



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:21 am BST

PDB ID : 6O6Z  
Title : Crystal structure of Csm6 H381A in complex with cA4 by cocrystallization of cA4 and Csm6  
Authors : Jia, N.; Patel, D.J.  
Deposited on : 2019-03-07  
Resolution : 2.10 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  
buster-report : 1.1.7 (2018)  
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

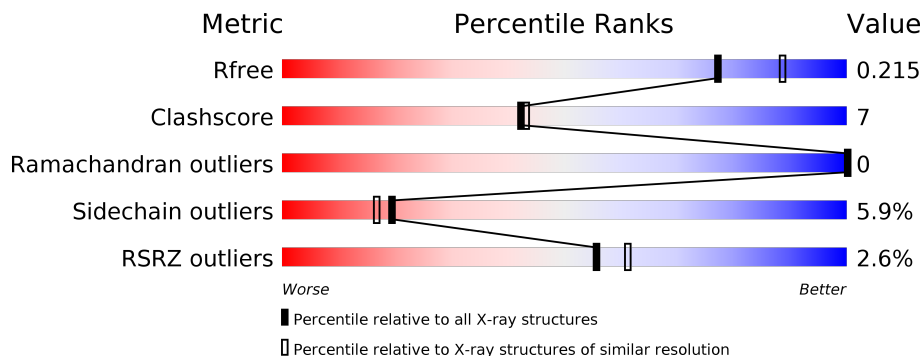
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	440	 2% 81% 16% ..
1	B	440	 3% 82% 14% ..

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 7098 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.

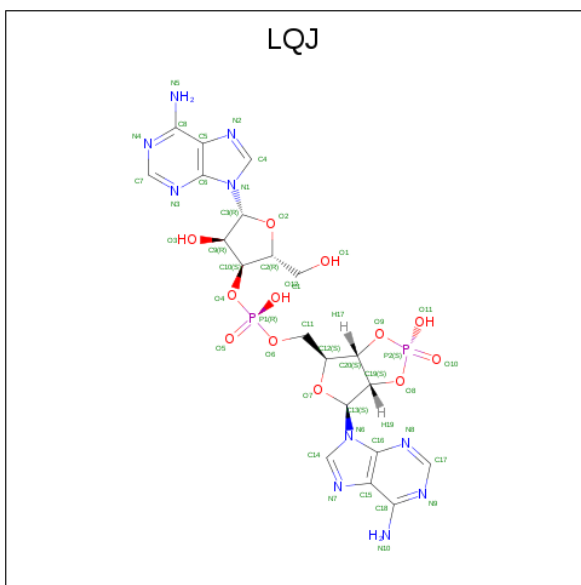
- Molecule 1 is a protein called Csm6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	431	Total	C	N	O	S	0	0	0
			3435	2211	582	634	8			
1	B	430	Total	C	N	O	S	0	0	0
			3425	2205	579	633	8			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP B6YWC3
A	0	GLY	-	expression tag	UNP B6YWC3
A	381	ALA	HIS	engineered mutation	UNP B6YWC3
A	433	HIS	-	expression tag	UNP B6YWC3
A	434	HIS	-	expression tag	UNP B6YWC3
A	435	HIS	-	expression tag	UNP B6YWC3
A	436	HIS	-	expression tag	UNP B6YWC3
A	437	HIS	-	expression tag	UNP B6YWC3
A	438	HIS	-	expression tag	UNP B6YWC3
B	-1	MET	-	initiating methionine	UNP B6YWC3
B	0	GLY	-	expression tag	UNP B6YWC3
B	381	ALA	HIS	engineered mutation	UNP B6YWC3
B	433	HIS	-	expression tag	UNP B6YWC3
B	434	HIS	-	expression tag	UNP B6YWC3
B	435	HIS	-	expression tag	UNP B6YWC3
B	436	HIS	-	expression tag	UNP B6YWC3
B	437	HIS	-	expression tag	UNP B6YWC3
B	438	HIS	-	expression tag	UNP B6YWC3

- Molecule 2 is 3'-O-[(R)-{[(2S,3aS,4S,6S,6aS)-6-(6-amino-9H-purin-9-yl)-2-hydroxy-2-oxotetrahydro-2H-2lambda 5-furo[3,4-d][1,3,2]dioxaphosphol-4-yl]methoxy}(hydroxy)phosphoryl]adenosine (three-letter code: LQJ) (formula: C<sub>20</sub>H<sub>24</sub>N<sub>10</sub>O<sub>12</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	A	1	Total	C	N	O	P	0	0
			44	20	10	12	2		
2	B	1	Total	C	N	O	P	0	0
			44	20	10	12	2		

