



# Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 05:27 am BST

PDB ID : 4LEV  
Title : Structure of human cGAS  
Authors : Li, P.  
Deposited on : 2013-06-26  
Resolution : 1.95 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

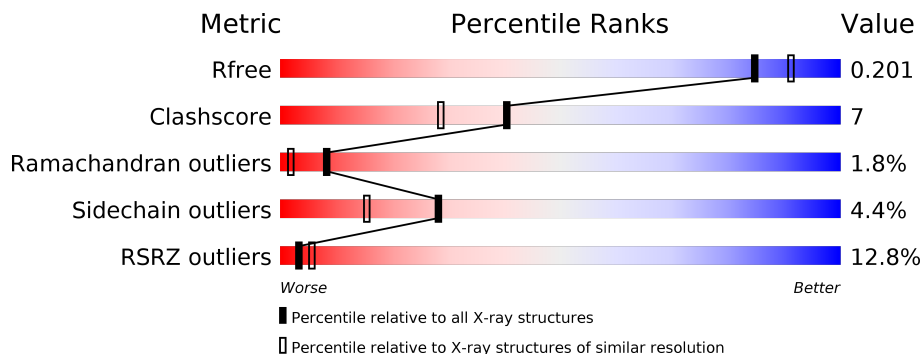
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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	369	
1	B	369	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6567 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 Cyclic GMP-AMP synthase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	369	3057	1955	525	561	10	6	0	3	0
1	B	369	3057	1954	525	561	10	7	0	3	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	154	PHE	-	EXPRESSION TAG	UNP Q8N884
A	155	GLU	-	EXPRESSION TAG	UNP Q8N884
A	156	LEU	-	EXPRESSION TAG	UNP Q8N884
B	154	PHE	-	EXPRESSION TAG	UNP Q8N884
B	155	GLU	-	EXPRESSION TAG	UNP Q8N884
B	156	LEU	-	EXPRESSION TAG	UNP Q8N884

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

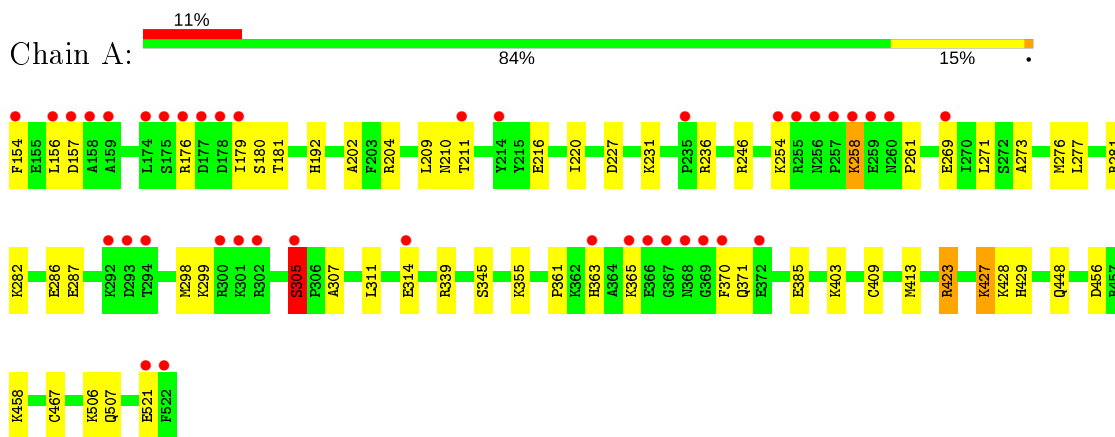
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	241	Total	O	0	0
			241	241		
3	B	210	Total	O	0	0
			210	210		

