



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:26 AM UTC

PDB ID : 5LMK / pdb_00005lmk
Title : Structure of phopsho-CDK2-cyclin A in complex with an ATP-competitive inhibitor
Authors : Echalier, A.
Deposited on : 2016-08-01
Resolution : 2.40 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

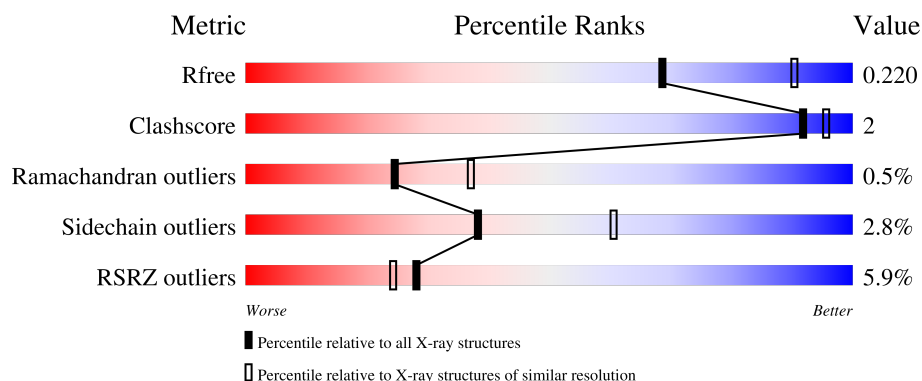
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	299	4% 93% 6%
2	B	258	2% 92% 8%
2	D	258	9% 91% 7%
3	C	297	7% 84% 6% 10%

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 9043 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
1	A	297	Total	C	N	O	P	S	0	1	0
			2399	1558	410	422	1	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P24941

- Molecule 2 is a protein called Cyclin-A2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	258	Total	C	N	O	S		0	3	0
			2104	1365	343	384	12				
2	D	257	Total	C	N	O	S		0	0	0
			2076	1345	338	382	11				

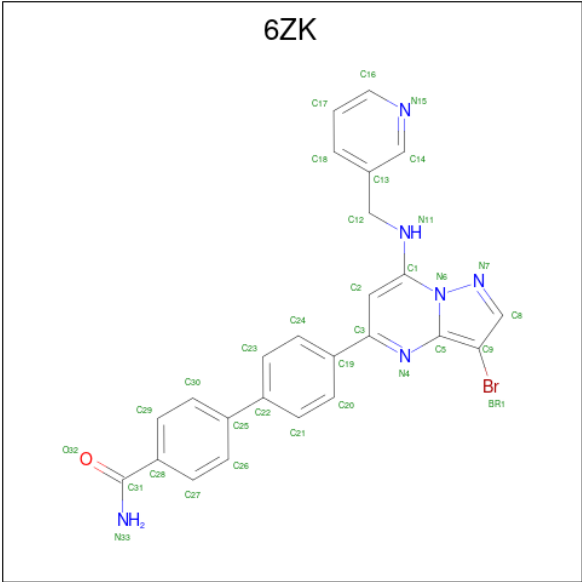
- Molecule 3 is a protein called Cyclin-dependent kinase 2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	266	Total	C	N	O	P	S	0	3	0
			2155	1396	371	380	1	7			

There is a discrepancy between the modelled and reference sequences:

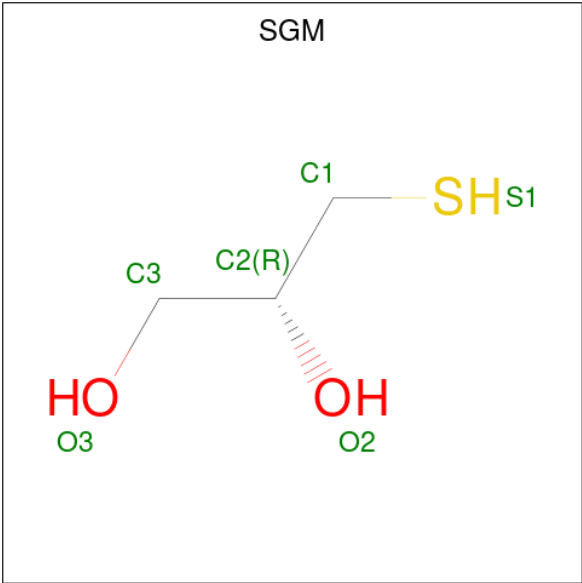
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP P24941

