



# Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 07:51 am BST

PDB ID : 5OV3  
Title : Structure of the RbBP5 beta-propeller domain  
Authors : Mittal, A.; Zhang, Y.; Gamblin, S.J.; Wilson, J.R.  
Deposited on : 2017-08-27  
Resolution : 2.45 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](mailto: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.

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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 : 2.11  
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) : 2.11

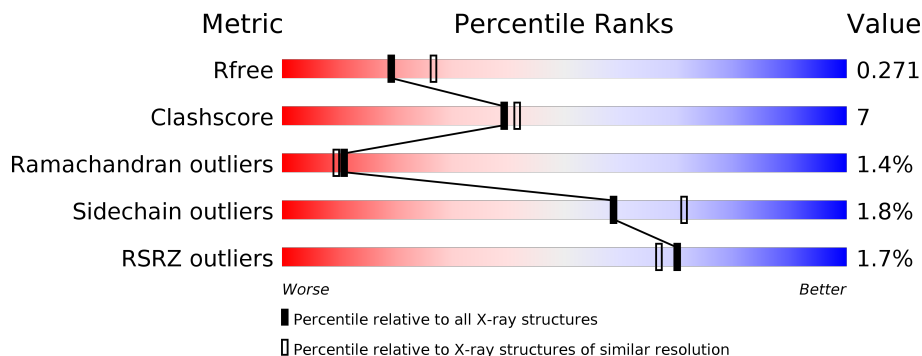
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.45 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	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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  $\geq 3$ , 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	381	 2% (red), 68% (green), 17% (yellow), 13% (grey)
1	B	381	 % (red), 69% (green), 13% (yellow), 17% (grey)
2	C	6	 83% (green), 17% (yellow)

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5153 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 Retinoblastoma-binding protein 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	330	Total	C	N	O	S	0	0	0
			2586	1634	452	485	15			
1	B	317	Total	C	N	O	S	0	0	0
			2462	1554	437	456	15			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q8BX09
A	1	PRO	-	expression tag	UNP Q8BX09
A	363	LYS	GLU	variant	UNP Q8BX09
B	0	GLY	-	expression tag	UNP Q8BX09
B	1	PRO	-	expression tag	UNP Q8BX09
B	363	LYS	GLU	variant	UNP Q8BX09

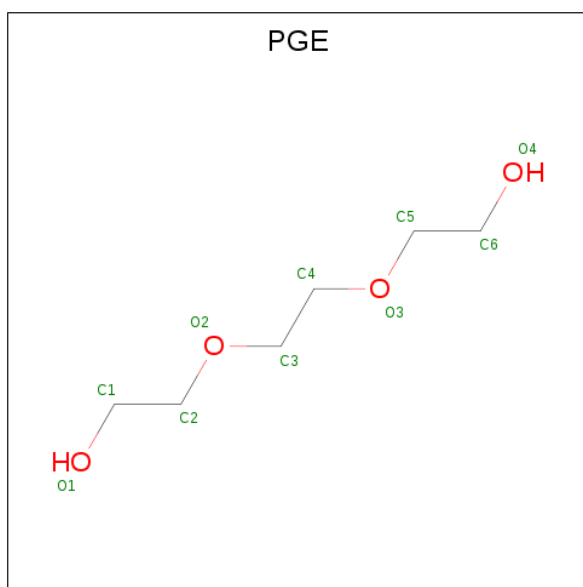
- Molecule 2 is a protein called Retinoblastoma-binding protein 5.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	6	Total	C	N	O	0	0	0
			45	27	8	10			

There is a discrepancy between the modelled and reference sequences:

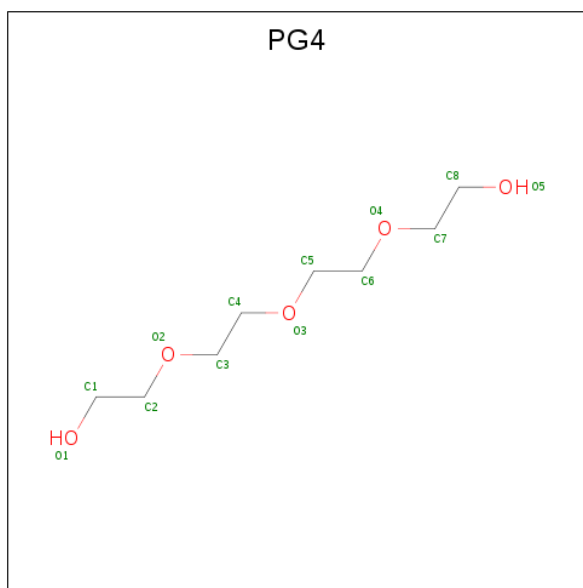
Chain	Residue	Modelled	Actual	Comment	Reference
C	363	LYS	GLU	variant	UNP Q8BX09

