



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 04:16 AM UTC

PDB ID : 4JPS / pdb_00004jps
Title : Co-crystal Structures of the Lipid Kinase PI3K alpha with Pan and Isoform Selective Inhibitors
Authors : Knapp, M.S.; Elling, R.A.
Deposited on : 2013-03-19
Resolution : 2.20 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

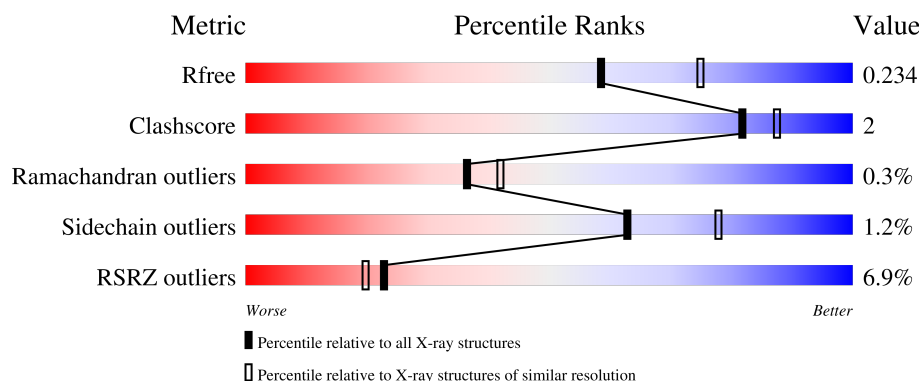
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	1074	5% 85% 8% 7%
2	B	293	9% 69% 8% 23%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10533 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 Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit alpha isoform.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	999	Total	C	N	O	S	0	6	0
			8111	5179	1393	1474	65			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	232	LYS	MET	engineered mutation	UNP P42336
A	233	LYS	LEU	engineered mutation	UNP P42336
A	1069	HIS	-	expression tag	UNP P42336
A	1070	HIS	-	expression tag	UNP P42336
A	1071	HIS	-	expression tag	UNP P42336
A	1072	HIS	-	expression tag	UNP P42336
A	1073	HIS	-	expression tag	UNP P42336
A	1074	HIS	-	expression tag	UNP P42336

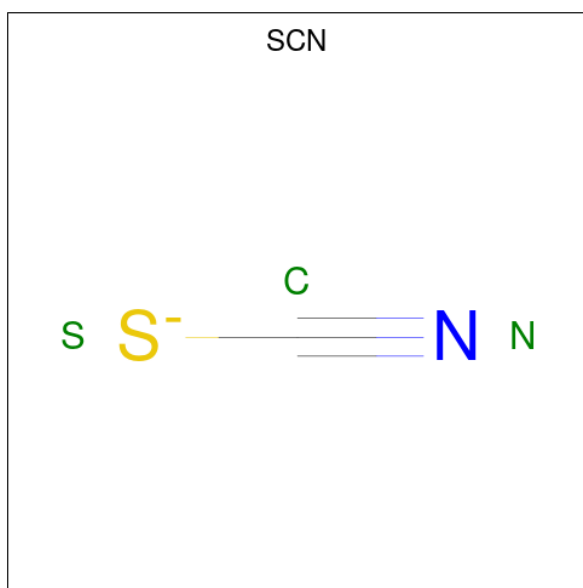
- Molecule 2 is a protein called Phosphatidylinositol 3-kinase regulatory subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	227	Total	C	N	O	S	0	1	0
			1883	1175	337	367	4			

There are 6 discrepancies between the modelled and reference sequences:

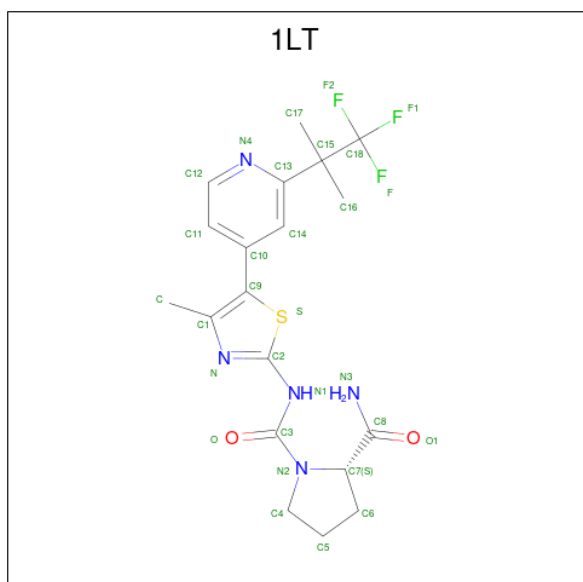
Chain	Residue	Modelled	Actual	Comment	Reference
B	301	MET	-	expression tag	UNP P27986
B	302	HIS	-	expression tag	UNP P27986
B	303	ASN	-	expression tag	UNP P27986
B	304	LEU	-	expression tag	UNP P27986
B	305	GLN	-	expression tag	UNP P27986
B	306	TYR	-	expression tag	UNP P27986

