



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 03:03 PM BST

PDB ID : 6C6K  
Title : Structural basis for preferential recognition of cap 0 RNA by a human IFIT1-IFIT3 protein complex  
Authors : Amarasinghe, G.K.; Leung, D.W.; Johnson, B.; Xu, W.  
Deposited on : 2018-01-18  
Resolution : 2.54 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.

---

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 : 2.13.1  
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) : 2.13.1

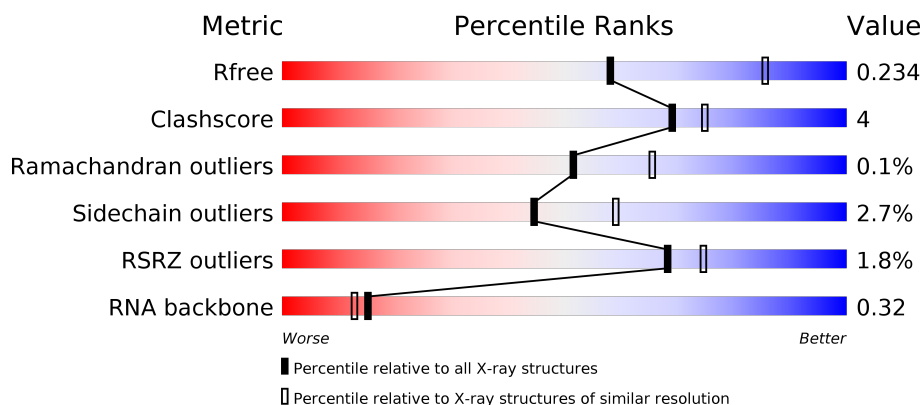
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.54 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)
RNA backbone	3102	1026 (2.88-2.20)

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	465	<div> <div>2%</div> <div>88%</div> <div>10%</div> <div>•</div> </div>
1	B	465	<div> <div>2%</div> <div>87%</div> <div>10%</div> <div>•</div> </div>
2	C	44	<div> <div>93%</div> <div>7%</div> </div>
2	D	44	<div> <div>2%</div> <div>93%</div> <div>7%</div> </div>

*Continued on next page...*

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	E	11	<div><div></div><div>9%</div><div>45%</div><div>27%</div><div>27%</div></div>
3	F	11	<div><div></div><div>45%</div><div>45%</div><div>9%</div></div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8993 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 Interferon-induced protein with tetratricopeptide repeats 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	456	Total	C	N	O	S	0	5	0
			3745	2386	648	690	21			
1	B	454	Total	C	N	O	S	0	2	0
			3717	2362	649	685	21			

- Molecule 2 is a protein called Interferon-induced protein with tetratricopeptide repeats 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	44	Total	C	N	O	S	0	0	0
			351	226	61	63	1			
2	D	44	Total	C	N	O	S	0	1	0
			357	230	62	64	1			

- Molecule 3 is a RNA chain called RNA (5'-R((M7G)\*P\*AP\*UP\*AP\*GP\*GP\*CP\*GP\*GP\*CP\*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	11	Total	C	N	O	P	0	0	0
			248	108	48	80	12			
3	F	11	Total	C	N	O	P	0	0	0
			248	108	48	80	12			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	E	3	Total	Mg	0	0
			3	3		

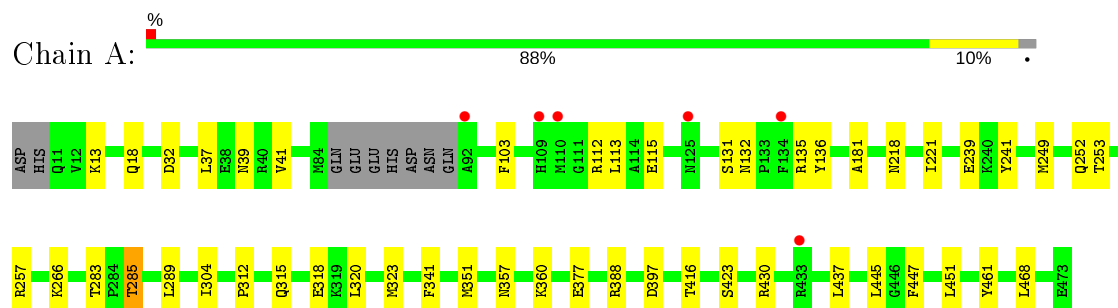
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	141	Total 141	O 141	0	0
5	B	139	Total 139	O 139	0	0
5	C	8	Total 8	O 8	0	0
5	D	6	Total 6	O 6	0	0
5	E	14	Total 14	O 14	0	0
5	F	15	Total 15	O 15	0	0

