



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 27, 2023 – 09:41 AM EDT

PDB ID : 3IK3
Title : AP24534, a Pan-BCR-ABL Inhibitor for Chronic Myeloid Leukemia, Potently Inhibits the T315I Mutant and Overcomes Mutation-Based Resistance
Authors : Zhou, T.
Deposited on : 2009-08-05
Resolution : 1.90 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35
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) : 2.35

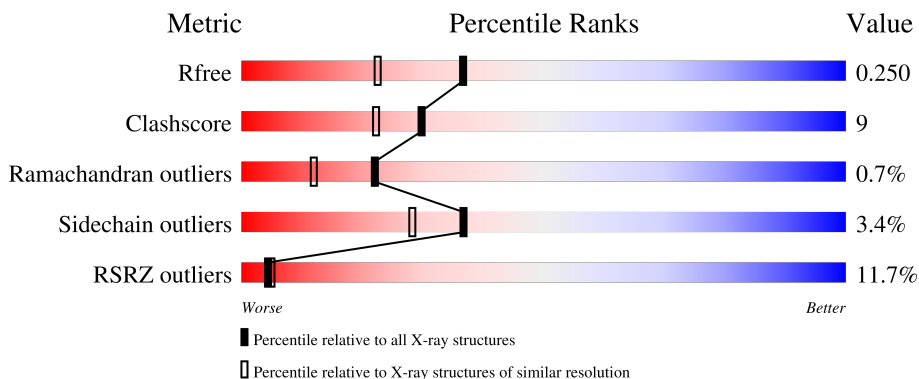
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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	288	 12% 82% 14% ..
1	B	288	 10% 80% 18% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5137 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.

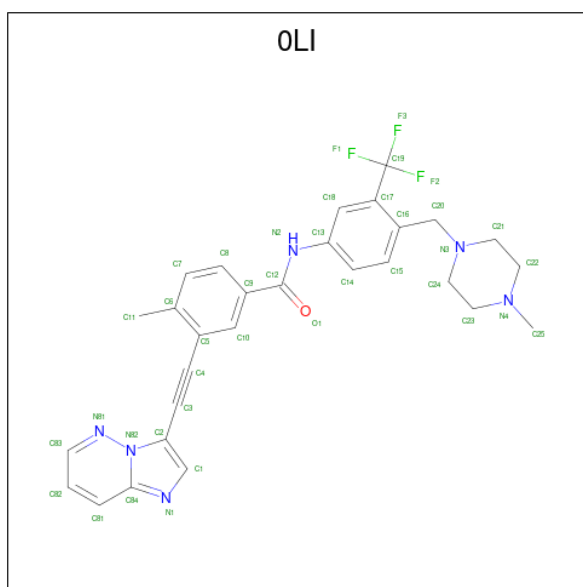
- Molecule 1 is a protein called Proto-oncogene tyrosine-protein kinase ABL1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	280	2277	1465	370	426	16	0	0	0
1	B	286	2323	1492	379	435	17	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	GLY	-	expression tag	UNP P00520
A	315	ILE	THR	engineered mutation	UNP P00520
A	514	GLY	-	expression tag	UNP P00520
A	515	THR	-	expression tag	UNP P00520
B	228	GLY	-	expression tag	UNP P00520
B	315	ILE	THR	engineered mutation	UNP P00520
B	514	GLY	-	expression tag	UNP P00520
B	515	THR	-	expression tag	UNP P00520

- Molecule 2 is 3-(imidazo[1,2-b]pyridazin-3-ylethynyl)-4-methyl-N-{4-[(4-methylpiperazin-1-yl)methyl]-3-(trifluoromethyl)phenyl}benzamide (three-letter code: OLI) (formula: C₂₉H₂₇F₃N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	A	1	Total	C	F	N	O	0	0
			39	29	3	6	1		
2	B	1	Total	C	F	N	O	0	0
			39	29	3	6	1		

