



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2024 – 01:49 PM EST

PDB ID : 2TS1
Title : STRUCTURE OF TYROSYL-T/RNA SYNTHETASE REFINED AT 2.3
ANGSTROMS RESOLUTION. INTERACTION OF THE ENZYME WITH
THE TYROSYL ADENYLATE INTERMEDIATE
Authors : Brick, P.; Bhat, T.N.; Blow, D.M.
Deposited on : 1989-06-29
Resolution : 2.30 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

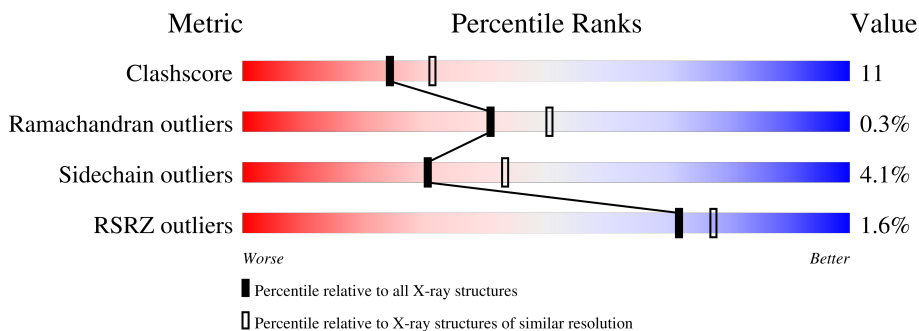
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	419	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2567 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 TYROSYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	317	2457	1568	425	457	7	0	0	0

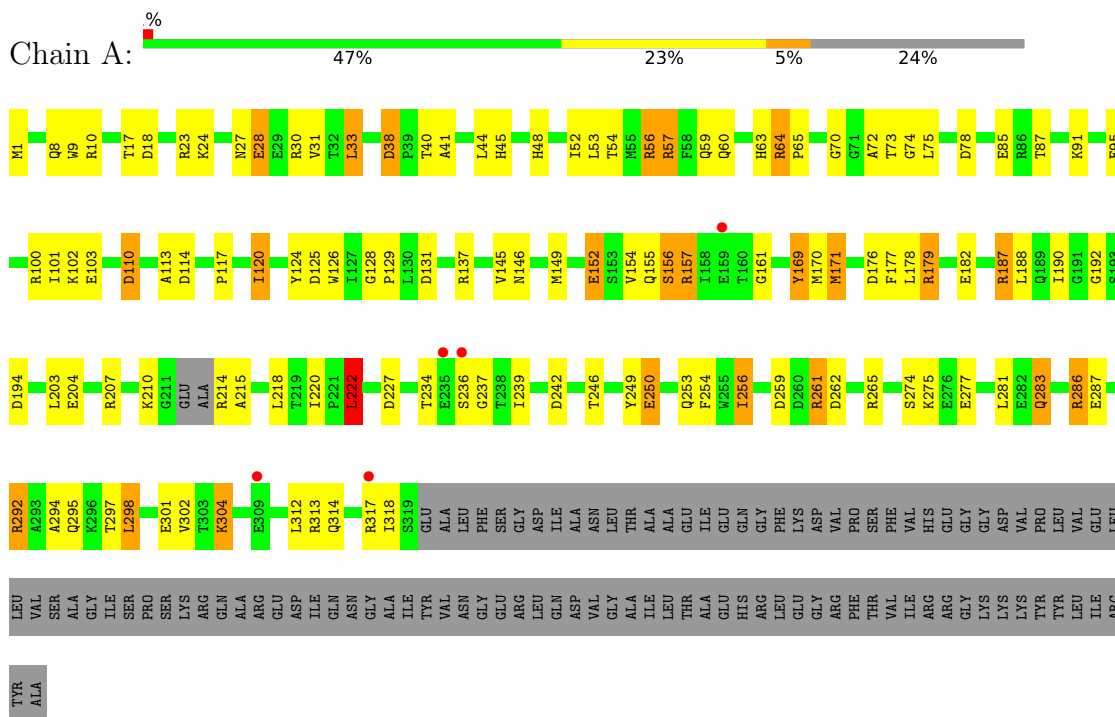
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	110	Total	O	0	0
			110	110		