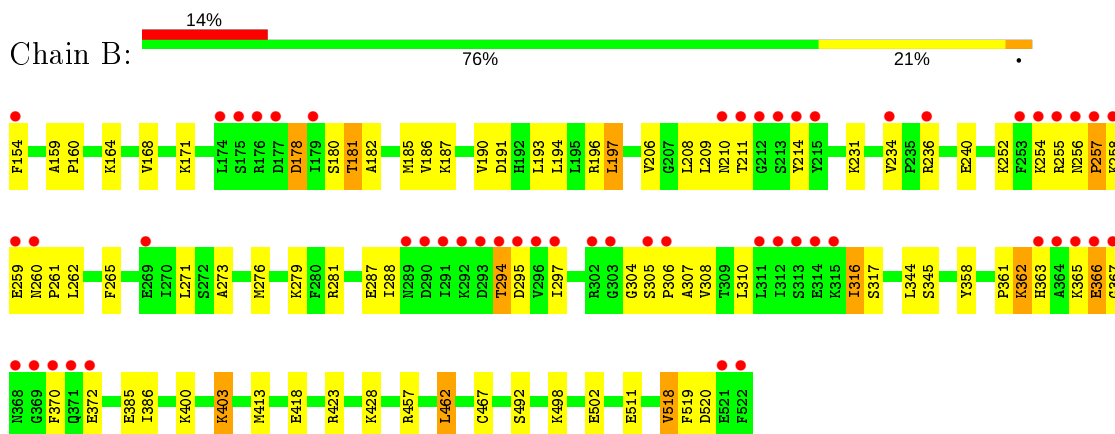
### 3 Residue-property plots [i](#)

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: Cyclic GMP-AMP synthase



- Molecule 1: Cyclic GMP-AMP synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	212.98Å 47.72Å 86.87Å 90.00° 113.89° 90.00°	Depositor
Resolution (Å)	38.45 – 1.95 38.45 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.8 (38.45-1.95) 93.5 (38.45-1.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 1.95Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.167 , 0.202 0.167 , 0.201	Depositor DCC
$R_{free}$ test set	1993 reflections (3.48%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.1	Xtrriage
Anisotropy	0.373	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 53.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.018 for -h-2*1,-k,l	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6567	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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:  
ZN

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	0.52	0/3114	0.63	1/4169 (0.0%)
1	B	0.50	0/3114	0.62	0/4168
All	All	0.51	0/6228	0.63	1/8337 (0.0%)

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	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	305	SER	C-N-CA	-5.47	99.04	122.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	305	SER	Peptide

### 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	3057	0	3088	36	2
1	B	3057	0	3086	52	1
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	241	0	0	9	0
3	B	210	0	0	5	0
All	All	6567	0	6174	88	2

