



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 01:30 AM UTC

PDB ID : 6C1O / pdb_00006c1o
Title : FGFR1 kinase domain complexed with FIIN-1
Authors : Kalyukina, M.; Yosaatmadja, Y.; Smaill, J.B.; Squire, C.J.
Deposited on : 2018-01-05
Resolution : 2.29 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

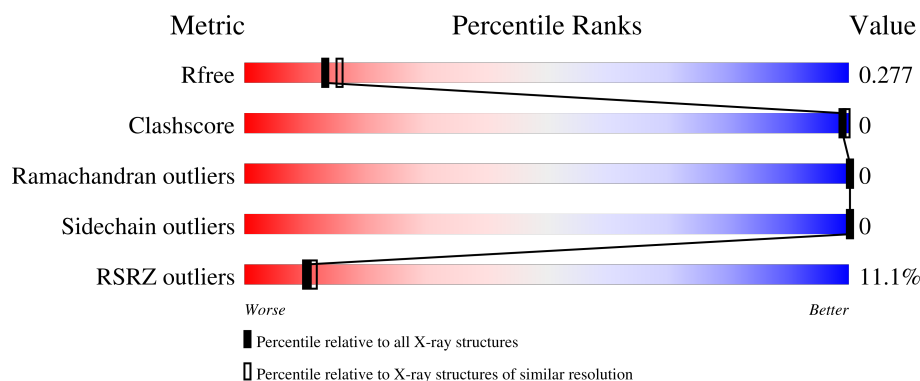
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	311	11% 89% 10%
1	B	311	8% 84% 16%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4171 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.

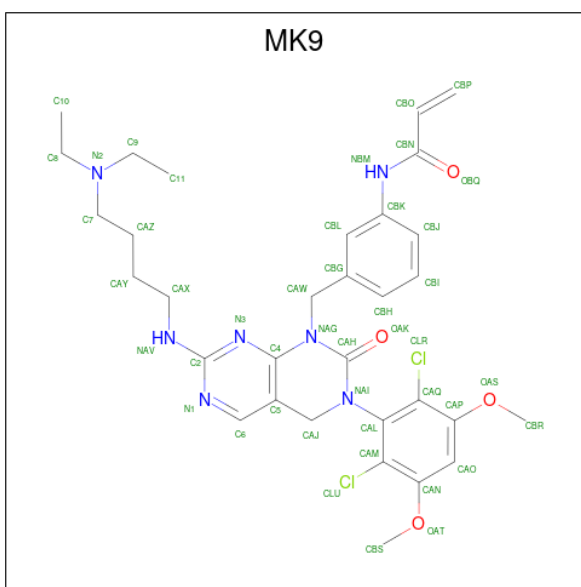
- Molecule 1 is a protein called Fibroblast growth factor receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	280	Total	C	N	O	S	0	1	0
			2122	1347	365	394	16			
1	B	262	Total	C	N	O	S	0	0	0
			1952	1240	336	360	16			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	455	GLY	-	expression tag	UNP P11362
A	456	ALA	-	expression tag	UNP P11362
A	457	MET	-	expression tag	UNP P11362
A	458	SER	-	expression tag	UNP P11362
A	488	ALA	CYS	engineered mutation	UNP P11362
A	584	SER	CYS	engineered mutation	UNP P11362
B	455	GLY	-	expression tag	UNP P11362
B	456	ALA	-	expression tag	UNP P11362
B	457	MET	-	expression tag	UNP P11362
B	458	SER	-	expression tag	UNP P11362
B	488	ALA	CYS	engineered mutation	UNP P11362
B	584	SER	CYS	engineered mutation	UNP P11362

- Molecule 2 is N-(3-{[3-(2,6-dichloro-3,5-dimethoxyphenyl)-7-{[4-(diethylamino)butyl]amino}-2-oxo-3,4-dihydropyrimido[4,5-d]pyrimidin-1(2H)-yl]methyl}phenyl)prop-2-enamide (CCD ID: MK9) (formula: C₃₂H₃₉Cl₂N₇O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 45	C 32	Cl 2	N 7	O 4	0	0
2	B	1	Total 33	C 23	Cl 2	N 5	O 3	0	0