- Molecule 3 is THIOCYANATE ION (CCD ID: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	S	0	0
			3	1	1	1		

- Molecule 4 is (2S)-N 1 -{4-methyl-5-[2-(1,1,1-trifluoro-2-methylpropan-2-yl)pyridin-4-yl]-1,3-thiazol-2-yl}pyrrolidine-1,2-dicarboxamide (CCD ID: 1LT) (formula: C₁₉H₂₂F₃N₅O₂S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	S	0	0
			30	19	3	5	2	1		

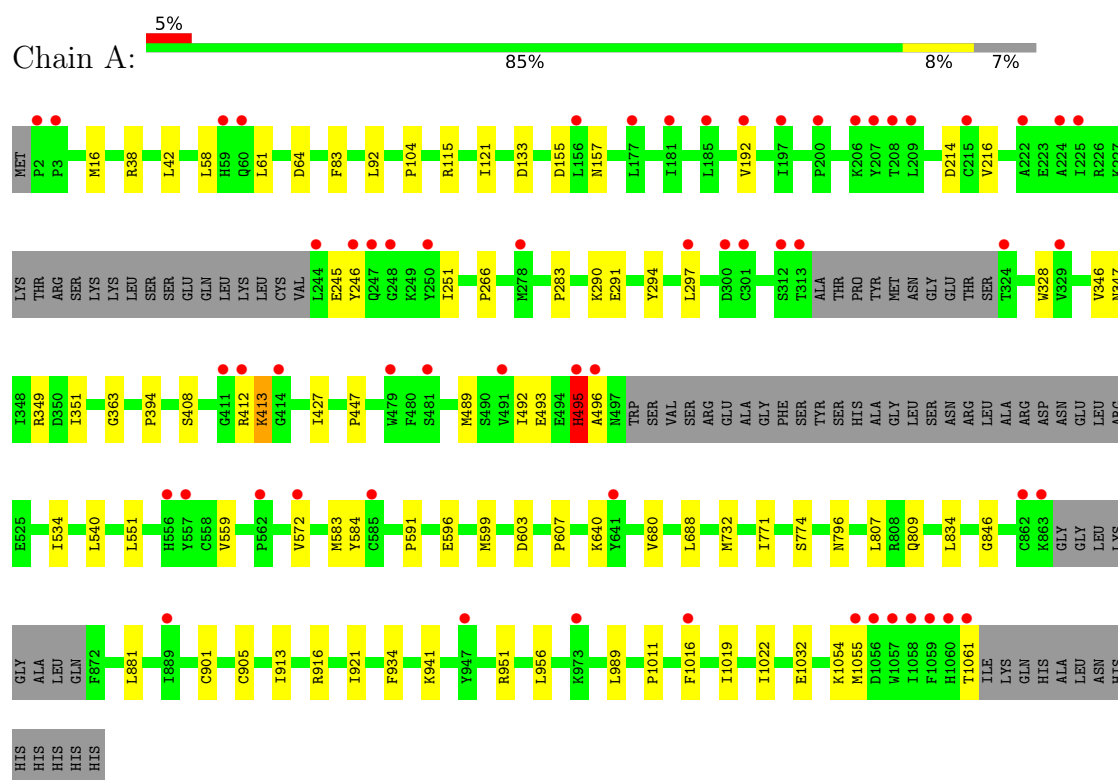
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	439	Total 439	O 439	0	0
5	B	67	Total 67	O 67	0	0