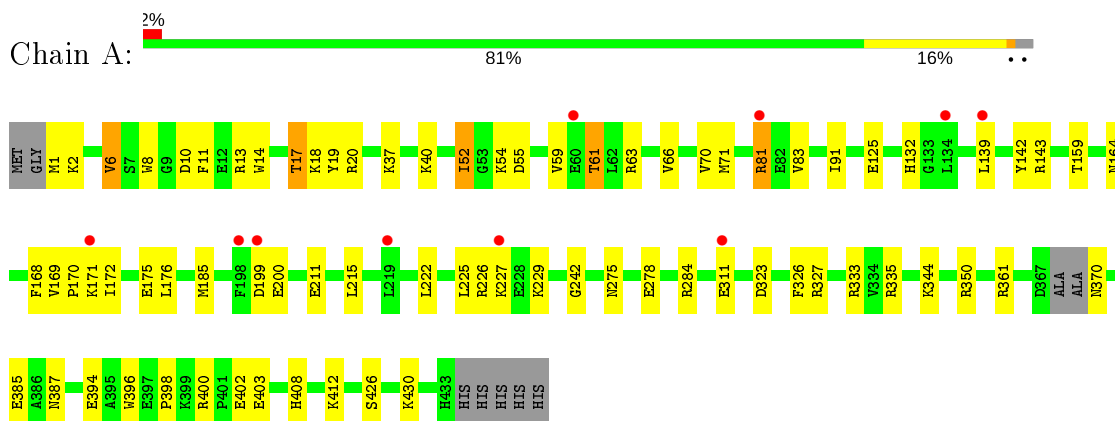
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	79	Total	O	0	0
			79	79		
3	B	71	Total	O	0	0
			71	71		

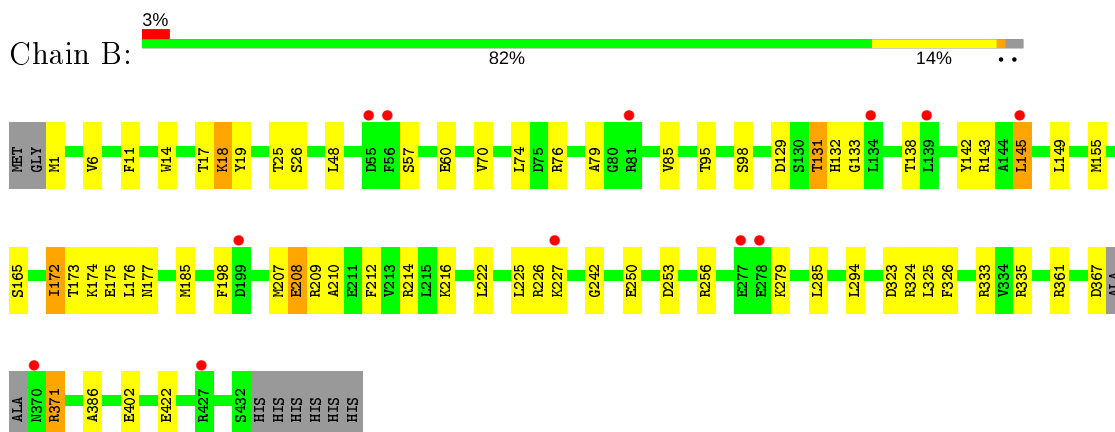
### 3 Residue-property plots [i](#)

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: Csm6



- Molecule 1: Csm6



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	54.29Å 115.94Å 159.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.35 – 2.10 48.30 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.2 (48.35-2.10) 97.9 (48.30-2.08)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.40 (at 2.08Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, $R_{free}$	0.192 , 0.222 0.198 , 0.215	Depositor DCC
$R_{free}$ test set	3064 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.6	Xtrriage
Anisotropy	0.422	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 36.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7098	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.66% 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: LQJ

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	0.74	0/3509	0.88	0/4747
1	B	0.71	0/3498	0.89	0/4732
All	All	0.73	0/7007	0.88	0/9479

There are no bond length outliers.

