



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:24 AM UTC

PDB ID : 4HEU / pdb_00004heu
Title : Crystal Structure of PDE10A with a biaryl ether inhibitor ((1-(3-(4-((1H-benzo[d]imidazol-2-yl)amino)phenoxy)pyridin-2-yl)piperidin-4-yl)methanol)
Authors : Chmait, S.; Jordan, S.
Deposited on : 2012-10-04
Resolution : 2.00 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

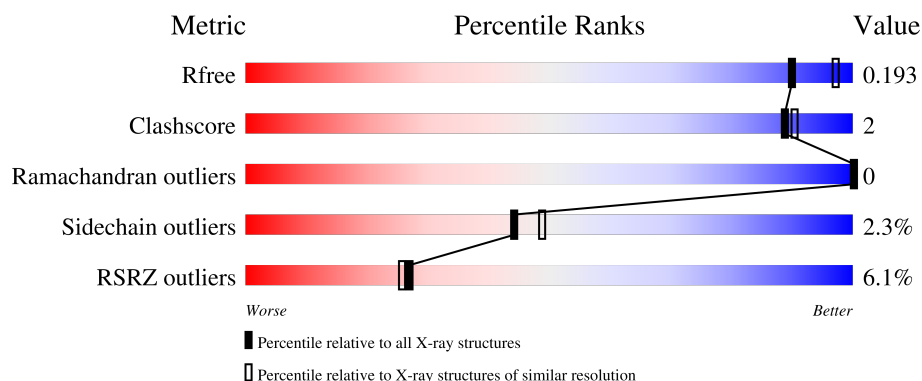
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	318	7% 94% 5%
1	B	318	5% 90% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	1004	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5607 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.

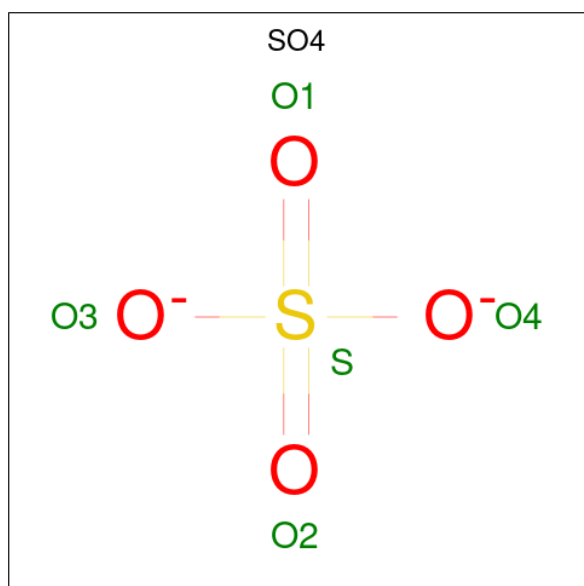
- Molecule 1 is a protein called cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2582	1650	439	470	23			
1	B	308	Total	C	N	O	S	0	0	0
			2499	1599	426	452	22			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

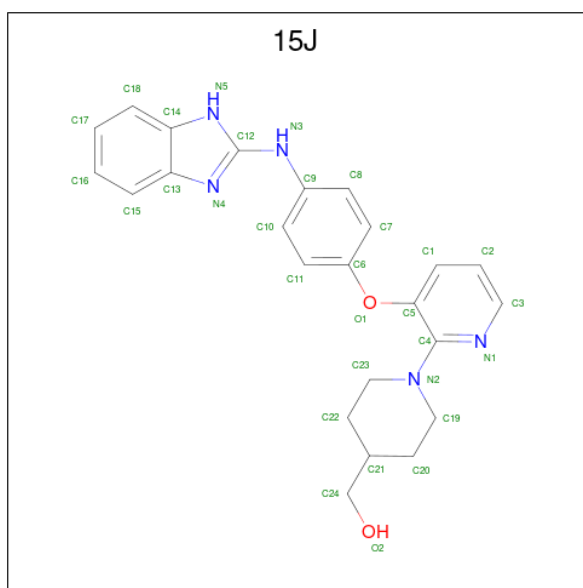
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is (1-{3-[4-(1H-benzimidazol-2-ylamino)phenoxy]pyridin-2-yl}piperidin-4-yl)methanol (CCD ID: 15J) (formula: C₂₄H₂₅N₅O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			31	24	5	2		
4	B	1	Total	C	N	O	0	0
			31	24	5	2		

