



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 05:53 pm BST

PDB ID : 3LS8
Title : Crystal structure of human PIK3C3 in complex with 3-[4-(4-Morpholinyl)thieno[3,2-d]pyrimidin-2-yl]-phenol
Authors : Tresaugues, L.; Welin, M.; Arrowsmith, C.H.; Berglund, H.; Bountra, C.; Collins, R.; Edwards, A.M.; Flodin, S.; Flores, A.; Graslund, S.; Hammarstrom, M.; Johansson, I.; Karlberg, T.; Kotenyova, T.; Kraulis, P.; Moche, M.; Nyman, T.; Persson, C.; Schuler, H.; Schutz, P.; Siponen, M.I.; Thorsell, A.G.; Van den Berg, S.; Wahlberg, E.; Weigelt, J.; Wisniewska, M.; Nordlund, P.; Structural Genomics Consortium (SGC)
Deposited on : 2010-02-12
Resolution : 2.25 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 following versions of software and data (see [references](#) ①) 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	:	2.11
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)	:	2.11

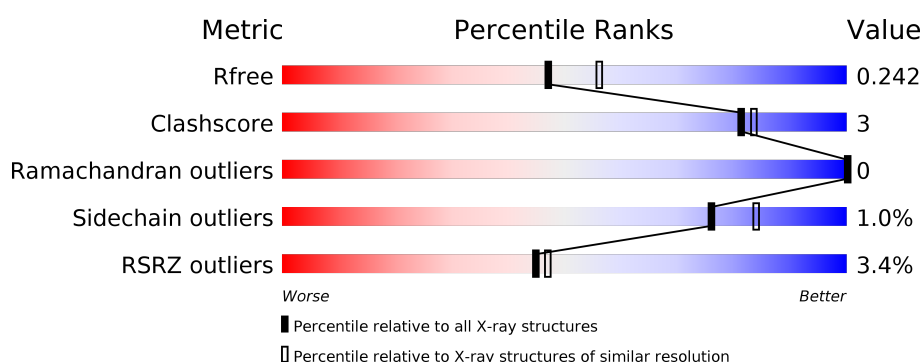
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 on 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	614	
1	B	614	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9027 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.

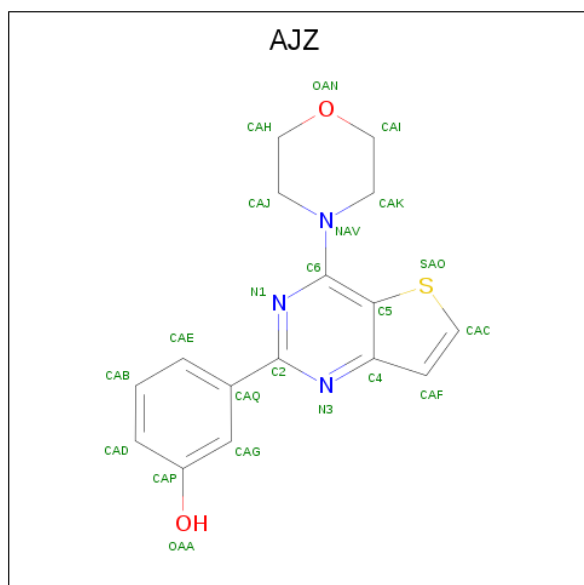
- Molecule 1 is a protein called Phosphatidylinositol 3-kinase catalytic subunit type 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	547	Total	C	N	O	S	0	5	0
			4443	2839	749	831	24			
1	B	522	Total	C	N	O	S	0	6	0
			4270	2735	719	790	26			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	266	SER	-	EXPRESSION TAG	UNP Q8NEB9
A	267	MET	-	EXPRESSION TAG	UNP Q8NEB9
B	266	SER	-	EXPRESSION TAG	UNP Q8NEB9
B	267	MET	-	EXPRESSION TAG	UNP Q8NEB9

- Molecule 2 is 3-(4-morpholin-4-ylthieno[3,2-d]pyrimidin-2-yl)phenol (three-letter code: AJZ) (formula: C₁₆H₁₅N₃O₂S).

