



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:24 AM UTC

PDB ID : 5EFT / pdb_00005eft
Title : Structural Basis for Specific Recognition of ssDNA by SRBSDV P9-1 Octamers
Authors : Li, X.Y.
Deposited on : 2015-10-26
Resolution : 2.50 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

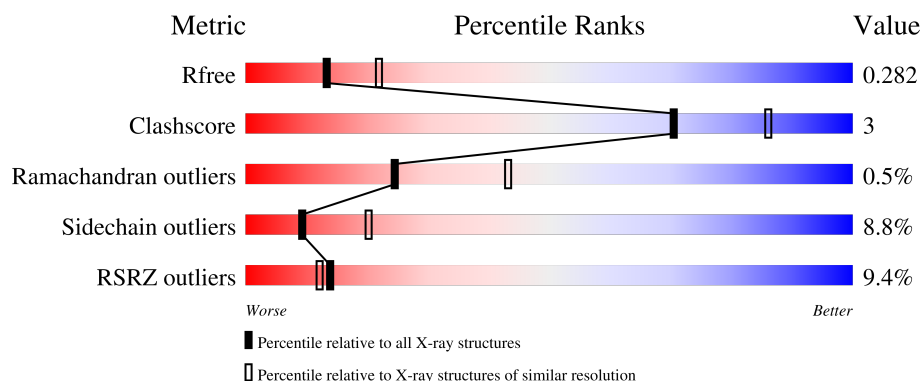
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	16	
1	C	16	
1	E	16	
1	G	16	
2	B	321	

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Mol	Chain	Length	Quality of chain
2	D	321	9% 68% 11% • 21%
2	F	321	12% 67% 11% • 21%
2	H	321	7% 66% 12% • 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8792 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 p9-1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	16	Total	C	N	O	0	0	0
			121	76	18	27			
1	C	16	Total	C	N	O	0	0	0
			121	76	18	27			
1	E	16	Total	C	N	O	0	0	0
			121	76	18	27			
1	G	16	Total	C	N	O	0	0	0
			121	76	18	27			

- Molecule 2 is a protein called p9-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	255	Total	C	N	O	S	0	0	0
			2077	1351	339	381	6			
2	D	255	Total	C	N	O	S	0	0	0
			2077	1351	339	381	6			
2	F	255	Total	C	N	O	S	0	0	0
			2077	1351	339	381	6			
2	H	255	Total	C	N	O	S	0	0	0
			2077	1351	339	381	6			

There are 4 discrepancies between the modelled and reference sequences:

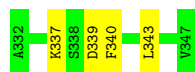
Chain	Residue	Modelled	Actual	Comment	Reference
B	162	THR	LYS	conflict	UNP B6SCH3
D	162	THR	LYS	conflict	UNP B6SCH3
F	162	THR	LYS	conflict	UNP B6SCH3
H	162	THR	LYS	conflict	UNP B6SCH3