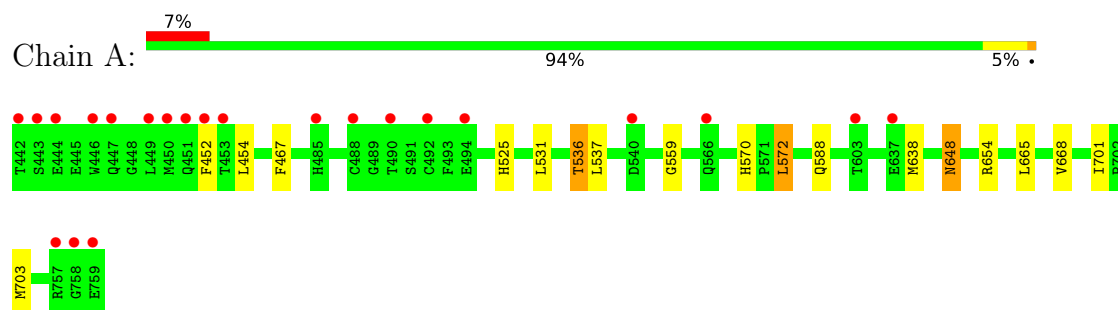
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	192	Total	O	0	0
			192	192		
5	B	208	Total	O	0	0
			208	208		

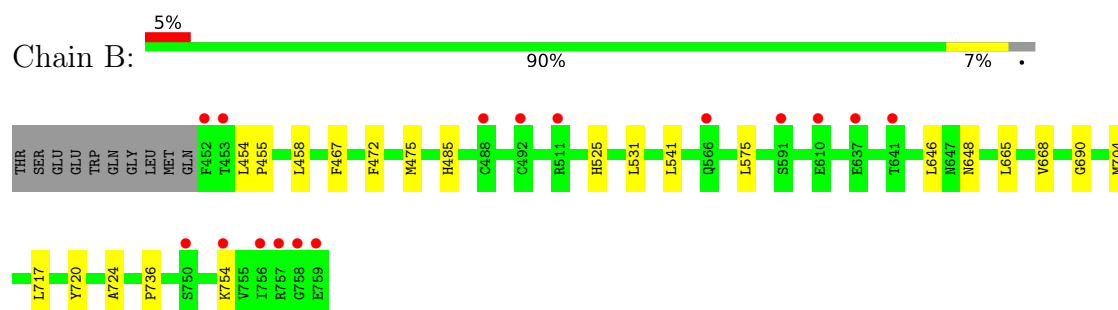
3 Residue-property plots [i](#)

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: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



- Molecule 1: cAMP and cAMP-inhibited cGMP 3',5'-cyclic phosphodiesterase 10A



4 Data and refinement statistics

Property	Value	Source
Space group	F 2 3	Depositor
Cell constants a, b, c, α , β , γ	252.86Å 252.86Å 252.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.66 – 2.00 48.66 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (48.66-2.00) 99.7 (48.66-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.186 , 0.199 (Not available) , 0.193	Depositor DCC
R_{free} test set	4524 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	32.3	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5607	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.98% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 15J, 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	0.69	0/2646	0.78	0/3583
1	B	0.69	1/2561 (0.0%)	0.78	1/3468 (0.0%)
All	All	0.69	1/5207 (0.0%)	0.78	1/7051 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	485	HIS	CG-ND1	-5.68	1.32	1.38

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	736	PRO	N-CA-C	5.07	116.88	110.70

