



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:11 AM UTC

PDB ID : 6JDG / pdb_00006jdg
Title : Complexed crystal structure of PaSSB with ssDNA dT20 at 2.39 angstrom resolution
Authors : Huang, Y.H.; Huang, C.Y.
Deposited on : 2019-02-01
Resolution : 2.39 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

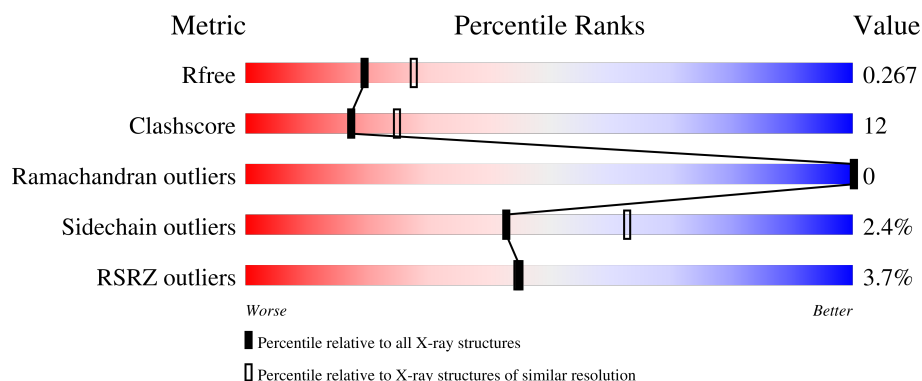
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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	7164 (2.40-2.36)
Clashscore	190562	7722 (2.40-2.36)
Ramachandran outliers	187476	7626 (2.40-2.36)
Sidechain outliers	187428	7627 (2.40-2.36)
RSRZ outliers	180081	7170 (2.40-2.36)

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	121	
1	B	121	
1	C	121	
1	D	121	
2	E	20	

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Mol	Chain	Length	Quality of chain
2	F	20	 20% 35% 45%
2	G	20	 30% 35% 35%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3724 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 Single-stranded DNA-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	101	Total	C	N	O	S	0	0	0
			797	494	148	153	2			
1	B	98	Total	C	N	O	S	0	0	0
			773	481	143	147	2			
1	C	97	Total	C	N	O	S	0	0	0
			764	476	141	145	2			
1	D	104	Total	C	N	O	S	0	0	0
			825	514	153	156	2			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	116	HIS	-	expression tag	UNP P40947
A	117	HIS	-	expression tag	UNP P40947
A	118	HIS	-	expression tag	UNP P40947
A	119	HIS	-	expression tag	UNP P40947
A	120	HIS	-	expression tag	UNP P40947
A	121	HIS	-	expression tag	UNP P40947
B	116	HIS	-	expression tag	UNP P40947
B	117	HIS	-	expression tag	UNP P40947
B	118	HIS	-	expression tag	UNP P40947
B	119	HIS	-	expression tag	UNP P40947
B	120	HIS	-	expression tag	UNP P40947
B	121	HIS	-	expression tag	UNP P40947
C	116	HIS	-	expression tag	UNP P40947
C	117	HIS	-	expression tag	UNP P40947
C	118	HIS	-	expression tag	UNP P40947
C	119	HIS	-	expression tag	UNP P40947
C	120	HIS	-	expression tag	UNP P40947
C	121	HIS	-	expression tag	UNP P40947
D	116	HIS	-	expression tag	UNP P40947
D	117	HIS	-	expression tag	UNP P40947
D	118	HIS	-	expression tag	UNP P40947

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Chain	Residue	Modelled	Actual	Comment	Reference
D	119	HIS	-	expression tag	UNP P40947
D	120	HIS	-	expression tag	UNP P40947
D	121	HIS	-	expression tag	UNP P40947

- Molecule 2 is a DNA chain called DNA (5'-D(P*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*TP*T)-3').

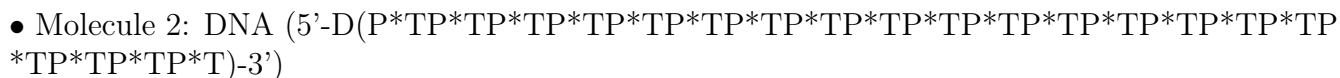
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	P	0	0	0
			60	30	6	21	3			
2	F	11	Total	C	N	O	P	0	0	0
			220	110	22	77	11			
2	G	13	Total	C	N	O	P	0	0	0
			260	130	26	91	13			

- Molecule 3 is water.

