



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:30 PM UTC

PDB ID : 8TSD / pdb_00008tsd
Title : Human PI3K p85alpha/p110alpha bound to RLY-2608
Authors : Holliday, M.; Tang, Y.; Bulku, A.; Wilbur, J.; Fraser, J.
Deposited on : 2023-08-11
Resolution : 2.70 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

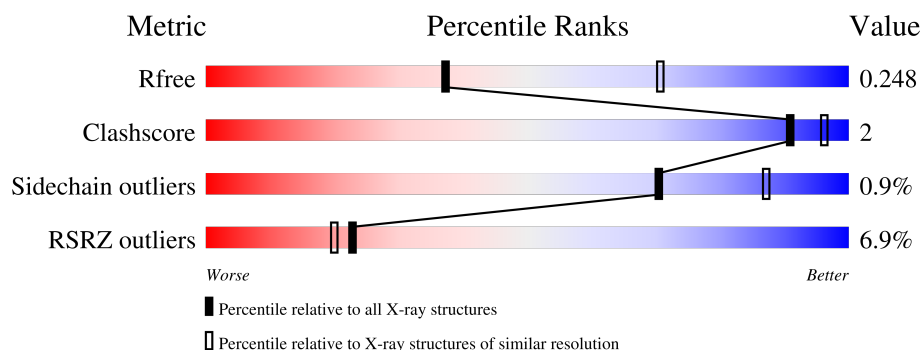
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	1063	7% 90% 5% 5%
2	B	300	4% 84% 6% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21060 atoms, of which 10525 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	1005	Total	C	H	N	O	S	47	0	0
			16437	5246	8237	1401	1485	68			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	GLY	-	expression tag	UNP P42336
A	-8	SER	-	expression tag	UNP P42336
A	-7	PRO	-	expression tag	UNP P42336
A	-6	GLY	-	expression tag	UNP P42336
A	-5	ILE	-	expression tag	UNP P42336
A	-4	SER	-	expression tag	UNP P42336
A	-3	GLY	-	expression tag	UNP P42336
A	-2	GLY	-	expression tag	UNP P42336
A	-1	GLY	-	expression tag	UNP P42336
A	0	GLY	-	expression tag	UNP P42336
A	1	GLY	-	expression tag	UNP P42336

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

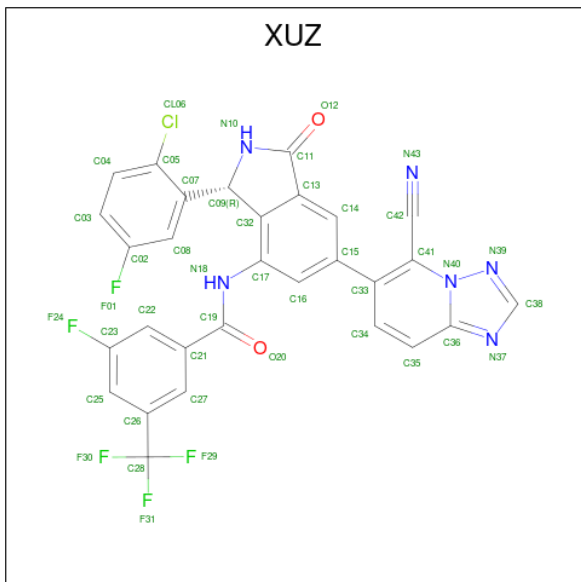
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	270	Total	C	H	N	O	S	63	0	0
			4566	1435	2274	409	442	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	316	GLY	-	expression tag	UNP P27986
B	317	PRO	-	expression tag	UNP P27986

- Molecule 3 is N-{(3R,6M)-3-(2-chloro-5-fluorophenyl)-6-[(4S)-5-cyano[1,2,4]triazolo[1,5-a]py

ridin-6-yl]-1-oxo-2,3-dihydro-1H-isoindol-4-yl}-3-fluoro-5-(trifluoromethyl)benzamide (CCD ID: XUZ) (formula: C₂₉H₁₄ClF₅N₆O₂) (labeled as "Ligand of Interest" by depositor).