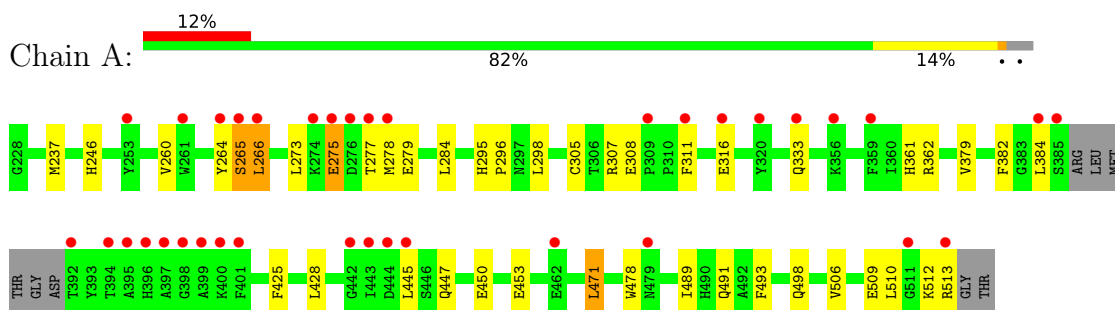
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	229	Total	O	0	0
			229	229		
3	B	230	Total	O	0	0
			230	230		

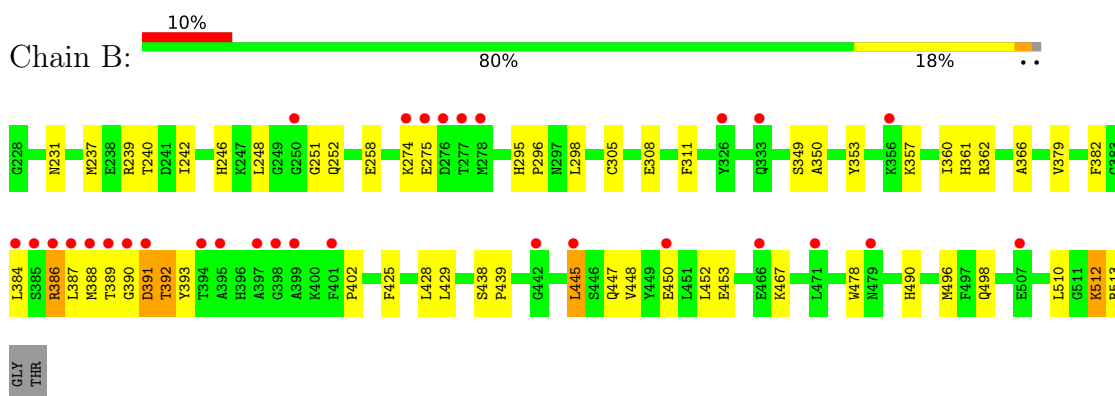
3 Residue-property plots [i](#)

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: Proto-oncogene tyrosine-protein kinase ABL1



- Molecule 1: Proto-oncogene tyrosine-protein kinase ABL1



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	43.91Å 131.45Å 60.20Å 90.00° 92.87° 90.00°	Depositor
Resolution (Å)	50.00 – 1.90 36.31 – 1.90	Depositor EDS
% Data completeness (in resolution range)	90.7 (50.00-1.90) 91.4 (36.31-1.90)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.94 (at 1.91Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.216 , 0.258 0.207 , 0.250	Depositor DCC
R_{free} test set	2536 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	13.5	Xtrriage
Anisotropy	0.829	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 55.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.166 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5137	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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: OLI

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	0.50	0/2334	0.60	0/3155
1	B	0.50	0/2381	0.61	0/3219
All	All	0.50	0/4715	0.61	0/6374

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	2277	0	2228	39	0
1	B	2323	0	2276	48	0
2	A	39	0	27	0	0
2	B	39	0	27	0	0
3	A	229	0	0	4	0
3	B	230	0	0	5	0
All	All	5137	0	4558	87	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 9.