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

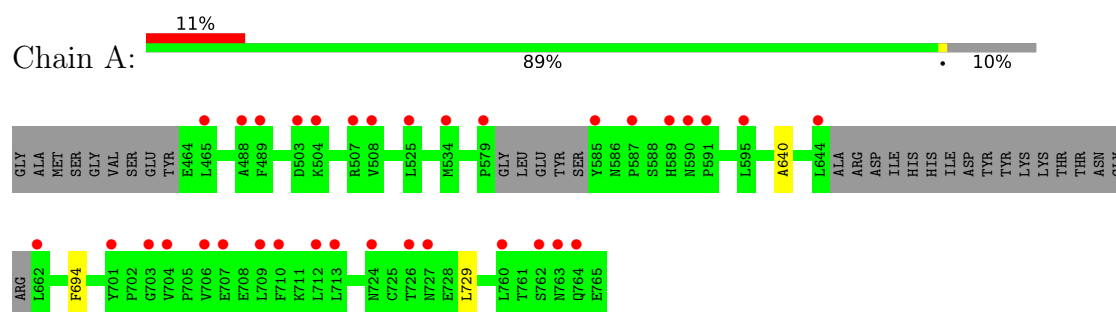
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total 7	O 7	0	0
4	B	7	Total 7	O 7	0	0

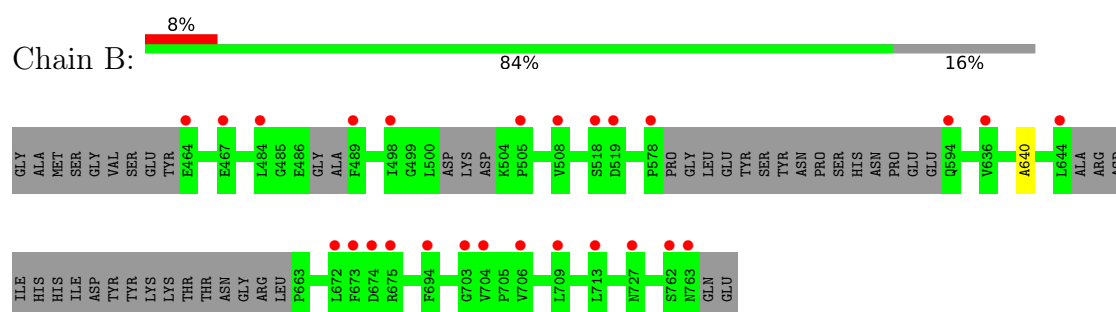
3 Residue-property plots

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: Fibroblast growth factor receptor 1



- Molecule 1: Fibroblast growth factor receptor 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	211.03Å 50.38Å 65.51Å 90.00° 106.24° 90.00°	Depositor
Resolution (Å)	101.30 – 2.29 101.30 – 2.29	Depositor EDS
% Data completeness (in resolution range)	97.6 (101.30-2.29) 97.9 (101.30-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.237 , 0.276 0.236 , 0.277	Depositor DCC
R_{free} test set	1573 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å ²)	52.1	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 30.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4171	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.81% 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: MK9, SO4

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	0.54	0/2165	0.76	0/2945
1	B	0.55	0/1989	0.79	0/2707
All	All	0.54	0/4154	0.78	0/5652

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

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	2122	0	2046	2	0
1	B	1952	0	1855	1	0
2	A	45	0	0	1	0
2	B	33	0	0	1	0
3	A	5	0	0	0	0
4	A	7	0	0	0	0
4	B	7	0	0	0	0
All	All	4171	0	3901	3	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 0.