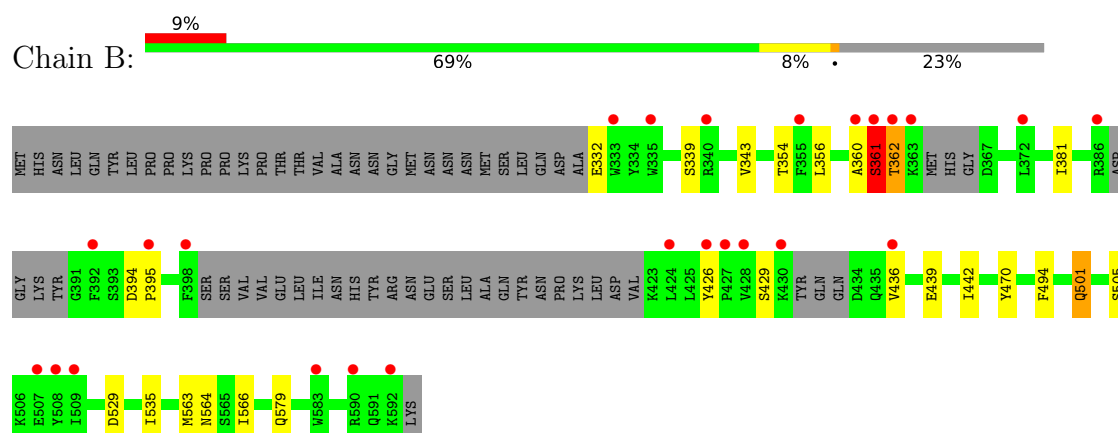
3 Residue-property plots

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: Phosphatidylinositol 4,5-bisphosphate 3-kinase catalytic subunit alpha isoform



- Molecule 2: Phosphatidylinositol 3-kinase regulatory subunit alpha



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.86Å 106.19Å 133.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.6 (50.00-2.20) 96.6 (50.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.00Å)	Xtriage
Refinement program	BUSTER-TNT, BUSTER 2.11.4	Depositor
R, R_{free}	0.205 , 0.228 0.208 , 0.234	Depositor DCC
R_{free} test set	4902 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	41.4	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10533	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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: SCN, 1LT

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	0.76	1/8299 (0.0%)	1.23	18/11236 (0.2%)
2	B	0.73	0/1909	1.40	7/2560 (0.3%)
All	All	0.76	1/10208 (0.0%)	1.26	25/13796 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	216	VAL	CA-C	5.30	1.58	1.52

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	361	SER	CA-C-N	8.20	137.21	121.54
2	B	361	SER	C-N-CA	8.20	137.21	121.54
1	A	495	HIS	CA-C-N	7.32	133.28	122.74
1	A	495	HIS	C-N-CA	7.32	133.28	122.74
1	A	603	ASP	CA-CB-CG	6.42	119.02	112.60
1	A	1054	LYS	CA-C-N	5.83	132.69	121.54
1	A	1054	LYS	C-N-CA	5.83	132.69	121.54
1	A	246	TYR	N-CA-C	5.80	117.27	111.07
1	A	245	GLU	CA-C-N	5.61	127.73	120.44
1	A	245	GLU	C-N-CA	5.61	127.73	120.44
1	A	796	ASN	CA-CB-CG	5.60	118.20	112.60
2	B	439	GLU	CA-C-N	5.60	128.24	120.63
2	B	439	GLU	C-N-CA	5.60	128.24	120.63
1	A	155	ASP	CA-C-N	5.52	127.62	120.44
1	A	155	ASP	C-N-CA	5.52	127.62	120.44
2	B	529	ASP	CA-CB-CG	5.51	118.11	112.60
1	A	157	ASN	CA-C-N	5.48	128.27	120.49
1	A	157	ASN	C-N-CA	5.48	128.27	120.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	64	ASP	CA-C-N	5.41	127.80	120.38
1	A	64	ASP	C-N-CA	5.41	127.80	120.38
1	A	133	ASP	CA-CB-CG	5.30	117.90	112.60
1	A	496	ALA	N-CA-C	5.18	118.89	112.58
1	A	934	PHE	CA-CB-CG	-5.04	108.76	113.80
2	B	339	SER	CA-C-N	5.03	127.03	120.28
2	B	339	SER	C-N-CA	5.03	127.03	120.28