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 (88) 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:164:LYS:NZ	3:B:836:HOH:O	2.08	0.85
1:A:458:LYS:NZ	3:A:875:HOH:O	2.16	0.79
1:B:418:GLU:OE2	3:B:811:HOH:O	2.03	0.75
1:A:236:ARG:HB2	1:A:254:LYS:HD2	1.69	0.73
1:A:365:LYS:HB3	1:A:370:PHE:HB3	1.72	0.71
1:A:271:LEU:HD21	1:A:276:MSE:HE1	1.73	0.70
1:B:236:ARG:HB2	1:B:254:LYS:HG3	1.74	0.70
1:A:204:ARG:HH21	1:A:261:PRO:HB3	1.60	0.67
1:A:286:GLU:OE1	3:A:869:HOH:O	2.12	0.67
1:B:294:THR:OG1	1:B:295:ASP:N	2.22	0.67
1:A:157:ASP:OD2	3:A:837:HOH:O	2.13	0.67
1:B:193:LEU:HD11	1:B:310:LEU:HD11	1.78	0.66
1:A:507:GLN:NE2	3:A:818:HOH:O	2.21	0.66
1:B:191:ASP:OD1	1:B:214:TYR:OH	2.14	0.65
1:B:255:ARG:HD2	1:B:259:GLU:OE2	1.98	0.64
1:A:273:ALA:HB1	1:A:361:PRO:HB3	1.83	0.60
1:B:365:LYS:HB3	1:B:370:PHE:HB3	1.83	0.60
1:B:164:LYS:HE3	1:B:520:ASP:HB3	1.83	0.60
1:A:448:GLN:NE2	3:A:913:HOH:O	2.32	0.60
1:B:305:SER:OG	1:B:361:PRO:O	2.19	0.60
1:A:209:LEU:O	1:A:211:THR:N	2.37	0.58
1:B:168:VAL:HG21	1:B:519:PHE:HB3	1.85	0.58
1:A:423:ARG:NH1	3:A:818:HOH:O	2.37	0.57
1:B:194:LEU:HD11	1:B:208:LEU:HG	1.88	0.55
1:B:186:VAL:HG23	1:B:316:ILE:HD11	1.90	0.54
1:A:202:ALA:HA	1:A:261:PRO:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:ILE:HD11	1:B:310:LEU:HD22	1.89	0.54
1:B:240:GLU:OE1	1:B:252:LYS:NZ	2.25	0.53
1:A:427:LYS:O	1:A:429:HIS:N	2.42	0.53
1:A:339:ARG:NH1	3:A:825:HOH:O	2.41	0.52
1:B:273:ALA:HB1	1:B:361:PRO:HB3	1.90	0.52
1:A:281:ARG:HH12	1:A:298:MSE:HE3	1.74	0.52
1:B:257:PRO:O	1:B:259:GLU:N	2.43	0.51
1:A:176:ARG:HH11	1:A:176:ARG:HG3	1.75	0.51
1:A:192:HIS:NE2	1:A:287:GLU:OE2	2.43	0.50
1:B:428:LYS:HA	3:B:839:HOH:O	2.12	0.50
1:A:299:LYS:O	3:A:762:HOH:O	2.20	0.49
1:B:187:LYS:HA	1:B:214:TYR:CE1	2.48	0.49
1:B:363:HIS:CE1	1:B:372:GLU:HG3	2.47	0.49
1:B:178:ASP:O	1:B:181:THR:HG23	2.13	0.48
1:B:413[A]:MSE:SE	1:B:467:CYS:HB3	2.63	0.48
1:A:409:CYS:O	1:A:413[A]:MSE:HG3	2.14	0.48
1:A:176:ARG:HH22	1:A:220:ILE:HA	1.79	0.48
1:B:265:PHE:HZ	1:B:279:LYS:HD3	1.79	0.47
1:B:498:LYS:HE3	1:B:502:GLU:OE2	2.13	0.47
1:B:281:ARG:HG3	1:B:308:VAL:HG22	1.97	0.47
1:A:236:ARG:HH11	1:A:236:ARG:HG2	1.79	0.47
1:B:385:GLU:HG2	3:B:767:HOH:O	2.15	0.47
1:A:281:ARG:NH1	1:A:298:MSE:HE3	2.31	0.46
1:B:403:LYS:HG2	1:B:457:ARG:HD2	1.97	0.46
1:A:423:ARG:NH2	1:A:521:GLU:OE1	2.50	0.45
1:B:260:ASN:C	1:B:262:LEU:H	2.19	0.45
1:A:236:ARG:HG2	1:A:236:ARG:NH1	2.32	0.45
1:B:305:SER:HB2	1:B:362:LYS:NZ	2.30	0.45
1:A:413[A]:MSE:SE	1:A:467:CYS:HB3	2.66	0.45
1:A:176:ARG:NH1	1:A:176:ARG:HG3	2.32	0.44
1:B:181:THR:OG1	1:B:182:ALA:N	2.50	0.44
1:B:260:ASN:O	1:B:262:LEU:N	2.50	0.44
1:A:282:LYS:O	1:A:286:GLU:HG3	2.18	0.44
1:A:258:LYS:H	1:A:258:LYS:HG3	1.42	0.43
1:B:181:THR:O	1:B:185[A]:MSE:HG3	2.17	0.43
1:B:209:LEU:HG	1:B:210:ASN:N	2.34	0.43
1:A:307:ALA:HB2	3:A:849:HOH:O	2.18	0.43
1:B:196:ARG:HE	1:B:287:GLU:CD	2.22	0.43
1:B:186:VAL:CG2	1:B:316:ILE:HD11	2.49	0.43
1:B:197:LEU:HA	1:B:197:LEU:HD12	1.85	0.43
1:A:355:LYS:HE3	1:A:385:GLU:OE2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:LEU:CD1	1:B:386:ILE:HD12	2.49	0.42
1:A:176:ARG:NH2	1:A:220:ILE:HD12	2.34	0.42
1:B:423:ARG:NH1	3:B:794:HOH:O	2.52	0.42
1:B:305:SER:OG	1:B:362:LYS:HG2	2.20	0.42
1:B:209:LEU:HD13	1:B:358:TYR:CZ	2.53	0.42
1:A:277:LEU:O	1:A:281:ARG:HG3	2.18	0.42
1:B:366:GLU:HB3	1:B:367:GLY:H	1.67	0.42
1:B:511:GLU:OE1	1:B:518:VAL:HB	2.20	0.42
1:A:403:LYS:HD2	1:A:403:LYS:HA	1.80	0.41
1:B:159:ALA:HA	1:B:160:PRO:HD3	1.90	0.41
1:B:271:LEU:HD21	1:B:276:MSE:HE1	2.01	0.41
1:B:462:LEU:HA	1:B:462:LEU:HD12	1.75	0.41
1:B:197:LEU:HB3	1:B:206:VAL:HG11	2.02	0.41
1:B:256:ASN:CB	1:B:257:PRO:HD2	2.51	0.41
1:A:179:ILE:O	1:A:181:THR:N	2.53	0.40
1:B:305:SER:HB2	1:B:362:LYS:HZ3	1.85	0.40
1:A:156:LEU:HA	1:A:156:LEU:HD23	1.94	0.40
1:B:190:VAL:HB	1:B:214:TYR:HE1	1.85	0.40
1:B:210:ASN:HD22	1:B:210:ASN:HA	1.70	0.40
1:B:344:LEU:HD11	1:B:386:ILE:HG23	2.03	0.40
1:B:168:VAL:CG2	1:B:519:PHE:HB3	2.51	0.40