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: p9-1

Chain A:  75% 25%



- Molecule 1: p9-1

Chain C:  6% 81% 19%



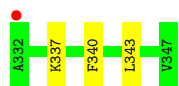
- Molecule 1: p9-1

Chain E:  6% 88% 12%



- Molecule 1: p9-1

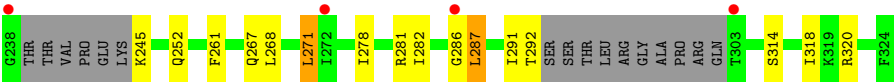
Chain G:  6% 81% 19%



- Molecule 2: p9-1

Chain B:  3% 67% 12% 21%





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	118.52Å 155.72Å 157.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.60 – 2.50 29.60 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (29.60-2.50) 99.5 (29.60-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.253 , 0.284 0.252 , 0.282	Depositor DCC
R_{free} test set	2566 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 50.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8792	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.61% 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	0.94	0/121	0.83	0/160
1	C	0.74	0/121	0.77	0/160
1	E	0.61	0/121	0.71	0/160
1	G	0.71	0/121	0.73	0/160
2	B	0.84	0/2115	0.87	1/2855 (0.0%)
2	D	0.72	0/2115	0.86	1/2855 (0.0%)
2	F	0.69	0/2115	0.83	1/2855 (0.0%)
2	H	0.63	0/2115	0.83	2/2855 (0.1%)
All	All	0.73	0/8944	0.84	5/12060 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	210	ASN	N-CA-C	5.70	118.97	108.94
2	H	261	PHE	N-CA-C	5.32	117.08	111.28
2	F	210	ASN	N-CA-C	5.16	118.02	108.94
2	H	210	ASN	N-CA-C	5.12	117.98	109.79
2	B	261	PHE	N-CA-C	5.03	117.42	111.33

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	121	0	122	5	0
1	C	121	0	122	2	0
1	E	121	0	122	1	0
1	G	121	0	122	2	0
2	B	2077	0	2075	19	0
2	D	2077	0	2075	18	0
2	F	2077	0	2075	13	2
2	H	2077	0	2075	14	2
All	All	8792	0	8788	60	2

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 3.

All (60) 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:269:SER:CB	2:D:74:ASN:HD22	1.97	0.76
2:B:269:SER:OG	2:D:74:ASN:ND2	2.18	0.75
2:B:128:VAL:HG12	2:B:162:THR:HG22	1.87	0.57
2:H:128:VAL:HG12	2:H:162:THR:HG22	1.88	0.56
2:D:128:VAL:HG12	2:D:162:THR:HG22	1.88	0.55
2:F:128:VAL:HG12	2:F:162:THR:HG22	1.88	0.54
1:A:339:ASP:OD2	2:D:6:ARG:NH1	2.40	0.53
2:D:117:LYS:HG2	2:D:318:ILE:HG12	1.90	0.52
2:B:117:LYS:HG2	2:B:318:ILE:HG12	1.91	0.52
2:H:117:LYS:HG2	2:H:318:ILE:HG12	1.91	0.52
2:B:269:SER:CB	2:D:74:ASN:ND2	2.68	0.52
2:F:117:LYS:HG2	2:F:318:ILE:HG12	1.92	0.51
2:H:286:GLY:O	2:H:287:LEU:HB2	2.12	0.50
2:F:56:LEU:HD21	2:F:116:ILE:HD11	1.94	0.49
2:B:286:GLY:O	2:B:287:LEU:HB2	2.12	0.49
2:H:268:LEU:HD12	2:H:271:LEU:HD22	1.95	0.48
2:H:56:LEU:HD21	2:H:116:ILE:HD11	1.95	0.48
2:B:268:LEU:HD12	2:B:271:LEU:HD22	1.96	0.48
2:B:269:SER:HB2	2:D:74:ASN:ND2	2.29	0.47
1:A:337:LYS:HE3	2:B:209:GLU:O	2.15	0.47
2:B:269:SER:HG	2:D:74:ASN:ND2	2.09	0.47
2:D:268:LEU:HD12	2:D:271:LEU:HD22	1.97	0.47
2:D:56:LEU:HD21	2:D:116:ILE:HD11	1.96	0.47
2:F:268:LEU:HD12	2:F:271:LEU:HD22	1.96	0.47
1:A:339:ASP:CG	2:D:6:ARG:HH12	2.22	0.47
1:A:339:ASP:HB2	2:D:6:ARG:HH22	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:286:GLY:O	2:D:287:LEU:HB2	2.13	0.47
2:F:286:GLY:O	2:F:287:LEU:HB2	2.14	0.46
2:H:222:LYS:NZ	2:H:320:ARG:O	2.49	0.46
2:F:117:LYS:HB2	2:F:122:THR:HB	1.98	0.46
1:G:337:LYS:HE3	2:H:209:GLU:O	2.16	0.45
2:D:222:LYS:NZ	2:D:320:ARG:O	2.49	0.45
2:H:117:LYS:HB2	2:H:122:THR:HB	1.98	0.45
2:B:56:LEU:HD21	2:B:116:ILE:HD11	1.98	0.44
2:B:117:LYS:HB2	2:B:122:THR:HB	1.99	0.44
2:B:222:LYS:NZ	2:B:320:ARG:O	2.50	0.44
2:D:117:LYS:HB2	2:D:122:THR:HB	1.98	0.44
2:F:278:ILE:O	2:F:282:ILE:HG12	2.18	0.44
2:D:278:ILE:O	2:D:282:ILE:HG12	2.18	0.44
2:F:222:LYS:NZ	2:F:320:ARG:O	2.51	0.44
2:H:268:LEU:HA	2:H:271:LEU:HD13	2.00	0.43
1:A:340:PHE:HB3	1:A:343:LEU:HD12	2.00	0.43
2:B:278:ILE:O	2:B:282:ILE:HG12	2.18	0.43
1:C:334:SER:HB3	2:H:281:ARG:NH2	2.33	0.43
2:B:268:LEU:HA	2:B:271:LEU:HD13	2.00	0.43
2:B:269:SER:HB2	2:D:74:ASN:HD22	1.81	0.43
2:B:46:PRO:HB3	2:B:214:TYR:CG	2.54	0.43
1:E:340:PHE:HB3	1:E:343:LEU:HD12	2.01	0.42
2:H:278:ILE:O	2:H:282:ILE:HG12	2.19	0.42
2:F:53:PHE:CD1	2:F:53:PHE:C	2.98	0.42
2:H:46:PRO:HB3	2:H:214:TYR:CG	2.54	0.42
1:C:340:PHE:HB3	1:C:343:LEU:HD12	2.01	0.42
2:F:268:LEU:HA	2:F:271:LEU:HD13	2.01	0.42
1:G:340:PHE:HB3	1:G:343:LEU:HD12	2.01	0.41
2:B:64:ASN:ND2	2:F:74:ASN:H	2.18	0.41
2:B:64:ASN:HD21	2:F:73:PHE:HA	1.86	0.41
2:H:53:PHE:CD1	2:H:53:PHE:C	2.98	0.41
2:D:268:LEU:HA	2:D:271:LEU:HD13	2.02	0.41
2:F:110:ARG:HD2	2:F:125:LEU:HG	2.03	0.40
2:H:110:ARG:HD2	2:H:125:LEU:HG	2.03	0.40