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	8111	0	8001	37	0
2	B	1883	0	1796	11	0
3	A	3	0	0	0	0
4	A	30	0	22	2	0
5	A	439	0	0	1	0
5	B	67	0	0	1	0
All	All	10533	0	9819	49	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 (49) 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:436:VAL:HB	2:B:579:GLN:HG2	1.68	0.75
1:A:121:ILE:HG12	1:A:688:LEU:HB3	1.71	0.71
2:B:394:ASP:HB3	2:B:395:PRO:HD3	1.76	0.68
1:A:807:LEU:HD12	1:A:846:GLY:HA3	1.81	0.62
2:B:332:GLU:N	2:B:429:SER:HG	1.99	0.60
1:A:42:LEU:HD21	1:A:92:LEU:HD11	1.85	0.58
1:A:58:LEU:HB3	1:A:61:LEU:HD12	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:732:MET:HE1	1:A:771:ILE:H	1.71	0.54
1:A:251:ILE:HG23	1:A:290:LYS:HG2	1.88	0.54
1:A:294:TYR:HA	1:A:297:LEU:HD12	1.89	0.54
1:A:916:ARG:HB3	1:A:921:ILE:HD11	1.90	0.53
2:B:563:MET:HA	2:B:566:ILE:HG22	1.89	0.53
1:A:16:MET:HB2	1:A:38:ARG:HG3	1.92	0.52
1:A:492:ILE:HG21	1:A:584:TYR:HD2	1.75	0.52
2:B:360:ALA:C	2:B:362:THR:H	2.19	0.51
2:B:494:PHE:HB3	2:B:535:ILE:HG12	1.94	0.50
1:A:351:ILE:HG21	1:A:408:SER:HB2	1.95	0.48
1:A:941:LYS:HD2	1:A:951:ARG:HD3	1.96	0.48
1:A:809[B]:GLN:HE21	1:A:1011:PRO:HD2	1.79	0.47
4:A:1102:1LT:H1	4:A:1102:1LT:H12	1.96	0.46
1:A:83:PHE:HE1	1:A:104:PRO:HB3	1.80	0.46
1:A:1019:ILE:O	1:A:1022:ILE:HG22	2.15	0.46
1:A:413:LYS:HA	1:A:413:LYS:HD2	1.77	0.46
2:B:354:THR:HA	2:B:426:TYR:O	2.16	0.46
1:A:774:SER:HB3	5:A:1555:HOH:O	2.15	0.45
1:A:328:TRP:HA	1:A:394:PRO:HB3	1.99	0.45
1:A:347:ASN:OD1	1:A:349:ARG:HG2	2.16	0.45
1:A:905[B]:CYS:SG	1:A:956:LEU:HD11	2.56	0.45
1:A:192:VAL:HG12	1:A:283:PRO:HG2	1.99	0.45
1:A:534:ILE:HG21	1:A:551:LEU:HD11	1.99	0.45
1:A:596:GLU:HA	1:A:599:MET:HE2	2.00	0.44
1:A:572:VAL:HG21	1:A:583:MET:HG2	2.00	0.44
1:A:346:VAL:HB	1:A:351:ILE:HD13	2.00	0.43
1:A:447:PRO:HB3	2:B:470:TYR:OH	2.19	0.43
2:B:343:VAL:HG22	2:B:356:LEU:HD11	2.01	0.43
2:B:501:GLN:O	2:B:505:SER:HB2	2.19	0.42
1:A:640:LYS:HE2	1:A:680:VAL:HG11	2.00	0.42
1:A:989:LEU:HD22	1:A:1032:GLU:HG2	2.01	0.42
1:A:540:LEU:HD21	1:A:1016:PHE:CD1	2.55	0.42
2:B:564:ASN:HA	5:B:665:HOH:O	2.20	0.41
1:A:363:GLY:N	1:A:607:PRO:HG3	2.35	0.41
1:A:881:LEU:HD12	1:A:881:LEU:HA	1.95	0.41
4:A:1102:1LT:O	4:A:1102:1LT:S	2.78	0.41
1:A:572:VAL:CG2	1:A:583:MET:HE3	2.51	0.41
1:A:214:ASP:HA	1:A:266:PRO:HB3	2.00	0.41
1:A:572:VAL:HG22	1:A:583:MET:HE3	2.03	0.41
1:A:559:VAL:HG13	1:A:591:PRO:HD3	2.03	0.41
1:A:291:GLU:HA	1:A:294:TYR:HB2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:489:MET:O	1:A:493:GLU:HG3	2.22	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	995/1074 (93%)	967 (97%)	26 (3%)	2 (0%)	43	51
2	B	218/293 (74%)	212 (97%)	4 (2%)	2 (1%)	14	14
All	All	1213/1367 (89%)	1179 (97%)	30 (2%)	4 (0%)	36	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	362	THR
1	A	495	HIS
2	B	361	SER
1	A	1055	MET