All (3) 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:A:640:ALA:HB1	2:A:801:MK9:CLU	2.42	0.56
1:B:640:ALA:HB1	2:B:801:MK9:CLU	2.48	0.50
1:A:694:PHE:CZ	1:A:729:LEU:HD13	2.56	0.41

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	275/311 (88%)	270 (98%)	5 (2%)	0	100	100
1	B	252/311 (81%)	248 (98%)	4 (2%)	0	100	100
All	All	527/622 (85%)	518 (98%)	9 (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	218/271 (80%)	218 (100%)	0	100	100
1	B	196/271 (72%)	196 (100%)	0	100	100
All	All	414/542 (76%)	414 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	724	ASN
1	A	743	GLN
1	B	635	ASN
1	B	679	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](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	SO4	A	802	-	4,4,4	0.43	0	6,6,6	0.09	0
2	MK9	A	801	-	48,48,48	1.09	4 (8%)	55,66,66	1.32	5 (9%)
2	MK9	B	801	-	36,36,48	1.17	4 (11%)	40,51,66	1.44	6 (15%)

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	MK9	A	801	-	-	4/30/46/46	0/4/4/4
2	MK9	B	801	-	-	3/15/31/46	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	MK9	CAH-NAG	4.16	1.43	1.37
2	B	801	MK9	C6-C5	3.97	1.45	1.37
2	B	801	MK9	CAH-NAG	3.86	1.42	1.37
2	A	801	MK9	C6-C5	3.81	1.45	1.37
2	A	801	MK9	CAH-NAI	3.11	1.41	1.37
2	B	801	MK9	CAH-NAI	2.46	1.40	1.37
2	B	801	MK9	CAW-NAG	2.12	1.50	1.46
2	A	801	MK9	CAW-NAG	2.04	1.50	1.46

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	MK9	C4-NAG-CAH	-6.27	116.58	122.23
2	B	801	MK9	C4-NAG-CAH	-5.31	117.44	122.23
2	A	801	MK9	CAJ-NAI-CAH	-2.98	117.56	122.53
2	B	801	MK9	CAW-NAG-CAH	-2.89	114.97	116.75
2	B	801	MK9	CAJ-NAI-CAH	-2.60	118.20	122.53
2	A	801	MK9	C6-C5-C4	2.47	118.53	115.51
2	B	801	MK9	OAT-CAN-CAM	2.39	118.06	115.44
2	B	801	MK9	C6-C5-C4	2.33	118.36	115.51
2	A	801	MK9	C5-C6-N1	-2.20	120.25	123.83
2	B	801	MK9	C5-C6-N1	-2.05	120.49	123.83
2	A	801	MK9	OAT-CAN-CAM	2.01	117.64	115.44

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	MK9	CAM-CAN-OAT-CBS
2	B	801	MK9	CAM-CAN-OAT-CBS
2	A	801	MK9	CAO-CAN-OAT-CBS
2	B	801	MK9	CAO-CAN-OAT-CBS
2	A	801	MK9	N2-C7-CAZ-CAY