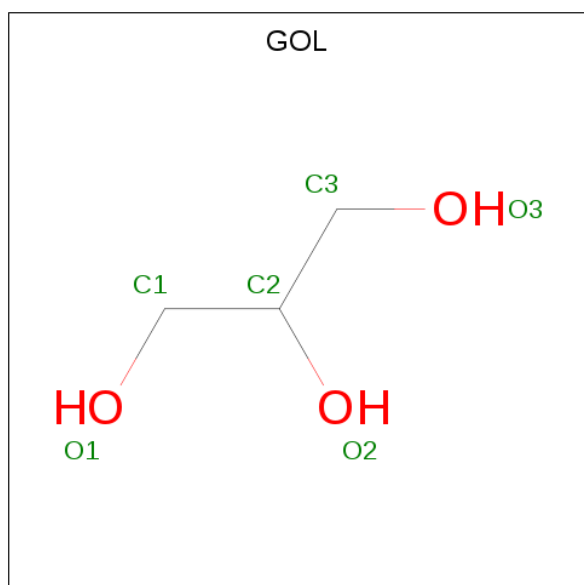


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			22	16	3	2	1		
2	B	1	Total	C	N	O	S	0	0
			22	16	3	2	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cl	0	0
			2	2		
3	A	2	Total	Cl	0	0
			2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	182	Total	O	0	0
			182	182		
5	B	78	Total	O	0	0
			78	78		

- Molecule 1: Phosphatidylinositol 3-kinase catalytic subunit type 3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.69Å 141.15Å 163.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	70.90 – 2.25 70.90 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.3 (70.90-2.25) 99.3 (70.90-2.25)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.209 , 0.251 0.204 , 0.242	Depositor DCC
R_{free} test set	3491 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	30.8	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9027	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 2.88% 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: GOL, AJZ, CL

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	0.47	0/4546	0.60	4/6147 (0.1%)
1	B	0.38	0/4372	0.53	0/5906
All	All	0.43	0/8918	0.57	4/12053 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	810	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	A	705	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	A	810	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	705	ARG	NE-CZ-NH1	5.02	122.81	120.30

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	4443	0	4474	21	0
1	B	4270	0	4329	22	0
2	A	22	0	15	3	0
2	B	22	0	15	3	0
3	A	2	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
4	A	6	0	8	0	0
5	A	182	0	0	0	0
5	B	78	0	0	0	0
All	All	9027	0	8841	46	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 3.

