



# Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 06:50 am BST

PDB ID : 6GUB  
Title : CDK2/CyclinA in complex with Flavopiridol  
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Deposited on : 2018-06-19  
Resolution : 2.52 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

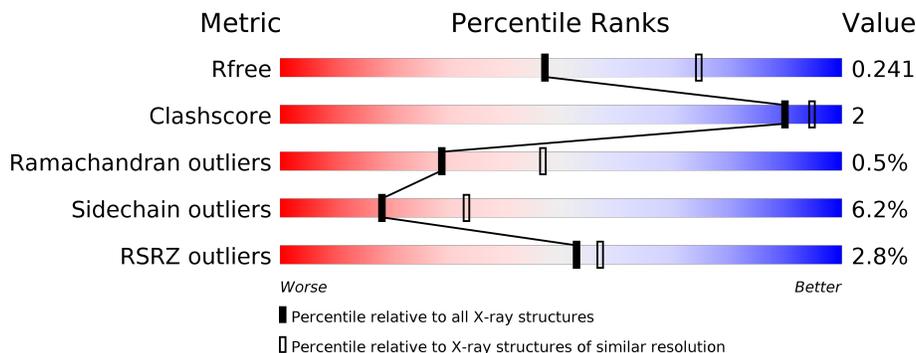
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.52 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	5743 (2.54-2.50)
Clashscore	141614	6463 (2.54-2.50)
Ramachandran outliers	138981	6335 (2.54-2.50)
Sidechain outliers	138945	6337 (2.54-2.50)
RSRZ outliers	127900	5630 (2.54-2.50)

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	302	
1	C	302	
2	B	268	
2	D	268	

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 9046 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 Cyclin-dependent kinase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	N	O	P				S
1	A	294	Total	C	N	O	P	S	0	0	0
			2364	1537	401	417	1	8			
1	C	292	Total	C	N	O	P	S	0	0	0
			2342	1519	394	420	1	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P24941
A	-2	PRO	-	expression tag	UNP P24941
A	-1	GLY	-	expression tag	UNP P24941
A	0	SER	-	expression tag	UNP P24941
C	-3	GLY	-	expression tag	UNP P24941
C	-2	PRO	-	expression tag	UNP P24941
C	-1	GLY	-	expression tag	UNP P24941
C	0	SER	-	expression tag	UNP P24941

- Molecule 2 is a protein called Cyclin-A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	261	Total	C	N	O	S	0	0	0
			2106	1364	343	389	10			
2	D	261	Total	C	N	O	S	0	0	0
			2106	1364	343	389	10			

There are 14 discrepancies between the modelled and reference sequences:

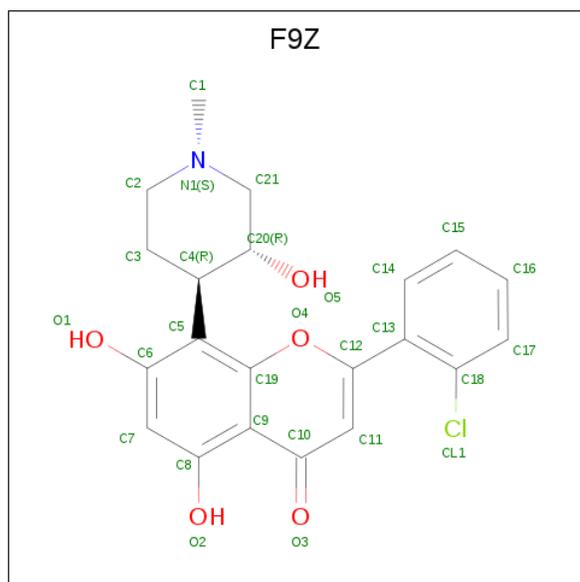
Chain	Residue	Modelled	Actual	Comment	Reference
B	171	GLY	-	expression tag	UNP P30274
B	433	HIS	-	expression tag	UNP P30274
B	434	HIS	-	expression tag	UNP P30274
B	435	HIS	-	expression tag	UNP P30274