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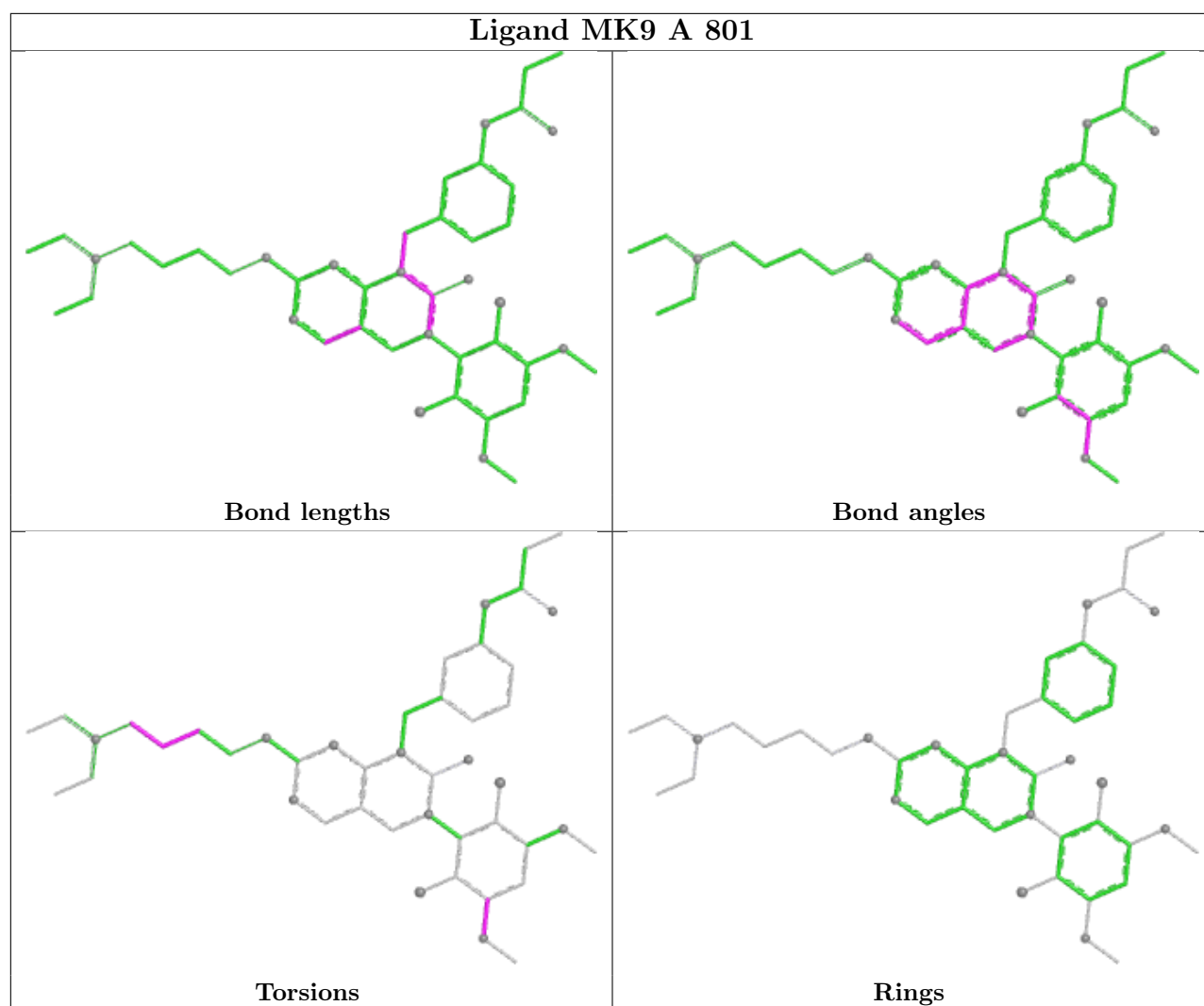
Mol	Chain	Res	Type	Atoms
2	A	801	MK9	CAX-CAY-CAZ-C7
2	B	801	MK9	CBG-CAW-NAG-CAH

