ASP	-	expression tag	UNP Q9RPH0

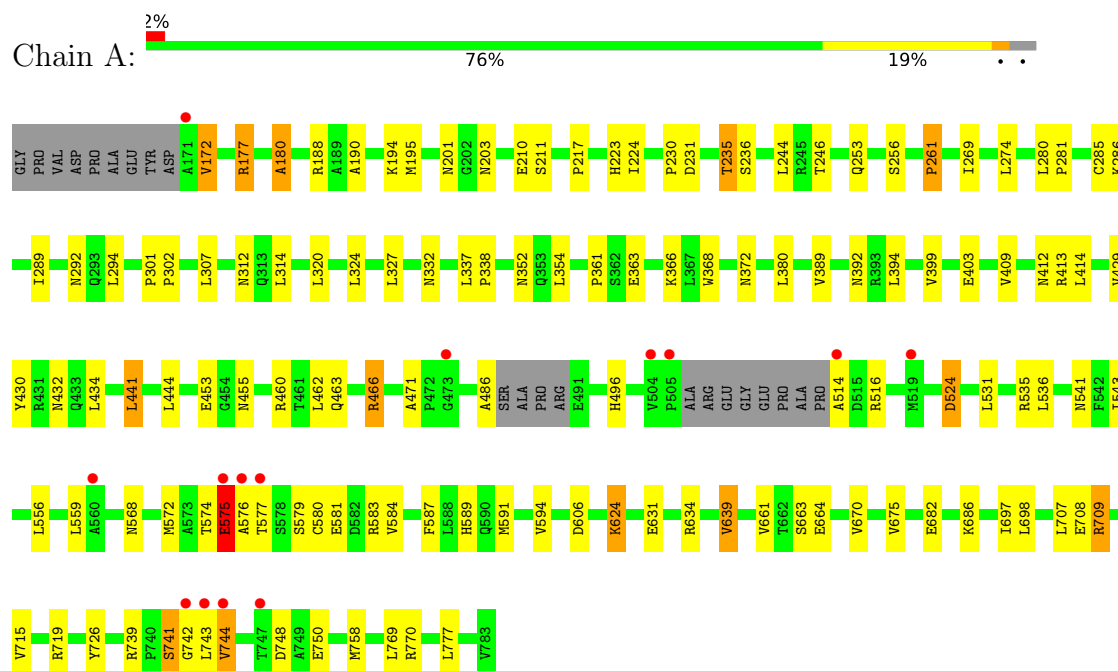
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	406	Total	O	0	0
			406	406		

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: SspH2 (Leucine-rich repeat protein)



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	76.89Å 76.89Å 281.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 50.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.1 (50.00-1.90) 98.1 (50.00-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.4.0063	Depositor
R, R_{free}	0.219 , 0.264 (Not available) , 0.270	Depositor DCC
R_{free} test set	3369 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5069	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 4.28% 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	1.19	4/4753 (0.1%)	1.22	26/6468 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	224	ILE	CA-CB	6.60	1.61	1.54
1	A	429	VAL	CA-CB	6.57	1.63	1.54
1	A	726	TYR	C-O	-6.44	1.16	1.24
1	A	399	VAL	CA-CB	6.21	1.60	1.54

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	663	SER	N-CA-C	8.24	121.38	111.82
1	A	575	GLU	N-CA-C	-7.96	97.30	109.95
1	A	639	VAL	CB-CA-C	-7.79	97.23	110.71
1	A	261	PRO	CA-C-N	7.25	127.88	119.47
1	A	261	PRO	C-N-CA	7.25	127.88	119.47
1	A	172	VAL	N-CA-C	-7.01	103.36	111.00
1	A	471	ALA	CA-C-N	-6.97	112.75	120.14
1	A	471	ALA	C-N-CA	-6.97	112.75	120.14
1	A	739	ARG	CA-C-N	-6.85	111.75	119.28
1	A	739	ARG	C-N-CA	-6.85	111.75	119.28
1	A	380	LEU	CA-C-N	6.60	126.63	119.90
1	A	380	LEU	C-N-CA	6.60	126.63	119.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	709	ARG	NE-CZ-NH2	-6.18	113.63	119.20
1	A	180	ALA	N-CA-C	6.16	118.55	109.50
1	A	337	LEU	CA-C-N	-6.07	114.38	120.21
1	A	337	LEU	C-N-CA	-6.07	114.38	120.21
1	A	180	ALA	CA-C-N	-5.99	113.79	119.90
1	A	180	ALA	C-N-CA	-5.99	113.79	119.90
1	A	580	CYS	CA-C-N	5.58	128.07	120.54
1	A	580	CYS	C-N-CA	5.58	128.07	120.54
1	A	709	ARG	N-CA-C	5.47	117.32	111.36
1	A	581	GLU	N-CA-CB	-5.38	101.92	109.94
1	A	758	MET	CA-C-N	5.33	127.69	120.44
1	A	758	MET	C-N-CA	5.33	127.69	120.44
1	A	543	ILE	N-CA-C	5.25	115.87	110.36
1	A	235	THR	N-CA-CB	-5.03	104.22	110.90

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	574	THR	Peptide
1	A	575	GLU	Peptide
1	A	741	SER	Peptide

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	4663	0	4711	81	0
2	A	406	0	0	13	0
All	All	5069	0	4711	81	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 9.