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	3435	0	3486	48	0
1	B	3425	0	3479	56	0
2	A	44	0	0	2	0
2	B	44	0	0	0	0
3	A	79	0	0	0	0
3	B	71	0	0	0	0
All	All	7098	0	6965	96	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 (96) 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:129:ASP:OD2	1:B:131:THR:HG22	1.75	0.85
1:B:129:ASP:OD2	1:B:131:THR:CG2	2.27	0.83
1:A:400:ARG:NH2	1:A:403:GLU:HG3	1.96	0.80
1:B:371:ARG:HH21	1:B:371:ARG:HG3	1.48	0.77
1:B:208:GLU:OE1	1:B:209:ARG:N	2.19	0.76
1:A:142:TYR:OH	1:B:143:ARG:HD3	1.93	0.68
1:B:253:ASP:OD2	1:B:256:ARG:NH1	2.26	0.68
1:A:172:ILE:HG12	1:B:95:THR:HG22	1.75	0.67
1:B:133:GLY:HA3	1:B:138:THR:OG1	1.95	0.67
1:A:199:ASP:OD1	1:A:200:GLU:CG	2.44	0.66
1:A:14:TRP:CZ2	1:A:132:HIS:HB3	2.32	0.65
1:B:210:ALA:O	1:B:214:ARG:HG3	1.99	0.63
1:A:199:ASP:OD1	1:A:200:GLU:HG3	2.00	0.62
1:B:207:MET:HE3	1:B:212:PHE:HB2	1.81	0.61
1:A:143:ARG:HD2	1:B:142:TYR:OH	2.03	0.58
1:B:129:ASP:OD2	1:B:131:THR:HG23	2.03	0.58
1:A:199:ASP:OD1	1:A:200:GLU:HG2	2.04	0.57
1:A:350:ARG:NH2	1:A:402:GLU:OE2	2.38	0.56
1:B:207:MET:HE3	1:B:212:PHE:CA	2.35	0.56
1:B:207:MET:HE3	1:B:212:PHE:CB	2.35	0.56
1:B:76:ARG:HG2	1:B:76:ARG:HH21	1.71	0.56
1:B:207:MET:CE	1:B:212:PHE:HA	2.36	0.56
1:B:14:TRP:CZ2	1:B:132:HIS:HB3	2.41	0.56
1:B:207:MET:CE	1:B:212:PHE:HD1	2.18	0.56
1:B:129:ASP:OD1	1:B:131:THR:HG23	2.06	0.55
1:B:172:ILE:O	1:B:172:ILE:HD12	2.06	0.55
1:A:396:TRP:O	1:A:398:PRO:HD3	2.05	0.55
1:A:323:ASP:OD2	1:A:335:ARG:NH1	2.40	0.54
1:B:207:MET:HE3	1:B:212:PHE:HA	1.89	0.54
1:A:394:GLU:OE2	1:A:408:HIS:HD2	1.90	0.53
1:B:11:PHE:HA	1:B:14:TRP:CE2	2.44	0.53
1:A:19:TYR:HA	1:A:176:LEU:O	2.09	0.53
1:B:129:ASP:CG	1:B:131:THR:HG23	2.29	0.53
1:B:74:LEU:HD22	1:B:79:ALA:CB	2.38	0.52
1:B:371:ARG:NH2	1:B:371:ARG:HG3	2.19	0.52
1:B:242:GLY:HA2	1:B:326:PHE:CE1	2.44	0.51
1:A:361:ARG:HH21	1:A:387:ASN:HD22	1.58	0.51
1:A:52:ILE:HG22	1:A:66:VAL:HG22	1.93	0.51
1:A:169:VAL:HG12	1:A:172:ILE:HD13	1.94	0.50
1:A:6:VAL:HG13	1:A:8:TRP:CD1	2.47	0.50
1:B:18:LYS:HA	1:B:26:SER:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:ARG:HG2	1:B:333:ARG:HD3	1.94	0.50
1:A:402:GLU:CD	1:A:402:GLU:H	2.15	0.50
1:B:18:LYS:HG3	1:B:175:GLU:HG3	1.93	0.50
1:A:225:LEU:O	1:A:229:LYS:HB2	2.13	0.49
1:A:172:ILE:CD1	1:B:95:THR:HG22	2.43	0.48
1:A:211:GLU:OE1	1:A:284:ARG:NH2	2.45	0.48
1:B:198:PHE:CZ	1:B:226:ARG:HA	2.48	0.48