All (87) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:386:ARG:HG2	1:B:386:ARG:HH21	1.19	1.05
1:B:390:GLY:O	1:B:391:ASP:HB2	1.71	0.88
1:B:386:ARG:HH21	1:B:386:ARG:CG	1.91	0.82
1:A:260:VAL:HG13	1:A:265:SER:O	1.84	0.77
1:B:386:ARG:HG2	1:B:386:ARG:NH2	1.98	0.76
1:B:467:LYS:HD3	1:B:496:MET:SD	2.29	0.72
1:B:510:LEU:HD23	1:B:513:ARG:NH1	2.06	0.70
1:A:445:LEU:H	1:A:445:LEU:HD12	1.56	0.70
1:A:506:VAL:O	1:A:510:LEU:HG	1.93	0.69
1:B:510:LEU:HD23	1:B:513:ARG:HH11	1.57	0.69
1:A:277:THR:HG22	1:A:279:GLU:H	1.57	0.68
1:A:361:HIS:O	1:A:362:ARG:HB2	1.94	0.67
1:A:471:LEU:HD21	1:A:489:ILE:HG23	1.76	0.67
1:A:471:LEU:HD22	1:A:493:PHE:CE2	2.31	0.65
1:A:512:LYS:O	1:A:513:ARG:CB	2.45	0.64
1:B:362:ARG:HD2	3:B:590:HOH:O	1.97	0.63
1:B:445:LEU:HD12	1:B:445:LEU:H	1.65	0.61
1:A:512:LYS:O	1:A:513:ARG:HB2	2.01	0.60
1:A:512:LYS:O	1:A:513:ARG:HG3	2.02	0.60
1:A:237:MET:HB2	3:A:566:HOH:O	2.02	0.59
1:A:445:LEU:HD12	1:A:445:LEU:N	2.17	0.59
1:B:298:LEU:HD23	1:B:379:VAL:HB	1.85	0.58
1:B:445:LEU:H	1:B:445:LEU:CD1	2.16	0.58
1:A:277:THR:HG22	1:A:279:GLU:N	2.17	0.58
1:B:361:HIS:O	1:B:362:ARG:HB2	2.03	0.58
1:A:277:THR:HG22	1:A:278:MET:N	2.20	0.56
1:B:251:GLY:O	1:B:387:LEU:HD21	2.06	0.56
1:A:445:LEU:H	1:A:445:LEU:CD1	2.17	0.56
1:A:275:GLU:HG2	3:A:577:HOH:O	2.05	0.56
1:B:386:ARG:CG	1:B:386:ARG:NH2	2.61	0.55
1:A:512:LYS:O	1:A:513:ARG:CG	2.56	0.54
1:B:248:LEU:HD21	1:B:258:GLU:HB2	1.87	0.54
1:A:471:LEU:CD2	1:A:489:ILE:HG23	2.37	0.53
1:B:447:GLN:HG2	3:B:572:HOH:O	2.08	0.53
1:B:448:VAL:O	1:B:452:LEU:HG	2.08	0.53
1:B:240:THR:HB	3:B:91:HOH:O	2.09	0.53
1:B:445:LEU:HD12	1:B:445:LEU:N	2.23	0.52
1:B:513:ARG:NH2	3:B:546:HOH:O	2.41	0.52
1:A:298:LEU:HD23	1:A:379:VAL:HB	1.92	0.52
1:B:382:PHE:HB2	1:B:384:LEU:HG	1.92	0.51
1:B:246:HIS:HE1	3:B:53:HOH:O	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:CYS:O	1:B:311:PHE:HA	2.11	0.51
1:B:390:GLY:O	1:B:391:ASP:CB	2.52	0.50
1:B:353:TYR:CE2	1:B:357:LYS:HD3	2.47	0.49
1:A:237:MET:CB	3:A:566:HOH:O	2.60	0.49
1:B:295:HIS:CG	1:B:296:PRO:HD2	2.48	0.49
1:B:353:TYR:CZ	1:B:357:LYS:HD3	2.47	0.49
1:B:512:LYS:N	1:B:512:LYS:HD2	2.27	0.48
1:B:392:THR:O	1:B:402:PRO:HA	2.14	0.47
1:B:389:THR:C	1:B:391:ASP:H	2.17	0.47
1:A:277:THR:CG2	1:A:278:MET:N	2.77	0.47
1:A:471:LEU:CD2	1:A:493:PHE:HE2	2.28	0.46
1:B:389:THR:OG1	1:B:390:GLY:N	2.49	0.46
1:B:240:THR:O	1:B:240:THR:HG22	2.15	0.46
1:B:391:ASP:O	1:B:393:TYR:N	2.47	0.46
1:B:239:ARG:HH22	1:B:308:GLU:HG3	1.81	0.45
1:B:246:HIS:H	1:B:246:HIS:CD2	2.33	0.45
1:A:509:GLU:O	1:A:512:LYS:O	2.35	0.45
1:B:237:MET:HE1	1:B:242:ILE:HD11	1.99	0.45
1:B:298:LEU:HG	1:B:350:ALA:HB1	1.98	0.45
1:B:239:ARG:NH2	1:B:308:GLU:OE2	2.50	0.45
1:A:265:SER:O	1:A:266:LEU:O	2.35	0.44
1:A:453:GLU:HA	1:A:478:TRP:CZ3	2.53	0.43
1:A:471:LEU:CD2	1:A:493:PHE:CE2	3.01	0.43
1:A:264:TYR:O	1:A:265:SER:HB2	2.19	0.43
1:A:275:GLU:HG2	1:A:275:GLU:H	1.53	0.43
1:B:349:SER:HB2	1:B:490:HIS:CE1	2.54	0.43
1:B:425:PHE:CE2	1:B:429:LEU:HD11	2.54	0.43
1:A:450:GLU:HG3	3:A:59:HOH:O	2.20	0.42
1:A:498:GLN:OE1	1:A:498:GLN:HA	2.20	0.42
1:A:265:SER:O	1:A:266:LEU:C	2.58	0.41
1:A:295:HIS:CG	1:A:296:PRO:HD2	2.55	0.41
1:B:389:THR:O	1:B:391:ASP:N	2.49	0.41
1:B:453:GLU:HA	1:B:478:TRP:CZ3	2.56	0.41
1:A:305:CYS:O	1:A:311:PHE:HA	2.20	0.41
1:A:509:GLU:OE2	1:A:509:GLU:HA	2.20	0.41
1:A:382:PHE:HB2	1:A:384:LEU:HG	2.02	0.41
1:B:252:GLN:C	1:B:387:LEU:HD21	2.42	0.41
1:B:360:ILE:HG22	1:B:362:ARG:HG3	2.02	0.41
1:B:438:SER:HA	1:B:439:PRO:HD3	1.92	0.41
1:A:425:PHE:CE2	1:A:489:ILE:HG21	2.55	0.41
1:B:498:GLN:OE1	1:B:498:GLN:HA	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:HIS:H	1:A:246:HIS:CD2	2.39	0.40
1:B:366:ALA:N	1:B:428:LEU:HD13	2.35	0.40
1:B:384:LEU:C	1:B:386:ARG:H	2.25	0.40
1:A:273:LEU:HD11	1:A:277:THR:HG21	2.04	0.40
1:A:307:ARG:HH11	1:A:307:ARG:HD2	1.74	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	276/288 (96%)	263 (95%)	11 (4%)	2 (1%)	22	12
1	B	284/288 (99%)	267 (94%)	15 (5%)	2 (1%)	22	12
All	All	560/576 (97%)	530 (95%)	26 (5%)	4 (1%)	22	12

