



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2019 – 12:10 PM EST

PDB ID : 6E4P
Title : Structure of the T. brucei RRM domain in complex with RNA
Authors : Schumacher, M.A.
Deposited on : 2018-07-18
Resolution : 1.95 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

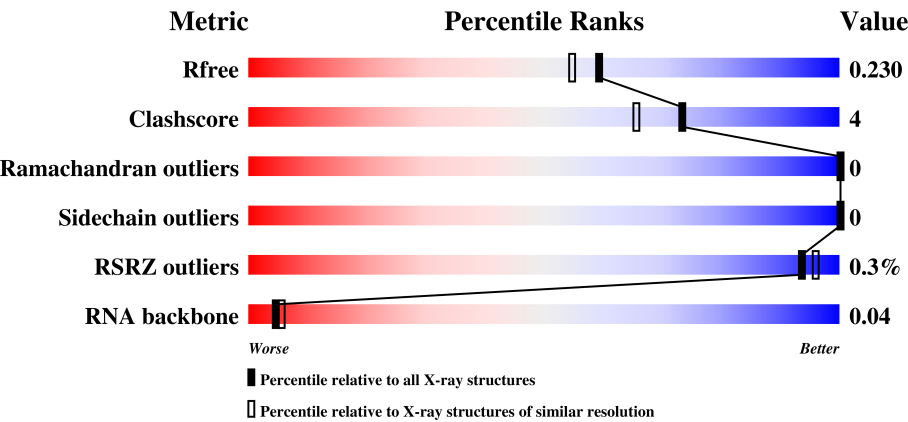
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	rb-20031633
Percentile statistics	:	20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac	:	5.8.0158
CCP4	:	7.0 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20031633

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	111664	3622 (1.96-1.92)
Clashscore	122126	3795 (1.96-1.92)
Ramachandran outliers	120053	3757 (1.96-1.92)
Sidechain outliers	120020	3757 (1.96-1.92)
RSRZ outliers	108989	3554 (1.96-1.92)
RNA backbone	2636	1093 (2.60-1.30)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	J	4	<div><div></div><div>50%25%25%</div></div>
1	K	4	<div><div></div><div>50%25%25%</div></div>
2	A	71	<div><div></div><div>96%..</div></div>
2	B	71	<div><div></div><div>85%13%. </div></div>

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Mol	Chain	Length	Quality of chain
2	C	71	<div><div></div><div>90%</div><div>6%</div><div>.</div></div>
2	D	71	<div>%<div><div></div><div>93%</div><div>7%</div></div></div>
2	E	71	<div><div></div><div>86%</div><div>13%</div><div>.</div></div>
2	F	71	<div>%<div><div></div><div>93%</div><div>7%</div></div></div>
2	G	71	<div><div></div><div>94%</div><div>.</div><div>.</div></div>
2	H	71	<div><div></div><div>97%</div><div>.</div></div>
2	I	71	<div><div></div><div>92%</div><div>8%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5319 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 RNA chain called RNA (5'-R(P*UP*UP*UP*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	J	3	Total	C	N	O	P	0	0	0
			61	27	6	25	3			
1	K	3	Total	C	N	O	P	0	0	0
			61	27	6	25	3			