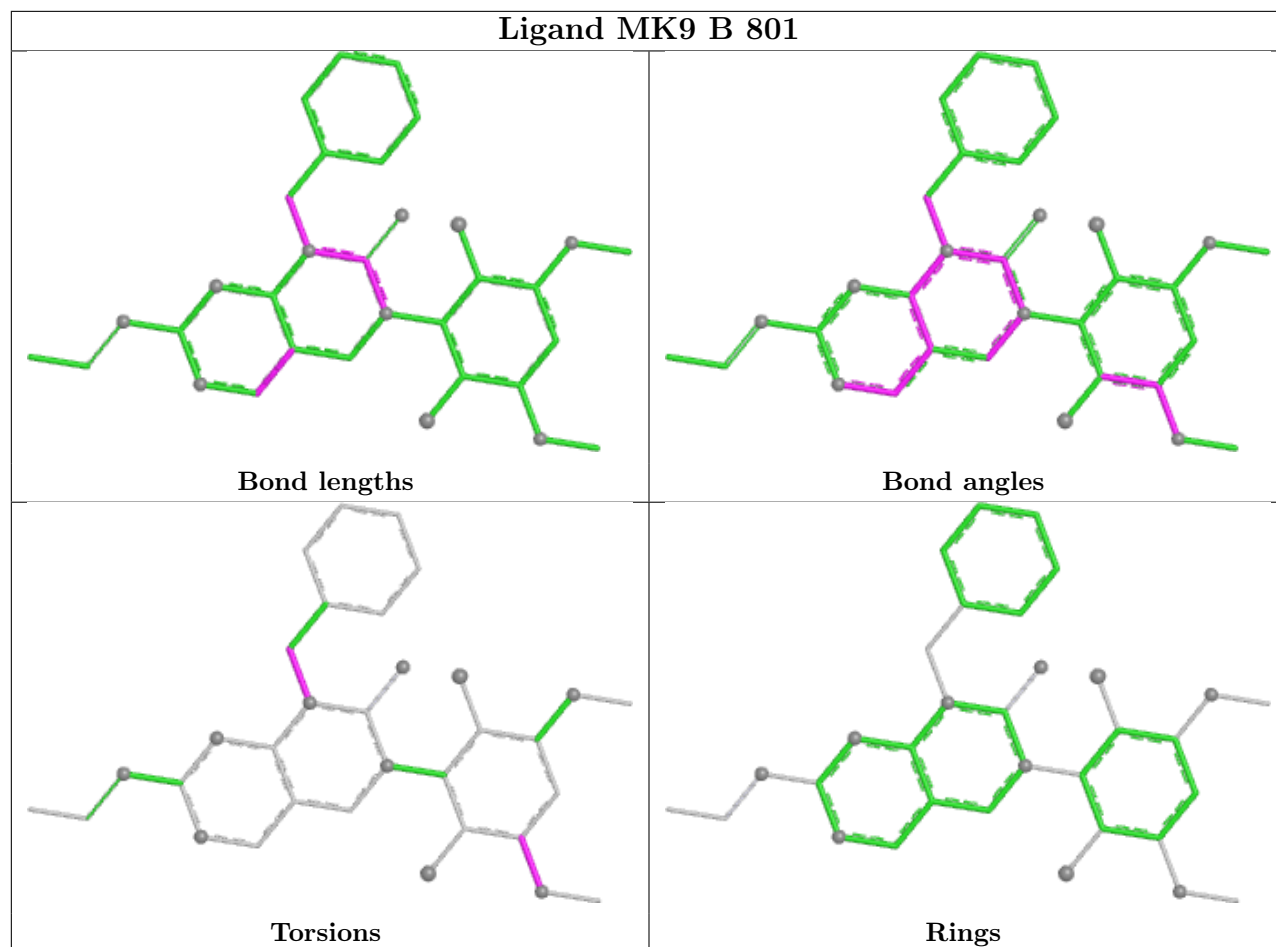
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	MK9	1	0
2	B	801	MK9	1	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	280/311 (90%)	0.75	34 (12%) 8 9	28, 62, 101, 111	1 (0%)
1	B	262/311 (84%)	0.75	26 (9%) 13 14	43, 64, 91, 106	0
All	All	542/622 (87%)	0.75	60 (11%) 10 11	28, 63, 97, 111	1 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	703	GLY	4.9
1	A	489	PHE	4.9
1	B	489	PHE	4.7
1	A	703	GLY	4.4
1	A	763	ASN	4.3
1	A	706	VAL	4.2
1	B	694	PHE	3.9
1	A	662	LEU	3.8
1	A	585	TYR	3.6
1	A	579	PRO	3.6
1	A	710	PHE	3.5
1	B	762	SER	3.5
1	B	464	GLU	3.3
1	B	763	ASN	3.2
1	B	519	ASP	3.1
1	B	644	LEU	3.1
1	A	712	LEU	2.9
1	A	508	VAL	2.9
1	B	578	PRO	2.8
1	A	727	ASN	2.8
1	B	518	SER	2.8
1	A	534	MET	2.8
1	B	706	VAL	2.8
1	B	673	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	726	THR	2.7
1	A	525	LEU	2.7
1	A	644	LEU	2.7
1	A	764	GLN	2.6
1	B	704	VAL	2.6
1	A	591	PRO	2.6
1	B	505	PRO	2.6
1	B	672	LEU	2.6
1	A	707	GLU	2.6
1	B	498	ILE	2.6
1	A	504	LYS	2.5
1	A	701	TYR	2.5
1	B	674	ASP	2.5
1	B	709	LEU	2.4
1	B	727	ASN	2.4
1	B	508	VAL	2.4
1	B	467	GLU	2.4
1	A	465	LEU	2.3
1	A	704	VAL	2.3
1	A	590	ASN	2.3
1	A	507	ARG	2.3
1	B	675	ARG	2.2
1	B	636	VAL	2.2
1	A	724	ASN	2.2
1	A	709	LEU	2.2
1	A	589	HIS	2.2
1	A	762	SER	2.2
1	B	484	LEU	2.2
1	A	488	ALA	2.2
1	A	595	LEU	2.1
1	A	760	LEU	2.1
1	A	503	ASP	2.1
1	A	713	LEU	2.1
1	A	587	PRO	2.0
1	B	594	GLN	2.0
1	B	713	LEU	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