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	894/980 (91%)	885 (99%)	9 (1%)	68	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	B	197/272 (72%)	193 (98%)	4 (2%)	48 64
All	All	1091/1252 (87%)	1078 (99%)	13 (1%)	63 78

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

Mol	Chain	Res	Type
1	A	115	ARG
1	A	412	ARG
1	A	413	LYS
1	A	427	ILE
1	A	495	HIS
1	A	834	LEU
1	A	901	CYS
1	A	913	ILE
1	A	1061	THR
2	B	361	SER
2	B	381	ILE
2	B	442	ILE
2	B	501	GLN

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	137	GLN
1	A	160	HIS
1	A	384	ASN
1	A	388	ASN
1	A	530	GLN
1	A	630	GLN
1	A	701	HIS
1	A	714	ASN
1	A	796	ASN

5.3.3 RNA ⓘ

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](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SCN	A	1101	-	1,2,2	0.74	0	0,1,1	-	-
4	1LT	A	1102	-	31,32,32	0.85	1 (3%)	44,49,49	0.82	2 (4%)

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	1LT	A	1102	-	-	5/29/41/41	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1102	1LT	C18-C15	4.44	1.63	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1102	1LT	C10-C9-C1	2.75	134.55	130.81
4	A	1102	1LT	C1-C9-S	-2.20	108.57	109.91

There are no chirality outliers.

All (5) torsion outliers are listed below:

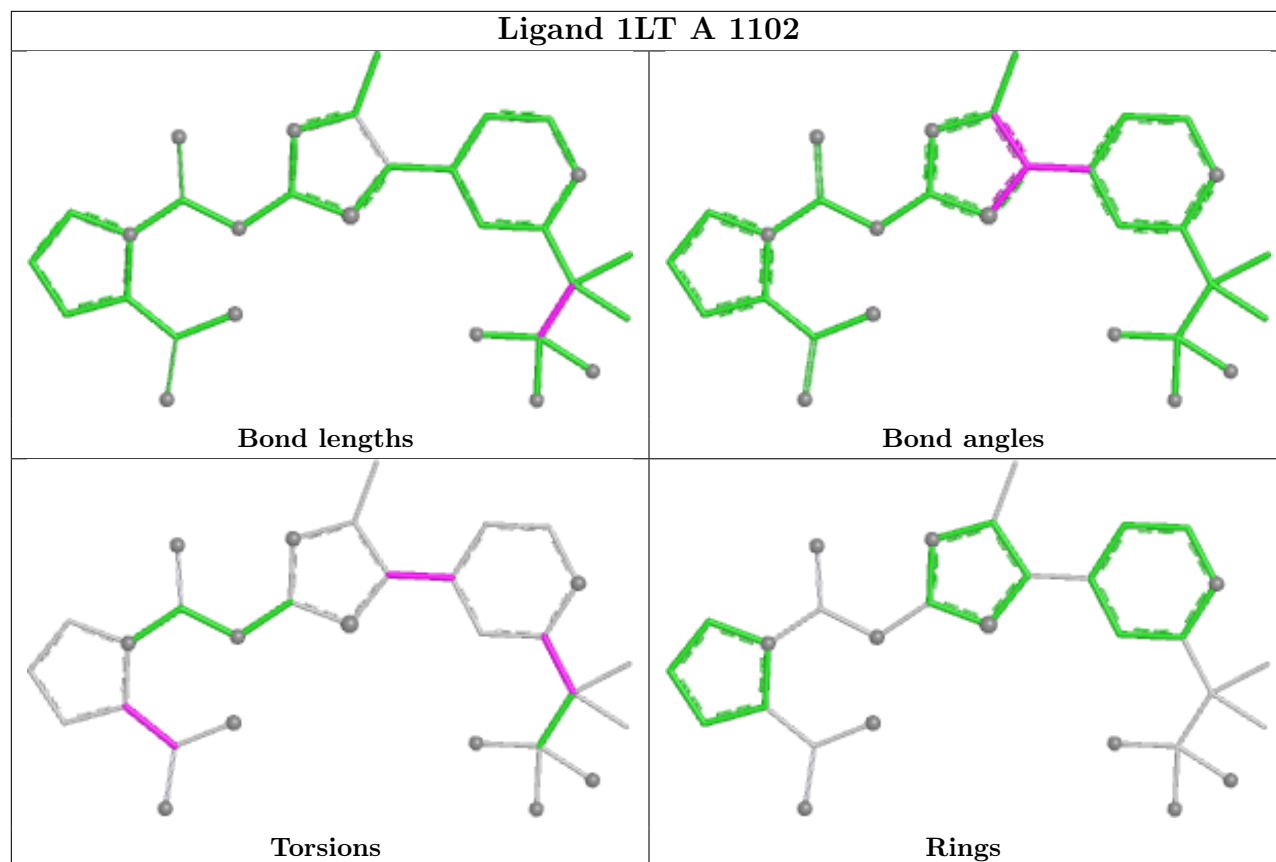
Mol	Chain	Res	Type	Atoms
4	A	1102	1LT	N2-C7-C8-O1
4	A	1102	1LT	N2-C7-C8-N3
4	A	1102	1LT	C14-C13-C15-C17
4	A	1102	1LT	C11-C10-C9-C1
4	A	1102	1LT	C14-C10-C9-C1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1102	1LT	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	999/1074 (93%)	0.37	59 (5%) 28 25	15, 47, 76, 112	6 (0%)
2	B	227/293 (77%)	0.83	25 (11%) 10 8	25, 63, 111, 138	1 (0%)
All	All	1226/1367 (89%)	0.45	84 (6%) 23 20	15, 49, 89, 138	7 (0%)