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	2582	0	2543	11	0
1	B	2499	0	2470	8	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	30	0	0	1	0
3	B	30	0	0	0	0
4	A	31	0	25	1	0
4	B	31	0	25	1	0
5	A	192	0	0	1	0
5	B	208	0	0	0	0
All	All	5607	0	5063	19	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 (19) 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:536:THR:HG23	1:B:646:LEU:HG	1.72	0.71
1:B:472:PHE:HB2	1:B:475:MET:HE2	1.72	0.71
1:A:570:HIS:CE1	1:A:572:LEU:HD23	2.46	0.51
1:B:690:GLY:HA3	1:B:704:MET:O	2.12	0.50
1:B:467:PHE:HB3	1:B:525:HIS:CE1	2.49	0.48
1:B:720:TYR:HA	1:B:724:ALA:HB3	1.96	0.48
1:A:638:MET:HE1	1:A:654:ARG:HD2	1.97	0.46
1:A:648:ASN:HD22	1:A:648:ASN:C	2.23	0.46
1:A:536:THR:CG2	1:B:646:LEU:HG	2.43	0.45
1:A:701:ILE:HD11	1:A:703:MET:HE2	1.98	0.45
1:A:467:PHE:HB3	1:A:525:HIS:CE1	2.52	0.45
1:A:559:GLY:HA3	3:A:1005:SO4:O2	2.18	0.44
1:A:536:THR:HG22	1:A:537:LEU:HD23	1.99	0.43
1:A:588:GLN:NE2	5:A:1232:HOH:O	2.52	0.42
1:B:454:LEU:HA	1:B:455:PRO:HD3	1.94	0.42
1:B:665:LEU:O	1:B:668:VAL:HG22	2.19	0.42
1:A:665:LEU:O	1:A:668:VAL:HG22	2.20	0.42
4:B:809:15J:H17	4:B:809:15J:N5	2.35	0.41
4:A:1009:15J:N5	4:A:1009:15J:H17	2.37	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	316/318 (99%)	313 (99%)	3 (1%)	0	100	100
1	B	306/318 (96%)	304 (99%)	2 (1%)	0	100	100
All	All	622/636 (98%)	617 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	287/287 (100%)	281 (98%)	6 (2%)	47	52
1	B	278/287 (97%)	271 (98%)	7 (2%)	42	45
All	All	565/574 (98%)	552 (98%)	13 (2%)	44	49

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	452	PHE
1	A	454	LEU
1	A	531	LEU
1	A	536	THR
1	A	572	LEU
1	A	648	ASN
1	B	458	LEU
1	B	531	LEU

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Mol	Chain	Res	Type
1	B	541	LEU
1	B	575	LEU
1	B	648	ASN
1	B	717	LEU
1	B	754	LYS

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

Mol	Chain	Res	Type
1	A	648	ASN
1	A	733	GLN
1	B	583	GLN
1	B	588	GLN
1	B	611	GLN
1	B	634	GLN
1	B	640	GLN
1	B	648	ASN
1	B	733	GLN

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](#)

Of 18 ligands modelled in this entry, 4 are monoatomic - leaving 14 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
3	SO4	B	804	-	4,4,4	0.40	0	6,6,6	0.11	0
3	SO4	B	803	-	4,4,4	0.42	0	6,6,6	0.07	0
3	SO4	A	1003	-	4,4,4	0.42	0	6,6,6	0.11	0
3	SO4	A	1005	-	4,4,4	0.39	0	6,6,6	0.08	0
3	SO4	A	1008	-	4,4,4	0.43	0	6,6,6	0.09	0
3	SO4	A	1004	-	4,4,4	0.44	0	6,6,6	3.26	5 (83%)
4	15J	A	1009	-	35,35,35	1.21	4 (11%)	44,48,48	1.32	5 (11%)
3	SO4	B	807	-	4,4,4	0.49	0	6,6,6	3.28	3 (50%)
3	SO4	B	808	-	4,4,4	0.41	0	6,6,6	0.07	0
3	SO4	A	1006	-	4,4,4	0.52	0	6,6,6	3.34	3 (50%)
3	SO4	B	805	-	4,4,4	0.43	0	6,6,6	0.06	0
3	SO4	A	1007	-	4,4,4	0.42	0	6,6,6	0.09	0
4	15J	B	809	-	35,35,35	1.25	4 (11%)	44,48,48	1.30	5 (11%)
3	SO4	B	806	-	4,4,4	0.43	0	6,6,6	0.09	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
4	15J	B	809	-	-	2/14/24/24	0/5/5/5
4	15J	A	1009	-	-	2/14/24/24	0/5/5/5