- Molecule 4 is 4-[4-[3-bromanyl-7-(pyridin-3-ylmethylamino)pyrazolo[1,5-a]pyrimidin-5-yl]phenyl]benzamide (CCD ID: 6ZK) (formula: C₂₅H₁₉BrN₆O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	Br	C	N	O	0	0
			33	1	25	6	1		
4	C	1	Total	Br	C	N	O	0	0
			33	1	25	6	1		

- Molecule 5 is MONOTHIOGLYCEROL (CCD ID: SGM) (formula: C₃H₈O₂S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	O	S	0	0
			6	3	2	1		
5	D	1	Total	C	O	S	0	0
			6	3	2	1		

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total 1	Mg 1	0	0

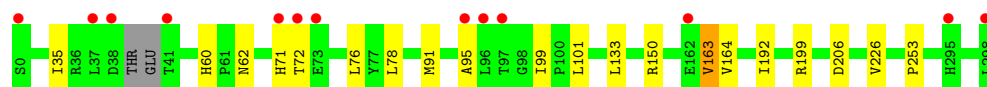
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	85	Total 85	O 85	0	0
7	B	73	Total 73	O 73	0	0
7	C	43	Total 43	O 43	0	0
7	D	29	Total 29	O 29	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.

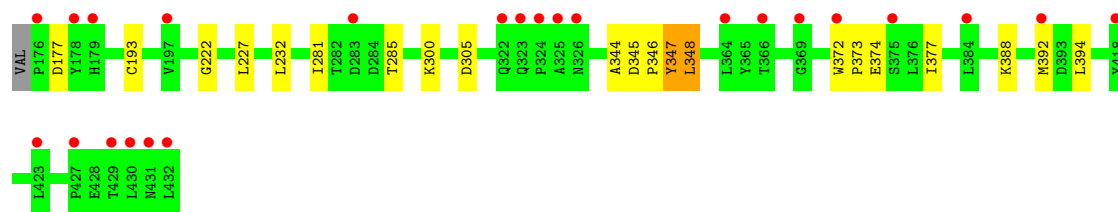
- Molecule 1: Cyclin-dependent kinase 2



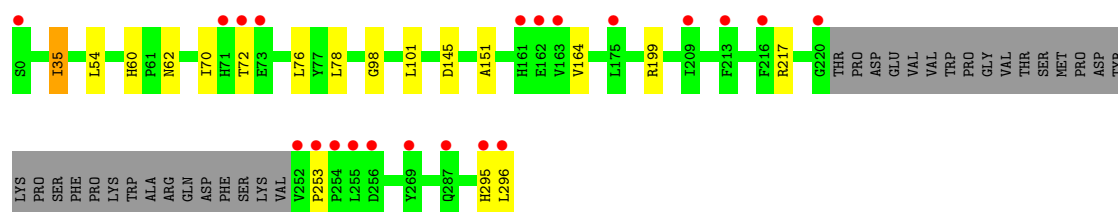
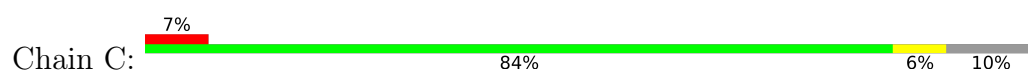
- Molecule 2: Cyclin-A2



- Molecule 2: Cyclin-A2



- Molecule 3: Cyclin-dependent kinase 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.47Å 134.40Å 149.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	99.94 – 2.40 99.94 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.1 (99.94-2.40) 91.3 (99.94-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0131	Depositor
R, R_{free}	0.212 , 0.249 (Not available) , 0.220	Depositor DCC
R_{free} test set	2862 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.299	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9043	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: TPO, 6ZK, SGM, MG