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	386	ARG	4.5
1	A	244	LEU	4.2
1	A	207	TYR	4.2
1	A	1058	ILE	4.1
1	A	246	TYR	3.9
1	A	3	PRO	3.8
2	B	361	SER	3.7
2	B	362	THR	3.7
2	B	430	LYS	3.6
1	A	200	PRO	3.5
2	B	392	PHE	3.5
1	A	224	ALA	3.4
1	A	1057	TRP	3.3
2	B	398	PHE	3.3
1	A	156	LEU	3.3
1	A	1059	PHE	3.3
1	A	1061	THR	3.2
1	A	247	GLN	3.1
1	A	414	GLY	3.1
2	B	335	TRP	3.0
2	B	426	TYR	3.0
2	B	592	LYS	3.0
2	B	363	LYS	3.0
1	A	557	TYR	2.8

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Mol	Chain	Res	Type	RSRZ
2	B	424	LEU	2.8
2	B	508	TYR	2.8
2	B	428	VAL	2.7
1	A	1056	ASP	2.7
1	A	181	ILE	2.7
1	A	215	CYS	2.7
1	A	1055	MET	2.7
2	B	333	TRP	2.7
1	A	572	VAL	2.7
1	A	496	ALA	2.6
1	A	329	VAL	2.6
1	A	863	LYS	2.6
1	A	324	THR	2.5
1	A	225	ILE	2.5
1	A	300	ASP	2.5
2	B	395	PRO	2.5
1	A	313	THR	2.4
1	A	177	LEU	2.4
2	B	340	ARG	2.4
1	A	2	PRO	2.4
2	B	355	PHE	2.4
1	A	209	LEU	2.4
1	A	248	GLY	2.4
1	A	491	VAL	2.4
1	A	411	GLY	2.3
1	A	973	LYS	2.3
2	B	360	ALA	2.3
2	B	372	LEU	2.3
1	A	479	TRP	2.3
2	B	427	PRO	2.3
1	A	301	CYS	2.3
1	A	278	MET	2.3
1	A	208	THR	2.3
1	A	312	SER	2.3
1	A	1060	HIS	2.3
2	B	436	VAL	2.3
1	A	585	CYS	2.2
1	A	556	HIS	2.2
1	A	947	TYR	2.2
2	B	583	TRP	2.2
1	A	206	LYS	2.2
1	A	185	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	641	TYR	2.2
1	A	197	ILE	2.2
1	A	862	CYS	2.2
1	A	250	TYR	2.1
1	A	1016	PHE	2.1
1	A	297	LEU	2.1
2	B	509	ILE	2.1
1	A	192	VAL	2.1
1	A	222	ALA	2.1
1	A	481	SER	2.1
1	A	412	ARG	2.1
1	A	59	HIS	2.1
1	A	60	GLN	2.1
1	A	495	HIS	2.1
2	B	507	GLU	2.0
1	A	889	ILE	2.0
1	A	562	PRO	2.0
2	B	590	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.