All (46) 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:802[B]:CYS:SG	1:B:863:ILE:HD11	2.29	0.72
1:A:741:VAL:HG12	1:A:744:ARG:HD3	1.70	0.72
1:B:741:VAL:HG12	1:B:744:ARG:HD3	1.74	0.69
1:A:305:THR:HG21	1:A:839:VAL:HG11	1.74	0.68
1:B:733[B]:CYS:SG	1:B:779:LEU:HD21	2.37	0.65
2:A:1000:AJZ:SAO	2:A:1000:AJZ:HAJ	2.37	0.65
1:A:749:LEU:HD11	1:A:757:LEU:HD21	1.78	0.64
2:A:1000:AJZ:CAJ	2:A:1000:AJZ:SAO	2.93	0.57
1:B:555:LEU:HD21	1:B:573:LEU:HD12	1.88	0.56
1:A:313:GLN:HA	1:A:336[A]:CYS:SG	2.47	0.54
1:A:768:ARG:HD2	1:A:833:LEU:HD12	1.89	0.54
1:A:664:ASP:OD2	1:A:666:LYS:NZ	2.34	0.53
1:B:385:ARG:O	1:B:482:ARG:NH2	2.41	0.53
1:B:733[B]:CYS:SG	1:B:779:LEU:CD2	2.97	0.52
1:A:685:ILE:O	2:A:1000:AJZ:HAH	2.10	0.51
2:B:1004:AJZ:SAO	2:B:1004:AJZ:CAK	2.98	0.51
1:B:494:TYR:CZ	1:B:498:ILE:HG21	2.47	0.50
1:B:494:TYR:O	1:B:498:ILE:HG23	2.12	0.50
1:B:305:THR:HG21	1:B:839:VAL:HG11	1.93	0.49
1:A:566:ARG:HD3	1:B:695:ASP:OD1	2.12	0.48
1:B:330:LEU:HD13	1:B:355:TRP:CD2	2.48	0.48
1:A:341:LEU:HD22	1:A:343:GLN:HB2	1.95	0.47
1:A:389:ASP:OD1	1:A:482:ARG:NH1	2.41	0.47
2:B:1004:AJZ:HAK	2:B:1004:AJZ:SAO	2.54	0.47
1:B:498:ILE:C	1:B:498:ILE:HD12	2.35	0.47
1:A:528:LEU:HD23	1:A:535:VAL:HG13	1.98	0.45
1:A:321:TYR:CE1	1:A:351:LEU:HD21	2.52	0.45
1:A:742:GLY:O	1:A:744:ARG:HD2	2.17	0.44
1:B:535:VAL:HA	1:B:538:MET:HE3	1.99	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:685:ILE:O	2:B:1004:AJZ:HAI	2.18	0.44
1:B:559:VAL:O	1:B:562:GLU:HB3	2.18	0.44
1:A:516:MET:O	1:A:520:VAL:HG23	2.18	0.44
1:A:701:GLN:OE1	1:A:705:ARG:NH2	2.51	0.43
1:A:305:THR:HG21	1:A:839:VAL:CG1	2.43	0.43
1:A:528:LEU:HA	1:A:535:VAL:CG1	2.49	0.43
1:B:308:LEU:HA	1:B:312:GLU:OE1	2.19	0.43
1:B:395:TYR:CE1	1:B:815:LEU:HD11	2.55	0.42
1:A:320:ARG:O	1:A:324:THR:HG23	2.20	0.42
1:B:535:VAL:HA	1:B:538:MET:CE	2.49	0.42
1:B:555:LEU:CD2	1:B:573:LEU:HD12	2.50	0.41
1:A:556:MET:HE2	1:A:637[B]:HIS:ND1	2.34	0.41
1:B:693:VAL:HG13	1:B:703:PHE:HB2	2.03	0.41
1:A:810:ARG:NH2	1:A:847:ARG:O	2.53	0.41
1:B:672:VAL:HG22	1:B:682:MET:HG2	2.03	0.41
1:A:555:LEU:O	1:A:559:VAL:HG23	2.21	0.41
1:B:676:SER:OG	1:B:677:THR:N	2.55	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	546/614 (89%)	539 (99%)	7 (1%)	0	100	100
1	B	524/614 (85%)	517 (99%)	7 (1%)	0	100	100
All	All	1070/1228 (87%)	1056 (99%)	14 (1%)	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	497/555 (90%)	490 (99%)	7 (1%)	67	76
1	B	478/555 (86%)	475 (99%)	3 (1%)	86	91
All	All	975/1110 (88%)	965 (99%)	10 (1%)	76	84

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	338	ASN
1	A	341	LEU
1	A	361	GLU
1	A	388	ASP
1	A	502	GLU
1	A	539	ARG
1	A	644	ASP
1	B	676	SER
1	B	694	LEU
1	B	827	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	526	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
2	AJZ	A	1000	-	23,25,25	1.03	1 (4%)	26,35,35	2.59	5 (19%)
4	GOL	A	1003	-	5,5,5	0.34	0	5,5,5	0.40	0
2	AJZ	B	1004	-	23,25,25	1.11	1 (4%)	26,35,35	2.76	7 (26%)

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	AJZ	A	1000	-	-	2/8/16/16	0/4/4/4
4	GOL	A	1003	-	-	2/4/4/4	-
2	AJZ	B	1004	-	-	2/8/16/16	0/4/4/4

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1004	AJZ	CAQ-C2	-3.46	1.39	1.48
2	A	1000	AJZ	CAQ-C2	-3.26	1.40	1.48

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1004	AJZ	C2-N3-C4	8.08	121.98	116.50
2	B	1004	AJZ	N3-C2-N1	-7.93	119.88	126.11