All (2) 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:314:GLU:O	1:A:506:LYS:NZ[4_655]	2.16	0.04
1:A:456:ASP:OD2	1:B:400:LYS:NZ[4_656]	2.17	0.03

## 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	370/369 (100%)	348 (94%)	17 (5%)	5 (1%)	11 3
1	B	370/369 (100%)	342 (92%)	20 (5%)	8 (2%)	6 1
All	All	740/738 (100%)	690 (93%)	37 (5%)	13 (2%)	8 2

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	428	LYS
1	B	257	PRO
1	B	258	LYS
1	B	307	ALA
1	A	180	SER
1	A	210	ASN
1	A	216	GLU
1	A	345	SER
1	B	345	SER
1	B	366	GLU
1	B	304	GLY
1	B	306	PRO
1	B	261	PRO

### 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	342/334 (102%)	330 (96%)	12 (4%)	36 24
1	B	342/334 (102%)	324 (95%)	18 (5%)	22 10
All	All	684/668 (102%)	654 (96%)	30 (4%)	28 15

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	154	PHE
1	A	227	ASP
1	A	231	LYS

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	246	ARG
1	A	258	LYS
1	A	269	GLU
1	A	305	SER
1	A	311	LEU
1	A	363	HIS
1	A	371	GLN
1	A	423	ARG
1	A	427	LYS
1	B	154	PHE
1	B	171	LYS
1	B	178	ASP
1	B	180	SER
1	B	181	THR
1	B	197	LEU
1	B	211	THR
1	B	231	LYS
1	B	234	VAL
1	B	294	THR
1	B	297	ILE
1	B	316	ILE
1	B	317	SER
1	B	362	LYS
1	B	403	LYS
1	B	462	LEU
1	B	492	SER
1	B	518	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	210	ASN
1	A	224	ASN
1	B	210	ASN

