



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:36 PM UTC

PDB ID : 4ACG / pdb_00004acg
Title : GSK3b in complex with inhibitor
Authors : Xue, Y.; Ormo, M.
Deposited on : 2011-12-15
Resolution : 2.60 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

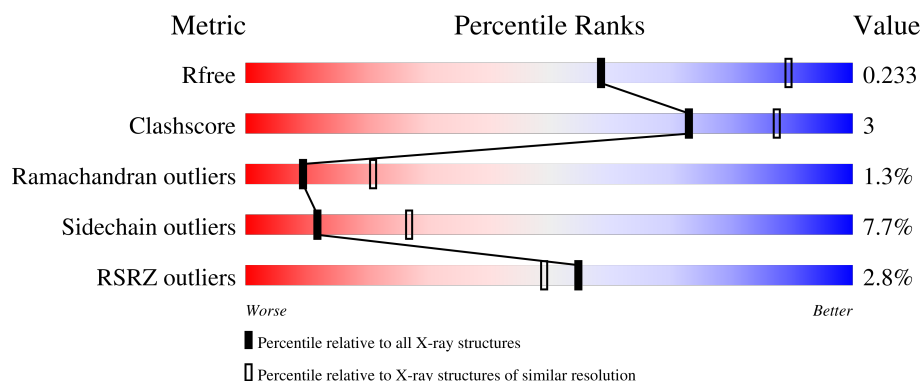
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	465	2% 65% 8% • 25%
1	B	465	3% 61% 12% • 25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6171 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 GLYCOGEN SYNTHASE KINASE-3 BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	351	Total	C	N	O	S	0	0	1
			2805	1804	482	508	11			
1	B	351	Total	C	N	O	S	0	0	1
			2805	1804	482	508	11			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-44	MET	-	expression tag	UNP P49841
A	-43	SER	-	expression tag	UNP P49841
A	-42	TYR	-	expression tag	UNP P49841
A	-41	TYR	-	expression tag	UNP P49841
A	-40	HIS	-	expression tag	UNP P49841
A	-39	HIS	-	expression tag	UNP P49841
A	-38	HIS	-	expression tag	UNP P49841
A	-37	HIS	-	expression tag	UNP P49841
A	-36	HIS	-	expression tag	UNP P49841
A	-35	HIS	-	expression tag	UNP P49841
A	-34	ASP	-	expression tag	UNP P49841
A	-33	TYR	-	expression tag	UNP P49841
A	-32	ASP	-	expression tag	UNP P49841
A	-31	ILE	-	expression tag	UNP P49841
A	-30	PRO	-	expression tag	UNP P49841
A	-29	THR	-	expression tag	UNP P49841
A	-28	THR	-	expression tag	UNP P49841
A	-27	GLU	-	expression tag	UNP P49841
A	-26	ASN	-	expression tag	UNP P49841
A	-25	LEU	-	expression tag	UNP P49841
A	-24	TYR	-	expression tag	UNP P49841
A	-23	PHE	-	expression tag	UNP P49841
A	-22	GLN	-	expression tag	UNP P49841
A	-21	GLY	-	expression tag	UNP P49841
A	-20	ALA	-	expression tag	UNP P49841

