



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 7, 2020 – 06:05 PM EDT

PDB ID : 6UBF
Title : Role of Beta-hairpin motifs in the DNA duplex opening by the Rad4/XPC nucleotide excision repair complex
Authors : Paul, D.; Min, J.H.
Deposited on : 2019-09-11
Resolution : 4.60 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
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
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)	:	2.14.6

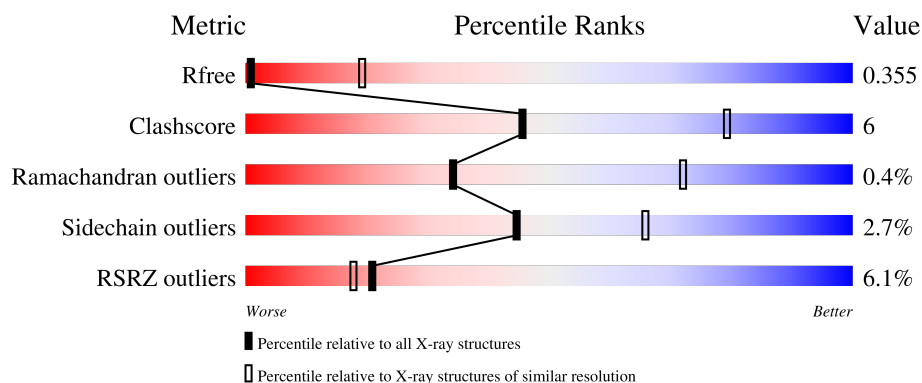
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



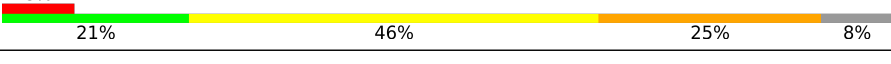

The reported resolution of this entry is 4.60 Å.

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}	130704	1062 (5.40-3.80)
Clashscore	141614	1130 (5.40-3.80)
Ramachandran outliers	138981	1074 (5.40-3.80)
Sidechain outliers	138945	1055 (5.40-3.80)
RSRZ outliers	127900	1113 (5.50-3.70)

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 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	531	
2	X	171	
3	W	24	
4	Y	24	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10147 atoms, of which 4913 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 DNA repair protein RAD4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	479	7950	2516	4004	705	699	26	0	0	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	95	GLY	-	expression tag	UNP P14736
A	96	SER	-	expression tag	UNP P14736
A	97	SER	-	expression tag	UNP P14736
A	98	ARG	-	expression tag	UNP P14736
A	99	ALA	-	expression tag	UNP P14736
A	100	MET	-	expression tag	UNP P14736
A	115	THR	LYS	conflict	UNP P14736
A	131	CYS	VAL	conflict	UNP P14736
A	132	SER	CYS	conflict	UNP P14736
A	223	GLU	VAL	conflict	UNP P14736
A	427	ARG	GLN	conflict	UNP P14736
A	?	-	GLU	deletion	UNP P14736
A	?	-	ARG	deletion	UNP P14736
A	?	-	GLY	deletion	UNP P14736
A	?	-	SER	deletion	UNP P14736
A	?	-	THR	deletion	UNP P14736
A	?	-	VAL	deletion	UNP P14736
A	?	-	LYS	deletion	UNP P14736

- Molecule 2 is a protein called UV excision repair protein RAD23.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
2	X	53	827	260	418	69	78	2	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	228	GLY	-	expression tag	UNP P32628
X	229	SER	-	expression tag	UNP P32628

- Molecule 3 is a DNA chain called DNA (5'-D(*TP*TP*GP*AP*CP*TP*CP*(G47)P*AP*CP*AP*TP*CP*CP*C*GP*CP*TP*AP*CP*AP*A)-3').

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
3	W	22	Total	C	H	N	O	P	S	0	0	0
			693	214	252	79	127	20	1			

- Molecule 4 is a DNA chain called DNA (5'-D(*AP*TP*TP*GP*TP*AP*GP*CP*GP*GP*GP*AP*TP*GP*TP*CP*GP*AP*GP*TP*CP*A)-3').