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Chain	Residue	Modelled	Actual	Comment	Reference
B	436	HIS	-	expression tag	UNP P30274
B	437	HIS	-	expression tag	UNP P30274
B	438	HIS	-	expression tag	UNP P30274
D	171	GLY	-	expression tag	UNP P30274
D	433	HIS	-	expression tag	UNP P30274
D	434	HIS	-	expression tag	UNP P30274
D	435	HIS	-	expression tag	UNP P30274
D	436	HIS	-	expression tag	UNP P30274
D	437	HIS	-	expression tag	UNP P30274
D	438	HIS	-	expression tag	UNP P30274

- Molecule 3 is 2-(2-chlorophenyl)-8-[(3 {R},4 {R})-1-methyl-3-oxidanyl-piperidin-4-yl]-5,7-bis(oxidanyl)chromen-4-one (three-letter code: F9Z) (formula: C<sub>21</sub>H<sub>20</sub>ClNO<sub>5</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Cl	N			O
3	A	1	28	21	1	1	5	0	0
3	C	1	28	21	1	1	5	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total	O	0	0
			33	33		

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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
4	B	17	Total O 17 17	0	0
4	C	11	Total O 11 11	0	0
4	D	11	Total O 11 11	0	0



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.18Å 136.22Å 149.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	101.04 – 2.52 75.18 – 2.52	Depositor EDS
% Data completeness (in resolution range)	100.0 (101.04-2.52) 100.0 (75.18-2.52)	Depositor EDS
$R_{merge}$	0.20	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 2.51Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.195 , 0.238 0.201 , 0.241	Depositor DCC
$R_{free}$ test set	2638 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.3	Xtrriage
Anisotropy	0.324	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.4	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	0.94	EDS
Total number of atoms	9046	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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: TPO, F9Z

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	0.73	0/2413	0.87	0/3272
1	C	0.69	0/2388	0.84	0/3238
2	B	0.67	0/2156	0.76	0/2932
2	D	0.66	0/2156	0.77	0/2932
All	All	0.69	0/9113	0.81	0/12374

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	2364	0	2412	11	0
1	C	2342	0	2381	17	0
2	B	2106	0	2125	1	0
2	D	2106	0	2125	4	0
3	A	28	0	0	0	0
3	C	28	0	0	0	0
4	A	33	0	0	1	0
4	B	17	0	0	0	0
4	C	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	11	0	0	0	0
All	All	9046	0	9043	32	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 2.