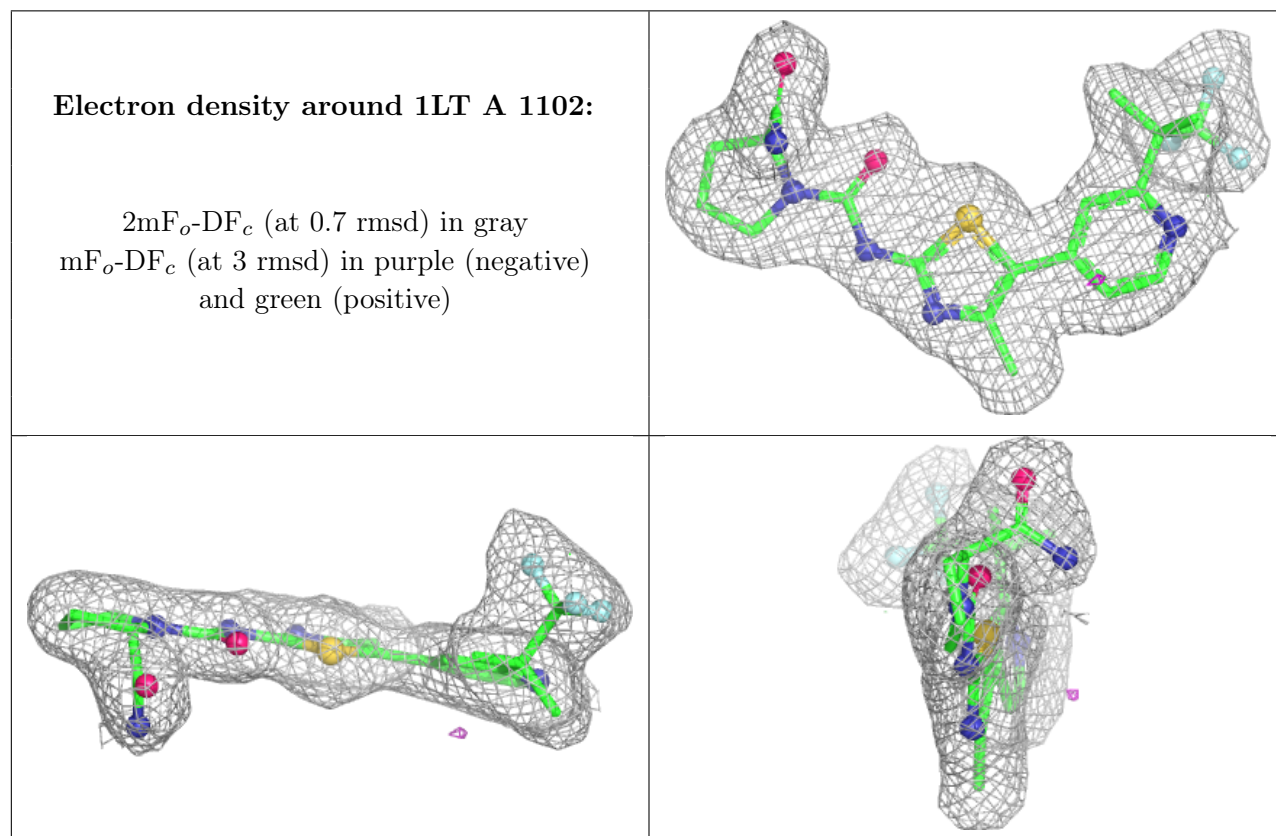
6.4 Ligands [i](#)

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	SCN	A	1101	3/3	0.91	0.15	71,71,72,73	0
4	1LT	A	1102	30/30	0.96	0.06	27,31,41,43	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.