Mol	Chain	Residues	Atoms							ZeroOcc	AltConf	Trace
4	Y	22	Total	C	H	N	O	P		0	0	0
			677	208	239	81	128	21				

- Molecule 4: DNA (5'-D(*AP*TP*TP*GP*TP*AP*GP*CP*GP*GP*GP*AP*TP*GP*TP*CP*GP*AP*GP*TP*CP*A)-3')

Chain Y:  63% 29% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	78.72Å 78.72Å 403.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.75 – 4.60 48.75 – 4.60	Depositor EDS
% Data completeness (in resolution range)	84.6 (48.75-4.60) 85.2 (48.75-4.60)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 4.64Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.314 , 0.356 0.316 , 0.355	Depositor DCC
R_{free} test set	309 reflections (4.67%)	wwPDB-VP
Wilson B-factor (Å ²)	185.7	Xtriage
Anisotropy	0.463	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 188.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	10147	wwPDB-VP
Average B, all atoms (Å ²)	279.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% 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

Bond lengths and bond angles in the following residue types are not validated in this section: G47

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	0.38	0/4030	0.74	3/5417 (0.1%)
2	X	0.34	0/415	0.83	0/565
3	W	1.51	8/463 (1.7%)	1.38	4/706 (0.6%)
4	Y	1.17	1/490 (0.2%)	1.15	0/754
All	All	0.66	9/5398 (0.2%)	0.87	7/7442 (0.1%)

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	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	W	22	DC	C1'-N1	13.56	1.66	1.49
4	Y	2	DT	C1'-N1	9.97	1.62	1.49
3	W	21	DA	C3'-O3'	-8.76	1.32	1.44
3	W	22	DC	C4'-O4'	7.49	1.52	1.45
3	W	12	DT	C1'-N1	6.48	1.57	1.49
3	W	23	DA	P-O5'	6.41	1.66	1.59
3	W	2	DT	C3'-O3'	-5.93	1.36	1.44
3	W	22	DC	N1-C6	5.90	1.40	1.37
3	W	3	DG	C3'-O3'	-5.43	1.36	1.44

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	W	14	DC	O4'-C1'-C2'	-5.36	101.61	105.90
1	A	129	ARG	NE-CZ-NH1	-5.30	117.65	120.30
3	W	22	DC	C6-N1-C2	-5.22	118.21	120.30
3	W	5	DC	OP1-P-OP2	5.17	127.36	119.60
1	A	243	PHE	CB-CG-CD2	-5.13	117.21	120.80
3	W	13	DC	P-O5'-C5'	-5.13	112.70	120.90
1	A	219	LYS	CA-CB-CG	5.07	124.55	113.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	243	PHE	Sidechain
1	A	245	THR	Mainchain

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	3946	4004	4002	48	0
2	X	409	418	418	7	0
3	W	441	252	254	14	1
4	Y	438	239	239	3	0
All	All	5234	4913	4913	63	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 6.