All (32) 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:C:247:ASP:CG	1:C:248:PHE:H	1.74	0.87
1:A:96:LEU:HD12	1:A:97:THR:HG23	1.65	0.77
2:D:289:LYS:O	2:D:293:ARG:HG3	1.91	0.70
1:C:247:ASP:CG	1:C:248:PHE:N	2.45	0.69
1:C:247:ASP:OD1	1:C:248:PHE:N	2.25	0.69
1:C:72:THR:HG22	1:C:74:ASN:H	1.59	0.66
1:C:237:LYS:HB3	1:C:238:PRO:CD	2.26	0.65
1:A:177:CYS:SG	1:A:233:MET:CE	2.87	0.62
1:C:95:ALA:HA	1:C:199:ARG:HD2	1.81	0.61
1:A:177:CYS:SG	1:A:233:MET:HE3	2.43	0.59
1:C:218:THR:HG23	1:C:251:VAL:HG11	1.85	0.59
1:C:237:LYS:HB3	1:C:238:PRO:HD2	1.84	0.58
2:D:289:LYS:O	2:D:293:ARG:CG	2.56	0.54
2:B:323:GLN:HA	2:B:323:GLN:HE21	1.73	0.53
2:D:407:GLN:OE1	2:D:410:ARG:HD3	2.08	0.53
1:C:198:THR:O	1:C:199:ARG:HB2	2.10	0.52
1:A:161:HIS:HB2	4:A:432:HOH:O	2.09	0.51
1:A:293:VAL:HB	1:A:294:PRO:HD2	1.93	0.50
1:C:197:VAL:HG11	1:C:252:VAL:HG12	1.92	0.50
1:A:71:HIS:CD2	1:A:76:LEU:HD13	2.48	0.49
1:C:198:THR:O	1:C:199:ARG:CB	2.61	0.49
1:A:96:LEU:HD12	1:A:96:LEU:C	2.33	0.49
1:C:72:THR:HB	1:C:75:LYS:O	2.15	0.47
1:C:129:LYS:HA	1:C:192:ILE:HD11	1.97	0.47
1:C:251:VAL:HG23	1:C:252:VAL:HG23	1.98	0.46
1:C:248:PHE:HA	1:C:251:VAL:HG22	1.99	0.44
1:A:247:ASP:OD1	1:A:248:PHE:N	2.50	0.44
1:C:53:SER:HB3	2:D:304:PHE:O	2.18	0.43
1:C:72:THR:HG22	1:C:74:ASN:N	2.32	0.43
1:A:177:CYS:SG	1:A:233:MET:HE2	2.59	0.43
1:A:129:LYS:HA	1:A:192:ILE:HD11	2.00	0.42
1:A:12:GLU:HG2	1:A:13:GLY:N	2.34	0.42

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	289/302 (96%)	276 (96%)	12 (4%)	1 (0%)	41	59
1	C	287/302 (95%)	272 (95%)	11 (4%)	4 (1%)	11	19
2	B	259/268 (97%)	254 (98%)	5 (2%)	0	100	100
2	D	259/268 (97%)	254 (98%)	5 (2%)	0	100	100
All	All	1094/1140 (96%)	1056 (96%)	33 (3%)	5 (0%)	29	47

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	237	LYS
1	C	239	SER
1	A	145	ASP
1	C	145	ASP
1	C	98	GLY

### 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	258/264 (98%)	242 (94%)	16 (6%)	18	33
1	C	257/264 (97%)	233 (91%)	24 (9%)	9	16
2	B	234/240 (98%)	222 (95%)	12 (5%)	24	43

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	D	234/240 (98%)	225 (96%)	9 (4%)	33 56
All	All	983/1008 (98%)	922 (94%)	61 (6%)	18 33

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	VAL
1	A	36	ARG
1	A	64	VAL
1	A	73	GLU
1	A	74	ASN
1	A	83	LEU
1	A	88	LYS
1	A	96	LEU
1	A	101	LEU
1	A	122	ARG
1	A	150	ARG
1	A	163	VAL
1	A	200	ARG
1	A	206	ASP
1	A	237	LYS
1	A	248	PHE
2	B	172	VAL
2	B	179	HIS
2	B	189	MET
2	B	196	LYS
2	B	199	TYR
2	B	201	LYS
2	B	245	SER
2	B	283	ASP
2	B	284	ASP
2	B	323	GLN
2	B	378	GLN
2	B	432	VAL
1	C	2	GLU
1	C	17	VAL
1	C	36	ARG
1	C	40	GLU
1	C	64	VAL
1	C	65	LYS
1	C	71	HIS
1	C	73	GLU

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	C	74	ASN
1	C	89	LYS
1	C	101	LEU
1	C	105	LYS
1	C	122	ARG
1	C	131	GLN
1	C	148	LEU
1	C	150	ARG
1	C	163	VAL
1	C	199	ARG
1	C	200	ARG
1	C	217	ARG
1	C	240	PHE
1	C	246	GLN
1	C	248	PHE
1	C	264	SER
2	D	172	VAL
2	D	189	MET
2	D	201	LYS
2	D	245	SER
2	D	283	ASP
2	D	289	LYS
2	D	391	LEU
2	D	400	ARG
2	D	432	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	71	HIS
1	A	85	GLN
1	A	295	HIS
2	B	183	HIS
2	B	254	GLN
2	B	317	GLN
2	B	323	GLN
2	B	395	HIS
1	C	85	GLN
2	D	183	HIS
2	D	254	GLN
2	D	395	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