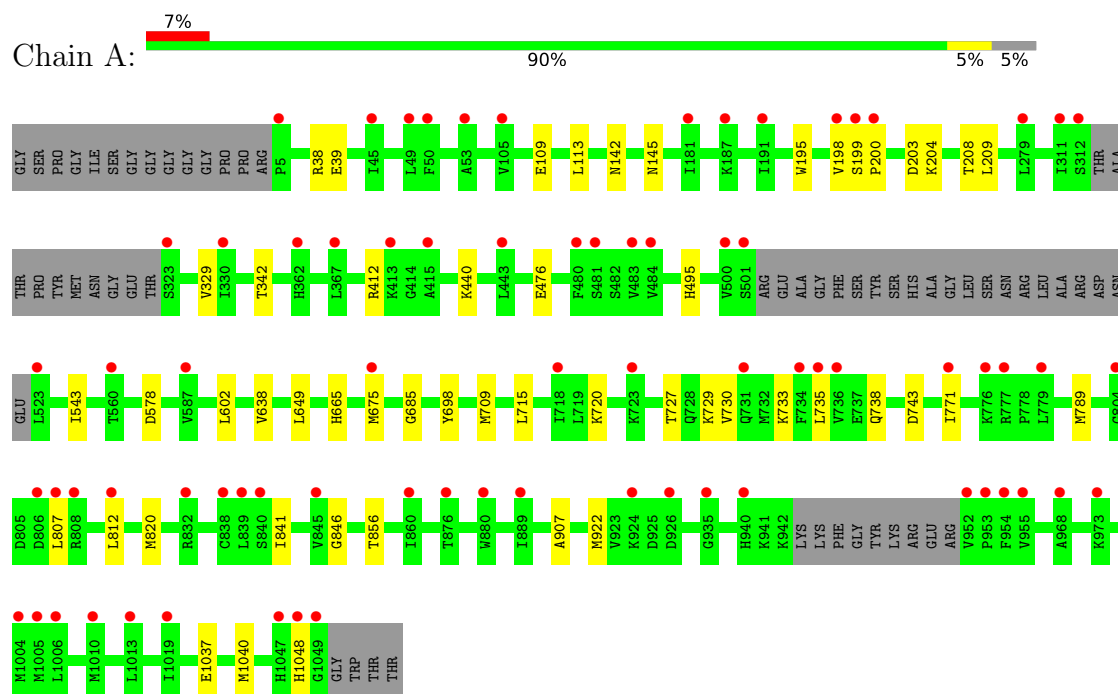


Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	Cl	F	H	N	O		
3	A	1	57	29	1	5	14	6	2	0	0

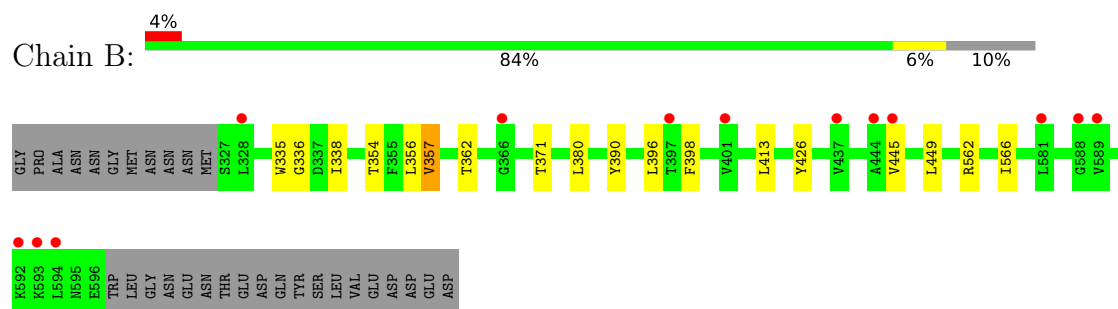
3 Residue-property plots [i](#)

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: 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	86.90Å 121.52Å 192.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	102.78 – 2.70 102.78 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (102.78-2.70) 100.0 (102.78-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.239 , 0.277 (Not available) , 0.248	Depositor DCC
R_{free} test set	2000 reflections (3.52%)	wwPDB-VP
Wilson B-factor (Å ²)	97.3	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 59.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21060	wwPDB-VP
Average B, all atoms (Å ²)	130.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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: XUZ

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	0.09	0/8382	0.24	0/11326
2	B	0.08	0/2330	0.25	0/3120
All	All	0.08	0/10712	0.25	0/14446