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	0.60	0/2451	0.97	2/3322 (0.1%)
2	B	0.59	0/2161	1.03	0/2933
2	D	0.56	0/2126	1.02	1/2886 (0.0%)
3	C	0.58	0/2201	0.95	1/2978 (0.0%)
All	All	0.58	0/8939	0.99	4/12119 (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
3	C	0	1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	253	PRO	N-CA-C	7.72	120.12	110.70
1	A	253	PRO	N-CA-C	6.91	119.13	110.70
1	A	163	VAL	N-CA-C	6.52	122.89	109.34
2	D	347	TYR	N-CA-C	5.21	119.26	113.01

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	70	ILE	Peptide

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	2399	0	2453	6	0
2	B	2104	0	2135	9	0
2	D	2076	0	2099	9	0
3	C	2155	0	2219	5	0
4	A	33	0	0	0	0
4	C	33	0	0	0	0
5	B	6	0	8	2	0
5	D	6	0	8	2	0
6	B	1	0	0	0	0
7	A	85	0	0	0	0
7	B	73	0	0	1	0
7	C	43	0	0	0	0
7	D	29	0	0	0	0
All	All	9043	0	8922	29	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 (29) 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)
2:B:193:CYS:SG	5:B:501:SGM:S1	2.53	1.02
2:D:305:ASP:HB3	5:D:501:SGM:H12	1.77	0.66
2:D:193:CYS:SG	5:D:501:SGM:S1	2.65	0.65
1:A:60:HIS:CD2	1:A:62:ASN:H	2.19	0.60
1:A:60:HIS:HD2	1:A:62:ASN:H	1.52	0.56
2:D:388:LYS:HE2	2:D:392:MET:HE1	1.89	0.54
1:A:71:HIS:CD2	1:A:76:LEU:HD13	2.46	0.51
3:C:98:GLY:HA2	3:C:199:ARG:HD3	1.92	0.51
2:D:374:GLU:HA	2:D:377:ILE:HD12	1.93	0.50
3:C:60:HIS:CD2	3:C:62:ASN:H	2.30	0.49
2:D:344:ALA:HB1	2:D:348:LEU:HD22	1.95	0.49
3:C:35:ILE:HD12	3:C:76:LEU:HB3	1.98	0.46
2:D:345:ASP:HA	2:D:346:PRO:HA	1.72	0.46
2:D:372:TRP:HA	2:D:373:PRO:HD3	1.82	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:229:ASN:HD22	2:B:334[A]:MET:HE3	1.81	0.45
3:C:54:LEU:HD21	3:C:151:ALA:HB2	1.97	0.45
3:C:60:HIS:HD2	3:C:62:ASN:H	1.63	0.45
2:B:192:LYS:HB3	5:B:501:SGM:H31	1.98	0.45
1:A:133:LEU:HD11	1:A:192:ILE:HD13	1.99	0.45
1:A:91:MET:HG2	1:A:99:ILE:HD11	1.99	0.44
2:B:345:ASP:HA	2:B:346:PRO:HA	1.75	0.43
2:B:344:ALA:HB1	2:B:348:LEU:HD22	1.99	0.43
2:D:347:TYR:OH	2:D:394:LEU:HA	2.19	0.43
1:A:95:ALA:HA	1:A:199:ARG:HE	1.84	0.43
2:B:227:LEU:HD12	2:B:261:MET:HE3	2.02	0.42
2:B:395:HIS:HE1	2:B:427:PRO:O	2.03	0.42
2:B:183:HIS:HD2	7:B:613:HOH:O	2.04	0.41
2:D:222:GLY:HA2	2:D:227:LEU:HD12	2.03	0.40
2:B:323:GLN:HA	2:B:324:PRO:HA	1.80	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	A	293/299 (98%)	286 (98%)	5 (2%)	2 (1%)	18	28
2	B	259/258 (100%)	257 (99%)	2 (1%)	0	100	100
2	D	255/258 (99%)	251 (98%)	4 (2%)	0	100	100
3	C	264/297 (89%)	254 (96%)	7 (3%)	3 (1%)	11	18
All	All	1071/1112 (96%)	1048 (98%)	18 (2%)	5 (0%)	24	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	163	VAL
3	C	164	VAL
1	A	164	VAL
3	C	295	HIS
3	C	145	ASP