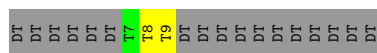
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total	O	0	0
			8	8		
3	B	10	Total	O	0	0
			10	10		
3	C	3	Total	O	0	0
			3	3		
3	D	4	Total	O	0	0
			4	4		

i

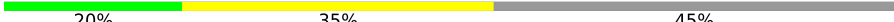
- Molecule 1: Single-stranded DNA-binding protein

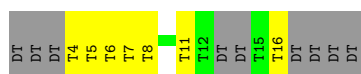


Chain E:  5% 10% 85%

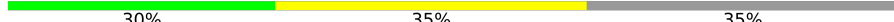


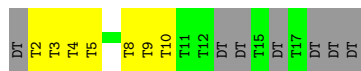
- Molecule 2: DNA (5'-D(P*T)-3')

Chain F:  20% 35% 45%



- Molecule 2: DNA (5'-D(P*T)-3')

Chain G:  30% 35% 35%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	60.04Å 60.04Å 129.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.25 – 2.39 29.25 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.25-2.39) 98.9 (29.25-2.39)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.39Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.213 , 0.264 0.217 , 0.267	Depositor DCC
R_{free} test set	1052 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	52.3	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.006 for -h,-k,l 0.057 for h,-h-k,-l 0.028 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3724	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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	0.50	0/808	0.62	0/1090
1	B	0.66	1/783 (0.1%)	0.75	2/1055 (0.2%)
1	C	0.42	0/774	0.57	0/1043
1	D	0.36	0/838	0.54	0/1130
2	E	0.52	0/65	0.73	0/98
2	F	0.40	0/240	0.87	0/366
2	G	0.41	0/283	0.67	0/430
All	All	0.48	1/3791 (0.0%)	0.65	2/5212 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	56	ARG	C-O	-5.60	1.17	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	94	GLN	CB-CA-C	-6.73	98.61	114.16
1	B	96	ARG	N-CA-C	-5.68	99.41	109.06

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	797	0	782	17	1
1	B	773	0	763	37	0
1	C	764	0	755	15	0
1	D	825	0	813	23	1
2	E	60	0	37	4	0
2	F	220	0	134	12	0
2	G	260	0	159	8	0
3	A	8	0	0	0	0
3	B	10	0	0	1	0
3	C	3	0	0	0	0
3	D	4	0	0	0	0
All	All	3724	0	3443	89	1

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 12.