1:A:172:ILE:CG1	1:B:95:THR:HG22	2.43	0.48
1:A:40:LYS:HE2	1:A:40:LYS:HA	1.94	0.48
1:A:11:PHE:HA	1:A:14:TRP:CE2	2.49	0.48
1:B:207:MET:HE1	1:B:212:PHE:HD1	1.79	0.48
1:B:361:ARG:HB3	1:B:386:ALA:HB1	1.95	0.47
1:A:275:ASN:OD1	1:A:278:GLU:OE1	2.33	0.47
1:A:17:THR:HG23	1:A:168:PHE:HE1	1.80	0.47
1:B:173:THR:HG21	1:B:176:LEU:HD21	1.97	0.46
1:A:333:ARG:CB	1:B:333:ARG:HD3	2.46	0.46
1:A:242:GLY:HA2	1:A:326:PHE:CZ	2.51	0.46
1:B:145:LEU:O	1:B:149:LEU:HG	2.17	0.45
1:A:11:PHE:HA	1:A:14:TRP:CD2	2.51	0.45
1:A:242:GLY:HA2	1:A:326:PHE:CE1	2.52	0.45
1:A:311:GLU:O	1:A:311:GLU:HG2	2.17	0.44
1:A:333:ARG:HB2	1:B:333:ARG:HD3	2.00	0.44
1:B:76:ARG:NH2	1:B:76:ARG:HG2	2.30	0.44
1:A:10:ASP:OD1	1:A:13:ARG:NH2	2.51	0.44
1:B:6:VAL:HG12	1:B:129:ASP:HB3	1.99	0.44
1:B:18:LYS:HD2	1:B:175:GLU:OE2	2.18	0.44
1:A:426:SER:O	1:A:430:LYS:HG3	2.18	0.44
1:A:55:ASP:O	1:A:61:THR:HG23	2.18	0.43
1:A:275:ASN:CG	1:A:278:GLU:OE1	2.57	0.43
1:A:18:LYS:HB3	1:A:175:GLU:HB2	2.01	0.43
1:B:250:GLU:OE2	1:B:422:GLU:OE2	2.37	0.42
1:A:400:ARG:CZ	1:A:403:GLU:HG3	2.46	0.42
1:B:323:ASP:OD2	1:B:335:ARG:NH1	2.41	0.42
1:A:81:ARG:O	1:A:83:VAL:HG12	2.20	0.42
1:B:165:SER:HA	1:B:177:ASN:O	2.20	0.42
1:B:207:MET:CE	1:B:212:PHE:CD1	3.01	0.42
1:A:394:GLU:OE2	1:A:408:HIS:CD2	2.72	0.42
1:A:52:ILE:HD12	2:A:501:LQJ:N9	2.35	0.42
1:B:155:MET:O	1:B:324:ARG:NH2	2.51	0.42
1:B:209:ARG:HB3	1:B:209:ARG:NH1	2.35	0.41
1:B:216:LYS:HG3	1:B:222:LEU:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:LYS:HA	1:A:125:GLU:O	2.20	0.41
1:B:48:LEU:CD1	1:B:70:VAL:HG21	2.51	0.41
1:A:70:VAL:HG12	1:A:71:MET:HE2	2.03	0.41
1:B:209:ARG:CZ	1:B:209:ARG:HB3	2.51	0.41
1:B:74:LEU:HD22	1:B:79:ALA:HB3	2.03	0.41
1:B:207:MET:CE	1:B:285:LEU:HD21	2.50	0.41
1:A:361:ARG:NH2	1:A:387:ASN:HD22	2.18	0.41
1:B:131:THR:HB	1:B:165:SER:H	1.86	0.41
2:A:501:LQJ:N2	1:B:19:TYR:OH	2.53	0.41
1:A:6:VAL:CG1	1:A:8:TRP:NE1	2.84	0.40
1:B:225:LEU:HD11	1:B:294:LEU:HD11	2.03	0.40
1:A:18:LYS:HB3	1:A:175:GLU:CB	2.51	0.40
1:A:159:THR:HG22	1:A:185:MET:SD	2.61	0.40
1:B:48:LEU:HD12	1:B:85:VAL:HG13	2.03	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	427/440 (97%)	417 (98%)	10 (2%)	0	100	100
1	B	426/440 (97%)	421 (99%)	5 (1%)	0	100	100
All	All	853/880 (97%)	838 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	371/377 (98%)	346 (93%)	25 (7%)	16	13
1	B	370/377 (98%)	351 (95%)	19 (5%)	24	22
All	All	741/754 (98%)	697 (94%)	44 (6%)	19	17