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1009	15J	C9-N3	-3.34	1.34	1.41
4	B	809	15J	C9-N3	-3.31	1.35	1.41
4	B	809	15J	C13-N4	-3.18	1.33	1.39
4	A	1009	15J	C13-N4	-3.12	1.33	1.39
4	B	809	15J	C14-N5	-2.93	1.33	1.38
4	A	1009	15J	C14-N5	-2.73	1.33	1.38
4	B	809	15J	C23-N2	2.17	1.50	1.46
4	A	1009	15J	C23-N2	2.13	1.50	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1004	SO4	O4-S-O2	-4.64	85.31	109.56
3	A	1006	SO4	O4-S-O2	-4.58	85.62	109.56
3	B	807	SO4	O4-S-O2	-4.55	85.78	109.56
3	A	1006	SO4	O4-S-O1	-4.53	85.86	109.56
3	A	1004	SO4	O3-S-O2	-4.51	85.96	109.56
3	B	807	SO4	O4-S-O1	-4.35	86.82	109.56
4	A	1009	15J	C19-N2-C23	4.05	120.67	111.57
3	A	1006	SO4	O4-S-O3	-4.02	86.35	108.54
3	B	807	SO4	O4-S-O3	-3.92	86.94	108.54
4	B	809	15J	C19-N2-C23	3.78	120.07	111.57
4	B	809	15J	C5-C4-N1	-3.33	119.62	124.58
4	B	809	15J	C3-N1-C4	3.30	122.52	115.05
3	A	1004	SO4	O2-S-O1	-3.26	86.12	109.06
4	A	1009	15J	C3-N1-C4	3.18	122.24	115.05
4	A	1009	15J	C5-C4-N1	-3.15	119.89	124.58
4	A	1009	15J	C13-N4-C12	2.85	107.56	104.21
4	B	809	15J	C13-N4-C12	2.70	107.38	104.21
4	A	1009	15J	N5-C12-N4	-2.31	110.37	113.73
4	B	809	15J	N5-C12-N4	-2.17	110.59	113.73
3	A	1004	SO4	O4-S-O3	2.10	120.10	108.54
3	A	1004	SO4	O4-S-O1	2.03	120.15	109.56

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1009	15J	C5-C4-N2-C19
4	B	809	15J	C5-C4-N2-C19
4	A	1009	15J	N1-C4-N2-C19
4	B	809	15J	N1-C4-N2-C19

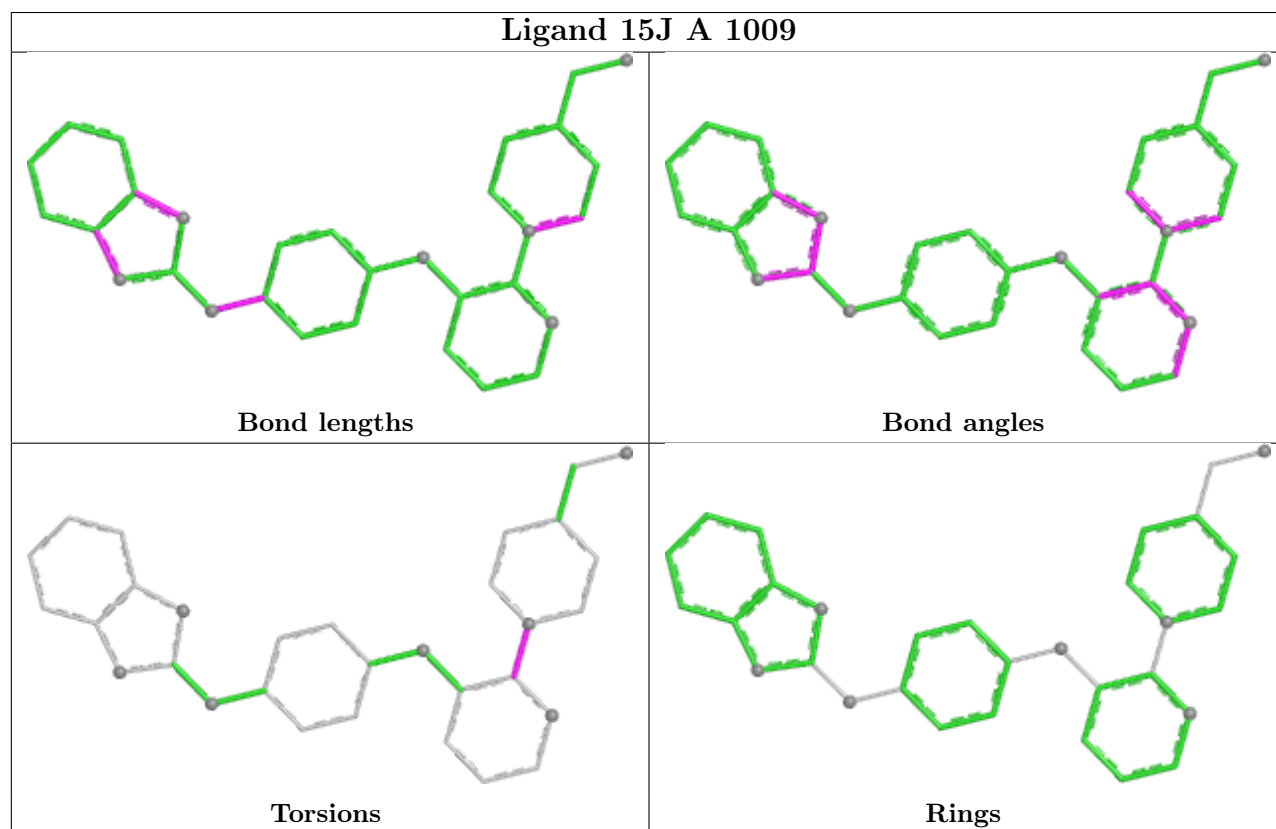
There are no ring outliers.