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	266	LEU
1	B	391	ASP
1	B	392	THR
1	A	265	SER

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	245/251 (98%)	236 (96%)	9 (4%)	34	25
1	B	250/251 (100%)	242 (97%)	8 (3%)	39	30
All	All	495/502 (99%)	478 (97%)	17 (3%)	37	28

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	275	GLU
1	A	284	LEU
1	A	308	GLU
1	A	316	GLU
1	A	333	GLN
1	A	428	LEU
1	A	447	GLN
1	A	471	LEU
1	A	491	GLN
1	B	231	ASN
1	B	274	LYS
1	B	275	GLU
1	B	386	ARG
1	B	388	MET
1	B	445	LEU
1	B	450	GLU
1	B	512	LYS

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

Mol	Chain	Res	Type
1	A	231	ASN
1	A	246	HIS
1	A	333	GLN
1	A	375	HIS
1	A	414	ASN
1	A	490	HIS
1	A	491	GLN
1	B	231	ASN
1	B	246	HIS
1	B	252	GLN
1	B	375	HIS
1	B	490	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 monosaccharides 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
2	0LI	B	2	-	40,43,43	1.86	12 (30%)	51,62,62	0.92	0
2	0LI	A	1	-	40,43,43	1.78	9 (22%)	51,62,62	0.93	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
2	0LI	B	2	-	-	2/20/33/33	0/5/5/5
2	0LI	A	1	-	-	2/20/33/33	0/5/5/5