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	6	VAL
1	A	17	THR
1	A	20	ARG
1	A	37	LYS
1	A	52	ILE
1	A	54	LYS
1	A	59	VAL
1	A	61	THR
1	A	63	ARG
1	A	81	ARG
1	A	91	ILE
1	A	139	LEU
1	A	164	ASN
1	A	170	PRO
1	A	171	LYS
1	A	215	LEU
1	A	222	LEU
1	A	226	ARG
1	A	227	LYS
1	A	327	ARG
1	A	344	LYS
1	A	370	ASN
1	A	385	GLU
1	A	412	LYS
1	B	1	MET
1	B	17	THR
1	B	18	LYS
1	B	25	THR
1	B	57	SER
1	B	60	GLU
1	B	98	SER
1	B	131	THR

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Mol	Chain	Res	Type
1	B	145	LEU
1	B	172	ILE
1	B	174	LYS
1	B	185	MET
1	B	208	GLU
1	B	227	LYS
1	B	279	LYS
1	B	325	LEU
1	B	367	ASP
1	B	371	ARG
1	B	402	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	HIS
1	A	370	ASN
1	A	377	ASN
1	A	387	ASN
1	A	408	HIS
1	B	112	HIS
1	B	156	ASN
1	B	164	ASN
1	B	224	ASN

### 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
2	LQJ	B	501	-	41,50,50	3.12	10 (24%)	44,78,78	2.89	10 (22%)
2	LQJ	A	501	-	41,50,50	3.25	14 (34%)	44,78,78	2.77	9 (20%)

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
2	LQJ	B	501	-	-	5/13/63/63	0/7/7/7
2	LQJ	A	501	-	-	1/13/63/63	0/7/7/7

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	LQJ	C17-N8	10.78	1.49	1.32
2	A	501	LQJ	C7-N3	9.69	1.47	1.32
2	B	501	LQJ	C7-N3	8.93	1.46	1.32
2	A	501	LQJ	C17-N8	8.77	1.46	1.32
2	A	501	LQJ	C17-N9	8.44	1.49	1.33
2	B	501	LQJ	C17-N9	8.42	1.49	1.33
2	A	501	LQJ	C7-N4	8.28	1.49	1.33
2	B	501	LQJ	C7-N4	7.65	1.48	1.33
2	A	501	LQJ	O7-C13	5.43	1.48	1.41
2	A	501	LQJ	C19-C20	4.44	1.62	1.53
2	A	501	LQJ	O2-C3	3.91	1.46	1.41
2	B	501	LQJ	C9-C3	-3.05	1.49	1.53
2	A	501	LQJ	C15-C16	-2.82	1.33	1.40
2	A	501	LQJ	C10-C2	2.77	1.60	1.52
2	B	501	LQJ	C5-C6	-2.49	1.34	1.40
2	B	501	LQJ	C18-C15	-2.46	1.34	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	LQJ	C8-C5	-2.42	1.34	1.43
2	A	501	LQJ	C15-N7	-2.18	1.31	1.39
2	B	501	LQJ	O6-C11	-2.14	1.36	1.44
2	A	501	LQJ	P1-O4	2.13	1.66	1.60
2	A	501	LQJ	C18-C15	-2.09	1.35	1.43
2	A	501	LQJ	C8-C5	-2.07	1.35	1.43
2	A	501	LQJ	O6-C11	-2.07	1.36	1.44
2	B	501	LQJ	C20-C12	2.07	1.58	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	LQJ	N8-C17-N9	-11.42	110.83	128.68
2	A	501	LQJ	N8-C17-N9	-10.40	112.43	128.68
2	B	501	LQJ	N3-C7-N4	-9.72	113.48	128.68
2	A	501	LQJ	N3-C7-N4	-8.92	114.73	128.68
2	B	501	LQJ	C13-N6-C16	5.95	137.10	126.64
2	A	501	LQJ	O7-C13-C19	-5.83	96.48	106.59
2	A	501	LQJ	O7-C12-C20	-5.21	93.70	104.87
2	B	501	LQJ	O7-C13-C19	-4.66	98.50	106.59
2	B	501	LQJ	O11-P2-O10	3.47	121.10	109.89
2	A	501	LQJ	O8-P2-O10	-3.16	107.42	115.76
2	A	501	LQJ	O7-C12-C11	-3.12	99.12	109.37
2	B	501	LQJ	C15-C18-N10	-2.89	115.95	120.35
2	A	501	LQJ	C13-N6-C16	2.89	131.72	126.64
2	A	501	LQJ	O11-P2-O10	2.86	119.11	109.89
2	B	501	LQJ	C20-C19-C13	-2.64	97.02	103.54
2	B	501	LQJ	O2-C3-C9	-2.59	103.14	106.93
2	A	501	LQJ	C16-C15-N7	-2.57	106.72	109.40
2	B	501	LQJ	O2-C2-C1	2.29	114.16	109.21
2	B	501	LQJ	O8-C19-C20	2.14	109.00	105.08

There are no chirality outliers.