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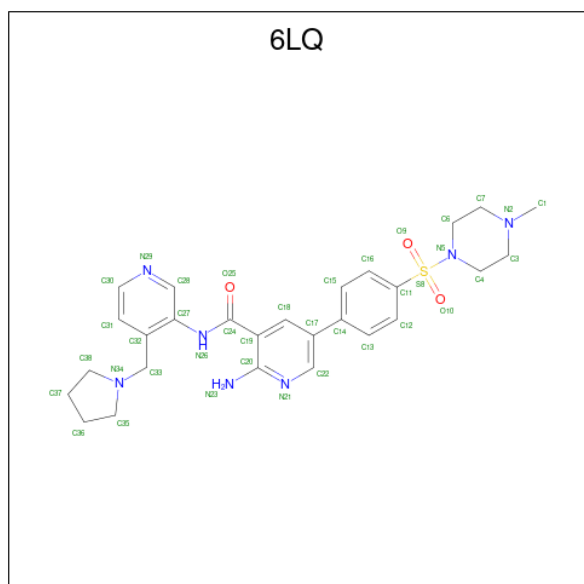
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P49841
A	-18	GLY	-	expression tag	UNP P49841
A	-17	SER	-	expression tag	UNP P49841
A	-16	SER	-	expression tag	UNP P49841
A	-15	HIS	-	expression tag	UNP P49841
A	-14	HIS	-	expression tag	UNP P49841
A	-13	HIS	-	expression tag	UNP P49841
A	-12	HIS	-	expression tag	UNP P49841
A	-11	HIS	-	expression tag	UNP P49841
A	-10	HIS	-	expression tag	UNP P49841
A	-9	SER	-	expression tag	UNP P49841
A	-8	SER	-	expression tag	UNP P49841
A	-7	GLY	-	expression tag	UNP P49841
A	-6	LEU	-	expression tag	UNP P49841
A	-5	VAL	-	expression tag	UNP P49841
A	-4	PRO	-	expression tag	UNP P49841
A	-3	ARG	-	expression tag	UNP P49841
A	-2	GLY	-	expression tag	UNP P49841
A	-1	SER	-	expression tag	UNP P49841
A	0	HIS	-	expression tag	UNP P49841
B	-44	MET	-	expression tag	UNP P49841
B	-43	SER	-	expression tag	UNP P49841
B	-42	TYR	-	expression tag	UNP P49841
B	-41	TYR	-	expression tag	UNP P49841
B	-40	HIS	-	expression tag	UNP P49841
B	-39	HIS	-	expression tag	UNP P49841
B	-38	HIS	-	expression tag	UNP P49841
B	-37	HIS	-	expression tag	UNP P49841
B	-36	HIS	-	expression tag	UNP P49841
B	-35	HIS	-	expression tag	UNP P49841
B	-34	ASP	-	expression tag	UNP P49841
B	-33	TYR	-	expression tag	UNP P49841
B	-32	ASP	-	expression tag	UNP P49841
B	-31	ILE	-	expression tag	UNP P49841
B	-30	PRO	-	expression tag	UNP P49841
B	-29	THR	-	expression tag	UNP P49841
B	-28	THR	-	expression tag	UNP P49841
B	-27	GLU	-	expression tag	UNP P49841
B	-26	ASN	-	expression tag	UNP P49841
B	-25	LEU	-	expression tag	UNP P49841
B	-24	TYR	-	expression tag	UNP P49841
B	-23	PHE	-	expression tag	UNP P49841

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-22	GLN	-	expression tag	UNP P49841
B	-21	GLY	-	expression tag	UNP P49841
B	-20	ALA	-	expression tag	UNP P49841
B	-19	MET	-	expression tag	UNP P49841
B	-18	GLY	-	expression tag	UNP P49841
B	-17	SER	-	expression tag	UNP P49841
B	-16	SER	-	expression tag	UNP P49841
B	-15	HIS	-	expression tag	UNP P49841
B	-14	HIS	-	expression tag	UNP P49841
B	-13	HIS	-	expression tag	UNP P49841
B	-12	HIS	-	expression tag	UNP P49841
B	-11	HIS	-	expression tag	UNP P49841
B	-10	HIS	-	expression tag	UNP P49841
B	-9	SER	-	expression tag	UNP P49841
B	-8	SER	-	expression tag	UNP P49841
B	-7	GLY	-	expression tag	UNP P49841
B	-6	LEU	-	expression tag	UNP P49841
B	-5	VAL	-	expression tag	UNP P49841
B	-4	PRO	-	expression tag	UNP P49841
B	-3	ARG	-	expression tag	UNP P49841
B	-2	GLY	-	expression tag	UNP P49841
B	-1	SER	-	expression tag	UNP P49841
B	0	HIS	-	expression tag	UNP P49841