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1000	AJZ	C2-N3-C4	7.80	121.79	116.50
2	A	1000	AJZ	N3-C2-N1	-7.45	120.25	126.11
2	A	1000	AJZ	CAK-NAV-CAJ	4.54	121.53	111.52
2	B	1004	AJZ	CAK-NAV-CAJ	4.44	121.32	111.52
2	B	1004	AJZ	OAN-CAH-CAJ	-4.43	102.04	111.80
2	A	1000	AJZ	CAF-C4-C5	3.33	113.05	110.62
2	B	1004	AJZ	CAF-C4-C5	3.30	113.03	110.62
2	A	1000	AJZ	OAN-CAI-CAK	-2.97	105.25	111.80
2	B	1004	AJZ	C6-N1-C2	2.30	121.56	116.17
2	B	1004	AJZ	CAQ-C2-N3	2.18	122.06	116.88

There are no chirality outliers.

All (6) torsion outliers are listed below:

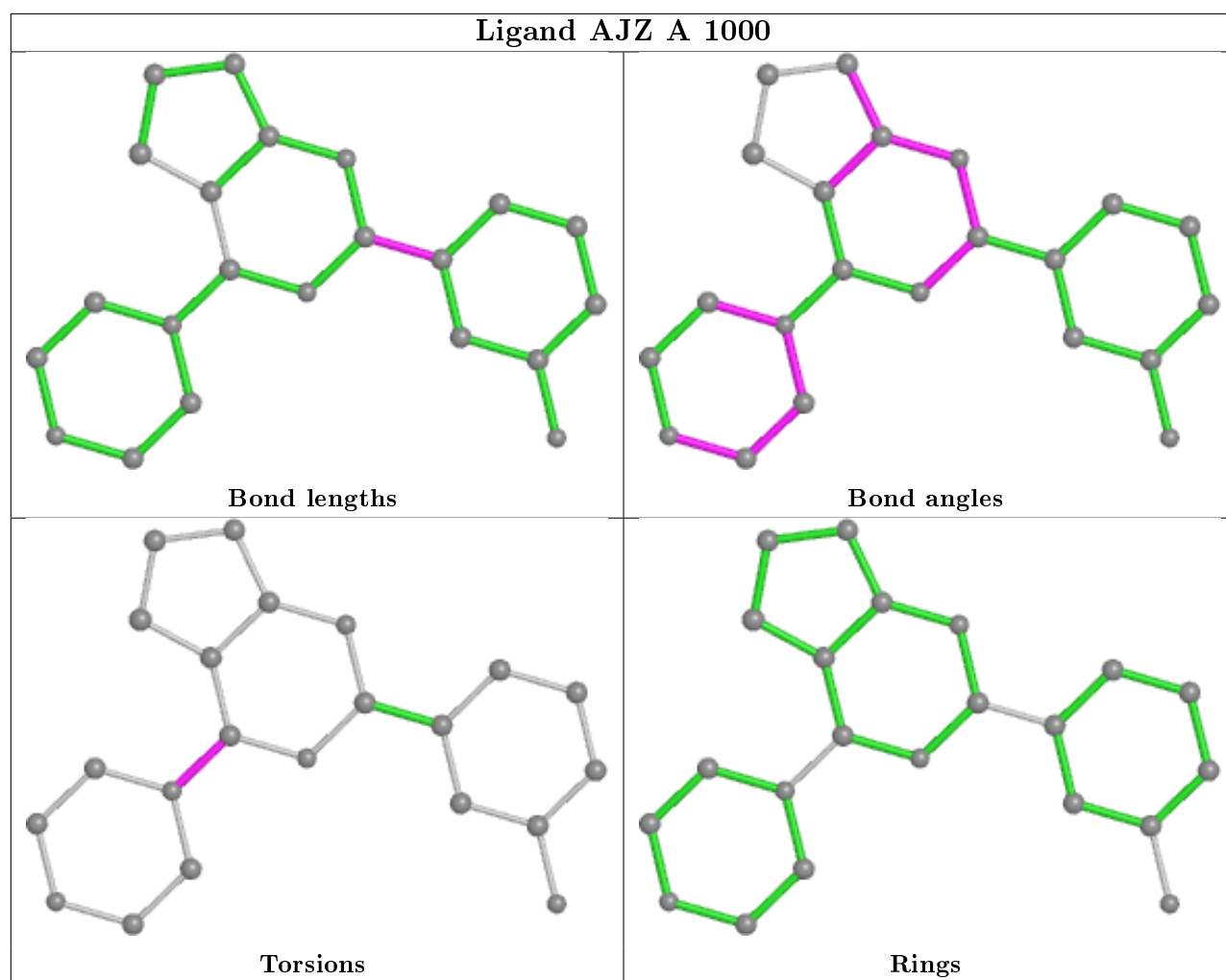
Mol	Chain	Res	Type	Atoms
2	A	1000	AJZ	N1-C6-NAV-CAJ
2	A	1000	AJZ	C5-C6-NAV-CAJ
2	B	1004	AJZ	N1-C6-NAV-CAK
2	B	1004	AJZ	C5-C6-NAV-CAK
4	A	1003	GOL	O1-C1-C2-C3
4	A	1003	GOL	O1-C1-C2-O2