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	0LI	C16-C17	4.36	1.46	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	OLI	C8-C7	4.20	1.46	1.38
2	B	2	OLI	C16-C17	4.14	1.45	1.40
2	A	1	OLI	C8-C9	3.70	1.45	1.39
2	B	2	OLI	C8-C7	3.60	1.45	1.38
2	B	2	OLI	C18-C13	3.51	1.45	1.39
2	B	2	OLI	C18-C17	3.14	1.44	1.39
2	A	1	OLI	C10-C9	3.05	1.43	1.39
2	B	2	OLI	C10-C9	3.02	1.43	1.39
2	B	2	OLI	C14-C15	2.94	1.44	1.38
2	B	2	OLI	C20-C16	2.82	1.56	1.51
2	A	1	OLI	C18-C17	2.80	1.43	1.39
2	B	2	OLI	C8-C9	2.80	1.44	1.39
2	A	1	OLI	C5-C6	2.72	1.45	1.40
2	A	1	OLI	C18-C13	2.67	1.43	1.39
2	B	2	OLI	C5-C6	2.39	1.44	1.40
2	A	1	OLI	C14-C15	2.27	1.42	1.38
2	B	2	OLI	C81-C84	2.15	1.43	1.40
2	A	1	OLI	C82-C83	2.10	1.43	1.37
2	B	2	OLI	C15-C16	2.03	1.43	1.39
2	B	2	OLI	C23-N4	2.00	1.50	1.46

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

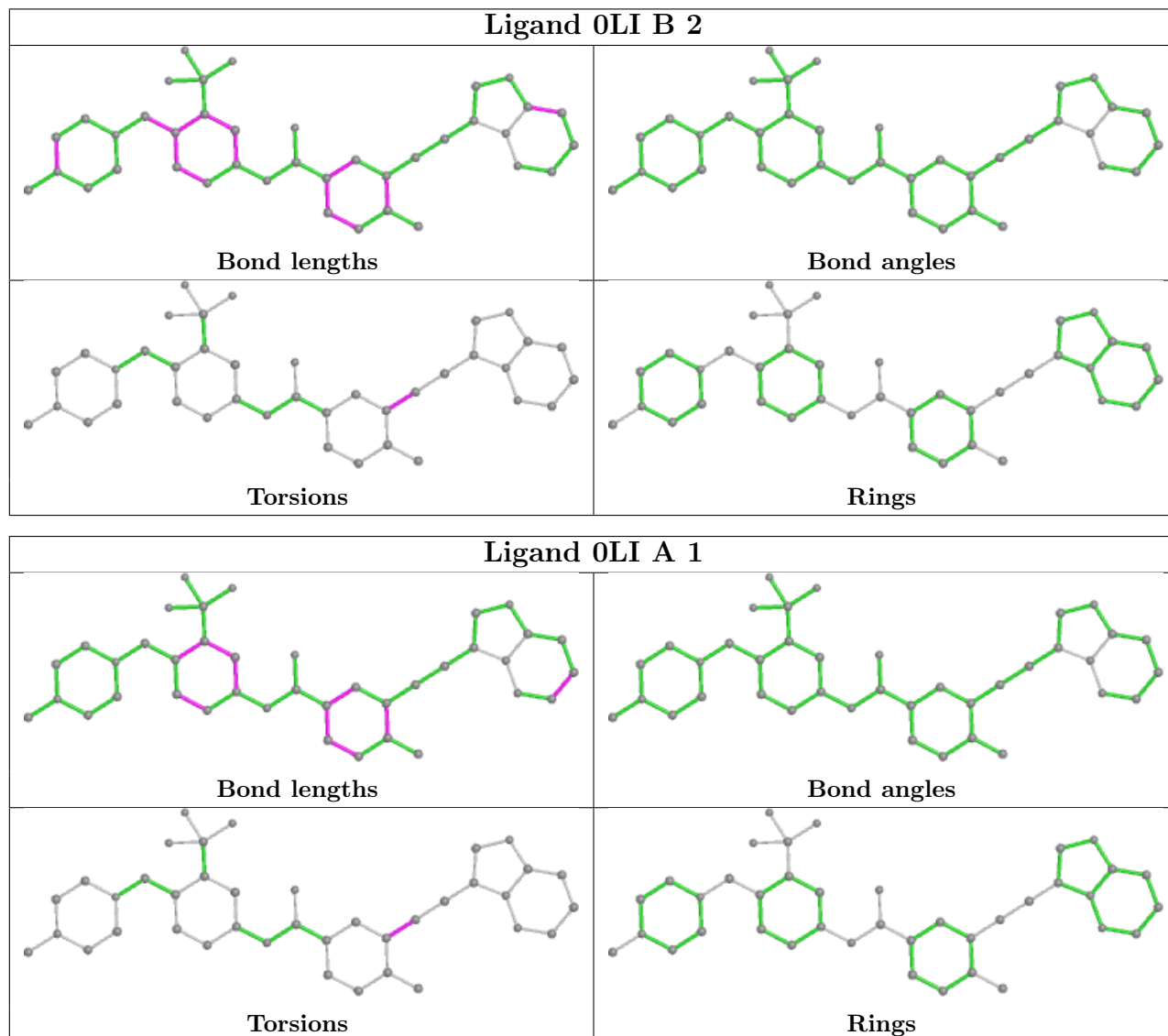
Mol	Chain	Res	Type	Atoms
2	A	1	OLI	C3-C4-C5-C6
2	B	2	OLI	C3-C4-C5-C6
2	A	1	OLI	C3-C4-C5-C10
2	B	2	OLI	C3-C4-C5-C10