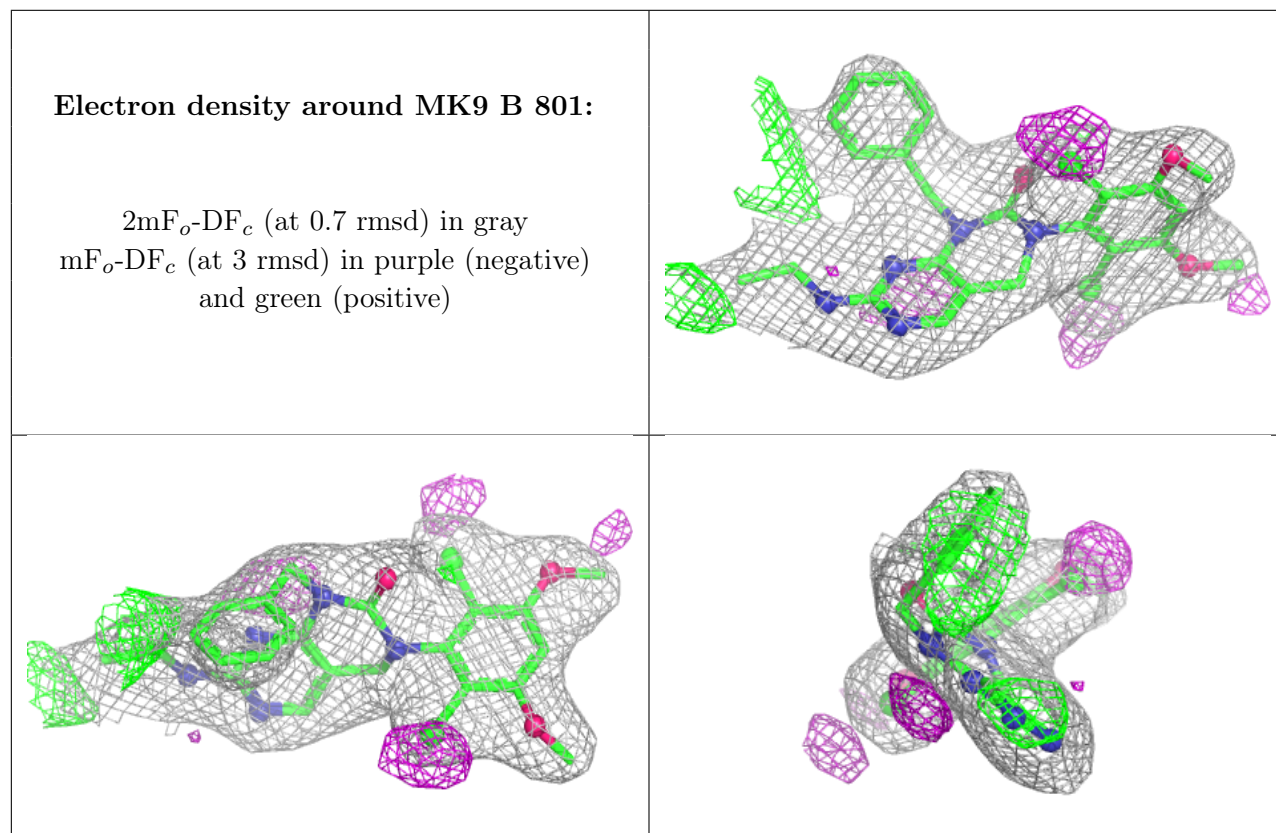
There are no oligosaccharides in this entry.