- Molecule 2 is 2-AMINO-5-{4-[(4-METHYLPIPERAZIN-1-YL)SULFONYL]PHENYL}-N-[4-(PYRROLIDIN-1-YLMETHYL)PYRIDIN-3-YL]PYRIDINE-3-CARBOXAMIDE (CCD ID: 6LQ) (formula: C₂₇H₃₃N₇O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			38	27	7	3	1		
2	B	1	Total	C	N	O	S	0	0
			38	27	7	3	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	224	Total	O	0	0
			224	224		
3	B	261	Total	O	0	0
			261	261		

- Molecule 1: GLYCOGEN SYNTHASE KINASE-3 BETA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.98Å 83.88Å 177.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.97 – 2.60 19.97 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (19.97-2.60) 98.8 (19.97-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.59Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.175 , 0.225 0.181 , 0.233	Depositor DCC
R_{free} test set	1520 reflections (4.01%)	wwPDB-VP
Wilson B-factor (Å ²)	41.2	Xtriage
Anisotropy	0.646	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 57.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.036 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6171	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 3.45% 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: 6LQ

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	0.85	1/2876 (0.0%)	1.32	15/3913 (0.4%)
1	B	0.88	1/2876 (0.0%)	1.31	16/3913 (0.4%)
All	All	0.86	2/5752 (0.0%)	1.31	31/7826 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	384	ILE	C-N	-5.46	1.25	1.33
1	B	384	ILE	C-N	-5.33	1.25	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	91	LYS	CA-C-N	7.99	131.33	120.38
1	B	91	LYS	C-N-CA	7.99	131.33	120.38
1	B	136	PRO	N-CA-C	7.09	123.22	113.65
1	A	133	ASP	CA-CB-CG	6.79	119.39	112.60
1	A	136	PRO	N-CA-C	6.74	122.75	113.65
1	B	133	ASP	CA-CB-CG	6.57	119.17	112.60
1	A	79	GLY	CA-C-N	6.32	129.68	120.71
1	A	79	GLY	C-N-CA	6.32	129.68	120.71
1	B	317	CYS	CA-C-N	6.24	128.93	120.38
1	B	317	CYS	C-N-CA	6.24	128.93	120.38
1	A	89	GLN	CA-C-N	6.13	129.11	120.28
1	A	89	GLN	C-N-CA	6.13	129.11	120.28
1	B	79	GLY	CA-C-N	6.01	129.25	120.71
1	B	79	GLY	C-N-CA	6.01	129.25	120.71
1	A	66	SER	CA-C-N	5.97	129.08	120.79
1	A	66	SER	C-N-CA	5.97	129.08	120.79
1	A	377	LEU	N-CA-C	5.93	117.74	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	GLY	CA-C-N	5.80	132.62	121.54
1	B	120	GLY	C-N-CA	5.80	132.62	121.54
1	B	377	LEU	N-CA-C	5.75	117.55	111.28
1	A	264	ASP	CA-CB-CG	5.58	118.18	112.60
1	A	371	PRO	N-CA-C	5.57	117.49	110.70
1	A	91	LYS	CA-C-N	5.26	127.33	120.28
1	A	91	LYS	C-N-CA	5.26	127.33	120.28
1	B	371	PRO	N-CA-C	5.22	117.06	110.70
1	A	368	SER	N-CA-C	5.15	118.34	111.75
1	B	58	ASP	CA-CB-CG	5.13	117.73	112.60
1	A	352	ASN	CA-CB-CG	5.11	117.71	112.60
1	B	264	ASP	CA-CB-CG	5.04	117.64	112.60
1	B	368	SER	N-CA-C	5.02	117.40	111.33
1	B	72	GLN	N-CA-C	-5.01	101.99	109.95

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	2805	0	2833	17	0
1	B	2805	0	2833	20	0
2	A	38	0	33	1	0
2	B	38	0	33	1	0
3	A	224	0	0	2	0
3	B	261	0	0	0	0
All	All	6171	0	5732	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 3.