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: TYROSYL-TRNA SYNTHETASE



4 Data and refinement statistics i

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	64.46Å 64.46Å 237.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.30 45.63 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.30) 97.5 (45.63-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 2.10Å)	Xtrriage
Refinement program	PROLSQ	Depositor
R, R_{free}	0.228 , (Not available) 0.236 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.8	Xtrriage
Anisotropy	0.195	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 86.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.055 for -h,-k,l	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	2567	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.58% 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	1.05	1/2505 (0.0%)	2.02	76/3389 (2.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	GLU	CB-CG	-5.96	1.40	1.52

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	ARG	NE-CZ-NH1	14.35	127.47	120.30
1	A	157	ARG	NE-CZ-NH2	-14.17	113.21	120.30
1	A	56	ARG	NE-CZ-NH2	-13.39	113.61	120.30
1	A	23	ARG	NE-CZ-NH2	-13.37	113.61	120.30
1	A	261	ARG	NE-CZ-NH1	12.01	126.30	120.30
1	A	157	ARG	NE-CZ-NH1	11.06	125.83	120.30
1	A	179	ARG	NE-CZ-NH1	10.86	125.73	120.30
1	A	23	ARG	NE-CZ-NH1	10.76	125.68	120.30
1	A	261	ARG	NE-CZ-NH2	-10.60	115.00	120.30
1	A	292	ARG	NE-CZ-NH1	10.46	125.53	120.30
1	A	64	ARG	CD-NE-CZ	10.35	138.10	123.60
1	A	110	ASP	CB-CG-OD1	10.33	127.59	118.30
1	A	10	ARG	NE-CZ-NH1	10.09	125.35	120.30
1	A	137	ARG	NE-CZ-NH2	9.88	125.24	120.30
1	A	313	ARG	NE-CZ-NH1	9.01	124.81	120.30
1	A	137	ARG	CD-NE-CZ	8.96	136.14	123.60
1	A	179	ARG	NE-CZ-NH2	-8.93	115.83	120.30
1	A	41	ALA	N-CA-CB	8.45	121.92	110.10
1	A	286	ARG	NE-CZ-NH2	-8.07	116.26	120.30
1	A	169	TYR	CB-CG-CD1	-7.96	116.22	121.00
1	A	56	ARG	NH1-CZ-NH2	7.77	127.94	119.40
1	A	259	ASP	CB-CG-OD2	7.66	125.20	118.30
1	A	313	ARG	CD-NE-CZ	7.66	134.32	123.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	261	ARG	CG-CD-NE	7.44	127.43	111.80
1	A	57	ARG	NE-CZ-NH2	-7.27	116.67	120.30
1	A	125	ASP	CB-CG-OD1	7.19	124.77	118.30
1	A	194	ASP	CB-CG-OD2	7.16	124.75	118.30
1	A	85	GLU	OE1-CD-OE2	7.15	131.88	123.30
1	A	28	GLU	CA-CB-CG	7.13	129.10	113.40
1	A	283	GLN	CB-CG-CD	7.03	129.89	111.60
1	A	38	ASP	CB-CG-OD1	-6.90	112.09	118.30
1	A	227	ASP	CB-CG-OD1	6.74	124.37	118.30
1	A	171	MET	CA-CB-CG	6.61	124.54	113.30
1	A	298	LEU	CB-CG-CD2	-6.57	99.83	111.00
1	A	261	ARG	CD-NE-CZ	6.55	132.77	123.60
1	A	57	ARG	NE-CZ-NH1	6.45	123.52	120.30
1	A	207	ARG	NE-CZ-NH1	6.44	123.52	120.30
1	A	187	ARG	CD-NE-CZ	-6.41	114.63	123.60
1	A	157	ARG	N-CA-C	6.36	128.16	111.00
1	A	131	ASP	CB-CG-OD1	6.34	124.01	118.30
1	A	30	ARG	N-CA-CB	6.32	121.97	110.60
1	A	214	ARG	N-CA-CB	-6.24	99.37	110.60
1	A	317	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	126	TRP	CA-C-O	-5.94	107.62	120.10
1	A	87	THR	N-CA-CB	5.91	121.54	110.30
1	A	286	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	A	18	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	101	ILE	CB-CA-C	5.82	123.24	111.60
1	A	182	GLU	OE1-CD-OE2	-5.79	116.36	123.30
1	A	283	GLN	CB-CA-C	5.77	121.94	110.40
1	A	249	TYR	CB-CG-CD1	-5.71	117.58	121.00
1	A	126	TRP	CB-CG-CD1	5.63	134.33	127.00
1	A	182	GLU	CG-CD-OE1	5.61	129.52	118.30
1	A	179	ARG	CD-NE-CZ	5.53	131.35	123.60
1	A	262	ASP	CB-CG-OD1	5.53	123.27	118.30
1	A	265	ARG	CA-CB-CG	5.43	125.35	113.40
1	A	204	GLU	CG-CD-OE1	5.40	129.11	118.30
1	A	103	GLU	CA-C-O	-5.40	108.77	120.10
1	A	110	ASP	CB-CG-OD2	-5.38	113.46	118.30
1	A	250	GLU	CA-CB-CG	5.36	125.18	113.40
1	A	317	ARG	CG-CD-NE	-5.35	100.56	111.80
1	A	222	LEU	CB-CA-C	5.33	120.33	110.20
1	A	237	GLY	N-CA-C	-5.32	99.80	113.10
1	A	176	ASP	CB-CG-OD1	-5.31	113.52	118.30
1	A	215	ALA	CB-CA-C	5.30	118.05	110.10