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	8200	8237	8234	27	0
2	B	2292	2274	2273	9	0
3	A	43	14	0	1	0
All	All	10535	10525	10507	36	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 (36) 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:356:LEU:HD23	2:B:357:VAL:N	2.12	0.65
1:A:198:VAL:HG22	1:A:200:PRO:HD2	1.78	0.64
1:A:812:LEU:HD23	3:A:1101:XUZ:F29	1.88	0.63
1:A:204:LYS:HD2	1:A:789:MET:HE3	1.81	0.62
1:A:208:THR:C	1:A:209:LEU:HD12	2.33	0.54
1:A:820:MET:HE1	1:A:907:ALA:HB1	1.91	0.52
1:A:543:ILE:HG23	1:A:543:ILE:O	2.11	0.51
1:A:665:HIS:ND1	1:A:698:TYR:OH	2.39	0.50
1:A:807:LEU:HD12	1:A:846:GLY:HA3	1.95	0.49
1:A:675:MET:HE1	1:A:685:GLY:N	2.28	0.48
1:A:142:ASN:O	1:A:145:ASN:ND2	2.47	0.47
2:B:562:ARG:O	2:B:566:ILE:HD13	2.16	0.45
1:A:638:VAL:HG23	1:A:649:LEU:HD21	1.99	0.45
2:B:354:THR:HG22	2:B:426:TYR:HB2	1.99	0.45
1:A:38:ARG:NH1	1:A:743:ASP:OD2	2.47	0.45
1:A:727:THR:OG1	1:A:730:VAL:HG23	2.17	0.44
1:A:39:GLU:OE1	1:A:39:GLU:N	2.49	0.44
1:A:440:LYS:HA	1:A:476:GLU:HA	1.98	0.44
1:A:495:HIS:NE2	1:A:578:ASP:OD1	2.40	0.44
1:A:715:LEU:HD22	1:A:738:GLN:HG2	1.99	0.44
1:A:109:GLU:N	1:A:109:GLU:OE1	2.51	0.43
1:A:856:THR:HG22	1:A:922:MET:HG2	2.01	0.43
1:A:199:SER:O	1:A:200:PRO:C	2.62	0.43
1:A:729:LYS:N	1:A:729:LYS:CD	2.82	0.43
1:A:412:ARG:NE	1:A:412:ARG:HA	2.35	0.42
1:A:735:LEU:HD22	1:A:771:ILE:HG23	2.01	0.42
2:B:390:TYR:N	2:B:398:PHE:O	2.53	0.42
1:A:195:TRP:CD2	1:A:789:MET:HE1	2.55	0.41
1:A:1037:GLU:HA	1:A:1040:MET:HE3	2.02	0.41
1:A:198:VAL:HG12	1:A:203:ASP:O	2.20	0.41
2:B:445:VAL:O	2:B:449:LEU:HB3	2.20	0.41
2:B:335:TRP:CD1	2:B:356:LEU:HD21	2.56	0.41
2:B:371:THR:HG22	2:B:380:LEU:HG	2.03	0.41
1:A:709:MET:HE2	1:A:841:ILE:HD13	2.03	0.41
2:B:362:THR:O	2:B:362:THR:HG22	2.22	0.40
2:B:336:GLY:O	2:B:338:ILE:N	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

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	921/963 (96%)	914 (99%)	7 (1%)	73	88
2	B	251/277 (91%)	248 (99%)	3 (1%)	63	84
All	All	1172/1240 (94%)	1162 (99%)	10 (1%)	70	87

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

Mol	Chain	Res	Type
1	A	113	LEU
1	A	329	VAL
1	A	342	THR
1	A	602	LEU
1	A	720	LYS
1	A	733	LYS
1	A	1048	HIS
2	B	357	VAL
2	B	396	LEU
2	B	413	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	HIS
1	A	554	HIS
1	A	731	GLN
1	A	822	ASN
2	B	453	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	XUZ	A	1101	-	46,48,48	2.18	17 (36%)	59,73,73	1.69	17 (28%)

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
3	XUZ	A	1101	-	-	3/22/36/36	0/6/6/6

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	XUZ	C11-N10	8.37	1.41	1.34
3	A	1101	XUZ	C13-C11	3.66	1.53	1.48
3	A	1101	XUZ	C38-N39	3.37	1.38	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1101	XUZ	C21-C19	3.26	1.57	1.50
3	A	1101	XUZ	C09-N10	3.13	1.49	1.46
3	A	1101	XUZ	C15-C33	3.00	1.53	1.48
3	A	1101	XUZ	C28-C26	2.79	1.55	1.49
3	A	1101	XUZ	C36-N37	2.69	1.37	1.33
3	A	1101	XUZ	C17-N18	2.59	1.46	1.41
3	A	1101	XUZ	C19-N18	2.54	1.43	1.35
3	A	1101	XUZ	C38-N37	2.48	1.40	1.36
3	A	1101	XUZ	C16-C17	2.45	1.43	1.39
3	A	1101	XUZ	C14-C15	2.35	1.43	1.39
3	A	1101	XUZ	C07-C09	2.35	1.55	1.52
3	A	1101	XUZ	C22-C23	2.32	1.41	1.37
3	A	1101	XUZ	C36-N40	2.30	1.42	1.39
3	A	1101	XUZ	C08-C02	2.22	1.41	1.37