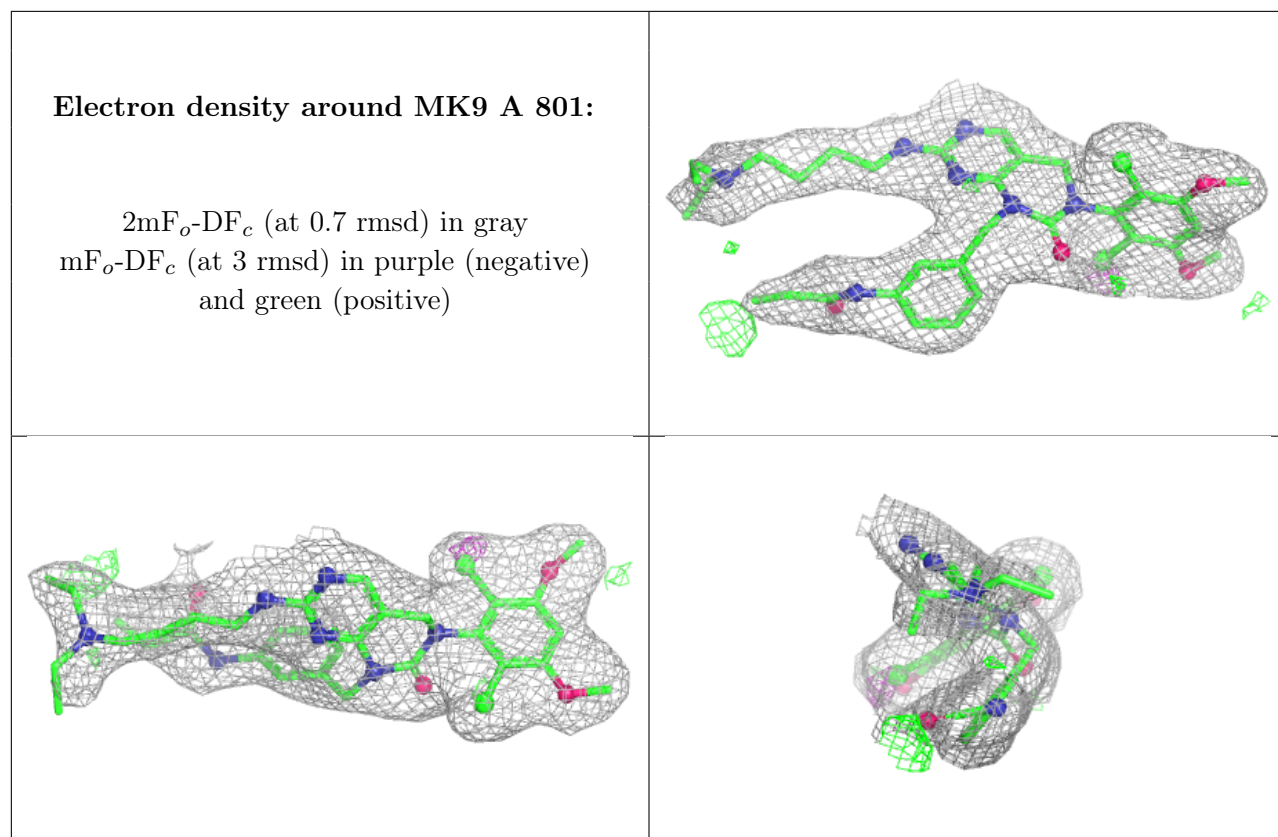
6.4 Ligands [i](#)

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
3	SO4	A	802	5/5	0.70	0.10	92,92,93,93	0
2	MK9	B	801	33/45	0.86	0.13	51,57,68,68	0
2	MK9	A	801	45/45	0.91	0.12	48,56,69,70	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.