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	A	262/263 (100%)	255 (97%)	7 (3%)	39	62
2	B	223/232 (96%)	215 (96%)	8 (4%)	31	52
2	D	231/232 (100%)	225 (97%)	6 (3%)	40	63
3	C	235/261 (90%)	229 (97%)	6 (3%)	40	63
All	All	951/988 (96%)	924 (97%)	27 (3%)	38	60

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	ILE
1	A	72	THR
1	A	78	LEU
1	A	101	LEU
1	A	150	ARG
1	A	206	ASP
1	A	226	VAL
2	B	177	ASP
2	B	196[A]	LYS
2	B	196[B]	LYS
2	B	232	LEU
2	B	292	LEU
2	B	348	LEU
2	B	392	MET
2	B	432	LEU
3	C	35	ILE
3	C	72	THR

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Mol	Chain	Res	Type
3	C	78	LEU
3	C	101	LEU
3	C	217	ARG
3	C	296	LEU
2	D	177	ASP
2	D	232	LEU
2	D	281	ILE
2	D	285	THR
2	D	300	LYS
2	D	348	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	ASN
1	A	60	HIS
1	A	268	HIS
2	B	228	GLN
2	B	237	ASN
2	B	254	GLN
2	B	296	HIS
2	B	323	GLN
2	B	395	HIS
2	B	403	GLN
3	C	3	ASN
3	C	60	HIS
3	C	62	ASN
3	C	84	HIS
3	C	119	HIS
3	C	268	HIS
2	D	179	HIS
2	D	228	GLN
2	D	254	GLN
2	D	312	ASN
2	D	317	GLN
2	D	322	GLN
2	D	370	GLN
2	D	395	HIS
2	D	415	ASN
2	D	419	HIS
2	D	431	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	$\# Z > 2$
3	TPO	C	160	3	8,10,11	0.81	0	10,14,16	1.08	0
1	TPO	A	160	1	8,10,11	0.82	0	10,14,16	1.04	0

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	TPO	C	160	3	-	0/9/11/13	-
1	TPO	A	160	1	-	1/9/11/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	160	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 for Mogul analysis.

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
5	SGM	B	501	-	5,5,5	0.41	0	5,5,5	0.56	0
4	6ZK	A	301	-	36,37,37	1.65	6 (16%)	45,52,52	1.80	12 (26%)
5	SGM	D	501	-	5,5,5	0.36	0	5,5,5	0.69	0
4	6ZK	C	301	-	36,37,37	1.63	7 (19%)	45,52,52	1.67	8 (17%)

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
5	SGM	B	501	-	-	1/4/4/4	-
4	6ZK	A	301	-	-	5/16/17/17	0/5/5/5
5	SGM	D	501	-	-	1/4/4/4	-
4	6ZK	C	301	-	-	5/16/17/17	0/5/5/5

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	6ZK	BR1-C9	5.19	2.00	1.88
4	C	301	6ZK	BR1-C9	4.87	1.99	1.88
4	C	301	6ZK	C5-N6	-3.79	1.33	1.38
4	A	301	6ZK	C5-C9	3.42	1.48	1.38
4	A	301	6ZK	C5-N6	-3.37	1.33	1.38
4	C	301	6ZK	C5-C9	3.25	1.47	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	301	6ZK	C8-N7	3.02	1.37	1.32
4	C	301	6ZK	C8-N7	2.96	1.37	1.32
4	A	301	6ZK	C3-N4	2.87	1.36	1.33
4	C	301	6ZK	C3-N4	2.85	1.36	1.33
4	A	301	6ZK	C1-N6	-2.29	1.35	1.37
4	C	301	6ZK	N6-N7	2.26	1.41	1.37
4	C	301	6ZK	C1-N6	-2.20	1.35	1.37