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, 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	280/288 (97%)	0.92	36 (12%) 3 3	8, 16, 35, 47	0
1	B	286/288 (99%)	0.89	30 (10%) 6 7	7, 16, 37, 53	0
All	All	566/576 (98%)	0.90	66 (11%) 4 5	7, 16, 36, 53	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	390	GLY	8.0
1	B	276	ASP	7.9
1	B	387	LEU	5.7
1	A	392	THR	5.3
1	B	386	ARG	5.1
1	B	278	MET	4.9
1	A	276	ASP	4.8
1	A	277	THR	4.8
1	B	391	ASP	4.7
1	B	389	THR	4.6
1	A	278	MET	4.2
1	A	401	PHE	4.1
1	A	397	ALA	4.0
1	B	398	GLY	4.0
1	A	384	LEU	4.0
1	A	333	GLN	3.9
1	A	398	GLY	3.6
1	A	399	ALA	3.6
1	A	445	LEU	3.5
1	B	275	GLU	3.4
1	A	274	LYS	3.4
1	B	507	GLU	3.2
1	A	395	ALA	3.2
1	A	385	SER	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	513	ARG	3.1
1	A	394	THR	3.1
1	B	442	GLY	3.0
1	B	445	LEU	2.9
1	B	333	GLN	2.9
1	B	399	ALA	2.9
1	A	511	GLY	2.9
1	A	443	ILE	2.8
1	A	265	SER	2.8
1	B	277	THR	2.8
1	A	266	LEU	2.8
1	A	261	TRP	2.8
1	A	444	ASP	2.8
1	A	479	ASN	2.8
1	A	316	GLU	2.7
1	B	274	LYS	2.7
1	B	466	GLU	2.7
1	B	401	PHE	2.6
1	A	275	GLU	2.6
1	A	442	GLY	2.6
1	B	384	LEU	2.5
1	B	385	SER	2.5
1	A	359	PHE	2.5
1	A	253	TYR	2.4
1	B	250	GLY	2.4
1	B	471	LEU	2.4
1	B	397	ALA	2.4
1	B	450	GLU	2.4
1	A	462	GLU	2.4
1	B	326	TYR	2.2
1	A	309	PRO	2.2
1	B	388	MET	2.1
1	A	311	PHE	2.1
1	A	320	TYR	2.1
1	B	356	LYS	2.1
1	B	394	THR	2.1
1	B	395	ALA	2.1
1	A	396	HIS	2.1
1	A	264	TYR	2.1
1	A	400	LYS	2.0
1	B	479	ASN	2.0
1	A	356	LYS	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 monosaccharides in this entry.