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)
1:B:352:ASN:HD21	1:B:354:ARG:HB2	1.67	0.58
1:B:337:HIS:HD2	1:B:339:PHE:H	1.52	0.57
1:A:352:ASN:HD21	1:A:354:ARG:HB2	1.73	0.53
1:A:352:ASN:HD22	1:A:354:ARG:H	1.57	0.52
1:B:90:ASP:C	1:B:92:ARG:H	2.17	0.52
1:A:135:VAL:HB	1:A:188:LEU:HB3	1.92	0.51
1:A:42:ALA:HB1	1:A:114:TYR:HB3	1.94	0.50
1:B:319:ARG:NH2	1:B:327:ALA:O	2.41	0.50
1:B:135:VAL:HB	1:B:188:LEU:HB3	1.92	0.49
1:B:42:ALA:HB1	1:B:114:TYR:HB3	1.95	0.49
1:B:135:VAL:O	2:B:1385:6LQ:H13	2.13	0.49
1:A:110:VAL:HG22	1:A:133:ASP:HB2	1.95	0.48
1:A:218:CYS:HB3	1:A:223:ARG:HG2	1.94	0.48
1:B:278:ARG:HG3	1:B:289:THR:HG23	1.97	0.47
1:B:218:CYS:HB3	1:B:223:ARG:HG2	1.98	0.45
1:A:185:GLN:O	2:A:1385:6LQ:H381	2.16	0.45
1:B:67:PHE:CE2	1:B:70:VAL:HG23	2.52	0.45
1:A:385:GLN:N	3:A:2224:HOH:O	2.50	0.44
1:A:123:LYS:C	1:A:125:GLU:H	2.26	0.44
1:A:252:LEU:HD11	1:A:256:ILE:HD12	2.00	0.44
1:A:162:MET:HE2	1:A:247:LEU:HB2	2.00	0.43
1:A:162:MET:HG3	1:A:247:LEU:HD13	2.00	0.43
1:B:252:LEU:HD11	1:B:256:ILE:HD12	2.00	0.43
1:A:352:ASN:ND2	1:A:354:ARG:HB2	2.33	0.43
1:A:220:ARG:HD3	1:B:220:ARG:HD3	2.00	0.43
1:B:162:MET:HG3	1:B:247:LEU:HD13	2.01	0.43
1:B:67:PHE:HE2	1:B:70:VAL:HG23	1.83	0.42
1:B:162:MET:HE2	1:B:247:LEU:HB2	2.00	0.42
1:A:151:GLN:HG3	3:A:2081:HOH:O	2.20	0.42
1:B:288:TYR:HA	1:B:291:PHE:HD2	1.84	0.42
1:A:155:VAL:HG12	1:A:159:LYS:HE3	2.02	0.42
1:B:155:VAL:HG12	1:B:159:LYS:HE3	2.02	0.42
1:B:137:GLU:OE2	1:B:141:ARG:NH1	2.54	0.41
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.96	0.40
1:B:266:LEU:HD12	1:B:266:LEU:HA	1.96	0.40
1:B:310:PRO:HA	1:B:311:PRO:HD3	2.00	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	349/465 (75%)	323 (93%)	22 (6%)	4 (1%)	11	25
1	B	349/465 (75%)	323 (93%)	21 (6%)	5 (1%)	9	19
All	All	698/930 (75%)	646 (93%)	43 (6%)	9 (1%)	9	21

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	121	GLU
1	A	77	ASP
1	A	120	GLY
1	A	221	TYR
1	B	221	TYR
1	B	295	GLN
1	A	295	GLN
1	B	120	GLY
1	B	384	ILE

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	312/404 (77%)	288 (92%)	24 (8%)	12	27
1	B	312/404 (77%)	288 (92%)	24 (8%)	12	27
All	All	624/808 (77%)	576 (92%)	48 (8%)	12	27

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LYS
1	A	64	ASN
1	A	89	GLN
1	A	91	LYS
1	A	99	GLN
1	A	110	VAL
1	A	121	GLU
1	A	122	LYS
1	A	147	SER
1	A	151	GLN
1	A	185	GLN
1	A	206	GLN
1	A	218	CYS
1	A	220	ARG
1	A	256	ILE
1	A	266	LEU
1	A	275	THR
1	A	289	THR
1	A	292	LYS
1	A	295	GLN
1	A	312	GLU
1	A	316	LEU
1	A	352	ASN
1	A	384	ILE
1	B	64	ASN
1	B	66	SER
1	B	89	GLN
1	B	110	VAL
1	B	111	ARG
1	B	125	GLU
1	B	126	VAL
1	B	141	ARG
1	B	147	SER
1	B	185	GLN
1	B	206	GLN
1	B	218	CYS
1	B	220	ARG
1	B	256	ILE
1	B	266	LEU
1	B	275	THR
1	B	283	GLU
1	B	290	GLU