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	6ZK	C8-N7-N6	5.68	107.43	103.43
4	C	301	6ZK	C8-N7-N6	5.65	107.40	103.43
4	A	301	6ZK	C26-C25-C22	-3.68	114.91	121.25
4	A	301	6ZK	C20-C19-C3	-3.20	117.28	120.87
4	C	301	6ZK	C19-C3-N4	3.09	120.23	116.94
4	C	301	6ZK	N6-C5-N4	3.08	125.56	122.15
4	C	301	6ZK	C2-C3-N4	-2.98	118.99	121.83
4	A	301	6ZK	C16-N15-C14	2.79	121.73	116.85
4	A	301	6ZK	C19-C3-N4	2.76	119.88	116.94
4	C	301	6ZK	C16-N15-C14	2.74	121.65	116.85
4	A	301	6ZK	N6-C5-N4	2.73	125.17	122.15
4	A	301	6ZK	C21-C22-C25	-2.71	116.57	121.25
4	A	301	6ZK	C2-C3-N4	-2.53	119.42	121.83
4	C	301	6ZK	C20-C19-C3	-2.36	118.23	120.87
4	C	301	6ZK	C30-C25-C22	-2.30	117.29	121.25
4	A	301	6ZK	C30-C25-C22	2.26	125.15	121.25
4	A	301	6ZK	C2-C1-N11	-2.15	121.23	127.46
4	C	301	6ZK	C2-C1-N6	2.08	117.47	114.85
4	A	301	6ZK	C2-C1-N6	2.05	117.44	114.85
4	A	301	6ZK	C13-C14-N15	-2.05	119.84	123.75

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	301	6ZK	C23-C22-C25-C26
4	C	301	6ZK	C21-C22-C25-C30
4	C	301	6ZK	C21-C22-C25-C26
4	C	301	6ZK	C23-C22-C25-C30
4	A	301	6ZK	C21-C22-C25-C26
4	A	301	6ZK	C21-C22-C25-C30

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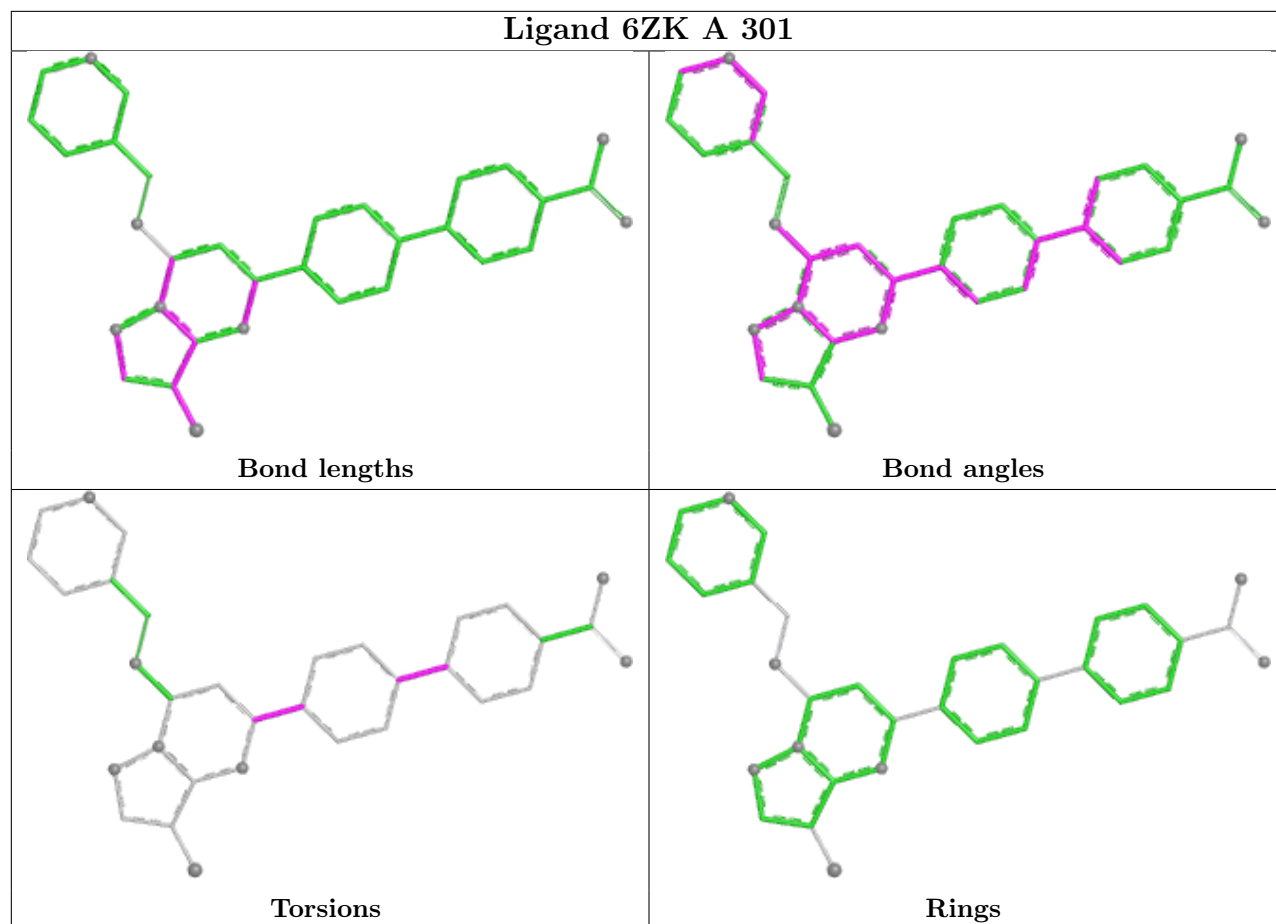
Mol	Chain	Res	Type	Atoms
4	A	301	6ZK	C23-C22-C25-C26
4	A	301	6ZK	C23-C22-C25-C30
5	B	501	SGM	S1-C1-C2-O2
5	D	501	SGM	S1-C1-C2-O2
4	C	301	6ZK	C29-C28-C31-N33
4	A	301	6ZK	C20-C19-C3-N4