All (6) torsion outliers are listed below:

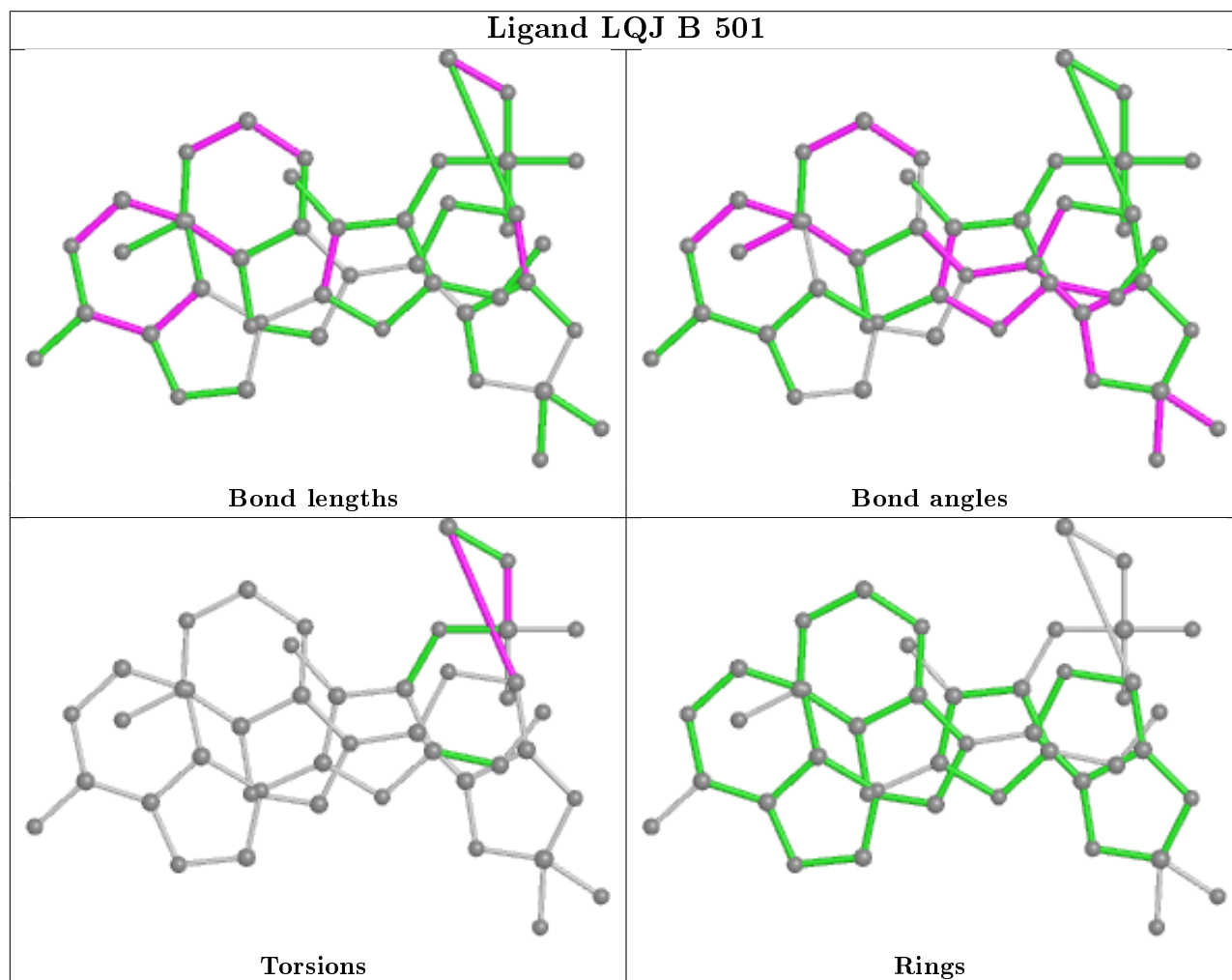
Mol	Chain	Res	Type	Atoms
2	B	501	LQJ	O6-C11-C12-O7
2	B	501	LQJ	O6-C11-C12-C20
2	B	501	LQJ	C11-O6-P1-O4
2	B	501	LQJ	C11-O6-P1-O5
2	B	501	LQJ	C11-O6-P1-O12
2	A	501	LQJ	O6-C11-C12-O7