2 non-standard protein/DNA/RNA residues 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
1	TPO	A	160	1	8,10,11	0.92	0	10,14,16	1.55	2 (20%)
1	TPO	C	160	1	8,10,11	0.94	0	10,14,16	1.50	1 (10%)

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
1	TPO	A	160	1	-	0/9/11/13	-
1	TPO	C	160	1	-	0/9/11/13	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	160	TPO	P-OG1-CB	-3.38	112.98	123.21
1	A	160	TPO	P-OG1-CB	-2.93	114.35	123.21
1	A	160	TPO	O-C-CA	-2.46	118.33	124.78

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	F9Z	C	301	-	25,31,31	1.33	2 (8%)	37,46,46	1.60	6 (16%)
3	F9Z	A	301	-	25,31,31	1.32	2 (8%)	37,46,46	1.56	5 (13%)

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	F9Z	C	301	-	-	4/7/21/21	0/4/4/4
3	F9Z	A	301	-	-	4/7/21/21	0/4/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	301	F9Z	C11-C10	4.21	1.46	1.37
3	C	301	F9Z	C6-C5	4.15	1.42	1.38
3	A	301	F9Z	C11-C10	4.13	1.45	1.37
3	A	301	F9Z	C6-C5	3.85	1.41	1.38

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	301	F9Z	C11-C10-C9	-5.25	116.78	123.16
3	A	301	F9Z	C11-C10-C9	-4.63	117.53	123.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	301	F9Z	C3-C4-C20	-4.38	103.52	109.54
3	A	301	F9Z	C3-C4-C20	-4.25	103.70	109.54
3	A	301	F9Z	C20-C21-N1	-3.39	104.78	110.74
3	C	301	F9Z	C20-C21-N1	-3.10	105.28	110.74
3	C	301	F9Z	O4-C19-C5	3.03	119.77	116.17
3	A	301	F9Z	O4-C19-C5	2.97	119.70	116.17
3	A	301	F9Z	C3-C2-N1	2.92	115.14	111.22
3	C	301	F9Z	C3-C2-N1	2.55	114.64	111.22
3	C	301	F9Z	C2-N1-C21	2.10	112.08	110.16

There are no chirality outliers.