All (63) 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:X:266:ARG:NH1	2:X:305:LEU:O	2.16	0.78
1:A:414:ARG:NH1	1:A:419:ASP:OD1	2.22	0.72
1:A:377:ARG:NH2	1:A:384:ASN:OD1	2.29	0.65
1:A:230:TRP:NE1	1:A:290:ASP:O	2.26	0.63
1:A:244:LYS:O	2:X:272:ASN:ND2	2.23	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:CYS:O	1:A:296:ILE:HD11	2.00	0.62
2:X:260:GLU:N	2:X:260:GLU:OE1	2.33	0.60
1:A:367:ASP:OD1	1:A:371:GLY:N	2.33	0.60
2:X:289:GLN:N	2:X:289:GLN:OE1	2.36	0.59
1:A:465:GLU:N	1:A:465:GLU:OE1	2.36	0.59
1:A:457:GLN:NE2	1:A:489:ASP:OD1	2.33	0.59
1:A:492:SER:N	1:A:495:GLN:OE1	2.33	0.56
1:A:211:GLN:NE2	1:A:332:PRO:O	2.39	0.55
4:Y:13:DG:H2''	4:Y:14:DA:C8	2.42	0.55
1:A:590:PHE:CD2	1:A:612:ILE:HD12	2.41	0.55
1:A:137:ARG:NE	1:A:294:MET:SD	2.80	0.55
3:W:9:DA:H2''	3:W:10:DC:C6	2.42	0.54
1:A:438:VAL:HG23	3:W:4:DA:OP1	2.09	0.53
1:A:464:LYS:HB2	1:A:465:GLU:OE1	2.09	0.52
1:A:495:GLN:NE2	3:W:15:DC:OP1	2.42	0.52
1:A:273:LEU:O	1:A:278:VAL:HG22	2.10	0.51
1:A:134:ASN:ND2	3:W:11:DA:H4'	2.26	0.51
4:Y:3:DT:H1'	4:Y:4:DG:C8	2.47	0.50
1:A:588:VAL:HG21	1:A:618:LEU:HD13	1.93	0.50
3:W:9:DA:H2''	3:W:10:DC:H6	1.75	0.50
1:A:377:ARG:HD2	1:A:428:ARG:HB3	1.93	0.49
3:W:10:DC:OP1	3:W:10:DC:H4'	2.13	0.49
1:A:237:ALA:HB2	2:X:281:GLU:HG2	1.93	0.49
1:A:513:ILE:HG12	1:A:529:GLU:O	2.13	0.48
3:W:14:DC:H2''	3:W:15:DC:C6	2.49	0.47
1:A:221:ASP:O	1:A:222:ASN:C	2.51	0.47
4:Y:18:DC:H2''	4:Y:19:DG:C8	2.48	0.47
1:A:217:THR:O	1:A:217:THR:HG23	2.15	0.47
1:A:137:ARG:HG2	1:A:294:MET:SD	2.55	0.46
1:A:233:ILE:O	1:A:234:GLU:C	2.55	0.46
1:A:130:ASN:OD1	1:A:296:ILE:HB	2.16	0.45
1:A:321:ASP:OD2	1:A:324:SER:OG	2.32	0.45
1:A:576:ASN:HB3	1:A:579:ALA:HB2	1.99	0.45
1:A:331:ASP:O	1:A:336:LYS:HA	2.16	0.44
1:A:342:ARG:O	1:A:428:ARG:NH1	2.51	0.44
1:A:434:ILE:H	1:A:434:ILE:HD12	1.82	0.44
1:A:158:ARG:NH1	1:A:267:GLN:OE1	2.44	0.44
1:A:438:VAL:HG13	1:A:449:LEU:HD11	2.00	0.44
1:A:168:SER:HA	1:A:275:ALA:HB2	2.00	0.43
1:A:227:MET:HB2	1:A:287:GLN:HB3	1.99	0.43
3:W:2:DT:H2''	3:W:3:DG:C8	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:W:11:DA:H2''	3:W:12:DT:H5'	2.01	0.42
1:A:129:ARG:CZ	3:W:9:DA:H4'	2.49	0.42
1:A:495:GLN:HG2	3:W:15:DC:OP2	2.19	0.42
1:A:373:ARG:NH1	1:A:418:ASP:HA	2.34	0.42
1:A:384:ASN:O	1:A:445:PRO:HD2	2.19	0.42
1:A:228:ARG:NH1	2:X:277:ALA:HB3	2.34	0.42
3:W:13:DC:H2''	3:W:14:DC:C6	2.55	0.42
1:A:355:CYS:HA	1:A:358:ASN:O	2.20	0.41
1:A:266:VAL:HG11	1:A:316:TRP:HA	2.02	0.41
1:A:134:ASN:HB2	3:W:11:DA:OP1	2.21	0.41
1:A:471:VAL:CG2	1:A:479:LEU:HG	2.51	0.41
2:X:287:TYR:O	2:X:290:LEU:HB3	2.20	0.41
1:A:137:ARG:CG	1:A:294:MET:SD	3.09	0.41
1:A:134:ASN:OD1	1:A:137:ARG:HD2	2.21	0.41
1:A:377:ARG:HB2	1:A:383:MET:SD	2.61	0.41
1:A:385:SER:HB2	1:A:440:ASP:HB3	2.04	0.41
3:W:23:DA:H2''	3:W:24:DA:C8	2.56	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)
3:W:1:DT:O5'	3:W:24:DA:O3'[1_565]	1.75	0.45