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	254	PHE	CB-CG-CD1	-5.28	117.11	120.80
1	A	194	ASP	O-C-N	5.24	131.08	122.70
1	A	126	TRP	O-C-N	5.22	131.06	122.70
1	A	313	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	28	GLU	CG-CD-OE2	-5.16	107.97	118.30
1	A	40	THR	CA-CB-OG1	-5.16	98.16	109.00
1	A	120	ILE	N-CA-CB	5.16	122.66	110.80
1	A	192	GLY	C-N-CA	5.14	134.54	121.70
1	A	152	GLU	CG-CD-OE2	-5.13	108.04	118.30
1	A	74	GLY	O-C-N	-5.09	114.56	122.70
1	A	157	ARG	CB-CA-C	-5.03	100.34	110.40

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	2457	0	2385	53	0
2	A	110	0	0	4	0
All	All	2567	0	2385	53	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 11.

All (53) 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:45:HIS:H	1:A:48:HIS:HD2	1.17	0.88
1:A:234:THR:HG22	1:A:236:SER:H	1.47	0.79
1:A:45:HIS:H	1:A:48:HIS:CD2	2.06	0.70
1:A:283:GLN:HG3	1:A:286:ARG:NH1	2.12	0.64
1:A:222:LEU:HD23	2:A:493:HOH:O	1.97	0.64
1:A:72:ALA:O	1:A:75:LEU:HB2	1.98	0.63
1:A:292:ARG:NH1	1:A:295:GLN:HG2	2.13	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:VAL:O	1:A:149:MET:HG2	2.00	0.61
1:A:54:THR:HG21	1:A:220:ILE:HD11	1.81	0.61
1:A:31:VAL:O	1:A:63:HIS:HB3	2.00	0.60
1:A:48:HIS:O	1:A:52:ILE:HG13	2.00	0.60
1:A:100:ARG:NH2	1:A:242:ASP:OD1	2.32	0.60
1:A:253:GLN:HA	1:A:256:ILE:HG22	1.83	0.60
1:A:110:ASP:OD2	1:A:113:ALA:HB2	2.01	0.60
1:A:24:LYS:NZ	1:A:28:GLU:OE2	2.35	0.59
1:A:64:ARG:HD2	2:A:485:HOH:O	2.04	0.57
1:A:292:ARG:HH11	1:A:295:GLN:HG2	1.71	0.56
1:A:314:GLN:O	1:A:318:ILE:HG13	2.06	0.55
1:A:152:GLU:O	1:A:156:SER:HB2	2.07	0.55
1:A:157:ARG:O	1:A:161:GLY:N	2.39	0.54
1:A:33:LEU:HA	1:A:188:LEU:O	2.08	0.53
1:A:9:TRP:CZ2	1:A:275:LYS:HG3	2.43	0.53
1:A:239:ILE:HG23	1:A:246:THR:HG21	1.91	0.53
1:A:297:THR:O	1:A:301:GLU:HG2	2.09	0.53
1:A:178:LEU:HD11	1:A:210:LYS:HE2	1.90	0.53
1:A:256:ILE:O	1:A:292:ARG:NH1	2.42	0.53
1:A:234:THR:HG22	1:A:236:SER:N	2.19	0.53
1:A:75:LEU:HD21	1:A:91:LYS:HA	1.92	0.52
1:A:91:LYS:O	1:A:95:GLU:HB2	2.11	0.50
