



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:43 PM UTC

PDB ID : 4IQ6 / pdb_00004iq6
Title : Gsk-3beta with inhibitor 6-chloro-N-cyclohexyl-4-(1H-pyrrolo[2,3-b]pyridin-3-yl)pyridin-2-amine
Authors : Tong, Y.; Stewart, K.D.; Florjancic, A.S.; Harlan, J.E.; Merta, P.J.; Przytulinska, M.; Soni, N.; Swinger, K.S.; Zhu, H.; Johnson, E.F.; Shoemaker, A.R.; Penning, T.D.
Deposited on : 2013-01-10
Resolution : 3.12 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](#) ①) 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

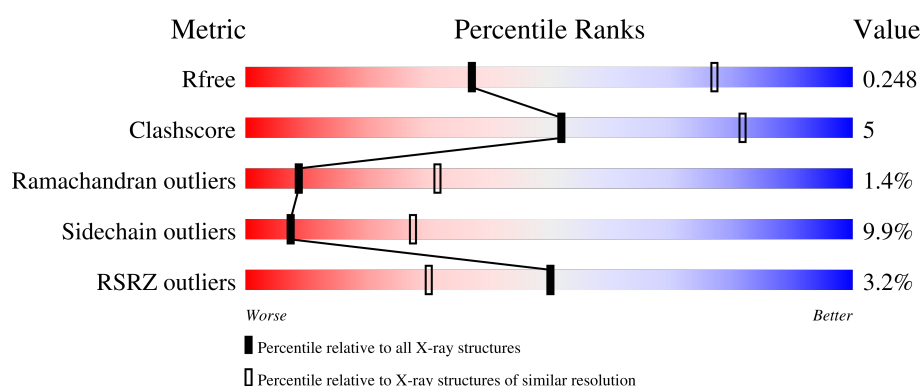
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.12 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	1816 (3.14-3.10)
Clashscore	190562	1906 (3.14-3.10)
Ramachandran outliers	187476	1802 (3.14-3.10)
Sidechain outliers	187428	1802 (3.14-3.10)
RSRZ outliers	180081	1816 (3.14-3.10)

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	426	4% 60% 14% • 24%
1	B	426	% 60% 16% • 22%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5164 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.

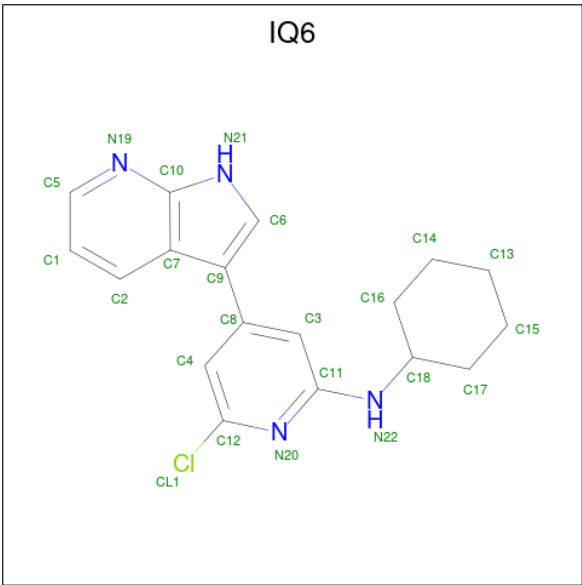
- Molecule 1 is a protein called Glycogen synthase kinase-3 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	324	Total	C	N	O	S	0	0	0
			2522	1625	426	460	11			
1	B	332	Total	C	N	O	S	0	0	0
			2596	1672	440	473	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	GLY	-	expression tag	UNP P49841
A	-4	ALA	-	expression tag	UNP P49841
A	-3	MET	-	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	-5	GLY	-	expression tag	UNP P49841
B	-4	ALA	-	expression tag	UNP P49841
B	-3	MET	-	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 6-chloro-N-cyclohexyl-4-(1H-pyrrolo[2,3-b]pyridin-3-yl)pyridin-2-amine (CCD ID: IQ6) (formula: C₁₈H₁₉ClN₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	0	0
			23	18	1	4		
2	B	1	Total	C	Cl	N	0	0
			23	18	1	4		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.83Å 84.99Å 178.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	89.03 – 3.12 89.03 – 3.12	Depositor EDS
% Data completeness (in resolution range)	99.9 (89.03-3.12) 100.0 (89.03-3.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 3.13Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.186 , 0.227 0.201 , 0.248	Depositor DCC
R_{free} test set	1183 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.821	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 70.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.031 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5164	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% 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: IQ6

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	0.88	1/2583 (0.0%)	1.39	18/3521 (0.5%)
1	B	0.87	2/2660 (0.1%)	1.37	18/3621 (0.5%)
All	All	0.88	3/5243 (0.1%)	1.38	36/7142 (0.5%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	284	MET	SD-CE	-9.00	1.57	1.79
1	B	284	MET	SD-CE	-8.03	1.59	1.79
1	B	377	LEU	C-N	5.92	1.38	1.33