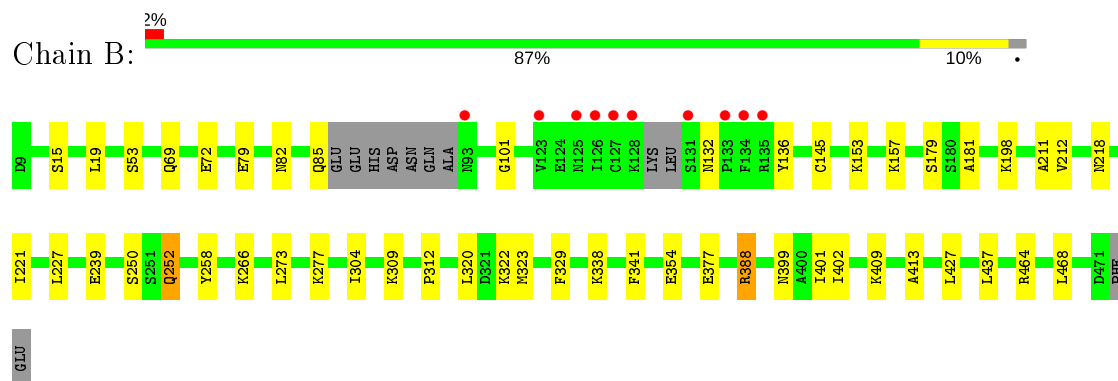
### 3 Residue-property plots [i](#)

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.

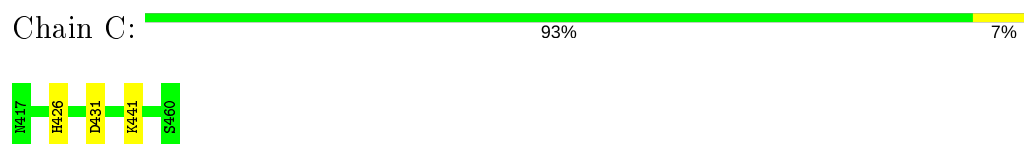
- Molecule 1: Interferon-induced protein with tetratricopeptide repeats 1



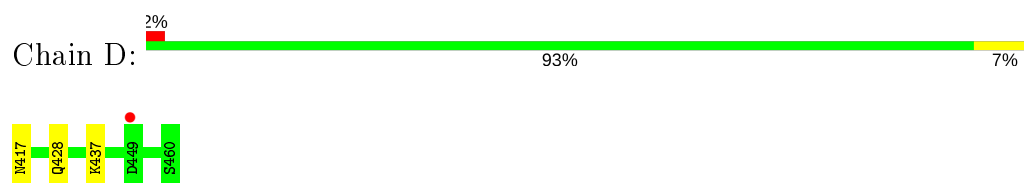
- Molecule 1: Interferon-induced protein with tetratricopeptide repeats 1



- Molecule 2: Interferon-induced protein with tetratricopeptide repeats 3



- Molecule 2: Interferon-induced protein with tetratricopeptide repeats 3



- Molecule 3: RNA (5'-R((M7G)\*P\*AP\*UP\*AP\*GP\*GP\*CP\*GP\*GP\*CP\*G)-3')



- Molecule 3: RNA (5'-R((M7G)\*P\*AP\*UP\*AP\*GP\*GP\*CP\*GP\*GP\*CP\*G)-3')



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.96 Å 80.48 Å 88.05 Å 79.82° 78.81° 90.04°	Depositor
Resolution (Å)	84.96 – 2.54 48.07 – 2.54	Depositor EDS
% Data completeness (in resolution range)	85.6 (84.96-2.54) 85.6 (48.07-2.54)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.50 (at 2.54 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.169 , 0.237 0.175 , 0.234	Depositor DCC
$R_{free}$ test set	1947 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.3	Xtriage
Anisotropy	0.111	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 51.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8993	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.84% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, M7G

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	0.70	0/3830	0.82	1/5145 (0.0%)
1	B	0.68	0/3792	0.81	1/5091 (0.0%)
2	C	0.62	0/358	0.75	0/481
2	D	0.59	0/367	0.72	0/493
3	E	0.51	0/245	0.90	0/381
3	F	0.54	0/245	0.94	0/381
All	All	0.68	0/8837	0.82	2/11972 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	351	MET	CG-SD-CE	-5.63	91.19	100.20
1	B	464	ARG	NE-CZ-NH1	5.47	123.03	120.30

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	3745	0	3781	35	1
1	B	3717	0	3736	27	1
2	C	351	0	351	1	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	357	0	359	2	0
3	E	248	0	126	6	0
3	F	248	0	126	4	0
4	B	1	0	0	0	0
4	E	3	0	0	0	0
5	A	141	0	0	4	0
5	B	139	0	0	3	0
5	C	8	0	0	0	0
5	D	6	0	0	1	0
5	E	14	0	0	0	0
5	F	15	0	0	0	0
All	All	8993	0	8479	64	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 4.