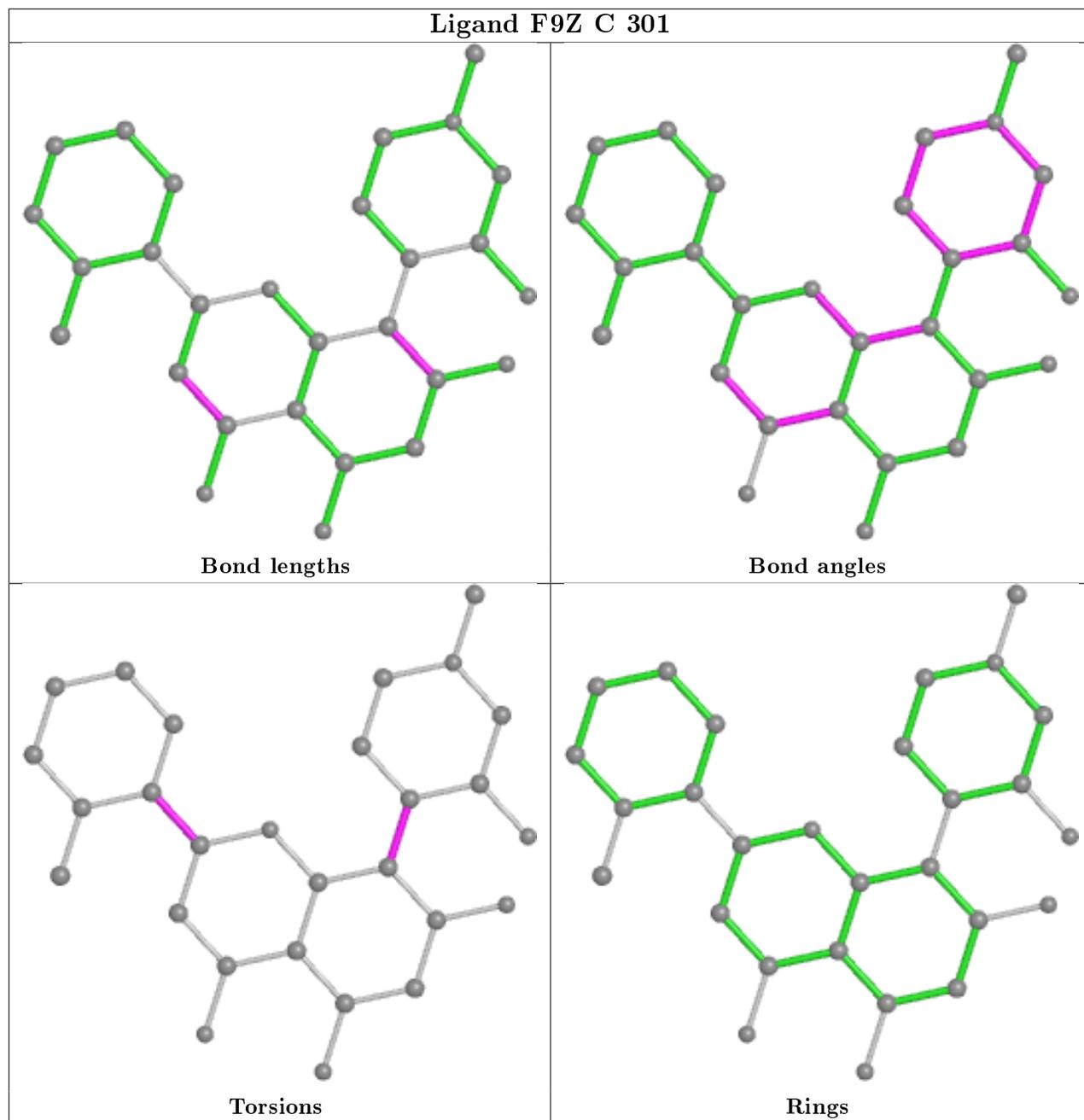
All (8) torsion outliers are listed below:

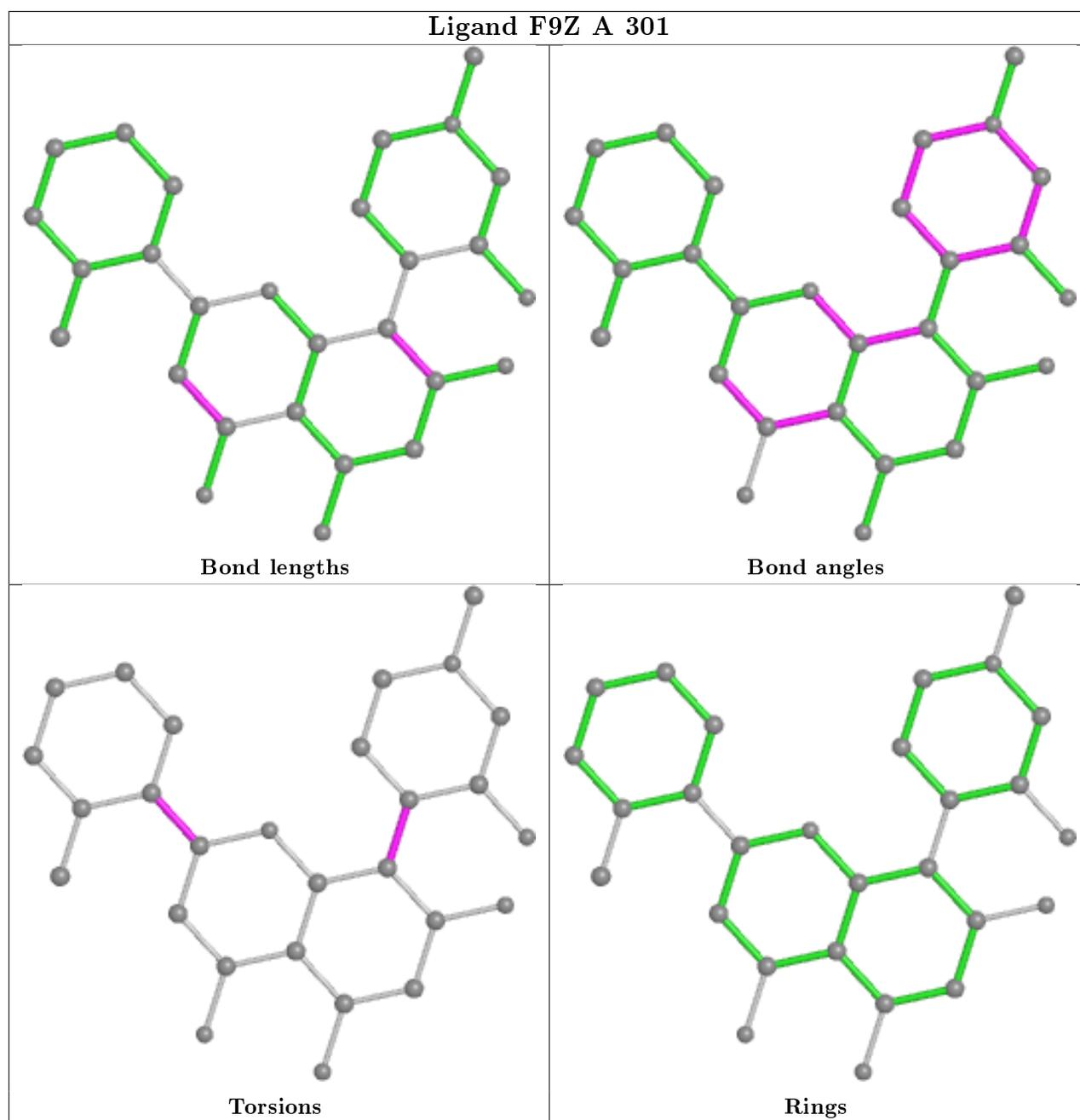
Mol	Chain	Res	Type	Atoms
3	C	301	F9Z	C20-C4-C5-C6
3	A	301	F9Z	C20-C4-C5-C6
3	C	301	F9Z	O4-C12-C13-C14
3	A	301	F9Z	O4-C12-C13-C14
3	C	301	F9Z	C11-C12-C13-C18
3	A	301	F9Z	C11-C12-C13-C18
3	C	301	F9Z	C20-C4-C5-C19
3	A	301	F9Z	C20-C4-C5-C19

There are no ring outliers.

No monomer is involved in short contacts.