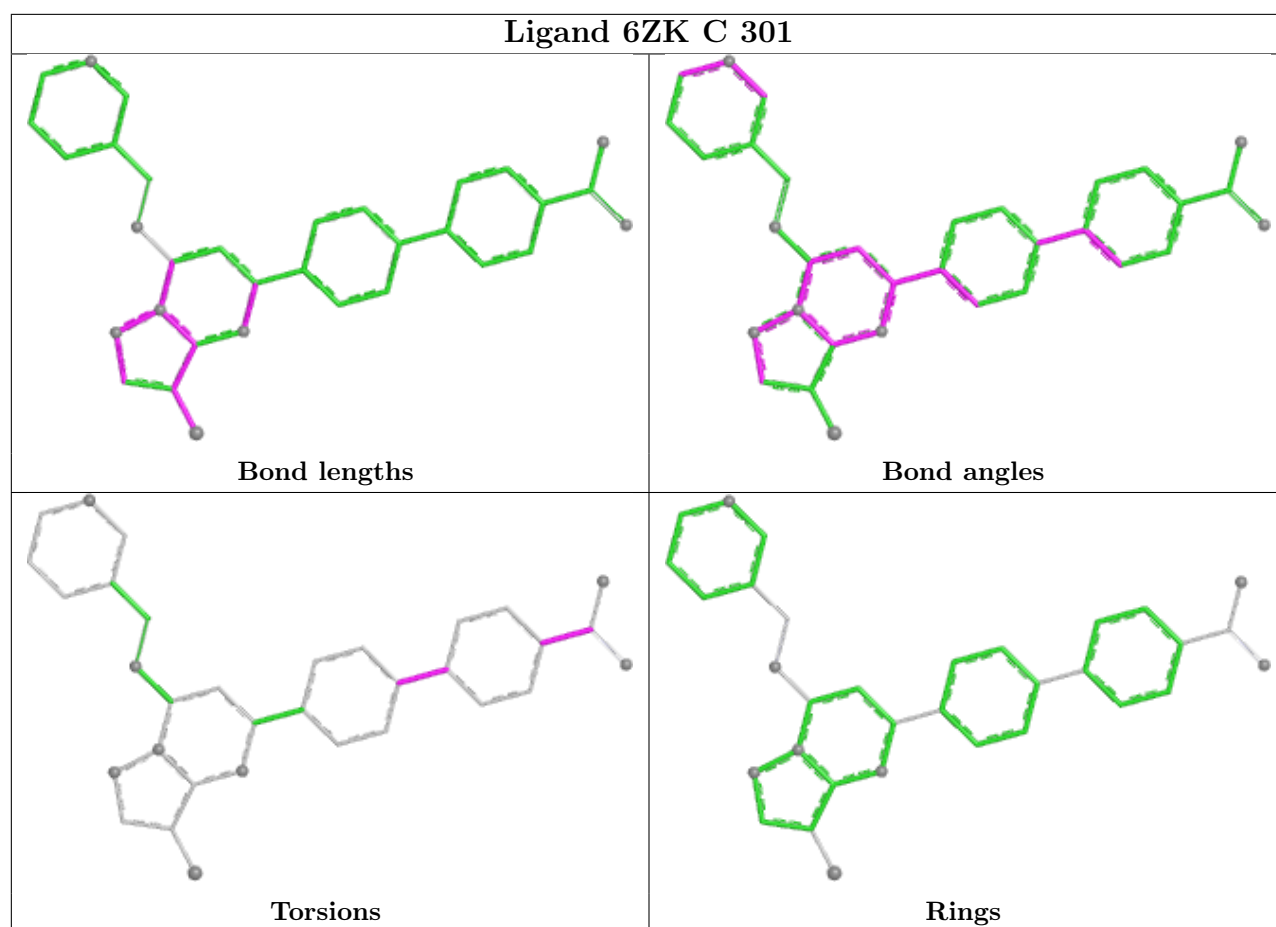
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	501	SGM	2	0
5	D	501	SGM	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, 95th 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 ²)	Q<0.9
1	A	296/299 (98%)	-0.04	13 (4%) 39 34	19, 28, 61, 85	1 (0%)
2	B	258/258 (100%)	-0.10	5 (1%) 66 62	14, 31, 53, 75	3 (1%)
2	D	257/258 (99%)	0.64	24 (9%) 14 11	25, 45, 84, 99	0
3	C	265/297 (89%)	0.56	21 (7%) 18 15	21, 44, 69, 83	3 (1%)
All	All	1076/1112 (96%)	0.26	63 (5%) 28 24	14, 37, 72, 99	7 (0%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	72	THR	6.2
2	D	176	PRO	4.8
1	A	298	LEU	4.6
3	C	252	VAL	4.2
1	A	38	ASP	4.2
2	D	324	PRO	4.1
1	A	96	LEU	4.1
1	A	97	THR	4.0
2	B	175	VAL	3.9
2	D	430	LEU	3.9
3	C	71	HIS	3.9
2	D	432	LEU	3.8
3	C	296	LEU	3.8
1	A	71	HIS	3.6
3	C	0	SER	3.4
2	B	176	PRO	3.3
1	A	41	THR	3.3
1	A	95	ALA	3.2
2	D	418	TYR	3.1
1	A	73	GLU	3.1
2	D	429	THR	3.0