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	469/531 (88%)	446 (95%)	21 (4%)	2 (0%)	34	72
2	X	51/171 (30%)	46 (90%)	5 (10%)	0	100	100
All	All	520/702 (74%)	492 (95%)	26 (5%)	2 (0%)	34	72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	247	LYS
1	A	238	ASN

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	432/474 (91%)	419 (97%)	13 (3%)	41	63
2	X	46/129 (36%)	46 (100%)	0	100	100
All	All	478/603 (79%)	465 (97%)	13 (3%)	44	66

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	ARG
1	A	143	PHE
1	A	147	TYR
1	A	221	ASP
1	A	243	PHE
1	A	325	LYS
1	A	345	SER
1	A	377	ARG
1	A	441	LEU
1	A	456	THR
1	A	465	GLU
1	A	539	LEU
1	A	555	THR

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

Mol	Chain	Res	Type
1	A	499	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	G47	W	8	1,3,4	20,27,28	4.74	10 (50%)	21,38,41	1.79	5 (23%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	G47	W	8	1,3,4	-	4/7/25/26	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	W	8	G47	C4-N3	9.57	1.50	1.35
3	W	8	G47	O4'-C1'	8.44	1.61	1.42
3	W	8	G47	C2-N2	8.32	1.47	1.34
3	W	8	G47	C6-C5	6.70	1.52	1.41
3	W	8	G47	O4'-C4'	-6.64	1.30	1.45
3	W	8	G47	C6-N1	6.57	1.44	1.33
3	W	8	G47	C2'-C1'	-6.24	1.34	1.52
3	W	8	G47	C2-N3	3.51	1.45	1.34
3	W	8	G47	O3'-C3'	-3.37	1.36	1.43
3	W	8	G47	C2-N1	3.24	1.44	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	W	8	G47	C2-N3-C4	4.37	120.23	115.28
3	W	8	G47	N3-C2-N1	-4.03	119.86	126.23
3	W	8	G47	C6-N1-C2	2.65	119.93	115.18
3	W	8	G47	C5-C6-N1	-2.65	119.81	123.43
3	W	8	G47	N2-C2-N3	2.09	120.81	117.19

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	W	8	G47	N2-C6A-C7A-SG
3	W	8	G47	C3'-C4'-C5'-O5'
3	W	8	G47	O4'-C4'-C5'-O5'
3	W	8	G47	C7A-C6A-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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

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	479/531 (90%)	0.33	33 (6%) 16 14	217, 247, 283, 323	0
2	X	53/171 (30%)	-0.27	0 100 100	232, 248, 258, 262	0
3	W	21/24 (87%)	0.35	2 (9%) 8 8	262, 287, 322, 343	0
4	Y	22/24 (91%)	0.11	0 100 100	269, 284, 313, 318	0
All	All	575/750 (76%)	0.27	35 (6%) 21 18	217, 250, 303, 343	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	459	VAL	5.4
1	A	526	GLU	4.5
1	A	458	ILE	4.0
1	A	238	ASN	3.3
1	A	516	THR	3.3
1	A	525	GLU	3.1
1	A	489	ASP	3.0
1	A	488	ALA	2.9
1	A	455	GLN	2.8
1	A	490	LEU	2.8
1	A	460	LYS	2.8
1	A	482	TYR	2.7
1	A	589	GLU	2.5
1	A	132	SER	2.5
1	A	463	CYS	2.3
1	A	464	LYS	2.3
1	A	551	ILE	2.3
1	A	588	VAL	2.3
1	A	493	ALA	2.3
1	A	468	TYR	2.2
1	A	568	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	552	THR	2.2
1	A	543	PRO	2.2
1	A	483	ALA	2.2
1	A	456	THR	2.2
1	A	527	GLU	2.1
3	W	13	DC	2.1
1	A	342	ARG	2.1
1	A	461	PRO	2.1
1	A	301	ASN	2.1
1	A	341	VAL	2.0
1	A	340	GLN	2.0
1	A	450	GLU	2.0
1	A	453	ILE	2.0
3	W	15	DC	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	G47	W	8	25/26	0.83	0.31	266,273,328,331	0

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.