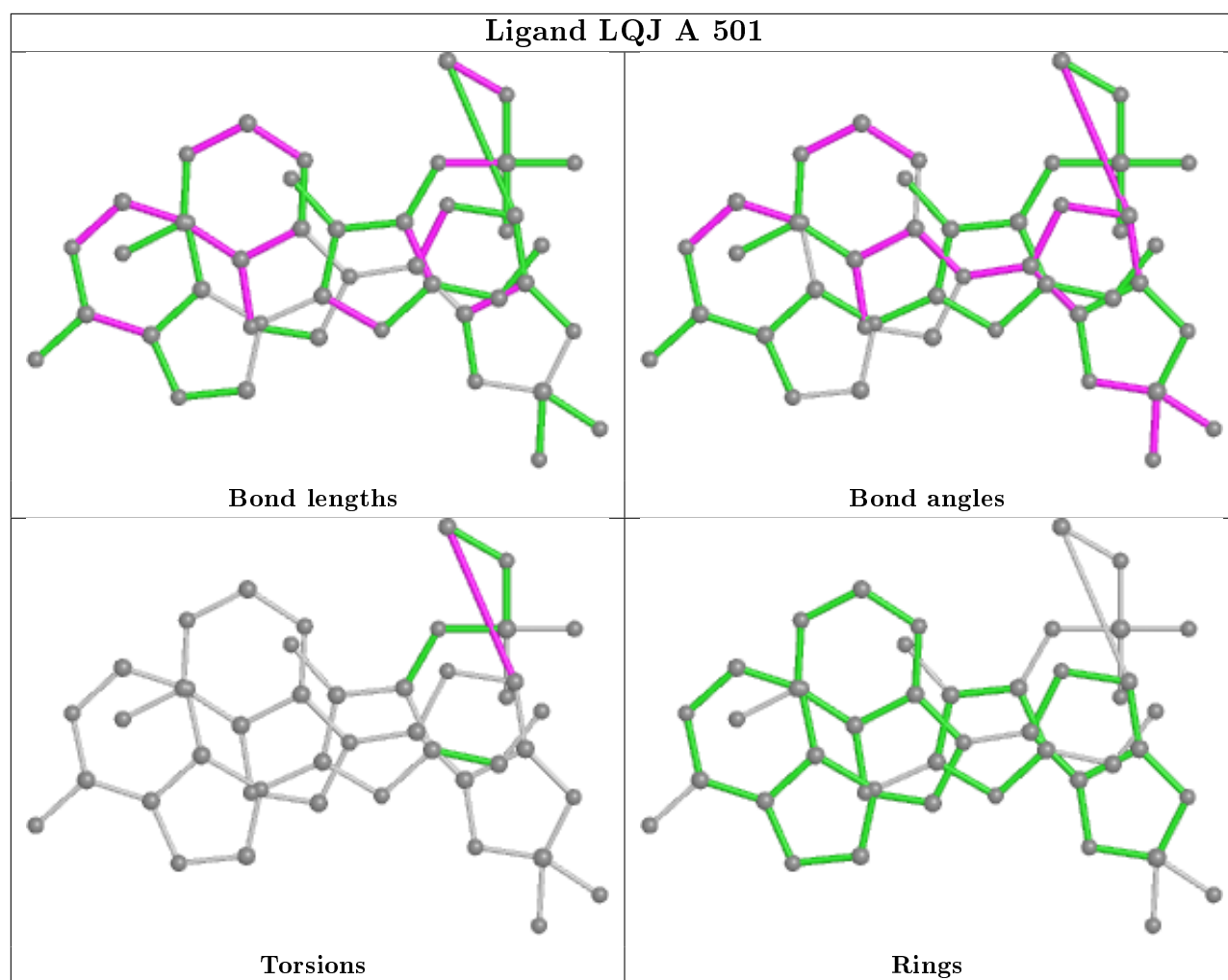
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	LQJ	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 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	431/440 (97%)	-0.12	10 (2%) 60 65	25, 38, 71, 107	0
1	B	430/440 (97%)	-0.07	12 (2%) 53 59	27, 40, 72, 97	0
All	All	861/880 (97%)	-0.09	22 (2%) 56 61	25, 39, 72, 107	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	81	ARG	4.6
1	A	81	ARG	4.2
1	A	139	LEU	4.2
1	B	370	ASN	4.0
1	B	199	ASP	3.8
1	A	199	ASP	3.7
1	A	311	GLU	3.3
1	B	134	LEU	3.2
1	A	60	GLU	2.9
1	A	134	LEU	2.9
1	A	227	LYS	2.5
1	B	139	LEU	2.5
1	A	171	LYS	2.5
1	B	56	PHE	2.5
1	B	277	GLU	2.4
1	B	145	LEU	2.4
1	A	219	LEU	2.3
1	B	227	LYS	2.3
1	B	55	ASP	2.2
1	B	427	ARG	2.2
1	A	198	PHE	2.1
1	B	278	GLU	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](#)

