



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:48 AM GMT

PDB ID : 2FDB  
Title : Crystal Structure of Fibroblast growth factor (FGF)8b in complex with FGF Receptor (FGFR) 2c  
Authors : Mohammadi, M.; Olsen, S.K.  
Deposited on : 2005-12-13  
Resolution : 2.28 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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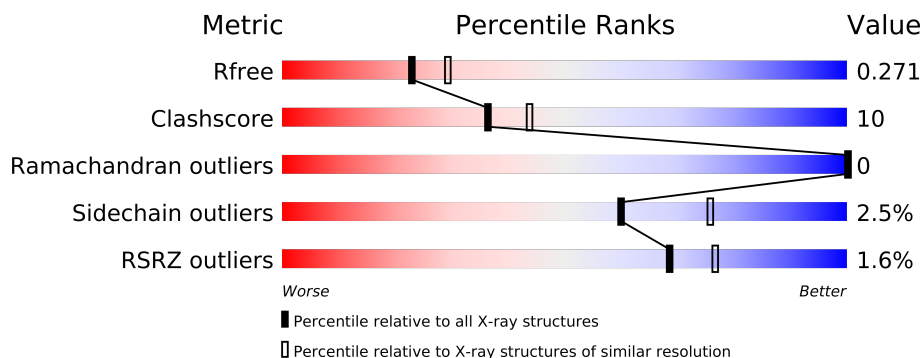
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.28 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))
$R_{free}$	66092	3861 (2.30-2.26)
Clashscore	79885	4801 (2.30-2.26)
Ramachandran outliers	78287	4729 (2.30-2.26)
Sidechain outliers	78261	4728 (2.30-2.26)
RSRZ outliers	66119	3864 (2.30-2.26)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	M	164	
1	N	164	
2	P	220	
2	R	220	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5345 atoms, of which 0 are hydrogen and 0 are deuterium.

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 fibroblast growth factor 8 isoform B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	M	149	Total	C	N	O	S	0	0	0
			1161	727	220	208	6			
1	N	148	Total	C	N	O	S	0	0	0
			1128	708	209	205	6			

- Molecule 2 is a protein called Fibroblast growth factor receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	197	Total	C	N	O	S	0	0	0
			1506	957	272	269	8			
2	R	191	Total	C	N	O	S	0	0	0
			1425	910	254	253	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	2253	ARG	PRO	ENGINEERED	UNP Q9UQH9
R	3253	ARG	PRO	ENGINEERED	UNP Q9UQH9

- Molecule 3 is water.

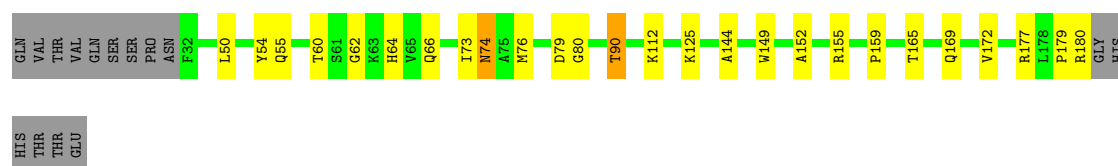
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	M	34	Total	O	0	0
			34	34		
3	N	21	Total	O	0	0
			21	21		
3	P	39	Total	O	0	0
			39	39		
3	R	31	Total	O	0	0
			31	31		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

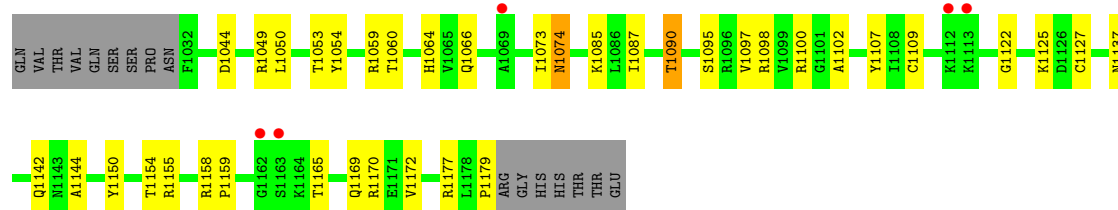
- Molecule 1: fibroblast growth factor 8 isoform B

Chain M: 



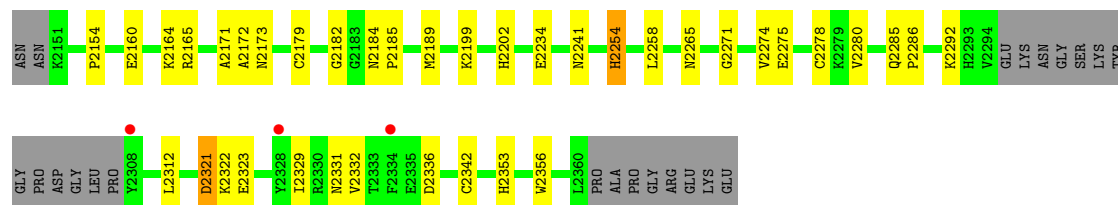
- Molecule 1: fibroblast growth factor 8 isoform B