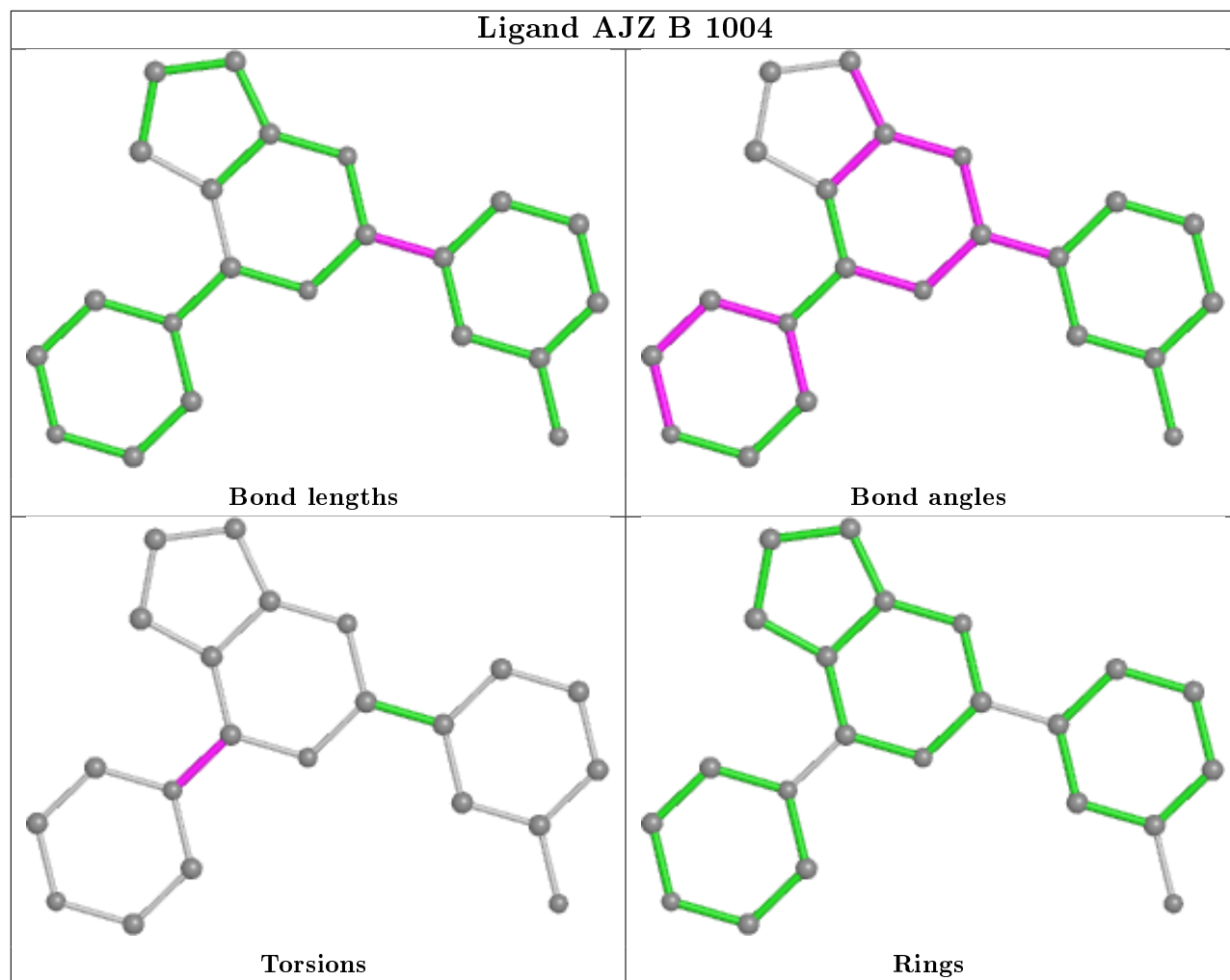
There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1000	AJZ	3	0
2	B	1004	AJZ	3	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	547/614 (89%)	-0.22	1 (0%) 95 96	13, 26, 45, 62	0
1	B	522/614 (85%)	0.22	35 (6%) 17 19	20, 37, 88, 135	0
All	All	1069/1228 (87%)	-0.00	36 (3%) 45 47	13, 31, 70, 135	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	317	TRP	5.8
1	B	471	LEU	5.4
1	B	300	VAL	5.2
1	B	315	LEU	4.7
1	B	322	TYR	4.6
1	B	341	LEU	4.1
1	B	336[A]	CYS	3.9
1	B	310	TYR	3.8
1	B	319	PHE	3.7
1	B	316	VAL	3.7
1	B	318	LYS	3.6
1	B	354	LYS	3.6
1	B	857	HIS	3.6
1	B	298	ILE	3.4
1	B	296	LEU	3.3
1	B	323	LEU	3.3
1	B	339	TRP	3.0
1	B	308	LEU	2.9
1	B	337	VAL	2.8
1	B	868	HIS	2.8
1	B	311	GLU	2.8
1	B	352	LEU	2.7
1	B	347	GLN	2.7
1	B	351	LEU	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	371	TYR	2.5
1	B	344	GLU	2.5
1	B	338	ASN	2.5
1	B	343	GLN	2.4
1	B	378	ARG	2.3
1	B	321	TYR	2.3
1	B	348	ALA	2.2
1	B	472	GLU	2.2
1	B	342	PRO	2.2
1	B	376	VAL	2.1
1	A	452	SER	2.0
1	B	309	THR	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 ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