All (64) 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:252:GLN:NE2	5:B:601:HOH:O	1.99	0.95
1:A:218:ASN:HD21	3:E:1:M7G:HM71	1.29	0.95
1:A:252[B]:GLN:HE21	1:A:252[B]:GLN:HA	1.28	0.95
1:A:252[B]:GLN:HA	1:A:252[B]:GLN:NE2	1.90	0.85
1:A:218:ASN:HD21	3:E:1:M7G:CM7	1.96	0.77
1:A:252[B]:GLN:HE21	1:A:252[B]:GLN:CA	1.96	0.74
1:B:218:ASN:HD21	3:F:1:M7G:HM71	1.54	0.73
1:A:388:ARG:NH2	3:E:9:G:N7	2.38	0.71
1:A:241:TYR:OH	5:A:501:HOH:O	2.16	0.62
1:A:257:ARG:NH2	1:A:283[B]:THR:HG21	2.15	0.62
1:B:145:CYS:SG	1:B:179:SER:OG	2.57	0.62
1:B:181:ALA:HB1	1:B:221:ILE:HD11	1.83	0.61
1:A:357:ASN:ND2	1:A:360:LYS:HG3	2.17	0.60
1:A:447:PHE:CE2	1:A:451:LEU:HD11	2.36	0.60
1:A:304:ILE:HD12	1:A:323:MET:HE1	1.85	0.58
1:A:218:ASN:ND2	3:E:1:M7G:HM71	2.11	0.57
1:A:397:ASP:OD2	1:A:430:ARG:NH2	2.37	0.56
1:B:212:VAL:HG23	1:B:221:ILE:HG22	1.87	0.56
1:A:257:ARG:HH21	1:A:283[B]:THR:HG21	1.71	0.55
1:A:304:ILE:HD12	1:A:323:MET:CE	2.37	0.55
1:A:257:ARG:HH21	1:A:283[A]:THR:HG21	1.72	0.54
1:B:69:GLN:HE21	1:B:72:GLU:HG3	1.73	0.53

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:LEU:HD13	1:B:101:GLY:HA2	1.89	0.53
1:B:239:GLU:OE1	1:B:266:LYS:HE3	2.09	0.53
1:A:181:ALA:HB1	1:A:221:ILE:HD11	1.91	0.52
1:A:252[A]:GLN:HE21	1:B:329:PHE:HE1	1.57	0.52
1:B:198:LYS:HD2	1:B:413:ALA:HB1	1.90	0.52
1:A:239:GLU:OE1	1:A:266:LYS:HE3	2.11	0.51
1:B:399:ASN:HA	1:B:402:ILE:HD12	1.93	0.50
1:B:211:ALA:HB1	1:B:221:ILE:HD13	1.94	0.50
1:A:283[A]:THR:HG21	5:A:563:HOH:O	2.12	0.50
1:B:401:ILE:HG13	1:B:427:LEU:HD23	1.93	0.50
1:B:212:VAL:CG2	1:B:221:ILE:HG22	2.43	0.49
1:B:377:GLU:HB2	5:B:681:HOH:O	2.12	0.49
1:A:132:ASN:HD22	1:A:132:ASN:N	2.10	0.49
1:A:39:ASN:HB3	1:A:285:THR:CG2	2.43	0.48
1:B:338:LYS:NZ	3:F:3:U:O4	2.47	0.48
1:B:82:ASN:ND2	5:B:612:HOH:O	2.45	0.47
1:A:112:ARG:NE	1:A:115:GLU:OE1	2.44	0.47
1:B:227:LEU:HD11	1:B:258:TYR:HB3	1.96	0.47
1:B:15:SER:HB3	1:B:136:TYR:CE2	2.51	0.47
1:A:388:ARG:HD3	5:A:548:HOH:O	2.15	0.46
1:A:285:THR:HB	5:A:565:HOH:O	2.15	0.46
1:A:18:GLN:HB2	1:A:136:TYR:HB2	1.98	0.45
1:A:257:ARG:HG2	1:A:289:LEU:HD21	2.00	0.44
2:D:417:ASN:N	5:D:502:HOH:O	2.50	0.44
1:A:249[A]:MET:SD	1:B:329:PHE:HB2	2.58	0.44
3:E:9:G:H1'	3:E:11:G:O4'	2.19	0.43
1:B:388:ARG:NH2	3:F:9:G:N7	2.66	0.43
1:A:312:PRO:HG3	1:A:320:LEU:HD22	2.00	0.42
2:C:426:HIS:HB3	2:C:431:ASP:HB2	2.01	0.42
1:B:312:PRO:HG2	1:B:320:LEU:HD22	2.01	0.42
1:B:79:GLU:HG2	2:D:428[B]:GLN:HG2	2.01	0.42
3:E:10:C:H2'	3:E:10:C:O2	2.19	0.42
1:A:37:LEU:O	1:A:41:VAL:HG23	2.20	0.41
1:A:377:GLU:OE1	1:A:416:THR:HG21	2.20	0.41
1:A:445:LEU:HD23	1:A:461:TYR:CE1	2.55	0.41
1:A:437:LEU:HD22	1:A:468:LEU:HD13	2.02	0.41
1:B:153:LYS:NZ	3:F:1:M7G:H5'2	2.35	0.41
1:B:304:ILE:HD12	1:B:323:MET:HE1	2.03	0.41
1:B:273:LEU:O	1:B:277:LYS:HG2	2.21	0.40
1:B:437:LEU:HD22	1:B:468:LEU:HD13	2.02	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:GLU:OE2	1:B:53:SER:OG[1_455]	2.18	0.02