Chain N: 



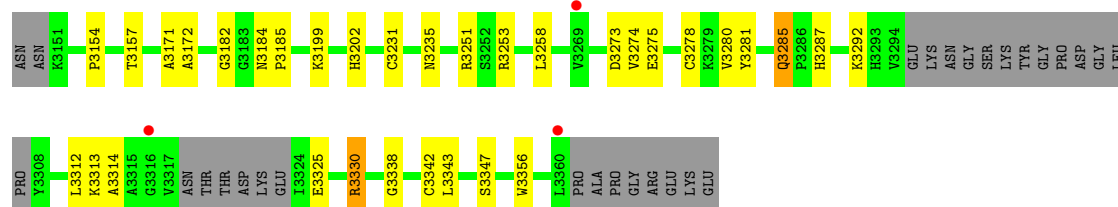
- Molecule 2: Fibroblast growth factor receptor 2

Chain P: 



- Molecule 2: Fibroblast growth factor receptor 2

Chain R: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	170.85Å 46.91Å 109.62Å 90.00° 91.66° 90.00°	Depositor
Resolution (Å)	25.00 – 2.28 29.95 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.8 (25.00-2.28) 97.6 (29.95-2.20)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.58 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.239 , 0.272 0.240 , 0.271	Depositor DCC
$R_{free}$ test set	6690 reflections (18.00%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.8	Xtriage
Anisotropy	0.174	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 28.0	EDS
Estimated twinning fraction	0.013 for -h,-k,l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 44825 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	5345	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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	M	0.37	0/1181	0.64	0/1592
1	N	0.35	0/1148	0.63	0/1553
2	P	0.37	0/1543	0.63	0/2099
2	R	0.34	0/1461	0.65	0/1993
All	All	0.36	0/5333	0.64	0/7237

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
2	R	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	R	3281	TYR	Sidechain

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	M	1161	0	1130	26	0
1	N	1128	0	1072	34	0
2	P	1506	0	1454	31	0
2	R	1425	0	1340	24	0
3	M	34	0	0	0	0
3	N	21	0	0	0	0
3	P	39	0	0	0	0
3	R	31	0	0	1	0
All	All	5345	0	4996	99	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 10.