All (2) 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:F:292:THR:OG1	2:H:183:GLU:OE2[8_475]	2.01	0.19
2:F:292:THR:OG1	2:H:183:GLU:CD[8_475]	2.18	0.02

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	14/16 (88%)	14 (100%)	0	0	100	100
1	C	14/16 (88%)	14 (100%)	0	0	100	100
1	E	14/16 (88%)	14 (100%)	0	0	100	100
1	G	14/16 (88%)	14 (100%)	0	0	100	100
2	B	243/321 (76%)	231 (95%)	11 (4%)	1 (0%)	30	49
2	D	243/321 (76%)	232 (96%)	9 (4%)	2 (1%)	16	31
2	F	243/321 (76%)	231 (95%)	11 (4%)	1 (0%)	30	49
2	H	243/321 (76%)	232 (96%)	10 (4%)	1 (0%)	30	49
All	All	1028/1348 (76%)	982 (96%)	41 (4%)	5 (0%)	24	43

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	287	LEU
2	D	287	LEU
2	F	287	LEU
2	H	287	LEU
2	D	69	CYS

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	15/15 (100%)	15 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	15/15 (100%)	15 (100%)	0	100	100
1	E	15/15 (100%)	15 (100%)	0	100	100
1	G	15/15 (100%)	15 (100%)	0	100	100
2	B	230/302 (76%)	208 (90%)	22 (10%)	8	17
2	D	230/302 (76%)	209 (91%)	21 (9%)	9	19
2	F	230/302 (76%)	208 (90%)	22 (10%)	8	17
2	H	230/302 (76%)	209 (91%)	21 (9%)	9	19
All	All	980/1268 (77%)	894 (91%)	86 (9%)	9	20