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Mol	Chain	Res	Type	RSRZ
3	C	175	LEU	3.0
2	B	177	ASP	2.9
3	C	213	PHE	2.9
3	C	161	HIS	2.8
2	D	364	LEU	2.8
1	A	295	HIS	2.8
1	A	0	SER	2.8
3	C	254	PRO	2.8
3	C	295	HIS	2.8
3	C	73	GLU	2.7
2	D	372	TRP	2.6
3	C	253	PRO	2.6
3	C	256	ASP	2.6
3	C	216	PHE	2.6
3	C	72	THR	2.6
2	D	283	ASP	2.5
3	C	269	TYR	2.5
3	C	209	ILE	2.5
1	A	37	LEU	2.4
3	C	220	GLY	2.4
2	D	423	LEU	2.4
2	D	179	HIS	2.4
3	C	163	VAL	2.4
2	D	326	ASN	2.4
1	A	162	GLU	2.3
2	D	366	THR	2.3
2	D	323	GLN	2.3
2	D	375	SER	2.3
2	D	369	GLY	2.3
2	D	392	MET	2.3
2	D	178	TYR	2.2
2	D	322	GLN	2.2
2	B	432	LEU	2.2
3	C	255	LEU	2.2
2	D	431	ASN	2.1
2	D	197	VAL	2.1
2	D	427	PRO	2.1
3	C	287	GLN	2.1
3	C	162	GLU	2.1
2	B	323	GLN	2.0
2	D	325	ALA	2.0
2	D	384	LEU	2.0