- Molecule 3 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula:  $C_8H_{18}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			13	8	5		

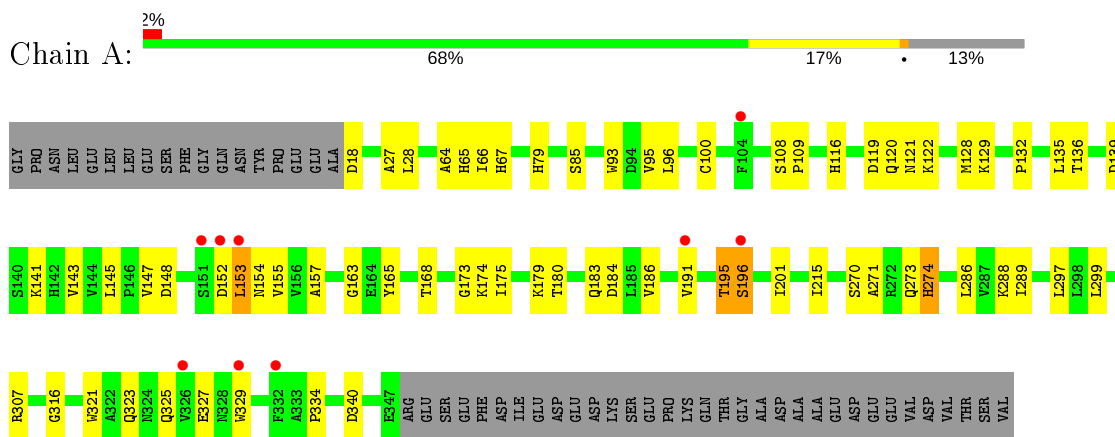
- Molecule 5 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	A	13	Total 13	O 13	0	0
5	B	24	Total 24	O 24	0	0

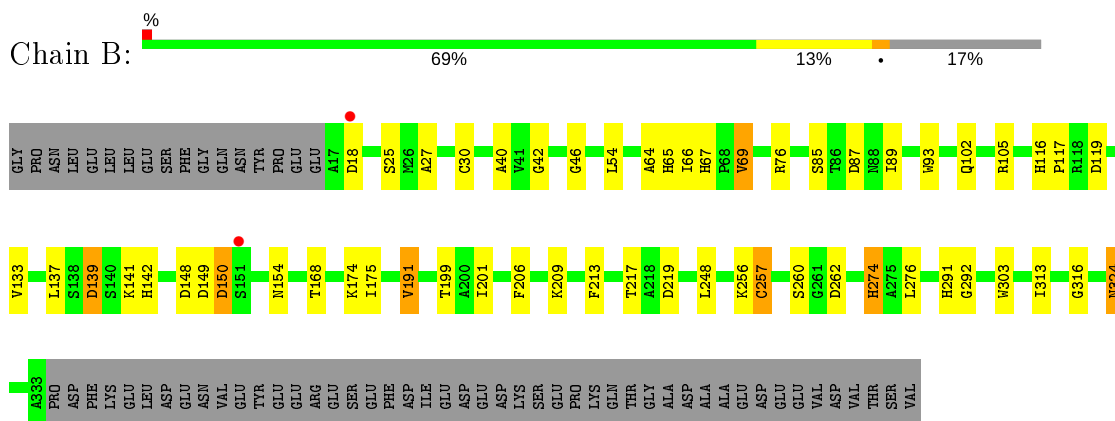
### 3 Residue-property plots [i](#)

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.

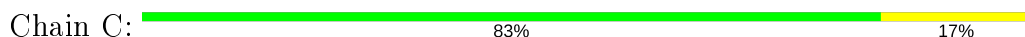
- Molecule 1: Retinoblastoma-binding protein 5



- Molecule 1: Retinoblastoma-binding protein 5



- Molecule 2: Retinoblastoma-binding protein 5



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.81Å 71.90Å 178.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.68 – 2.45 66.68 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.6 (66.68-2.45) 99.6 (66.68-2.45)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.62 (at 2.45Å)	Xtrriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, $R_{free}$	0.196 , 0.271 0.196 , 0.271	Depositor DCC
$R_{free}$ test set	1407 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	54.8	Xtrriage
Anisotropy	0.489	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 50.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5153	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, PGE

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	0.41	0/2642	0.64	0/3589
1	B	0.48	1/2514 (0.0%)	0.65	0/3416
2	C	0.48	0/45	0.63	0/59
All	All	0.44	1/5201 (0.0%)	0.65	0/7064

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	257	CYS	CB-SG	-5.97	1.72	1.81

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	A	2586	0	2565	44	0
1	B	2462	0	2461	35	0
2	C	45	0	43	1	0
3	A	10	0	14	2	0
4	B	13	0	18	0	0
5	A	13	0	0	0	0
5	B	24	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5153	0	5101	76	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 7.