- Molecule 2 is a protein called RNA-binding protein, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	70	Total	C	N	O	S	0	0	0
			543	339	97	105	2			
2	B	69	Total	C	N	O	S	0	0	0
			538	336	96	104	2			
2	C	68	Total	C	N	O	S	0	0	0
			528	330	93	103	2			
2	D	71	Total	C	N	O	S	0	0	0
			548	341	98	107	2			
2	E	70	Total	C	N	O	S	0	0	0
			538	333	97	106	2			
2	F	71	Total	C	N	O	S	0	0	0
			548	341	98	107	2			
2	G	70	Total	C	N	O	S	0	0	0
			544	339	97	106	2			
2	H	71	Total	C	N	O	S	0	0	0
			548	341	98	107	2			
2	I	71	Total	C	N	O	S	0	0	0
			548	341	98	107	2			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	199	GLY	-	expression tag	UNP Q389P7
A	200	SER	-	expression tag	UNP Q389P7
A	201	HIS	-	expression tag	UNP Q389P7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	202	MET	-	expression tag	UNP Q389P7
B	199	GLY	-	expression tag	UNP Q389P7
B	200	SER	-	expression tag	UNP Q389P7
B	201	HIS	-	expression tag	UNP Q389P7
B	202	MET	-	expression tag	UNP Q389P7
C	199	GLY	-	expression tag	UNP Q389P7
C	200	SER	-	expression tag	UNP Q389P7
C	201	HIS	-	expression tag	UNP Q389P7
C	202	MET	-	expression tag	UNP Q389P7
D	199	GLY	-	expression tag	UNP Q389P7
D	200	SER	-	expression tag	UNP Q389P7
D	201	HIS	-	expression tag	UNP Q389P7
D	202	MET	-	expression tag	UNP Q389P7
E	199	GLY	-	expression tag	UNP Q389P7
E	200	SER	-	expression tag	UNP Q389P7
E	201	HIS	-	expression tag	UNP Q389P7
E	202	MET	-	expression tag	UNP Q389P7
F	199	GLY	-	expression tag	UNP Q389P7
F	200	SER	-	expression tag	UNP Q389P7
F	201	HIS	-	expression tag	UNP Q389P7
F	202	MET	-	expression tag	UNP Q389P7
G	199	GLY	-	expression tag	UNP Q389P7
G	200	SER	-	expression tag	UNP Q389P7
G	201	HIS	-	expression tag	UNP Q389P7
G	202	MET	-	expression tag	UNP Q389P7
H	199	GLY	-	expression tag	UNP Q389P7
H	200	SER	-	expression tag	UNP Q389P7
H	201	HIS	-	expression tag	UNP Q389P7
H	202	MET	-	expression tag	UNP Q389P7
I	199	GLY	-	expression tag	UNP Q389P7
I	200	SER	-	expression tag	UNP Q389P7
I	201	HIS	-	expression tag	UNP Q389P7
I	202	MET	-	expression tag	UNP Q389P7

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	J	8	Total O 8 8	0	0
3	K	3	Total O 3 3	0	0
3	A	46	Total O 46 46	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	27	Total 27	O 27	0	0
3	C	30	Total 30	O 30	0	0
3	D	21	Total 21	O 21	0	0
3	E	27	Total 27	O 27	0	0
3	F	46	Total 46	O 46	0	0
3	G	33	Total 33	O 33	0	0
3	H	41	Total 41	O 41	0	0
3	I	32	Total 32	O 32	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

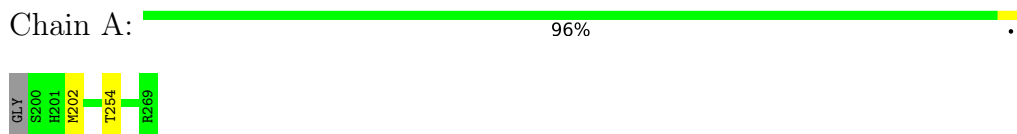
- Molecule 1: RNA (5'-R(P*UP*UP*UP*U)-3')



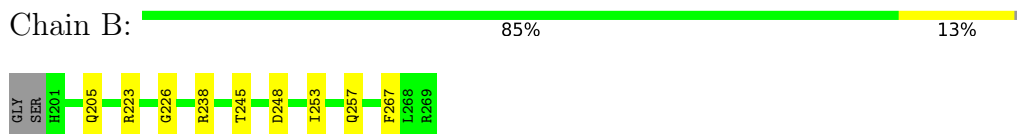
- Molecule 1: RNA (5'-R(P*UP*UP*UP*U)-3')



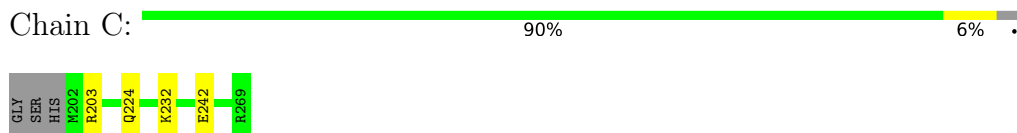
- Molecule 2: RNA-binding protein, putative



- Molecule 2: RNA-binding protein, putative



- Molecule 2: RNA-binding protein, putative



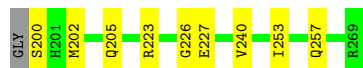
- Molecule 2: RNA-binding protein, putative