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	15	GLU
2	B	18	THR
2	B	43	LYS
2	B	56	LEU
2	B	71	CYS
2	B	95	ASN
2	B	125	LEU
2	B	167	MET
2	B	172	LEU
2	B	173	LEU
2	B	181	LEU
2	B	186	LEU
2	B	196	LEU
2	B	237	VAL
2	B	245	LYS
2	B	252	GLN
2	B	267	GLN
2	B	271	LEU
2	B	272	ILE
2	B	291	ILE
2	B	292	THR
2	B	314	SER
2	D	15	GLU
2	D	18	THR
2	D	43	LYS
2	D	56	LEU
2	D	71	CYS
2	D	95	ASN

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Mol	Chain	Res	Type
2	D	125	LEU
2	D	167	MET
2	D	172	LEU
2	D	173	LEU
2	D	181	LEU
2	D	186	LEU
2	D	196	LEU
2	D	237	VAL
2	D	245	LYS
2	D	252	GLN
2	D	267	GLN
2	D	271	LEU
2	D	291	ILE
2	D	292	THR
2	D	314	SER
2	F	15	GLU
2	F	18	THR
2	F	43	LYS
2	F	56	LEU
2	F	71	CYS
2	F	95	ASN
2	F	125	LEU
2	F	167	MET
2	F	172	LEU
2	F	173	LEU
2	F	181	LEU
2	F	186	LEU
2	F	196	LEU
2	F	237	VAL
2	F	245	LYS
2	F	252	GLN
2	F	267	GLN
2	F	271	LEU
2	F	272	ILE
2	F	291	ILE
2	F	292	THR
2	F	314	SER
2	H	15	GLU
2	H	18	THR
2	H	43	LYS
2	H	56	LEU
2	H	71	CYS

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Mol	Chain	Res	Type
2	H	95	ASN
2	H	125	LEU
2	H	167	MET
2	H	172	LEU
2	H	173	LEU
2	H	181	LEU
2	H	186	LEU
2	H	196	LEU
2	H	237	VAL
2	H	245	LYS
2	H	252	GLN
2	H	267	GLN
2	H	271	LEU
2	H	291	ILE
2	H	292	THR
2	H	314	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	64	ASN
2	B	80	ASN
2	B	107	ASN
2	B	194	ASN
2	B	217	GLN
2	B	248	HIS
2	B	252	GLN
2	B	317	ASN
2	D	64	ASN
2	D	80	ASN
2	D	107	ASN
2	D	194	ASN
2	D	217	GLN
2	D	248	HIS
2	D	252	GLN
2	D	267	GLN
2	D	317	ASN
2	F	64	ASN
2	F	80	ASN
2	F	107	ASN
2	F	194	ASN
2	F	217	GLN

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Mol	Chain	Res	Type
2	F	252	GLN
2	F	317	ASN
2	H	64	ASN
2	H	80	ASN
2	H	107	ASN
2	H	194	ASN
2	H	217	GLN
2	H	252	GLN
2	H	317	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	16/16 (100%)	-0.15	0 100 100	46, 57, 76, 95	0
1	C	16/16 (100%)	0.37	1 (6%) 26 23	51, 71, 93, 99	0
1	E	16/16 (100%)	0.91	1 (6%) 26 23	73, 88, 103, 113	0
1	G	16/16 (100%)	0.67	1 (6%) 26 23	73, 88, 108, 121	0
2	B	255/321 (79%)	0.16	10 (3%) 43 38	30, 59, 116, 148	0
2	D	255/321 (79%)	0.73	28 (10%) 10 8	45, 85, 152, 179	0
2	F	255/321 (79%)	0.92	40 (15%) 5 4	40, 90, 177, 202	0
2	H	255/321 (79%)	0.73	21 (8%) 17 15	47, 99, 155, 231	0
All	All	1084/1348 (80%)	0.62	102 (9%) 14 12	30, 84, 155, 231	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	17	LEU	7.7
2	H	17	LEU	6.5
2	D	4	LEU	5.3
2	D	271	LEU	5.0
2	F	72	GLY	4.9
2	B	4	LEU	4.6
2	F	6	ARG	4.5
2	F	4	LEU	4.4
2	F	65	PHE	4.3
2	F	290	ILE	4.0
2	D	130	LYS	3.9
2	H	18	THR	3.8
2	F	73	PHE	3.7
2	D	272	ILE	3.7
2	F	103	LEU	3.6
2	F	17	LEU	3.6