All (76) 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:199:THR:HG21	1:B:219:ASP:HB3	1.61	0.83
1:B:148:ASP:OD2	1:B:174:LYS:NZ	2.13	0.80
1:A:195:THR:OG1	1:A:196:SER:N	2.15	0.78
1:A:132:PRO:HB2	1:A:145:LEU:HD12	1.73	0.71
1:B:291:HIS:HB2	2:C:365:THR:HG22	1.73	0.69
1:B:139:ASP:OD1	1:B:141:LYS:HE2	1.93	0.68
1:A:157:ALA:HB2	1:A:168:THR:HG22	1.80	0.63
1:B:217:THR:OG1	1:B:219:ASP:OD1	2.16	0.60
1:A:334:PRO:HD3	1:B:25:SER:HB2	1.85	0.59
1:A:129:LYS:O	1:A:153:LEU:HD13	2.02	0.58
1:B:102:GLN:HE22	1:B:137:LEU:HB3	1.68	0.58
1:B:148:ASP:O	1:B:150:ASP:N	2.38	0.56
1:B:42:GLY:HA3	1:B:69:VAL:CG2	2.35	0.55
1:A:65:HIS:ND1	1:A:85:SER:OG	2.41	0.54
1:B:30:CYS:HB3	1:B:313:ILE:HD11	1.88	0.54
1:A:143:VAL:HG11	1:A:183:GLN:NE2	2.23	0.53
1:A:109:PRO:HG2	1:A:128:MET:HG3	1.89	0.53
1:A:96:LEU:HD22	1:A:96:LEU:H	1.73	0.53
1:B:76:ARG:HD2	1:B:117:PRO:O	2.10	0.52
1:A:18:ASP:HB2	1:A:321:TRP:CE3	2.44	0.51
1:A:270:SER:OG	1:A:273:GLN:HG2	2.11	0.51
1:B:67:HIS:ND1	1:B:87:ASP:OD1	2.44	0.51
1:B:274:HIS:CD2	1:B:292:GLY:HA3	2.46	0.51
1:B:168:THR:O	1:B:175:ILE:HA	2.11	0.50
1:A:163:GLY:O	1:A:180:THR:HG21	2.12	0.49
1:B:148:ASP:C	1:B:150:ASP:H	2.15	0.49
1:A:79:HIS:ND1	1:A:96:LEU:HD21	2.27	0.49
1:A:79:HIS:O	1:A:95:VAL:HG22	2.12	0.48
1:A:286:LEU:HD21	1:A:289:ILE:HD11	1.94	0.48
1:B:206:PHE:CD2	1:B:213:PHE:HB3	2.48	0.48
1:A:116:HIS:HB3	1:A:119:ASP:O	2.13	0.48
1:B:18:ASP:OD1	1:B:18:ASP:N	2.45	0.47
1:B:89:ILE:CD1	1:B:105:ARG:HG2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:HIS:HB3	1:B:119:ASP:O	2.15	0.47
1:B:133:VAL:CG1	1:B:142:HIS:HB2	2.46	0.46
1:A:323:GLN:HB2	3:A:401:PGE:H5	1.98	0.46
1:B:191:VAL:HG11	1:B:201:ILE:HD11	1.96	0.46
1:A:122:LYS:HG2	1:A:136:THR:HG22	1.98	0.46
1:B:66:ILE:HD12	1:B:66:ILE:H	1.81	0.46
1:A:325:GLN:OE1	1:A:325:GLN:N	2.50	0.45
1:A:201:ILE:HG23	1:A:215:ILE:HG23	1.99	0.45
1:A:143:VAL:HG11	1:A:183:GLN:HE21	1.82	0.44
1:A:96:LEU:HD22	1:A:96:LEU:N	2.32	0.44
1:A:152:ASP:HB3	1:A:155:VAL:HG23	1.98	0.44
1:A:121:ASN:O	1:A:136:THR:HA	2.18	0.44
1:A:93:TRP:CZ3	1:A:100:CYS:HB2	2.53	0.44
1:A:307:ARG:CZ	1:B:54:LEU:HD12	2.48	0.43
1:A:173:GLY:O	1:A:191:VAL:HG12	2.17	0.43
1:B:141:LYS:HE2	1:B:141:LYS:HB2	1.86	0.43
1:B:46:GLY:HA3	1:B:65:HIS:O	2.18	0.43
1:A:65:HIS:CG	1:A:85:SER:OG	2.72	0.43
1:A:168:THR:O	1:A:175:ILE:HA	2.18	0.43
1:A:139:ASP:O	1:A:141:LYS:N	2.42	0.43
1:A:152:ASP:O	1:A:154:ASN:N	2.51	0.43
1:A:340:ASP:HB3	1:B:324:ASN:OD1	2.19	0.43
1:A:271:ALA:HB1	1:A:297:LEU:CD1	2.49	0.42
1:B:27:ALA:O	1:B:316:GLY:HA2	2.19	0.42
1:A:27:ALA:O	1:A:316:GLY:HA2	2.20	0.42
1:A:66:ILE:HG23	1:A:67:HIS:HD2	1.85	0.42
1:A:165:TYR:CE1	1:A:179:LYS:HG3	2.55	0.42
1:A:271:ALA:HB1	1:A:297:LEU:HD11	2.01	0.42
1:B:256:LYS:NZ	5:B:501:HOH:O	2.50	0.42
1:B:65:HIS:CG	1:B:85:SER:HB3	2.54	0.42
1:B:64:ALA:HA	1:B:93:TRP:CH2	2.55	0.41
1:A:147:VAL:HG13	1:A:152:ASP:CG	2.41	0.41
1:A:28:LEU:HD12	1:A:299:LEU:HD13	2.02	0.41
1:B:65:HIS:ND1	1:B:85:SER:HB3	2.34	0.41
1:A:148:ASP:OD2	1:A:174:LYS:HG2	2.19	0.41
1:A:64:ALA:HA	1:A:93:TRP:CH2	2.55	0.41
1:B:260:SER:HA	1:B:303:TRP:CG	2.56	0.41
1:A:184:ASP:O	1:A:186:VAL:HG13	2.21	0.41
1:B:209:LYS:HG3	1:B:262:ASP:HA	2.03	0.41
1:B:30:CYS:HA	1:B:40:ALA:O	2.21	0.40
1:A:122:LYS:HA	1:A:135:LEU:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:LYS:NZ	3:A:401:PGE:O2	2.51	0.40
1:B:248:LEU:HA	1:B:248:LEU:HD23	1.95	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	328/381 (86%)	309 (94%)	12 (4%)	7 (2%)	7	5
1	B	315/381 (83%)	298 (95%)	15 (5%)	2 (1%)	25	29
2	C	4/6 (67%)	4 (100%)	0	0	100	100
All	All	647/768 (84%)	611 (94%)	27 (4%)	9 (1%)	11	9