All (81) 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:246:THR:HG22	2:A:967:HOH:O	1.45	1.16
1:A:576:ALA:HB2	1:A:587:PHE:HB2	1.46	0.95
1:A:742:GLY:HA3	1:A:743:LEU:HG	1.49	0.95
1:A:742:GLY:CA	1:A:743:LEU:HG	2.04	0.86
1:A:576:ALA:HB1	1:A:584:VAL:HA	1.60	0.84
1:A:414:LEU:H	1:A:432:ASN:HD22	1.31	0.78
1:A:352:ASN:HB2	1:A:372:ASN:HD21	1.50	0.74
1:A:631:GLU:OE2	1:A:634:ARG:NH1	2.20	0.74
1:A:434:LEU:H	1:A:455:ASN:HD22	1.35	0.74
1:A:413:ARG:HD2	2:A:975:HOH:O	1.87	0.74
1:A:496:HIS:HD2	2:A:898:HOH:O	1.72	0.72
1:A:177:ARG:CD	2:A:970:HOH:O	2.37	0.71
1:A:576:ALA:HB2	1:A:587:PHE:CB	2.20	0.70
1:A:392:ASN:HB2	1:A:412:ASN:HD21	1.58	0.69
1:A:432:ASN:HB2	1:A:455:ASN:HD21	1.59	0.68
1:A:177:ARG:HD3	2:A:970:HOH:O	1.95	0.67
1:A:741:SER:HB2	1:A:742:GLY:HA3	1.78	0.66
1:A:742:GLY:HA3	1:A:743:LEU:CG	2.24	0.66
1:A:514:ALA:N	2:A:971:HOH:O	2.30	0.64
1:A:741:SER:CB	1:A:742:GLY:HA3	2.28	0.63
1:A:403:GLU:HB2	2:A:122:HOH:O	1.99	0.63
1:A:675:VAL:HG23	2:A:1000:HOH:O	1.97	0.63
1:A:535:ARG:NH2	1:A:579:SER:O	2.34	0.61
1:A:312:ASN:HB2	1:A:332:ASN:HD21	1.64	0.60
1:A:177:ARG:HD2	2:A:970:HOH:O	1.98	0.59
1:A:180:ALA:HB2	1:A:188:ARG:HD2	1.85	0.58
1:A:587:PHE:O	1:A:591:MET:HG3	2.03	0.58
1:A:332:ASN:HB2	1:A:352:ASN:HD21	1.68	0.58
1:A:742:GLY:HA2	1:A:743:LEU:HG	1.85	0.57
1:A:572:MET:O	1:A:575:GLU:O	2.22	0.56
1:A:432:ASN:CB	1:A:455:ASN:HD21	2.18	0.55
1:A:576:ALA:CB	1:A:584:VAL:HA	2.35	0.55
1:A:606:ASP:OD2	1:A:709:ARG:HD2	2.07	0.55
1:A:412:ASN:HB2	1:A:432:ASN:HD21	1.71	0.54
1:A:363:GLU:HG3	2:A:1018:HOH:O	2.09	0.52
1:A:244:LEU:HB3	1:A:261:PRO:HG2	1.91	0.51
1:A:589:HIS:CE1	1:A:661:VAL:HG13	2.46	0.51
1:A:392:ASN:CB	1:A:412:ASN:HD21	2.22	0.51
1:A:541:ASN:HD21	1:A:664:GLU:H	1.60	0.50
1:A:314:LEU:H	1:A:332:ASN:HD22	1.60	0.49
1:A:210:GLU:HG3	1:A:230:PRO:HB2	1.94	0.49
1:A:707:LEU:HD13	1:A:769:LEU:HD22	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:715:VAL:O	1:A:719:ARG:HG2	2.14	0.47
1:A:414:LEU:H	1:A:432:ASN:ND2	2.04	0.47
1:A:285:CYS:O	1:A:286:LYS:HD3	2.15	0.47
1:A:294:LEU:H	1:A:312:ASN:HD22	1.63	0.47
1:A:466:ARG:HE	1:A:466:ARG:HB2	1.33	0.46
1:A:320:LEU:HG	1:A:338:PRO:HG2	1.97	0.46
1:A:568:ASN:O	1:A:572:MET:HG3	2.15	0.46
1:A:575:GLU:H	1:A:575:GLU:HG2	1.67	0.46
1:A:195:MET:HE1	1:A:217:PRO:HG3	1.98	0.45
1:A:486:ALA:C	2:A:968:HOH:O	2.58	0.45
1:A:190:ALA:O	1:A:194:LYS:HG3	2.17	0.45
1:A:579:SER:HB3	1:A:583:ARG:HB3	1.98	0.45
1:A:430:TYR:HA	1:A:453:GLU:O	2.16	0.45
1:A:274:LEU:HD12	1:A:274:LEU:HA	1.56	0.45
1:A:269:ILE:CG2	1:A:289:ILE:HG22	2.48	0.44
1:A:708:GLU:OE1	1:A:719:ARG:NH2	2.47	0.44
1:A:307:LEU:HD23	1:A:327:LEU:HD13	2.00	0.43
1:A:698:LEU:HD21	1:A:770:ARG:HA	2.00	0.43
1:A:223:HIS:CD2	1:A:223:HIS:H	2.34	0.43
1:A:354:LEU:H	1:A:372:ASN:HD22	1.64	0.43
1:A:572:MET:SD	1:A:594:VAL:HG21	2.58	0.43
1:A:434:LEU:H	1:A:455:ASN:ND2	2.09	0.43
1:A:741:SER:CB	1:A:742:GLY:CA	2.97	0.43
1:A:203:ASN:O	1:A:223:HIS:HE1	2.02	0.42
1:A:361:PRO:HB2	2:A:966:HOH:O	2.19	0.42
1:A:394:LEU:H	1:A:412:ASN:HD22	1.66	0.42
1:A:441:LEU:HD12	1:A:441:LEU:O	2.20	0.42
1:A:744:VAL:O	1:A:750:GLU:OE1	2.38	0.42
1:A:389:VAL:CG2	1:A:409:VAL:HG12	2.50	0.41
1:A:201:ASN:OD1	1:A:203:ASN:HB2	2.20	0.41
1:A:280:LEU:HA	1:A:281:PRO:HD3	1.88	0.41
1:A:682:GLU:OE2	1:A:686:LYS:HE3	2.21	0.41
1:A:624:LYS:HD3	2:A:877:HOH:O	2.20	0.41
1:A:301:PRO:HA	1:A:302:PRO:HD3	1.98	0.41
1:A:256:SER:HA	1:A:274:LEU:HD11	2.02	0.41
1:A:292:ASN:H	1:A:312:ASN:HD21	1.68	0.41
1:A:460:ARG:HH11	1:A:463:GLN:CD	2.29	0.40
1:A:366:LYS:HE3	1:A:368:TRP:CD2	2.56	0.40
1:A:556:LEU:HD23	1:A:556:LEU:HA	1.91	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	459/622 (74%)	436 (95%)	22 (5%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	524	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	516/531 (97%)	490 (95%)	26 (5%)	22	14