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Mol	Chain	Res	Type	RSRZ
2	H	19	ILE	3.6
2	D	238	GLY	3.5
2	D	161	GLU	3.4
2	F	105	LEU	3.3
2	F	9	PHE	3.3
2	D	273	SER	3.2
2	B	133	VAL	3.2
2	F	133	VAL	3.2
1	G	332	ALA	3.2
2	H	163	LYS	3.2
2	D	131	THR	3.1
2	B	162	THR	3.0
2	H	105	LEU	3.0
2	F	107	ASN	3.0
2	D	106	ALA	2.9
2	H	161	GLU	2.9
2	H	133	VAL	2.9
2	D	160	GLN	2.9
2	D	128	VAL	2.8
2	B	167	MET	2.8
2	F	19	ILE	2.8
2	F	165	VAL	2.8
2	H	303	THR	2.8
2	D	108	TYR	2.7
2	H	77	HIS	2.7
2	F	276	THR	2.7
2	H	286	GLY	2.7
1	E	332	ALA	2.7
2	F	324	PHE	2.6
2	F	175	ARG	2.6
2	D	133	VAL	2.6
2	H	160	GLN	2.6
2	D	209	GLU	2.5
1	C	332	ALA	2.5
2	F	58	TYR	2.5
2	D	45	ILE	2.5
2	H	272	ILE	2.5
2	F	245	LYS	2.5
2	H	238	GLY	2.5
2	H	78	LEU	2.5
2	H	165	VAL	2.5
2	D	107	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
2	F	10	GLY	2.4
2	D	19	ILE	2.4
2	D	129	LEU	2.4
2	H	70	TYR	2.4
2	D	18	THR	2.3
2	F	180	SER	2.3
2	F	81	LEU	2.3
2	F	173	LEU	2.3
2	F	291	ILE	2.3
2	F	70	TYR	2.3
2	F	176	GLU	2.3
2	D	324	PHE	2.3
2	D	286	GLY	2.3
2	H	181	LEU	2.3
2	F	164	PHE	2.3
2	B	132	VAL	2.3
2	F	238	GLY	2.3
2	H	4	LEU	2.3
2	F	5	GLU	2.3
2	B	128	VAL	2.3
2	F	68	THR	2.2
2	H	125	LEU	2.2
2	F	14	ILE	2.2
2	D	132	VAL	2.2
2	B	160	GLN	2.2
2	F	12	TYR	2.2
2	F	85	PHE	2.2
2	D	103	LEU	2.2
2	D	105	LEU	2.2
2	F	75	TYR	2.2
2	H	128	VAL	2.1
2	F	318	ILE	2.1
2	F	179	GLU	2.1
2	D	163	LYS	2.1
2	F	7	ARG	2.1
2	D	169	LEU	2.1
2	H	103	LEU	2.1
2	F	303	THR	2.1
2	B	268	LEU	2.0
2	D	303	THR	2.0
2	F	45	ILE	2.0
2	F	131	THR	2.0

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Mol	Chain	Res	Type	RSRZ
2	B	238	GLY	2.0
2	B	131	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.