6.4 Ligands [i](#)

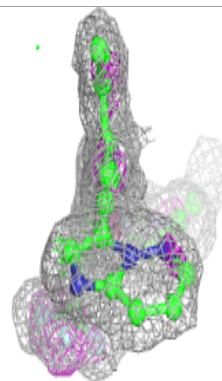
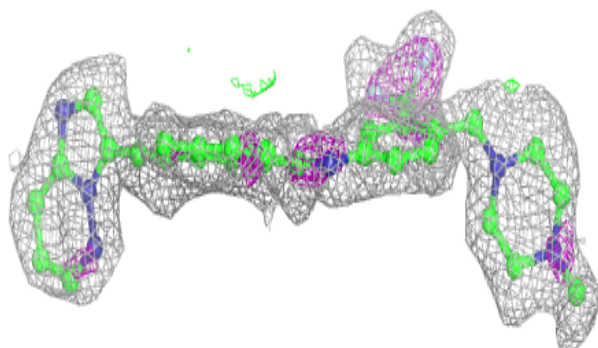
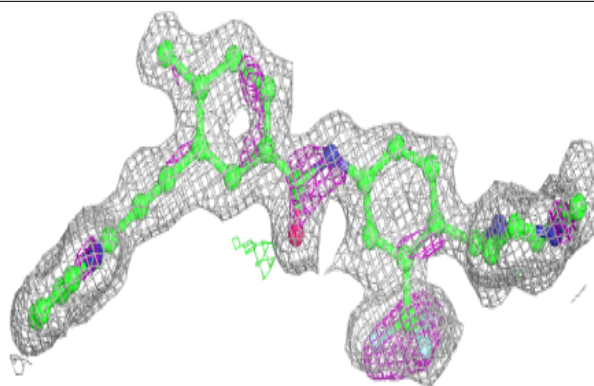
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	OLI	A	1	39/39	0.88	0.15	7,12,17,18	0
2	OLI	B	2	39/39	0.88	0.15	8,12,16,18	0

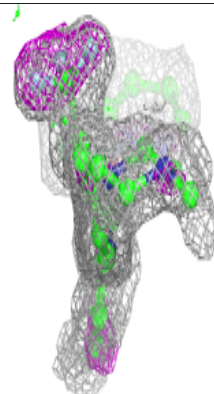
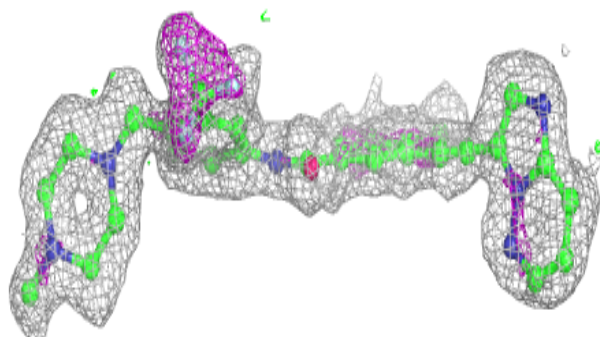
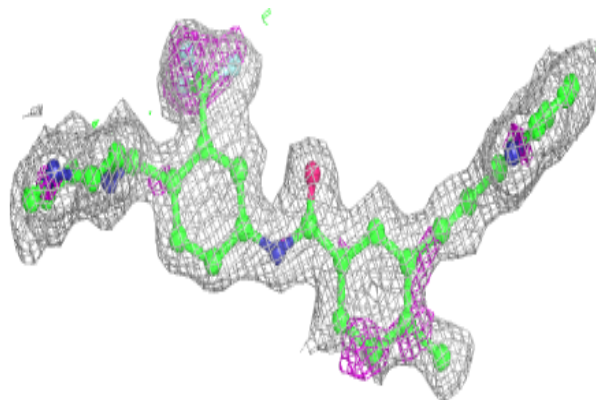
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 0LI A 1:

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

**Electron density around 0LI B 2:**

$2mF_o-DF_c$ (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.