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Mol	Chain	Res	Type
1	B	295	GLN
1	B	303	LYS
1	B	312	GLU
1	B	316	LEU
1	B	318	SER
1	B	350	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	129	ASN
1	A	352	ASN
1	B	46	GLN
1	B	337	HIS
1	B	352	ASN
1	B	381	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 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
2	6LQ	B	1385	-	42,42,42	0.94	2 (4%)	57,60,60	2.34	15 (26%)
2	6LQ	A	1385	-	42,42,42	0.81	0	57,60,60	2.34	13 (22%)

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	6LQ	B	1385	-	-	3/28/45/45	0/5/5/5
2	6LQ	A	1385	-	-	3/28/45/45	0/5/5/5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1385	6LQ	S8-N5	2.09	1.66	1.63
2	B	1385	6LQ	C20-N23	2.07	1.39	1.34

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1385	6LQ	C6-N5-C4	9.79	123.37	112.12
2	B	1385	6LQ	C6-N5-C4	9.08	122.55	112.12
2	B	1385	6LQ	C20-C19-C24	7.99	123.56	120.66
2	A	1385	6LQ	C20-C19-C24	7.17	123.26	120.66
2	A	1385	6LQ	C38-N34-C35	5.35	110.36	104.04
2	A	1385	6LQ	O10-S8-C11	-4.50	102.50	108.10
2	B	1385	6LQ	C22-N21-C20	4.36	122.86	118.64
2	A	1385	6LQ	C7-C6-N5	3.96	113.17	108.97
2	A	1385	6LQ	C3-C4-N5	3.78	112.98	108.97
2	B	1385	6LQ	C38-N34-C35	3.68	108.39	104.04
2	B	1385	6LQ	O10-S8-N5	3.64	110.11	106.69
2	B	1385	6LQ	C33-C32-C27	3.57	124.35	121.09
2	A	1385	6LQ	C22-N21-C20	3.54	122.06	118.64
2	B	1385	6LQ	C30-N29-C28	3.52	123.02	116.85
2	B	1385	6LQ	C7-C6-N5	3.33	112.50	108.97
2	B	1385	6LQ	C3-C4-N5	3.02	112.17	108.97
2	B	1385	6LQ	O10-S8-C11	-2.90	104.50	108.10
2	B	1385	6LQ	C16-C11-S8	2.84	122.56	119.73
2	A	1385	6LQ	C30-N29-C28	2.74	121.65	116.85
2	A	1385	6LQ	O10-S8-N5	2.64	109.17	106.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1385	6LQ	C11-S8-N5	-2.51	104.25	107.28
2	B	1385	6LQ	C4-C3-N2	-2.21	107.30	110.86
2	A	1385	6LQ	O9-S8-N5	2.19	108.75	106.69
2	A	1385	6LQ	C33-C32-C27	2.15	123.06	121.09
2	B	1385	6LQ	C31-C30-N29	-2.15	119.93	123.60
2	B	1385	6LQ	C17-C22-N21	-2.10	120.94	124.28
2	A	1385	6LQ	C18-C17-C22	2.02	119.27	117.10
2	A	1385	6LQ	O10-S8-O9	2.02	122.74	119.59

There are no chirality outliers.