All (99) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:60:THR:HG22	1:M:155:ARG:HA	1.44	0.96
2:P:2265:ASN:ND2	2:P:2356:TRP:H	1.69	0.91
1:M:55:GLN:HE22	1:M:80:GLY:H	1.21	0.88
2:R:3199:LYS:H	2:R:3202:HIS:CD2	1.92	0.87
1:N:1066:GLN:HE21	1:N:1074:ASN:HD21	1.24	0.86
2:P:2199:LYS:H	2:P:2202:HIS:CD2	1.94	0.84
2:R:3258:LEU:HD22	2:R:3280:VAL:HG22	1.60	0.84
1:N:1060:THR:HG22	1:N:1155:ARG:HA	1.62	0.80
2:P:2258:LEU:HD22	2:P:2280:VAL:HG22	1.65	0.77
1:N:1170:ARG:HG2	1:N:1170:ARG:HH11	1.49	0.77
2:P:2265:ASN:HD22	2:P:2356:TRP:H	1.32	0.75
2:P:2199:LYS:H	2:P:2202:HIS:HD2	1.37	0.73
1:N:1109:CYS:HG	1:N:1127:CYS:HG	0.73	0.72
1:M:60:THR:HG22	1:M:155:ARG:CA	2.20	0.71
1:M:90:THR:HG22	2:P:2285:GLN:HE21	1.54	0.71
1:M:66:GLN:HE21	1:M:74:ASN:HD21	1.34	0.71
1:N:1169:GLN:O	1:N:1172:VAL:HG12	1.92	0.69
1:N:1064:HIS:CE1	1:N:1177:ARG:HH11	2.11	0.69
1:M:50:LEU:HD11	2:P:2323:GLU:HG3	1.74	0.69
1:N:1137:ASN:O	2:R:3251:ARG:HG3	1.92	0.69
1:M:79:ASP:OD2	1:M:180:ARG:HD3	1.94	0.68
2:P:2254:HIS:HE1	2:R:3253:ARG:H	1.44	0.65
1:N:1064:HIS:HE1	1:N:1177:ARG:HH11	1.45	0.64
1:N:1050:LEU:HD22	2:R:3285:GLN:HE22	1.63	0.64
1:M:64:HIS:HE1	1:M:177:ARG:HH11	1.47	0.62
1:N:1150:TYR:HB3	1:N:1172:VAL:HG23	1.84	0.59
2:P:2254:HIS:CE1	2:R:3253:ARG:H	2.20	0.59
1:N:1064:HIS:HE1	1:N:1177:ARG:HD3	1.68	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:1125:LYS:HE2	1:N:1144:ALA:O	2.05	0.56
1:M:74:ASN:HD22	1:M:74:ASN:C	2.10	0.55
2:R:3278:CYS:CB	2:R:3342:CYS:HG	2.21	0.54
1:N:1060:THR:HG22	1:N:1155:ARG:CA	2.34	0.54
1:M:62:GLY:HA3	2:P:2164:LYS:HG3	1.89	0.54
1:M:50:LEU:CD1	2:P:2323:GLU:HG3	2.38	0.54
1:N:1050:LEU:HD22	2:R:3285:GLN:NE2	2.25	0.52
1:N:1049:ARG:HD2	1:N:1090:THR:O	2.09	0.52
2:R:3313:LYS:HG2	2:R:3314:ALA:O	2.10	0.52
2:P:2258:LEU:HD13	2:P:2278:CYS:SG	2.50	0.52
2:P:2278:CYS:HG	2:P:2342:CYS:HG	0.54	0.51
1:M:55:GLN:NE2	1:M:80:GLY:H	1.99	0.51
1:N:1074:ASN:C	1:N:1074:ASN:HD22	2.13	0.51
1:N:1100:ARG:HD2	1:N:1107:TYR:CZ	2.45	0.51
2:R:3274:VAL:HG22	2:R:3275:GLU:N	2.26	0.50
1:N:1170:ARG:HG2	1:N:1170:ARG:NH1	2.21	0.50
2:R:3157:THR:O	2:R:3157:THR:HG22	2.11	0.50
2:P:2171:ALA:O	2:P:2172:ALA:HB3	2.12	0.50
2:P:2271:GLY:HA2	2:P:2331:ASN:ND2	2.26	0.50
2:P:2271:GLY:HA2	2:P:2331:ASN:HD22	1.78	0.49
1:M:112:LYS:O	1:M:149:TRP:HH2	1.94	0.49
1:N:1090:THR:HA	1:N:1097:VAL:HG12	1.95	0.49
2:P:2321:ASP:O	2:P:2322:LYS:HB3	2.13	0.49
1:N:1142:GLN:HB2	1:N:1150:TYR:CE1	2.49	0.48
2:P:2179:CYS:HB2	2:P:2189:MET:CE	2.43	0.48
1:M:64:HIS:CE1	1:M:177:ARG:HH11	2.29	0.48
1:M:165:THR:HA	1:M:169:GLN:OE1	2.14	0.48
2:R:3292:LYS:NZ	2:R:3338:GLY:HA3	2.29	0.47
1:N:1172:VAL:O	1:N:1172:VAL:HG22	2.14	0.47
2:P:2154:PRO:HA	2:P:2182:GLY:O	2.15	0.47
2:P:2258:LEU:HB2	2:P:2353:HIS:CE1	2.51	0.46
1:N:1150:TYR:HB3	1:N:1172:VAL:CG2	2.43	0.46
2:P:2292:LYS:HB2	2:P:2312:LEU:HD11	1.97	0.46
1:N:1054:TYR:HA	1:N:1179:PRO:HG3	1.98	0.45
2:P:2154:PRO:HG2	2:P:2234:GLU:HA	1.98	0.45
2:R:3171:ALA:O	2:R:3172:ALA:HB3	2.15	0.45
1:M:125:LYS:HE2	1:M:144:ALA:O	2.17	0.45
1:M:74:ASN:HD22	1:M:76:MET:H	1.64	0.45
1:N:1165:THR:HA	1:N:1169:GLN:OE1	2.16	0.45
1:M:74:ASN:ND2	1:M:76:MET:H	2.14	0.45
1:M:90:THR:CG2	2:P:2285:GLN:HE21	2.28	0.44
1:N:1095:SER:HB2	2:R:3347:SER:HB3	1.99	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:N:1044:ASP:OD2	2:R:3287:HIS:HE1	2.00	0.44
1:N:1098:ARG:NE	1:N:1122:GLY:O	2.49	0.44
1:M:54:TYR:HA	1:M:179:PRO:HG3	2.00	0.44
1:M:50:LEU:HD11	2:P:2323:GLU:CG	2.46	0.44
2:R:3154:PRO:HA	2:R:3182:GLY:O	2.18	0.44
1:M:74:ASN:ND2	1:M:74:ASN:C	2.71	0.44
1:M:155:ARG:NH2	2:P:2160:GLU:O	2.51	0.43
1:M:90:THR:HG21	2:P:2286:PRO:O	2.18	0.43
1:N:1090:THR:HG22	2:R:3285:GLN:NE2	2.33	0.43
2:P:2179:CYS:HB2	2:P:2189:MET:HE3	2.00	0.43
2:R:3325:GLU:HB3	3:R:4104:HOH:O	2.18	0.43
2:P:2274:VAL:HG22	2:P:2275:GLU:N	2.34	0.43
2:P:2332:VAL:HG13	2:P:2336:ASP:HB2	2.01	0.42
1:N:1073:ILE:O	1:N:1159:PRO:HG3	2.19	0.42
2:R:3273:ASP:OD2	2:R:3330:ARG:HG2	2.19	0.42
2:P:2184:ASN:HA	2:P:2185:PRO:C	2.39	0.42
2:R:3231:CYS:SG	2:R:3231:CYS:O	2.77	0.41
1:N:1054:TYR:HB2	1:N:1177:ARG:O	2.20	0.41
1:N:1154:THR:HG23	1:N:1158:ARG:O	2.20	0.41
2:R:3154:PRO:HD3	2:R:3235:ASN:CG	2.40	0.41
2:R:3199:LYS:H	2:R:3202:HIS:HD2	1.59	0.41
2:R:3292:LYS:HB2	2:R:3312:LEU:HD11	2.03	0.41
1:N:1059:ARG:CZ	1:N:1170:ARG:NH1	2.84	0.41
1:N:1087:ILE:HG13	1:N:1102:ALA:HB2	2.02	0.41
2:R:3184:ASN:HA	2:R:3185:PRO:C	2.41	0.41
2:P:2274:VAL:HG12	2:P:2329:ILE:HB	2.03	0.41
1:M:152:ALA:HB2	1:M:172:VAL:HG12	2.02	0.41
1:N:1053:THR:HG21	1:N:1085:LYS:HE2	2.02	0.41
1:M:73:ILE:O	1:M:159:PRO:HG3	2.22	0.40