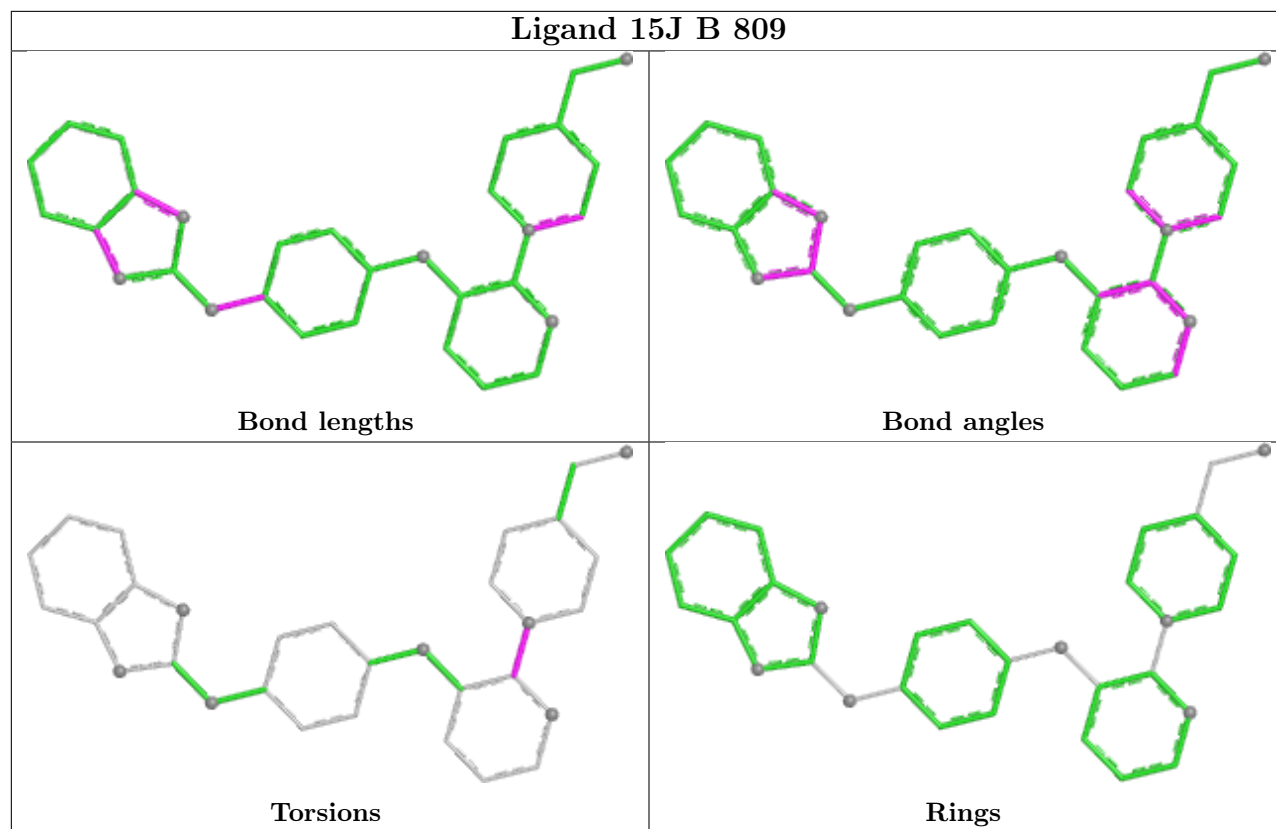
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1005	SO4	1	0
4	A	1009	15J	1	0
4	B	809	15J	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	318/318 (100%)	0.36	22 (6%)	23 21	24, 33, 51, 74	4 (1%)
1	B	308/318 (96%)	0.22	16 (5%)	33 31	24, 33, 47, 73	3 (0%)
All	All	626/636 (98%)	0.29	38 (6%)	27 26	24, 33, 48, 74	7 (1%)