All (6) torsion outliers are listed below:

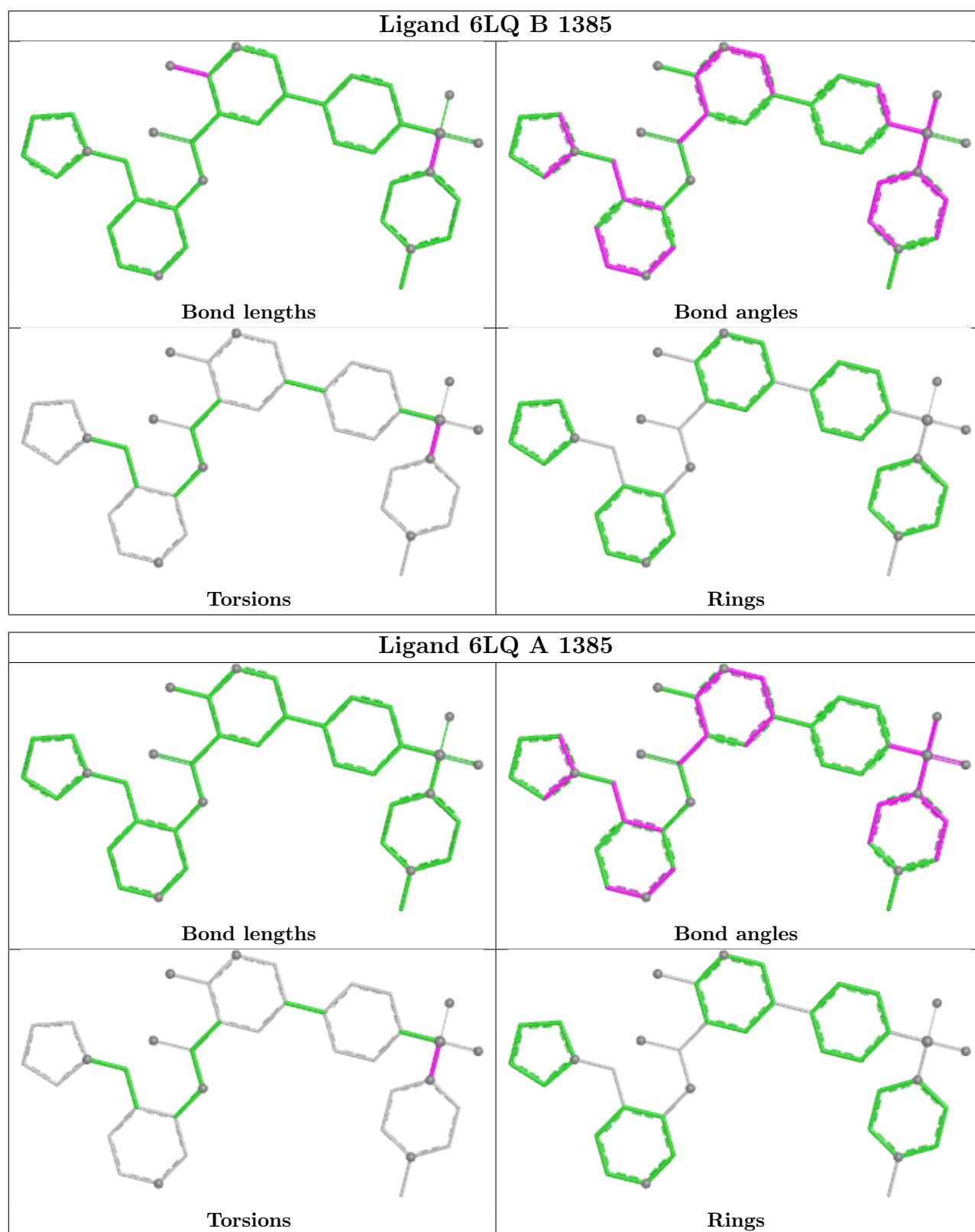
Mol	Chain	Res	Type	Atoms
2	B	1385	6LQ	C4-N5-S8-O10
2	A	1385	6LQ	C4-N5-S8-O10
2	B	1385	6LQ	C4-N5-S8-O9
2	A	1385	6LQ	C6-N5-S8-O9
2	B	1385	6LQ	C4-N5-S8-C11
2	A	1385	6LQ	C4-N5-S8-O9