There are no symmetry-related clashes.

## 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	M	147/164 (90%)	142 (97%)	5 (3%)	0	100	100
1	N	146/164 (89%)	141 (97%)	5 (3%)	0	100	100
2	P	193/220 (88%)	185 (96%)	8 (4%)	0	100	100
2	R	185/220 (84%)	177 (96%)	8 (4%)	0	100	100
All	All	671/768 (87%)	645 (96%)	26 (4%)	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	M	116/143 (81%)	114 (98%)	2 (2%)	73	86
1	N	110/143 (77%)	108 (98%)	2 (2%)	71	85
2	P	154/187 (82%)	149 (97%)	5 (3%)	51	66
2	R	140/187 (75%)	136 (97%)	4 (3%)	55	69
All	All	520/660 (79%)	507 (98%)	13 (2%)	60	76

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

Mol	Chain	Res	Type
1	M	74	ASN
1	M	90	THR
2	P	2165	ARG
2	P	2173	ASN
2	P	2241	ASN
2	P	2254	HIS
2	P	2321	ASP
1	N	1074	ASN
1	N	1090	THR
2	R	3285	GLN
2	R	3330	ARG
2	R	3343	LEU
2	R	3356	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	M	55	GLN
1	M	64	HIS
1	M	74	ASN
2	P	2202	HIS
2	P	2241	ASN
2	P	2265	ASN
2	P	2285	GLN
2	P	2287	HIS
2	P	2331	ASN
1	N	1064	HIS
1	N	1074	ASN
2	R	3202	HIS
2	R	3245	HIS
2	R	3285	GLN
2	R	3287	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

### 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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, 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	M	149/164 (90%)	-0.14	0 100 100	7, 21, 33, 38	0
1	N	148/164 (90%)	0.06	5 (3%) 43 52	6, 23, 34, 40	0
2	P	197/220 (89%)	-0.09	3 (1%) 70 78	6, 20, 39, 47	0
2	R	191/220 (86%)	-0.14	3 (1%) 68 77	7, 21, 38, 47	0
All	All	685/768 (89%)	-0.08	11 (1%) 68 77	6, 22, 38, 47	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	1163	SER	4.6
1	N	1113	LYS	4.3
1	N	1162	GLY	3.6
2	P	2334	PHE	3.1
2	P	2308	TYR	3.0
2	P	2328	TYR	2.7
2	R	3316	GLY	2.6
2	R	3269	VAL	2.6
1	N	1069	ALA	2.4
2	R	3360	LEU	2.2
1	N	1112	LYS	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.