All (89) 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:B:88:TRP:HD1	1:B:96:ARG:O	1.36	1.07
1:B:88:TRP:CD1	1:B:96:ARG:O	2.07	1.06
1:A:84:ARG:NH2	1:A:100:GLU:OE1	2.07	0.88
1:D:8:VAL:HG21	1:D:83:LEU:HD13	1.58	0.85
1:B:113:GLY:HA2	1:C:113:GLY:HA2	1.58	0.84
1:D:106:ASN:HA	2:F:16:DT:H1'	1.67	0.77
1:D:21:ARG:HH21	1:D:23:MET:HE1	1.51	0.76
1:D:40:TRP:H	1:D:40:TRP:CD1	2.01	0.74
1:C:90:GLY:CA	1:C:96:ARG:HH12	2.02	0.72
1:B:88:TRP:NE1	1:B:96:ARG:HB2	2.04	0.72
1:B:105:ILE:HG13	2:F:16:DT:H3'	1.71	0.72
2:F:16:DT:OP1	2:F:16:DT:O3'	2.05	0.70
1:A:3:ARG:HD2	1:C:110:GLN:OE1	1.92	0.68
2:G:4:DT:H2''	2:G:5:DT:H71	1.79	0.64
1:B:105:ILE:O	2:F:16:DT:H2'	1.98	0.64
1:B:86:ARG:HD3	1:B:100:GLU:CD	2.23	0.63
2:G:2:DT:H5''	2:G:3:DT:C2	2.34	0.62
1:C:83:LEU:HD11	1:D:83:LEU:HD11	1.82	0.61
1:B:88:TRP:CD1	1:B:96:ARG:C	2.79	0.61
1:D:3:ARG:NH1	2:E:9:DT:O2	2.35	0.60
1:B:51:ARG:O	1:B:51:ARG:HG3	1.96	0.60
1:C:90:GLY:HA2	1:C:96:ARG:HH12	1.64	0.60
1:B:106:ASN:ND2	3:B:201:HOH:O	2.28	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:ARG:HG2	1:C:102:VAL:HG21	1.85	0.59
1:B:73:LYS:HE2	2:G:8:DT:H5''	1.85	0.59
1:B:88:TRP:CD1	1:B:88:TRP:H	2.23	0.57
2:G:9:DT:H2'	2:G:10:DT:H5''	1.87	0.57
1:C:73:LYS:HE3	2:E:8:DT:H5''	1.86	0.56
2:G:4:DT:H4'	2:G:4:DT:OP1	2.06	0.55
1:A:94:GLN:NE2	1:B:96:ARG:HH22	2.06	0.54
1:B:106:ASN:O	2:F:16:DT:C4	2.61	0.54
2:F:7:DT:H2''	2:F:8:DT:O5'	2.07	0.54
1:B:88:TRP:CD1	1:B:96:ARG:HB2	2.42	0.53
1:B:88:TRP:HE1	1:B:96:ARG:CB	2.22	0.53
1:B:86:ARG:HD3	1:B:100:GLU:OE1	2.10	0.52
1:A:7:LYS:NZ	1:A:80:GLU:OE2	2.42	0.52
1:D:3:ARG:HH12	2:E:8:DT:H72	1.74	0.52
1:A:88:TRP:CH2	2:F:4:DT:H2'	2.44	0.51
1:D:40:TRP:H	1:D:40:TRP:HD1	1.51	0.51
1:A:3:ARG:N	2:G:10:DT:O2	2.44	0.51
1:C:84:ARG:NH1	1:C:100:GLU:OE1	2.39	0.51
1:D:30:THR:HG21	1:D:64:ALA:O	2.11	0.51
1:D:3:ARG:NH1	2:E:9:DT:H1'	2.26	0.50
1:A:94:GLN:NE2	1:B:96:ARG:NH2	2.59	0.50
1:B:3:ARG:HD2	1:D:111:LEU:O	2.12	0.50
1:B:8:VAL:HG21	1:B:83:LEU:HG	1.94	0.49
1:B:88:TRP:HE1	1:B:96:ARG:HB2	1.75	0.49
2:F:6:DT:H2''	2:F:7:DT:C6	2.48	0.49
1:A:83:LEU:HD11	1:B:83:LEU:HD21	1.95	0.48
1:A:7:LYS:HE3	1:C:80:GLU:OE2	2.13	0.48
1:A:86:ARG:NH1	2:F:4:DT:O4	2.47	0.47
1:B:88:TRP:CD1	1:B:88:TRP:N	2.78	0.47
2:F:6:DT:H2'	2:F:6:DT:O2	2.15	0.47
1:A:59:PHE:HB2	1:A:64:ALA:HA	1.97	0.46
1:B:7:LYS:HE3	1:D:80:GLU:OE2	2.16	0.46
1:A:3:ARG:NH2	1:C:111:LEU:H	2.14	0.46
1:A:94:GLN:HE22	1:B:96:ARG:HH22	1.62	0.46
1:D:21:ARG:NH2	1:D:23:MET:HE1	2.27	0.45
1:B:95:ASP:C	1:B:96:ARG:HG2	2.41	0.45
1:A:23:MET:HE3	1:A:27:ASN:HB3	1.99	0.45
1:D:19:GLU:OE1	1:D:21:ARG:HG3	2.17	0.45
1:B:86:ARG:NH2	1:B:100:GLU:OE2	2.45	0.44
2:G:2:DT:H3'	2:G:3:DT:H5'	1.99	0.44
1:B:71:LEU:O	1:B:72:ARG:HD2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:35:ALA:HB2	1:D:54:TRP:CH2	2.53	0.44
1:C:84:ARG:HA	1:D:53:GLU:OE2	2.18	0.43
1:C:8:VAL:HG21	1:C:83:LEU:HG	1.99	0.43
1:B:84:ARG:NH1	1:B:84:ARG:HG3	2.33	0.43
1:B:84:ARG:HG3	1:B:84:ARG:HH11	1.84	0.43
1:A:13:ASN:C	1:A:34:LEU:HD12	2.43	0.43
1:B:3:ARG:HG3	1:D:110:GLN:HE21	1.82	0.43
2:F:4:DT:H2''	2:F:5:DT:OP2	2.17	0.43
1:B:88:TRP:NE1	1:B:96:ARG:CB	2.77	0.43
1:B:79:VAL:HG13	1:B:103:VAL:HG21	2.00	0.42
1:D:19:GLU:OE2	1:D:31:ASN:HB2	2.19	0.42
1:B:8:VAL:CG2	1:B:83:LEU:HG	2.49	0.42
1:D:70:TYR:CE1	2:F:11:DT:H2''	2.54	0.42
1:B:88:TRP:HD1	1:B:88:TRP:H	1.62	0.42
2:G:8:DT:H2'	2:G:9:DT:O4'	2.19	0.42
1:A:10:LEU:HB2	1:A:79:VAL:HG22	2.01	0.41
1:D:7:LYS:NZ	1:D:80:GLU:OE1	2.54	0.41
1:B:25:ASN:HD21	1:B:27:ASN:CG	2.28	0.41
1:B:3:ARG:HG3	1:D:110:GLN:NE2	2.35	0.41
1:D:19:GLU:O	1:D:30:THR:HA	2.21	0.41
1:C:13:ASN:HA	1:C:75:SER:O	2.21	0.41
1:C:72:ARG:HA	1:C:72:ARG:HD3	1.79	0.40
1:D:20:THR:HG23	1:D:30:THR:HG22	2.01	0.40
1:A:7:LYS:HE2	1:A:7:LYS:HB2	1.89	0.40
1:C:90:GLY:C	1:C:96:ARG:HH12	2.28	0.40