## 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	457/465 (98%)	439 (96%)	18 (4%)	0	100	100
1	B	450/465 (97%)	438 (97%)	11 (2%)	1 (0%)	47	60
2	C	42/44 (96%)	42 (100%)	0	0	100	100
2	D	43/44 (98%)	42 (98%)	1 (2%)	0	100	100
All	All	992/1018 (97%)	961 (97%)	30 (3%)	1 (0%)	51	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	250	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	402/406 (99%)	391 (97%)	11 (3%)	44	59
1	B	398/406 (98%)	388 (98%)	10 (2%)	47	62
2	C	36/36 (100%)	35 (97%)	1 (3%)	43	58

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	37/36 (103%)	36 (97%)	1 (3%)	44	59
All	All	873/884 (99%)	850 (97%)	23 (3%)	44	61

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LYS
1	A	32	ASP
1	A	103	PHE
1	A	113	LEU
1	A	131	SER
1	A	135	ARG
1	A	253	THR
1	A	285	THR
1	A	315	GLN
1	A	341	PHE
1	A	423	SER
1	B	85	GLN
1	B	132	ASN
1	B	157	LYS
1	B	252	GLN
1	B	309	LYS
1	B	322	LYS
1	B	341	PHE
1	B	354	GLU
1	B	388	ARG
1	B	409	LYS
2	C	441	LYS
2	D	437	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	GLN
1	A	125	ASN
1	A	132	ASN
1	A	315	GLN
1	B	385	HIS

### 5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	E	9/11 (81%)	5 (55%)	1 (11%)
3	F	9/11 (81%)	4 (44%)	0
All	All	18/22 (81%)	9 (50%)	1 (5%)

All (9) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	E	4	A
3	E	8	G
3	E	9	G
3	E	10	C
3	E	11	G
3	F	4	A
3	F	6	G
3	F	8	G
3	F	9	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
3	E	10	C

## 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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	A	456/465 (98%)	-0.15	6 (1%) 77 82	14, 36, 76, 127	0
1	B	454/465 (97%)	-0.23	10 (2%) 62 68	16, 36, 64, 90	0
2	C	44/44 (100%)	0.15	0 100 100	34, 46, 75, 84	0
2	D	44/44 (100%)	0.06	1 (2%) 60 67	31, 45, 65, 81	0
3	E	10/11 (90%)	0.08	1 (10%) 7 9	22, 68, 100, 100	0
3	F	10/11 (90%)	-0.46	0 100 100	23, 69, 95, 99	0
All	All	1018/1040 (97%)	-0.16	18 (1%) 68 74	14, 37, 75, 127	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	134	PHE	8.1
1	A	92	ALA	4.5
1	B	128	LYS	3.5
1	A	110	MET	3.1
1	B	126	ILE	3.0
1	B	125	ASN	2.6
1	B	133	PRO	2.6
1	B	134	PHE	2.4
1	B	135	ARG	2.4
1	B	131	SER	2.4
1	B	127	CYS	2.2
1	B	93	ASN	2.1
3	E	10	C	2.1
1	A	125	ASN	2.1
1	A	433	ARG	2.1
2	D	449	ASP	2.1
1	A	109	HIS	2.1
1	B	123	VAL	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 monosaccharides 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
4	MG	E	102	1/1	0.94	0.21	53,53,53,53	0
4	MG	E	103	1/1	0.94	0.34	65,65,65,65	0
4	MG	E	101	1/1	0.97	0.19	55,55,55,55	0
4	MG	B	501	1/1	0.99	0.29	49,49,49,49	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.