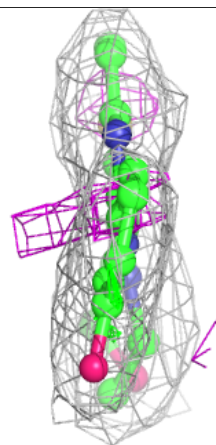
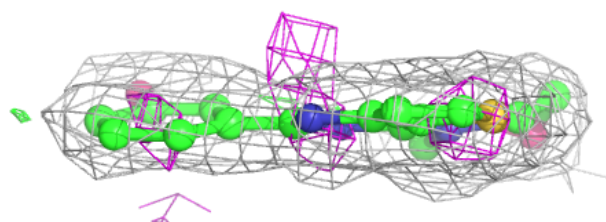
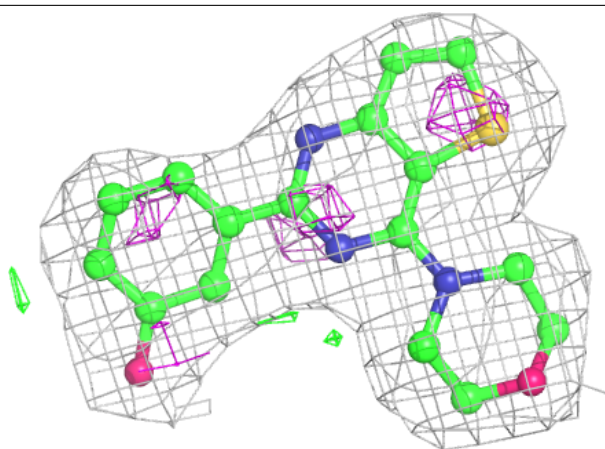
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