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	293/302 (97%)	-0.07	2 (0%) 87 89	21, 32, 69, 108	0
1	C	291/302 (96%)	0.39	25 (8%) 10 10	25, 46, 129, 189	0
2	B	261/268 (97%)	-0.07	1 (0%) 92 93	24, 40, 66, 93	0
2	D	261/268 (97%)	0.10	3 (1%) 80 83	26, 50, 88, 109	0
All	All	1106/1140 (97%)	0.09	31 (2%) 53 57	21, 42, 87, 189	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	236	TYR	10.5
1	C	240	PHE	7.3
1	C	222	PRO	7.1
1	C	246	GLN	6.5
1	C	238	PRO	6.5
1	C	295	HIS	6.0
1	C	296	LEU	5.4
1	C	247	ASP	5.4
1	C	227	TRP	5.3
1	A	96	LEU	4.5
1	C	237	LYS	4.2
1	C	248	PHE	4.2
2	D	419	HIS	3.9
1	C	228	PRO	3.8
1	C	71	HIS	3.7
1	C	249	SER	3.3
1	C	226	VAL	3.1
1	C	73	GLU	2.8
1	C	224	GLU	2.8
1	C	229	GLY	2.8
2	B	432	VAL	2.5

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	C	217	ARG	2.3
1	C	232	SER	2.3
1	C	233	MET	2.2
1	C	282	ALA	2.2
1	A	161	HIS	2.2
1	C	36	ARG	2.1
1	C	101	LEU	2.1
2	D	280	TYR	2.1
1	C	294	PRO	2.1
2	D	323	GLN	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	TPO	A	160	11/12	0.99	0.14	25,29,34,36	0
1	TPO	C	160	11/12	0.99	0.13	29,34,39,41	0

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

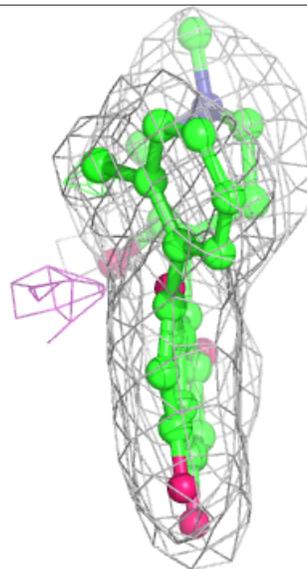
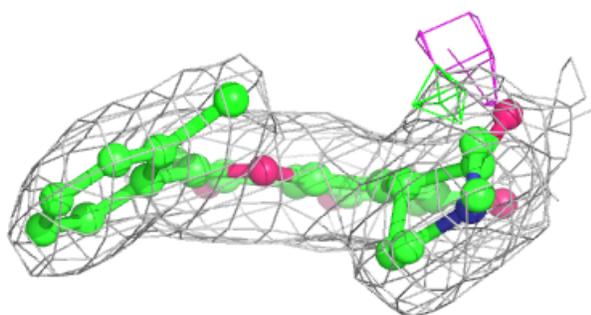
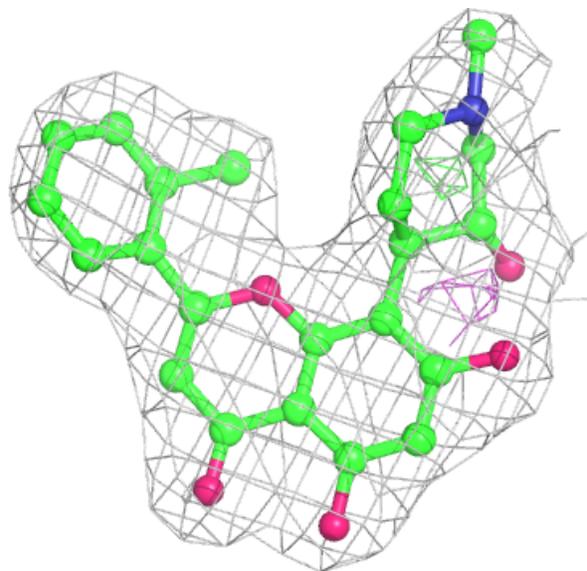
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