### 5.3.3 RNA

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 [i](#)

Of 2 ligands modelled in this entry, 2 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 [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, 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	364/369 (98%)	0.36	40 (10%) 5   9	18, 40, 118, 165	0
1	B	364/369 (98%)	0.61	53 (14%) 2   3	18, 46, 126, 174	0
All	All	728/738 (98%)	0.48	93 (12%) 3   6	18, 42, 123, 174	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	367	GLY	17.2
1	B	368	ASN	12.1
1	B	370	PHE	12.0
1	A	368	ASN	10.7
1	A	367	GLY	10.2
1	A	257	PRO	9.2
1	A	258	LYS	9.0
1	A	366	GLU	7.6
1	B	366	GLU	7.4
1	A	256	ASN	7.3
1	B	369	GLY	6.9
1	B	211	THR	6.9
1	B	522	PHE	6.8
1	A	154	PHE	6.6
1	A	369	GLY	6.6
1	A	156	LEU	6.5
1	B	292	LYS	6.4
1	B	214	TYR	6.3
1	A	370	PHE	6.2
1	B	294	THR	5.6
1	B	305	SER	5.5
1	B	303	GLY	5.5
1	B	256	ASN	5.4
1	A	255	ARG	5.3

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	214	TYR	5.2
1	B	255	ARG	5.2
1	B	254	LYS	5.2
1	B	297	ILE	5.1
1	A	211	THR	5.0
1	B	296	VAL	4.9
1	B	258	LYS	4.9
1	B	176	ARG	4.9
1	B	175	SER	4.8
1	B	215	TYR	4.7
1	B	291	ILE	4.7
1	B	212	GLY	4.5
1	B	260	ASN	4.5
1	A	176	ARG	4.5
1	B	315	LYS	4.4
1	B	521	GLU	4.4
1	B	179	ILE	4.3
1	A	305	SER	4.2
1	A	260	ASN	4.2
1	B	314	GLU	4.2
1	B	253	PHE	4.2
1	B	365	LYS	4.0
1	A	158	ALA	3.9
1	A	365	LYS	3.9
1	A	522	PHE	3.8
1	B	306	PRO	3.8
1	B	364	ALA	3.8
1	A	314	GLU	3.8
1	A	157	ASP	3.8
1	B	293	ASP	3.7
1	A	178	ASP	3.6
1	A	179	ILE	3.6
1	B	174	LEU	3.5
1	B	313	SER	3.5
1	A	292	LYS	3.4
1	A	259	GLU	3.4
1	A	174	LEU	3.4
1	B	154	PHE	3.3
1	A	254	LYS	3.2
1	B	289	ASN	3.2
1	B	213	SER	3.2
1	A	235	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	257	PRO	3.1
1	B	259	GLU	3.1
1	A	269	GLU	2.9
1	B	311	LEU	2.9
1	B	312	ILE	2.9
1	B	372	GLU	2.8
1	A	293	ASP	2.8
1	B	269	GLU	2.7
1	B	302	ARG	2.7
1	A	302	ARG	2.6
1	B	295	ASP	2.6
1	A	372	GLU	2.6
1	B	363	HIS	2.5
1	B	210	ASN	2.5
1	A	300	ARG	2.5
1	A	159	ALA	2.4
1	B	236	ARG	2.4
1	A	363	HIS	2.4
1	A	521	GLU	2.3
1	A	294	THR	2.3
1	A	177	ASP	2.3
1	A	301	LYS	2.2
1	B	371	GLN	2.2
1	B	234	VAL	2.2
1	B	290	ASP	2.2
1	B	177	ASP	2.0
1	A	175	SER	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 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(Å <sup>2</sup> )	Q<0.9
2	ZN	B	601	1/1	0.99	0.07	20,20,20,20	0
2	ZN	A	601	1/1	0.99	0.09	23,23,23,23	0

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.