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1385	6LQ	1	0
2	A	1385	6LQ	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	351/465 (75%)	-0.21	7 (1%) 65 60	26, 47, 110, 133	0
1	B	351/465 (75%)	-0.37	13 (3%) 45 39	23, 42, 104, 152	0
All	All	702/930 (75%)	-0.29	20 (2%) 55 49	23, 44, 105, 152	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	385	GLN	4.5
1	B	120	GLY	3.6
1	B	384	ILE	3.3
1	A	289	THR	3.3
1	A	286	PRO	3.2
1	A	296	ILE	3.0
1	A	294	PRO	2.9
1	B	93	PHE	2.8
1	B	91	LYS	2.6
1	B	122	LYS	2.4
1	B	119	SER	2.4
1	B	288	TYR	2.3
1	B	298	ALA	2.2
1	A	288	TYR	2.2
1	A	298	ALA	2.1
1	B	294	PRO	2.1
1	A	291	PHE	2.1
1	B	35	SER	2.1
1	B	289	THR	2.0
1	B	295	GLN	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](#)

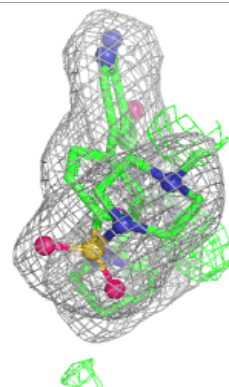
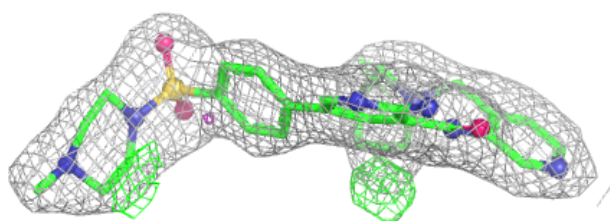
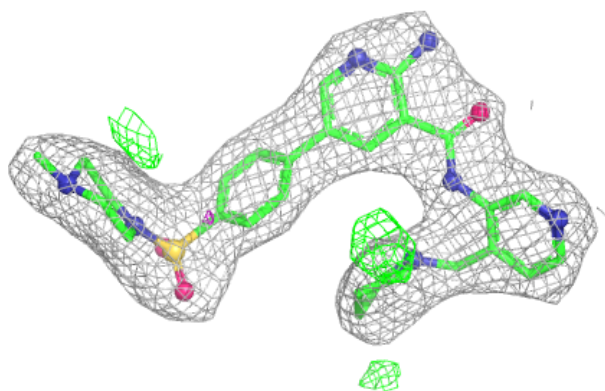
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	6LQ	A	1385	38/38	0.95	0.08	30,40,64,65	0
2	6LQ	B	1385	38/38	0.95	0.08	34,40,53,58	0

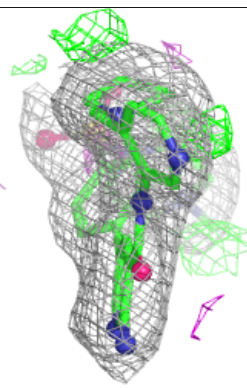
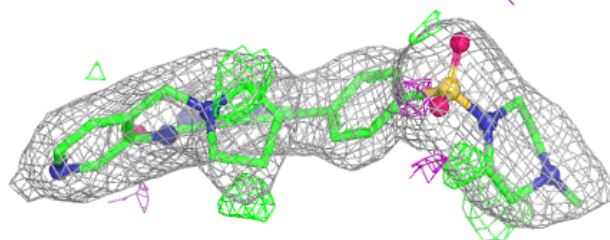
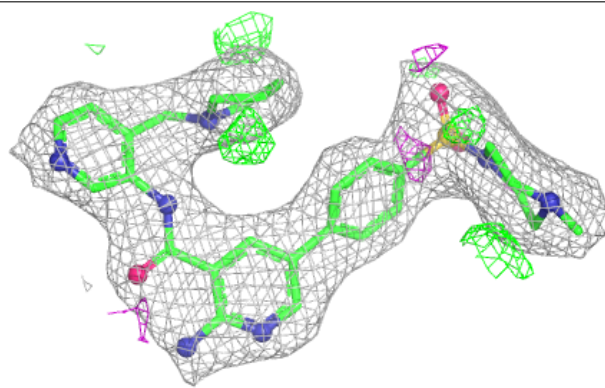
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 6LQ A 1385:

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

**Electron density around 6LQ B 1385:**

$2mF_o-DF_c$ (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.