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	VAL
1	A	177	ARG
1	A	211	SER
1	A	231	ASP
1	A	235	THR
1	A	236	SER
1	A	253	GLN
1	A	324	LEU
1	A	441	LEU
1	A	444	LEU

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Mol	Chain	Res	Type
1	A	462	LEU
1	A	466	ARG
1	A	516	ARG
1	A	524	ASP
1	A	531	LEU
1	A	536	LEU
1	A	559	LEU
1	A	575	GLU
1	A	577	THR
1	A	624	LYS
1	A	639	VAL
1	A	670	VAL
1	A	697	ILE
1	A	744	VAL
1	A	748	ASP
1	A	777	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	223	HIS
1	A	276	HIS
1	A	312	ASN
1	A	332	ASN
1	A	352	ASN
1	A	372	ASN
1	A	412	ASN
1	A	432	ASN
1	A	455	ASN
1	A	496	HIS
1	A	541	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	601/622 (96%)	0.09	14 (2%) 61 65	21, 33, 51, 66	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	576	ALA	6.7
1	A	575	GLU	4.1
1	A	171	ALA	3.7
1	A	743	LEU	3.5
1	A	742	GLY	3.1
1	A	747	THR	2.9
1	A	577	THR	2.8
1	A	744	VAL	2.8
1	A	514	ALA	2.7
1	A	505	PRO	2.6
1	A	560	ALA	2.4
1	A	519	MET	2.4
1	A	473	GLY	2.2
1	A	504	VAL	2.1

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.