- Molecule 2: RNA-binding protein, putative

Chain E: 86% 13%



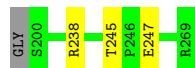
- Molecule 2: RNA-binding protein, putative

Chain F: 93% 7%



- Molecule 2: RNA-binding protein, putative

Chain G: 94%



- Molecule 2: RNA-binding protein, putative

Chain H: 97%



- Molecule 2: RNA-binding protein, putative

Chain I: 92% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.12Å 126.30Å 49.20Å 90.00° 118.57° 90.00°	Depositor
Resolution (Å)	35.66 – 1.95 35.66 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.66-1.95) 99.9 (35.66-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.38 (at 1.95Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.181 , 0.229 0.182 , 0.230	Depositor DCC
R_{free} test set	1995 reflections (5.21%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -h-l,k,h 0.000 for l,k,-h-l 0.023 for h,-k,-h-l 0.000 for -h-l,-k,l 0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5319	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 50.17 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.7250e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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	J	1.29	1/66 (1.5%)	0.83	0/98
1	K	1.31	1/66 (1.5%)	0.97	0/98
2	A	0.33	0/552	0.47	0/746
2	B	0.26	0/547	0.43	0/738
2	C	0.28	0/536	0.49	0/723
2	D	0.26	0/557	0.43	0/751
2	E	0.29	0/546	0.49	0/737
2	F	0.30	0/557	0.43	0/751
2	G	0.29	0/553	0.43	0/746
2	H	0.30	0/557	0.45	0/751
2	I	0.28	0/557	0.46	0/751
All	All	0.35	2/5094 (0.0%)	0.47	0/6890

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	K	1	U	OP3-P	-10.23	1.48	1.61
1	J	1	U	OP3-P	-10.06	1.49	1.61

There are no bond angle outliers.

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	J	61	0	31	0	0
1	K	61	0	31	1	0
2	A	543	0	531	2	0
2	B	538	0	526	9	0
2	C	528	0	519	6	0
2	D	548	0	534	3	0
2	E	538	0	524	7	0
2	F	548	0	534	4	0
2	G	544	0	531	2	0
2	H	548	0	534	1	0
2	I	548	0	534	4	0
3	A	46	0	0	0	0
3	B	27	0	0	0	0
3	C	30	0	0	0	0
3	D	21	0	0	0	0
3	E	27	0	0	1	0
3	F	46	0	0	0	0
3	G	33	0	0	1	0
3	H	41	0	0	0	0
3	I	32	0	0	0	0
3	J	8	0	0	0	0
3	K	3	0	0	0	0
All	All	5319	0	4829	35	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 (35) 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)
2:B:205:GLN:NE2	2:B:238:ARG:HH12	1.18	1.38
2:B:205:GLN:HE22	2:B:238:ARG:NH1	1.17	1.34
2:B:205:GLN:NE2	2:B:238:ARG:NH1	1.74	1.23
2:G:238:ARG:NH2	3:G:301:HOH:O	1.84	1.06
2:C:203:ARG:HH11	2:C:203:ARG:HG3	1.31	0.95
2:C:203:ARG:NH1	2:C:203:ARG:HG3	2.02	0.74
1:K:1:U:O2	1:K:1:U:H2'	1.89	0.73
2:B:205:GLN:NE2	2:B:238:ARG:HH11	1.87	0.72
2:I:244:VAL:HG13	2:I:245:THR:HG23	1.80	0.64
2:B:205:GLN:HE21	2:B:238:ARG:NH1	1.90	0.62
2:C:203:ARG:NH1	2:C:242:GLU:OE1	2.31	0.59
2:C:224:GLN:HA	2:I:232:LYS:HD3	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:203:ARG:NH1	2:C:203:ARG:CG	2.69	0.50
2:E:223:ARG:NH1	2:E:227:GLU:OE1	2.45	0.50
2:F:200:SER:HB3	2:F:246:PRO:HD3	1.94	0.49
2:D:206:VAL:HB	2:D:239:ALA:HB3	1.96	0.48
2:B:245:THR:OG1	2:B:248:ASP:OD2	2.28	0.48
2:F:200:SER:O	2:F:246:PRO:HD3	2.14	0.47
2:I:203:ARG:HD3	2:I:242:GLU:HB2	1.98	0.46
2:E:200:SER:C	2:E:202:MET:H	2.17	0.46
2:C:232:LYS:NZ	2:D:227:GLU:OE2	2.42	0.45
2:E:223:ARG:HA	2:E:226:GLY:O	2.16	0.45
2:F:219:LYS:HE3	2:F:223:ARG:HD2	1.98	0.45
2:B:223:ARG:HA	2:B:226:GLY:O	2.18	0.44
2:B:253:ILE:O	2:B:257:GLN:HB2	2.18	0.44
2:H:202:MET:HG3	2:H:246:PRO:HA	1.99	0.44
2:A:254:THR:HB	2:F:234:PHE:CZ	2.54	0.43
2:D:244:VAL:HG12	2:D:245:THR:HG23	2.01	0.43
2:E:200:SER:C	2:E:202:MET:N	2.71	0.42
2:A:202:MET:HG2	2:B:267:PHE:CD2	2.55	0.42
2:E:253:ILE:O	2:E:257:GLN:HB2	2.19	0.41
2:G:245:THR:HB	2:G:247:GLU:OE2	2.20	0.41
2:E:205:GLN:OE1	2:E:240:VAL:HG22	2.21	0.41
2:I:202:MET:HB2	2:I:242:GLU:HG3	2.02	0.41
2:E:223:ARG:HD2	3:E:304:HOH:O	2.20	0.41