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	THR
1	A	327	GLU
1	B	149	ASP
1	A	153	LEU
1	A	329	TRP
1	B	139	ASP
1	A	196	SER
1	A	274	HIS
1	A	108	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	288/332 (87%)	286 (99%)	2 (1%)	84	90
1	B	273/332 (82%)	265 (97%)	8 (3%)	42	53
2	C	5/5 (100%)	5 (100%)	0	100	100
All	All	566/669 (85%)	556 (98%)	10 (2%)	59	71

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

Mol	Chain	Res	Type
1	A	120	GLN
1	A	274	HIS
1	B	69	VAL
1	B	150	ASP
1	B	154	ASN
1	B	191	VAL
1	B	257	CYS
1	B	274	HIS
1	B	276	LEU
1	B	324	ASN

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

Mol	Chain	Res	Type
1	B	102	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 carbohydrates in this entry.

## 5.6 Ligand geometry

2 ligands are 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	PGE	A	401	-	9,9,9	0.27	0	8,8,8	0.29	0
4	PG4	B	401	-	12,12,12	0.56	0	11,11,11	0.39	0

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	PGE	A	401	-	-	1/7/7/7	-
4	PG4	B	401	-	-	2/10/10/10	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	401	PG4	O4-C7-C8-O5
4	B	401	PG4	O3-C5-C6-O4
3	A	401	PGE	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	PGE	2	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	330/381 (86%)	0.12	9 (2%) 54 50	39, 58, 95, 131	0
1	B	317/381 (83%)	0.10	2 (0%) 89 89	33, 53, 87, 120	0
2	C	6/6 (100%)	0.01	0 100 100	63, 64, 72, 75	0
All	All	653/768 (85%)	0.11	11 (1%) 70 67	33, 56, 91, 131	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	151	SER	6.4
1	B	151	SER	6.3
1	A	152	ASP	5.4
1	A	329	TRP	4.9
1	A	326	VAL	3.3
1	B	18	ASP	3.1
1	A	104	PHE	3.0
1	A	196	SER	2.6
1	A	332	PHE	2.4
1	A	153	LEU	2.3
1	A	191	VAL	2.1

### 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](#)

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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	PGE	A	401	10/10	0.95	0.16	54,58,68,68	0
4	PG4	B	401	13/13	0.95	0.14	44,49,65,66	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.