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1101	XUZ	O12-C11-N10	-4.77	121.88	125.97
3	A	1101	XUZ	C38-N39-N40	4.55	104.57	100.86
3	A	1101	XUZ	C09-N10-C11	-3.67	111.95	114.03
3	A	1101	XUZ	C03-C02-C08	-3.07	119.18	123.23
3	A	1101	XUZ	C13-C11-N10	2.94	107.79	106.51
3	A	1101	XUZ	C14-C13-C11	2.86	132.56	129.66
3	A	1101	XUZ	C21-C22-C23	2.80	121.36	118.48
3	A	1101	XUZ	C07-C08-C02	2.71	122.11	118.57
3	A	1101	XUZ	C25-C23-C22	-2.58	120.37	123.50
3	A	1101	XUZ	F01-C02-C08	2.49	121.82	118.28
3	A	1101	XUZ	C21-C19-N18	2.44	121.84	115.90
3	A	1101	XUZ	C14-C13-C32	-2.35	120.00	122.79
3	A	1101	XUZ	C16-C17-C32	-2.28	118.31	121.07
3	A	1101	XUZ	O20-C19-N18	-2.28	117.97	123.75
3	A	1101	XUZ	N39-C38-N37	-2.24	113.23	116.78
3	A	1101	XUZ	C27-C21-C22	-2.15	117.13	119.65
3	A	1101	XUZ	C16-C15-C14	-2.09	117.20	119.65

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1101	XUZ	C05-C07-C09-N10
3	A	1101	XUZ	C08-C07-C09-N10

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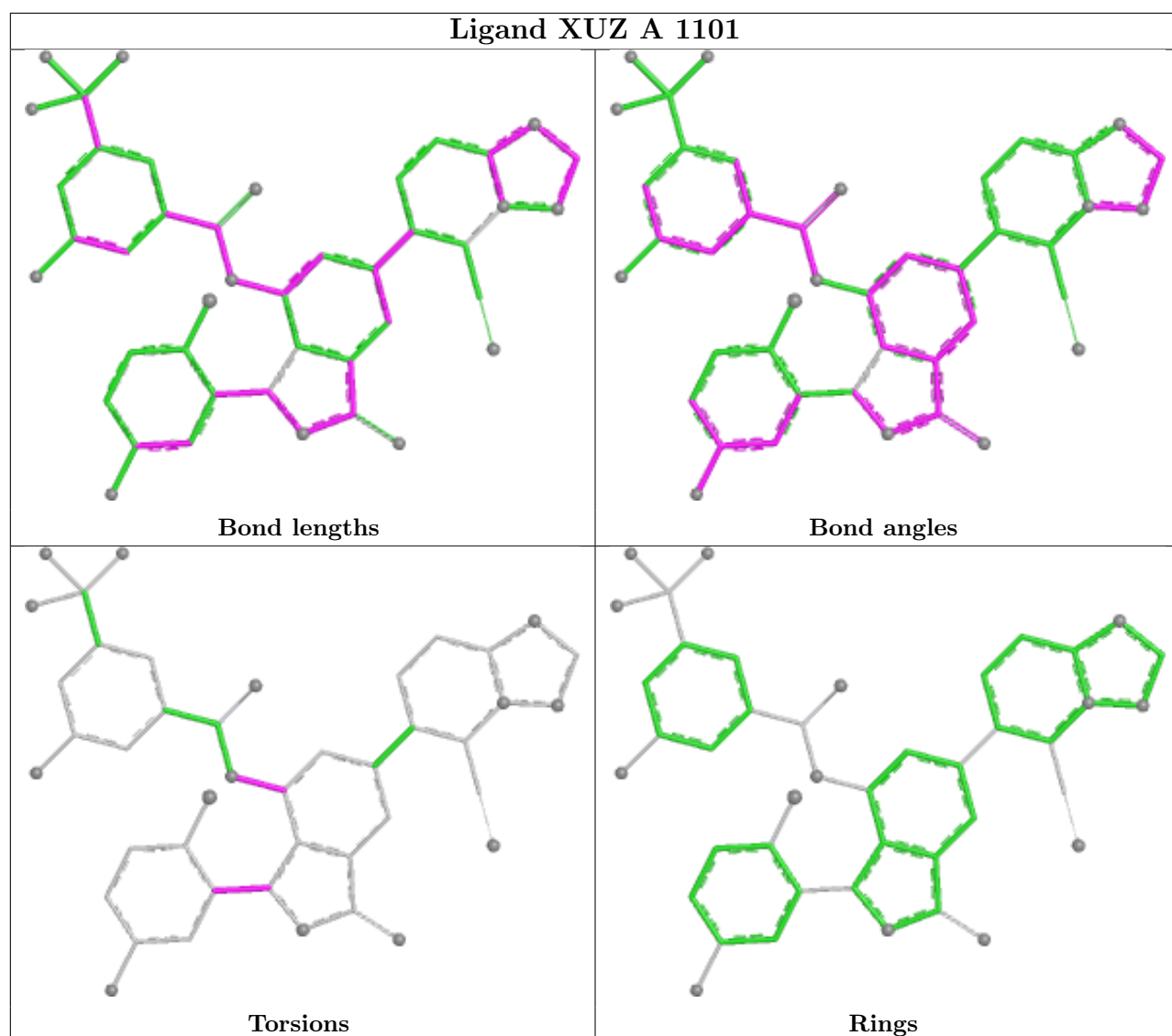
Mol	Chain	Res	Type	Atoms
3	A	1101	XUZ	C32-C17-N18-C19