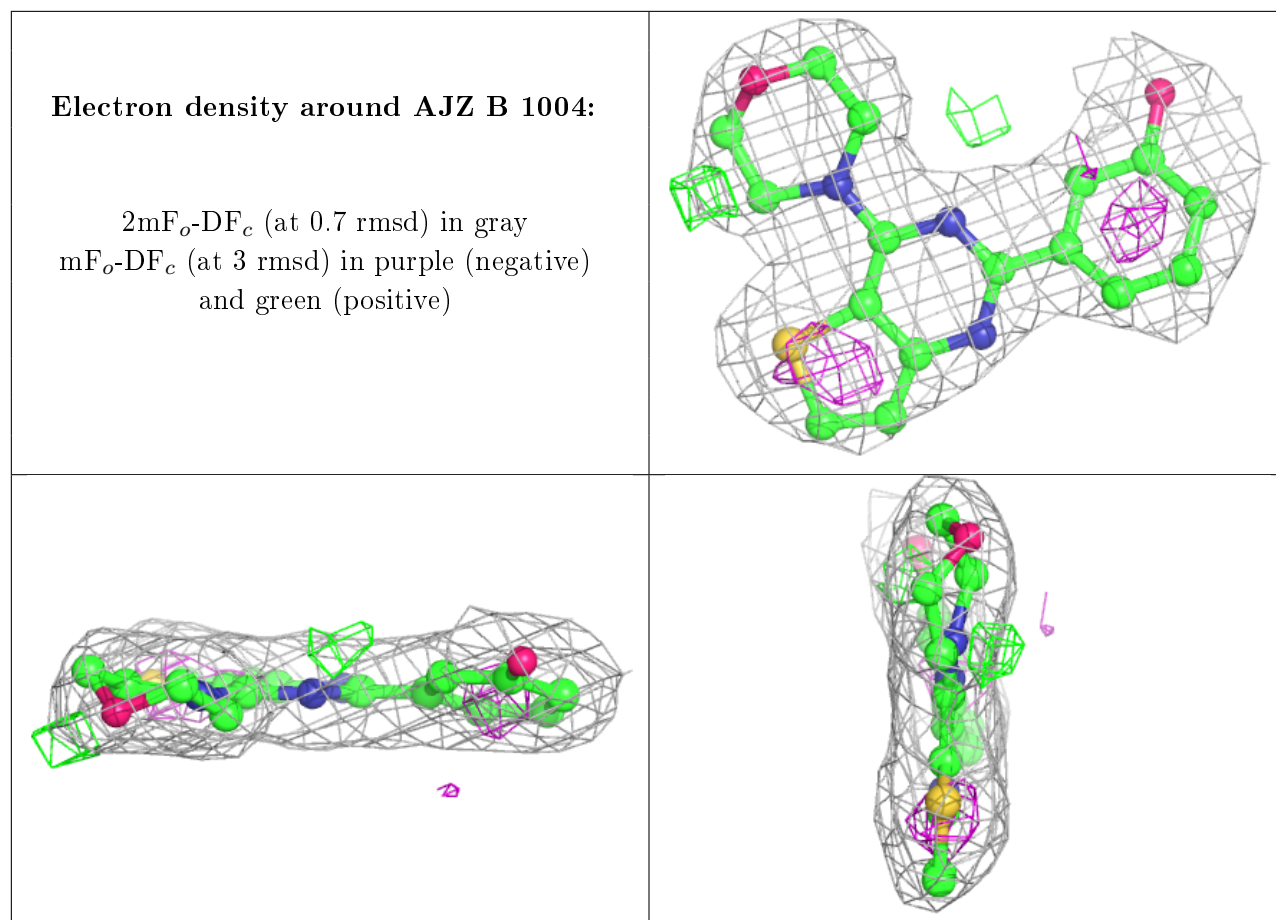
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CL	B	1005	1/1	0.69	0.12	73,73,73,73	0
4	GOL	A	1003	6/6	0.71	0.22	37,39,40,41	0
3	CL	A	1001	1/1	0.90	0.20	60,60,60,60	0
3	CL	B	1006	1/1	0.90	0.15	72,72,72,72	0
2	AJZ	A	1000	22/22	0.92	0.12	23,25,30,31	0
2	AJZ	B	1004	22/22	0.94	0.12	31,34,36,37	0
3	CL	A	1002	1/1	0.97	0.21	49,49,49,49	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 AJZ A 1000:

$2mF_o - DF_c$ (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.