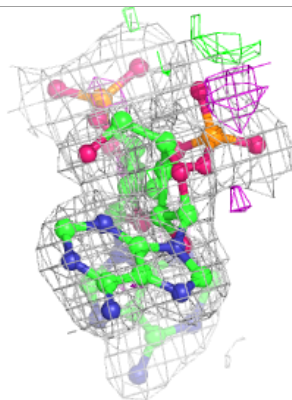
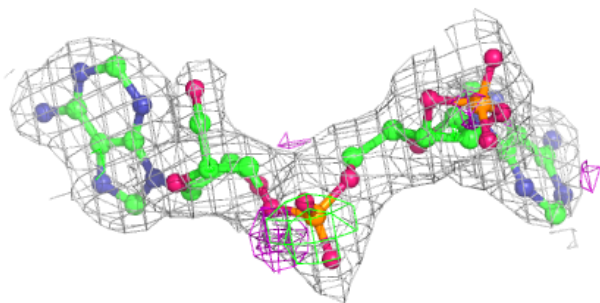
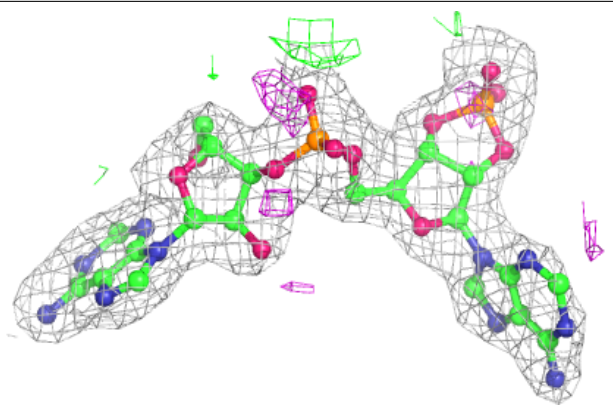
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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	LQJ	B	501	44/44	0.92	0.11	32,49,75,90	0
2	LQJ	A	501	44/44	0.95	0.10	29,41,51,59	0

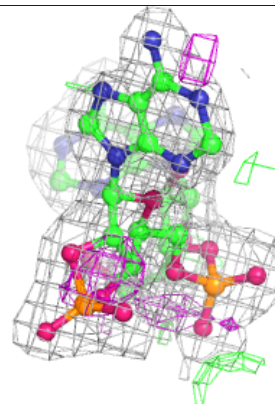
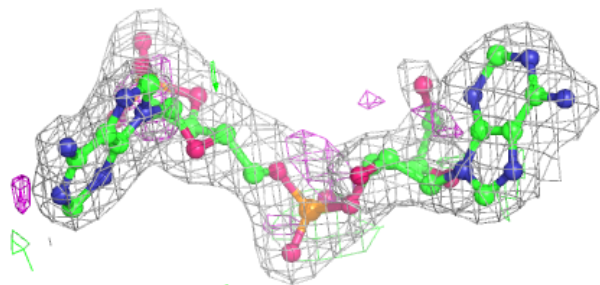
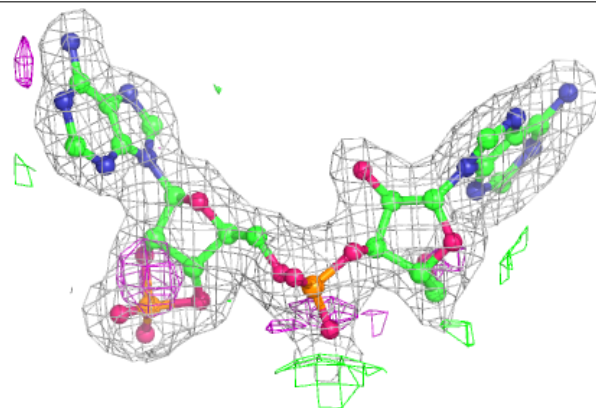
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around LQJ B 501:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around LQJ A 501:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers

There are no such residues in this entry.