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1101	XUZ	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	1005/1063 (94%)	0.42	75 (7%)	20 17	34, 115, 174, 227	3 (0%)
2	B	270/300 (90%)	0.54	13 (4%)	35 32	51, 161, 221, 260	5 (1%)
All	All	1275/1363 (93%)	0.45	88 (6%)	23 20	34, 121, 197, 260	8 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1049	GLY	6.2
1	A	734	PHE	5.1
1	A	311	ILE	4.5
1	A	1004	MET	4.2
1	A	731	GLN	4.1
1	A	523	LEU	3.7
1	A	838	CYS	3.7
1	A	198	VAL	3.7
1	A	953	PRO	3.7
1	A	367	LEU	3.6
1	A	481	SER	3.6
1	A	191	ILE	3.6
1	A	483	VAL	3.6
1	A	53	ALA	3.6
1	A	1010	MET	3.4
1	A	49	LEU	3.3
2	B	328	LEU	3.3
2	B	444	ALA	3.2
1	A	812	LEU	3.2
1	A	808	ARG	3.2
2	B	592	LYS	3.2
1	A	279	LEU	3.2
1	A	804	GLY	3.2
2	B	445	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	880	TRP	3.1
1	A	312	SER	3.1
1	A	777	ARG	3.1
1	A	1013	LEU	3.0
1	A	1048	HIS	2.9
1	A	952	VAL	2.9
1	A	779	LEU	2.9
1	A	840	SER	2.9
1	A	105	VAL	2.9
1	A	926	ASP	2.8
1	A	413	LYS	2.8
1	A	1005	MET	2.8
1	A	50	PHE	2.8
2	B	437	VAL	2.8
2	B	593	LYS	2.7
1	A	45	ILE	2.7
1	A	876	THR	2.7
1	A	954	PHE	2.7
1	A	1006	LEU	2.6
1	A	973	LYS	2.6
1	A	415	ALA	2.6
2	B	594	LEU	2.6
1	A	480	PHE	2.6
1	A	968	ALA	2.6
2	B	589	VAL	2.5
1	A	889	ILE	2.5
1	A	832	ARG	2.5
1	A	723	LYS	2.4
1	A	500	VAL	2.4
1	A	718	ILE	2.4
2	B	588	GLY	2.4
1	A	587	VAL	2.3
1	A	1047	HIS	2.3
2	B	397	THR	2.3
1	A	501	SER	2.3
1	A	736	VAL	2.3
2	B	401	VAL	2.3
1	A	735	LEU	2.3
1	A	807	LEU	2.3
1	A	839	LEU	2.3
1	A	181	ILE	2.3
1	A	771	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	935	GLY	2.3
1	A	955	VAL	2.3
1	A	187	LYS	2.2
1	A	675	MET	2.2
1	A	200	PRO	2.2
2	B	366	GLY	2.2
1	A	362	HIS	2.2
1	A	1019	ILE	2.2
1	A	560	THR	2.2
1	A	484	VAL	2.2
2	B	581	LEU	2.1
1	A	860	ILE	2.1
1	A	5	PRO	2.1
1	A	443	LEU	2.1
1	A	940	HIS	2.1
1	A	776	LYS	2.1
1	A	806	ASP	2.1
1	A	323	SER	2.1
1	A	199	SER	2.0
1	A	330	ILE	2.0
1	A	845	VAL	2.0
1	A	924	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 [\(i\)](#)