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	452	PHE	12.9
1	A	446	TRP	12.2
1	A	758	GLY	5.9
1	B	488	CYS	5.2
1	B	758	GLY	5.1
1	B	452	PHE	5.1
1	A	449	LEU	4.9
1	B	453	THR	4.7
1	A	450	MET	4.6
1	A	490	THR	3.9
1	A	451	GLN	3.7
1	A	492	CYS	3.5
1	B	641	THR	3.5
1	A	488	CYS	3.5
1	B	759	GLU	3.4
1	A	443	SER	3.2
1	A	759	GLU	3.2
1	B	610	GLU	3.2
1	B	492	CYS	3.2
1	A	540	ASP	3.2
1	A	494	GLU	3.2
1	A	444	GLU	3.1
1	A	453	THR	2.9
1	B	566	GLN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	637	GLU	2.9
1	A	442	THR	2.7
1	B	754	LYS	2.7
1	B	750	SER	2.6
1	A	603	THR	2.6
1	A	485	HIS	2.6
1	A	447	GLN	2.4
1	A	637	GLU	2.4
1	A	566	GLN	2.2
1	B	591	SER	2.1
1	B	756	ILE	2.1
1	A	757	ARG	2.1
1	B	511	ARG	2.0
1	B	757	ARG	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 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	B	808	5/5	0.81	0.24	38,40,40,40	5
3	SO4	B	804	5/5	0.82	0.25	45,46,47,47	5
3	SO4	A	1003	5/5	0.86	0.17	46,48,49,50	5
3	SO4	A	1005	5/5	0.86	0.22	31,31,32,32	5
3	SO4	B	803	5/5	0.87	0.24	36,36,36,37	5
3	SO4	A	1006	5/5	0.89	0.18	75,76,78,80	0
3	SO4	B	805	5/5	0.90	0.17	52,53,54,54	5
3	SO4	A	1008	5/5	0.90	0.15	61,62,65,65	5
3	SO4	A	1007	5/5	0.92	0.20	55,56,56,56	5