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, 95th 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
3	TPO	C	160	11/12	0.96	0.10	43,46,51,52	0
1	TPO	A	160	11/12	0.98	0.05	27,28,31,32	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

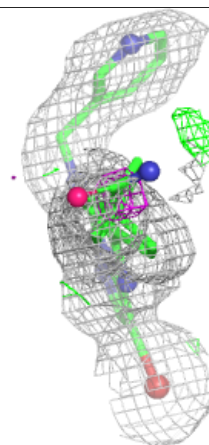
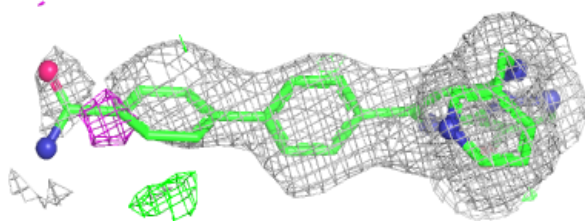
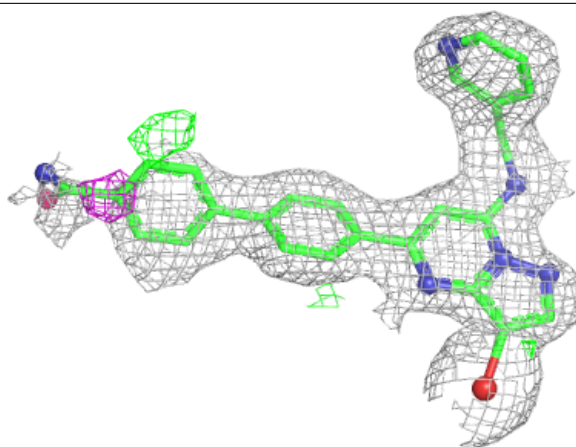
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
5	SGM	D	501	6/6	0.90	0.17	47,49,52,56	0
5	SGM	B	501	6/6	0.91	0.15	37,40,41,45	0
6	MG	B	502	1/1	0.94	0.20	48,48,48,48	0
4	6ZK	C	301	33/33	0.97	0.11	32,43,68,69	0
4	6ZK	A	301	33/33	0.98	0.08	25,33,44,45	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.

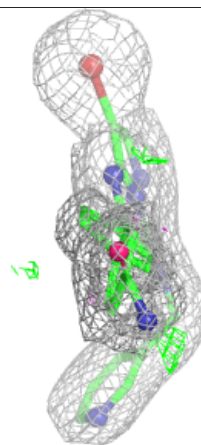
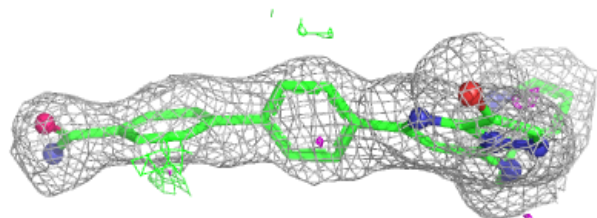
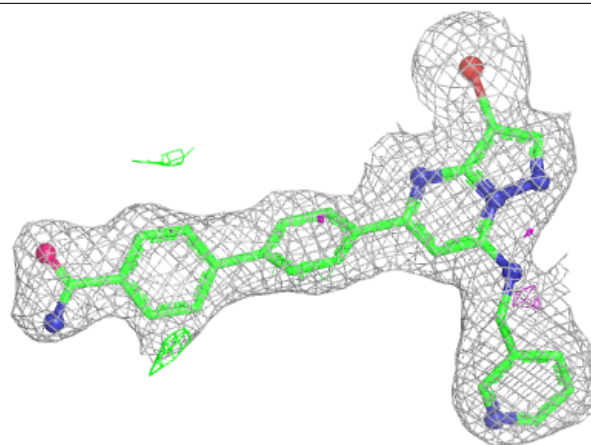
Electron density around 6ZK C 301:

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



Electron density around 6ZK A 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.