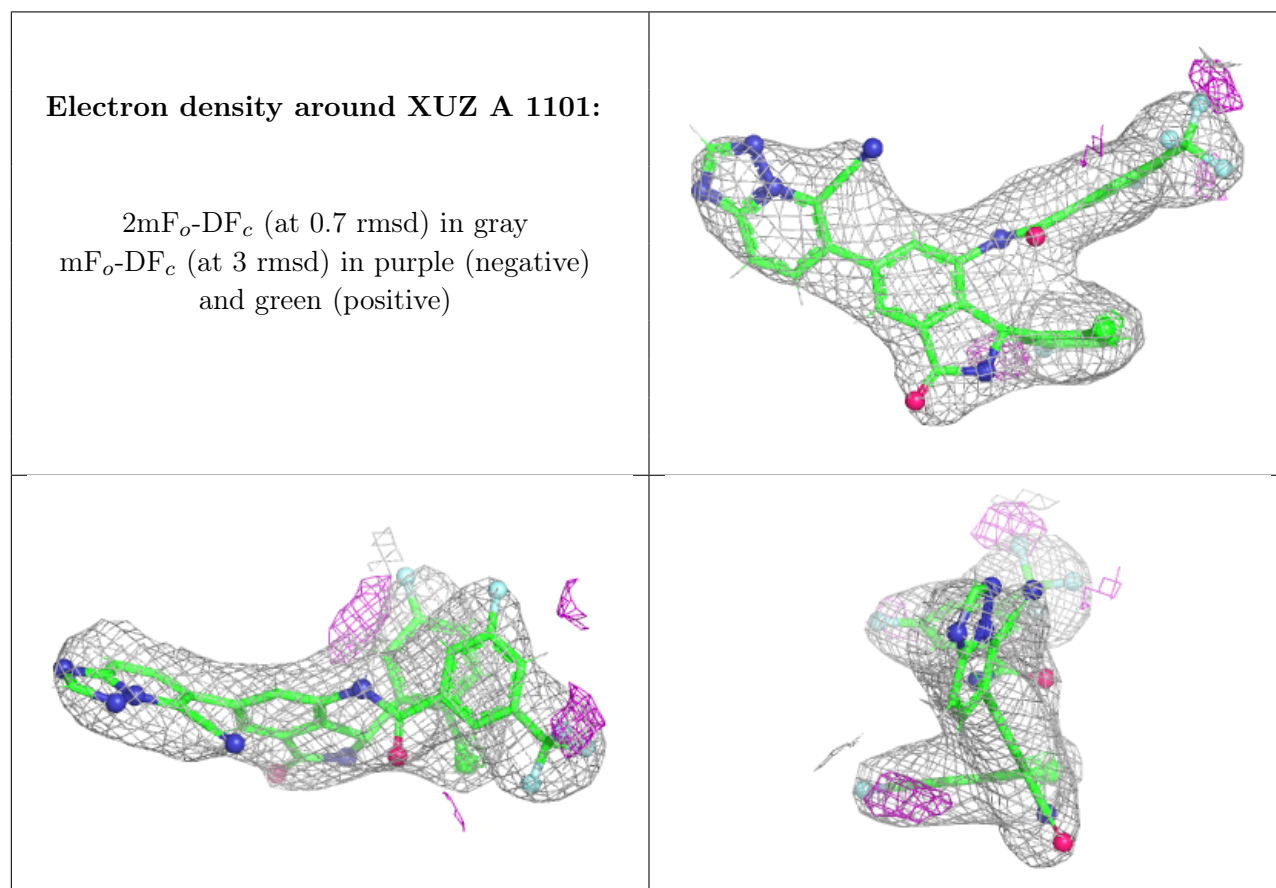
There are no oligosaccharides in this entry.

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	XUZ	A	1101	43/43	0.89	0.12	86,102,142,175	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 ⓘ

There are no such residues in this entry.