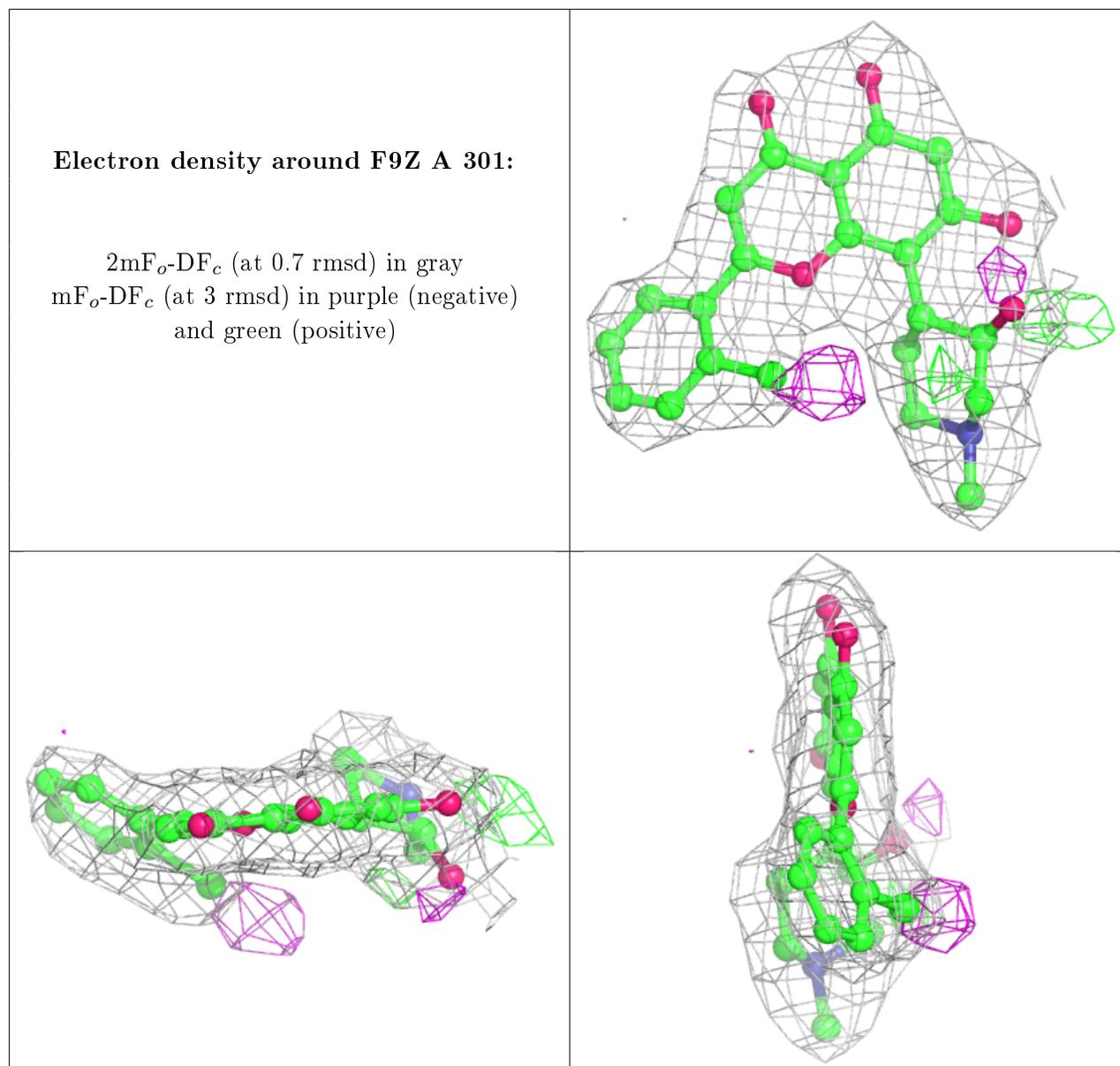
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	F9Z	C	301	28/28	0.95	0.16	26,34,49,71	0
3	F9Z	A	301	28/28	0.96	0.15	24,27,43,70	0

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 F9Z C 301:**

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





## 6.5 Other polymers [i](#)

There are no such residues in this entry.