1:A:281:LEU:HD22	1:A:294:ALA:HA	1.94	0.50
1:A:73:THR:HB	1:A:169:TYR:CE1	2.48	0.49
1:A:59:GLN:HB2	1:A:65:PRO:HG3	1.95	0.49
1:A:274:SER:N	1:A:277:GLU:OE2	2.23	0.49
1:A:53:LEU:O	1:A:56:ARG:HB3	2.15	0.47
1:A:56:ARG:O	1:A:60:GLN:HG3	2.15	0.47
1:A:1:MET:N	1:A:27:ASN:HD21	2.12	0.47
1:A:154:VAL:O	1:A:155:GLN:C	2.53	0.47
1:A:53:LEU:O	1:A:57:ARG:HG3	2.15	0.46
1:A:78:ASP:HB2	1:A:169:TYR:CZ	2.51	0.46
1:A:17:THR:HG21	1:A:203:LEU:CD1	2.46	0.45
1:A:102:LYS:HG3	1:A:120:ILE:HG21	1.98	0.45
1:A:44:LEU:HD13	1:A:52:ILE:HD11	1.98	0.45
1:A:170:MET:HG3	1:A:171:MET:HE2	1.99	0.45
1:A:190:ILE:HA	1:A:218:LEU:O	2.17	0.44
1:A:187:ARG:HG3	2:A:420:HOH:O	2.18	0.44
1:A:72:ALA:HB2	1:A:124:TYR:HA	2.01	0.43
1:A:298:LEU:O	1:A:302:VAL:HG23	2.19	0.43
1:A:38:ASP:OD2	1:A:70:GLY:HA3	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:TRP:CH2	1:A:275:LYS:HA	2.55	0.42
1:A:304:LYS:HG3	1:A:312:LEU:HD22	2.02	0.42
1:A:250:GLU:HG2	2:A:525:HOH:O	2.20	0.41
1:A:113:ALA:O	1:A:117:PRO:HB3	2.22	0.40
1:A:128:GLY:N	1:A:129:PRO:HD2	2.37	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	313/419 (75%)	304 (97%)	8 (3%)	1 (0%)	41 50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	SER

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	244/349 (70%)	234 (96%)	10 (4%)	30 43

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	GLN
1	A	33	LEU
1	A	114	ASP
1	A	146	ASN
1	A	177	PHE
1	A	179	ARG
1	A	222	LEU
1	A	256	ILE
1	A	261	ARG
1	A	304	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	48	HIS
1	A	60	GLN
1	A	89	ASN
1	A	155	GLN
1	A	283	GLN
1	A	314	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 monosaccharides 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	317/419 (75%)	-0.38	5 (1%) 72 77	2, 13, 33, 48	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	236	SER	3.4
1	A	317	ARG	2.6
1	A	309	GLU	2.6
1	A	159	GLU	2.3
1	A	235	GLU	2.2

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 monosaccharides 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.