All (1) 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)
1:A:84:ARG:NH1	1:D:91:GLN:O[1_655]	1.99	0.21

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	97/121 (80%)	95 (98%)	2 (2%)	0	100	100
1	B	92/121 (76%)	89 (97%)	3 (3%)	0	100	100
1	C	91/121 (75%)	89 (98%)	2 (2%)	0	100	100
1	D	100/121 (83%)	95 (95%)	5 (5%)	0	100	100
All	All	380/484 (78%)	368 (97%)	12 (3%)	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	
1	A	84/101 (83%)	83 (99%)	1 (1%)	63	79
1	B	82/101 (81%)	77 (94%)	5 (6%)	17	27
1	C	81/101 (80%)	81 (100%)	0	100	100
1	D	87/101 (86%)	85 (98%)	2 (2%)	44	64
All	All	334/404 (83%)	326 (98%)	8 (2%)	43	63

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	ARG
1	B	20	THR
1	B	51	ARG
1	B	56	ARG
1	B	94	GLN
1	B	100	GLU
1	D	30	THR
1	D	83	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	94	GLN
1	C	31	ASN
1	D	110	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	101/121 (83%)	0.11	0 100 100	29, 43, 74, 97	0
1	B	98/121 (80%)	0.37	2 (2%) 65 63	31, 48, 91, 101	0
1	C	97/121 (80%)	0.41	4 (4%) 41 41	30, 55, 84, 88	0
1	D	104/121 (85%)	0.74	10 (9%) 13 12	31, 59, 97, 109	0
2	E	3/20 (15%)	0.21	0 100 100	70, 70, 71, 74	0
2	F	11/20 (55%)	0.99	0 100 100	60, 92, 121, 122	0
2	G	13/20 (65%)	0.93	0 100 100	55, 85, 126, 127	0
All	All	427/544 (78%)	0.44	16 (3%) 45 45	29, 53, 96, 127	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	40	TRP	3.7
1	B	97	TYR	3.3
1	C	90	GLY	3.0
1	D	15	GLY	2.7
1	D	22	TYR	2.5
1	D	71	LEU	2.5
1	B	93	GLY	2.3
1	D	43	LYS	2.3
1	C	88	TRP	2.3
1	D	20	THR	2.3
1	C	62	ARG	2.1
1	C	21	ARG	2.1
1	D	32	ILE	2.1
1	D	70	TYR	2.1
1	D	29	VAL	2.1
1	D	66	ILE	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.