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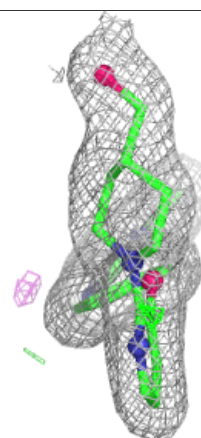
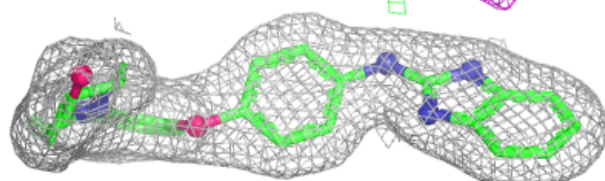
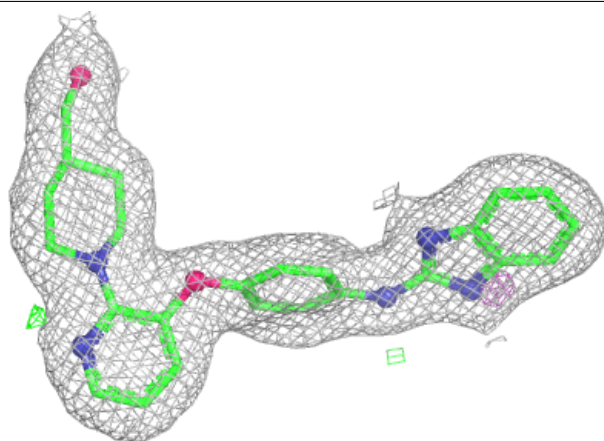
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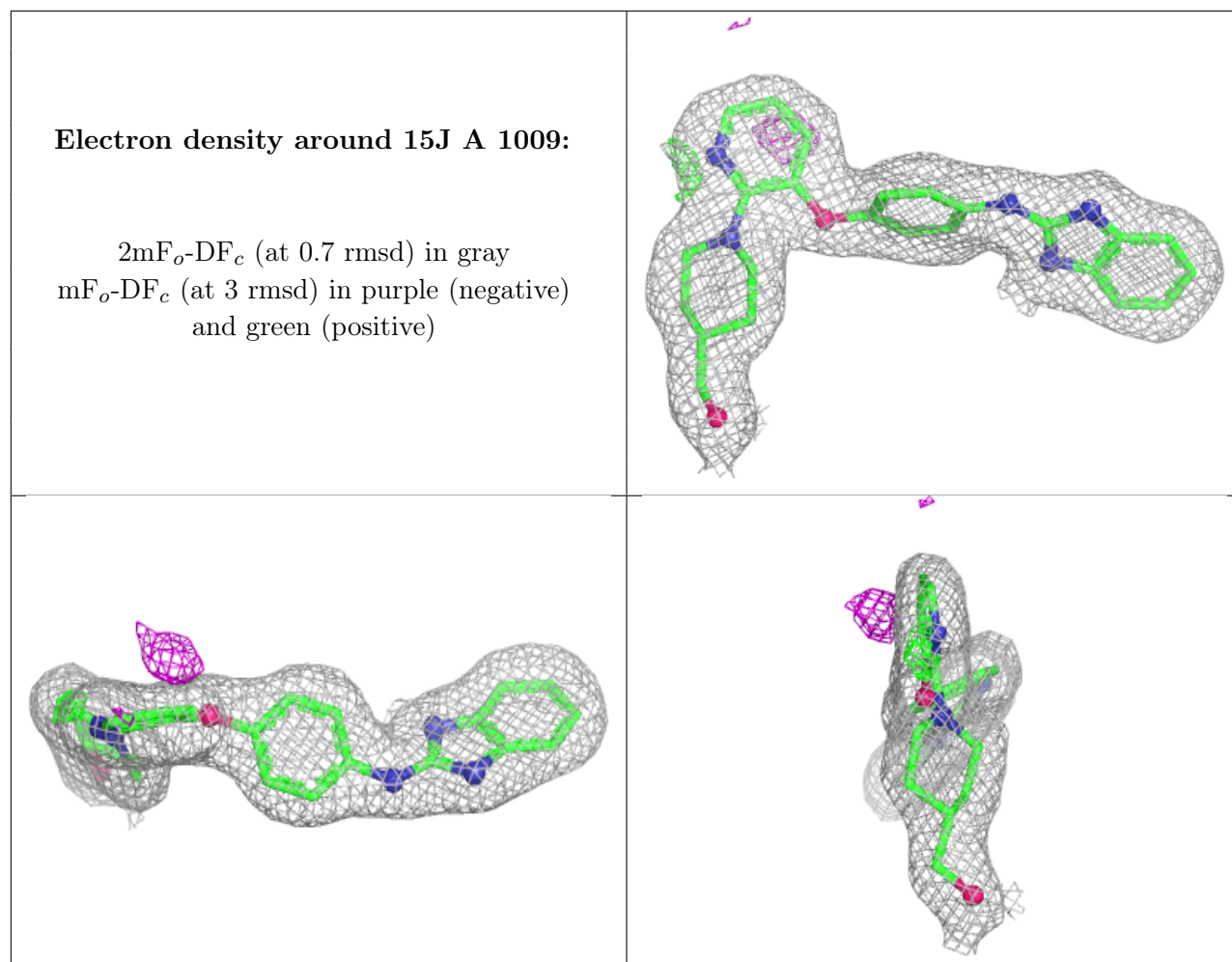
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	806	5/5	0.93	0.20	57,58,59,59	5
3	SO4	B	807	5/5	0.94	0.14	48,50,51,51	0
3	SO4	A	1004	5/5	0.94	0.09	53,54,55,55	0
4	15J	B	809	31/31	0.95	0.08	28,31,34,34	0
4	15J	A	1009	31/31	0.96	0.08	29,32,33,33	0
2	ZN	B	801	1/1	1.00	0.02	30,30,30,30	0
2	ZN	B	802	1/1	1.00	0.03	32,32,32,32	1
2	ZN	A	1001	1/1	1.00	0.02	28,28,28,28	0
2	ZN	A	1002	1/1	1.00	0.03	26,26,26,26	1

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 15J B 809:

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.