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	
2	A	68/71 (96%)	68 (100%)	0	0	100	100
2	B	67/71 (94%)	67 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	66/71 (93%)	66 (100%)	0	0	100	100
2	D	69/71 (97%)	69 (100%)	0	0	100	100
2	E	68/71 (96%)	68 (100%)	0	0	100	100
2	F	69/71 (97%)	68 (99%)	1 (1%)	0	100	100
2	G	68/71 (96%)	68 (100%)	0	0	100	100
2	H	69/71 (97%)	68 (99%)	1 (1%)	0	100	100
2	I	69/71 (97%)	69 (100%)	0	0	100	100
All	All	613/639 (96%)	611 (100%)	2 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	
2	A	58/58 (100%)	58 (100%)	0	100	100
2	B	57/58 (98%)	57 (100%)	0	100	100
2	C	56/58 (97%)	56 (100%)	0	100	100
2	D	58/58 (100%)	58 (100%)	0	100	100
2	E	57/58 (98%)	57 (100%)	0	100	100
2	F	58/58 (100%)	58 (100%)	0	100	100
2	G	58/58 (100%)	58 (100%)	0	100	100
2	H	58/58 (100%)	58 (100%)	0	100	100
2	I	58/58 (100%)	58 (100%)	0	100	100
All	All	518/522 (99%)	518 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	205	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	J	2/4 (50%)	0	0
1	K	2/4 (50%)	0	0
All	All	4/8 (50%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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 carbohydrates 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	J	3/4 (75%)	-0.06	0 100 100	30, 30, 36, 46	0
1	K	3/4 (75%)	0.73	0 100 100	50, 50, 52, 63	0
2	A	70/71 (98%)	-0.19	0 100 100	15, 22, 36, 54	0
2	B	69/71 (97%)	-0.08	0 100 100	20, 28, 43, 50	0
2	C	68/71 (95%)	-0.13	0 100 100	15, 27, 43, 57	0
2	D	71/71 (100%)	0.05	1 (1%) 75 81	19, 31, 46, 60	0
2	E	70/71 (98%)	0.03	0 100 100	20, 30, 47, 52	0
2	F	71/71 (100%)	-0.10	1 (1%) 75 81	16, 23, 39, 47	0
2	G	70/71 (98%)	-0.07	0 100 100	18, 29, 50, 58	0
2	H	71/71 (100%)	-0.14	0 100 100	17, 23, 38, 53	1 (1%)
2	I	71/71 (100%)	-0.31	0 100 100	18, 27, 41, 45	0
All	All	637/647 (98%)	-0.10	2 (0%) 93 96	15, 27, 46, 63	1 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	263	GLY	3.0
2	F	200	SER	2.4

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 carbohydrates 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.