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	370	ASN	CA-CB-CG	8.69	121.29	112.60
1	A	370	ASN	CA-CB-CG	8.61	121.21	112.60
1	B	371	PRO	N-CA-C	7.95	120.40	110.70
1	A	307	PRO	N-CA-C	7.90	124.76	113.47
1	A	371	PRO	N-CA-C	7.65	120.03	110.70
1	B	66	SER	CA-C-N	6.51	129.88	120.38
1	B	66	SER	C-N-CA	6.51	129.88	120.38
1	A	311	PRO	CA-C-N	6.28	128.70	120.28
1	A	311	PRO	C-N-CA	6.28	128.70	120.28
1	A	325	PRO	CA-C-N	6.26	129.19	120.29
1	A	325	PRO	C-N-CA	6.26	129.19	120.29
1	B	377	LEU	N-CA-C	6.26	118.90	111.33
1	A	66	SER	CA-C-N	6.20	129.44	120.38
1	A	66	SER	C-N-CA	6.20	129.44	120.38
1	B	325	PRO	CA-C-N	6.13	129.00	120.29
1	B	325	PRO	C-N-CA	6.13	129.00	120.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	293	PHE	CA-CB-CG	6.11	119.91	113.80
1	B	311	PRO	CA-C-N	6.09	128.44	120.28
1	B	311	PRO	C-N-CA	6.09	128.44	120.28
1	A	195	VAL	N-CA-CB	6.05	117.13	110.53
1	A	377	LEU	N-CA-C	6.02	118.33	111.11
1	A	172	ILE	N-CA-CB	5.94	117.90	110.47
1	A	190	ASP	N-CA-C	-5.93	100.99	109.24
1	B	190	ASP	N-CA-C	-5.83	100.17	109.04
1	B	172	ILE	N-CA-CB	5.82	117.74	110.47
1	B	133	ASP	CA-CB-CG	5.70	118.30	112.60
1	B	195	VAL	N-CA-CB	5.45	116.66	110.72
1	A	91	LYS	CA-C-N	5.42	128.95	120.82
1	A	91	LYS	C-N-CA	5.42	128.95	120.82
1	A	133	ASP	CA-CB-CG	5.41	118.00	112.60
1	B	91	LYS	CA-C-N	5.39	128.90	120.82
1	B	91	LYS	C-N-CA	5.39	128.90	120.82
1	A	77	ASP	CA-C-N	5.22	127.28	120.28
1	A	77	ASP	C-N-CA	5.22	127.28	120.28
1	B	294	PRO	N-CA-C	5.14	119.30	111.34
1	B	39	THR	CB-CA-C	5.05	118.22	110.14

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	2522	0	2498	23	0
1	B	2596	0	2544	24	0
2	A	23	0	19	4	0
2	B	23	0	19	2	0
All	All	5164	0	5080	52	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 5.

All (52) 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:337:HIS:HD2	1:B:339:PHE:H	1.33	0.77
1:A:266:LEU:O	1:A:270:ILE:HG12	1.94	0.66
1:A:284:MET:HE1	1:A:323:TYR:HD1	1.58	0.66
1:B:266:LEU:O	1:B:270:ILE:HG12	1.95	0.66
1:B:284:MET:HE1	1:B:323:TYR:HD1	1.61	0.65
1:A:284:MET:HE1	1:A:323:TYR:CD1	2.35	0.61
1:A:215:SER:O	1:A:223:ARG:NH2	2.34	0.61
1:A:283:GLU:HG3	1:A:324:THR:HG23	1.83	0.60
1:B:215:SER:O	1:B:223:ARG:NH2	2.34	0.59
1:A:332:LEU:HD11	1:A:367:LEU:HA	1.88	0.56
1:A:162:MET:HG3	1:A:247:LEU:HD13	1.89	0.54
1:B:332:LEU:HD11	1:B:367:LEU:HA	1.89	0.53
1:B:284:MET:HE1	1:B:323:TYR:CD1	2.43	0.52
2:A:501:IQ6:H3	2:A:501:IQ6:H13	1.92	0.52
2:A:501:IQ6:H13	2:A:501:IQ6:H2	1.91	0.52
1:B:257:PHE:HE1	1:B:272:VAL:HG13	1.75	0.51
1:A:96:ARG:CZ	1:A:205:LYS:HG3	2.41	0.50
1:A:106:HIS:HB3	1:A:109:ILE:HD12	1.93	0.50
2:A:501:IQ6:H3	2:A:501:IQ6:H2	1.93	0.50
1:B:162:MET:HG3	1:B:247:LEU:HD13	1.93	0.50
1:A:257:PHE:HE1	1:A:272:VAL:HG13	1.76	0.50
1:B:193:THR:OG1	1:B:195:VAL:HG13	2.11	0.49
1:B:273:LEU:HB3	1:B:323:TYR:CE2	2.48	0.49
1:B:96:ARG:CZ	1:B:205:LYS:HG3	2.42	0.49
1:A:273:LEU:HB3	1:A:323:TYR:CE2	2.47	0.48
1:A:144:ARG:HG2	1:A:148:ARG:HE	1.79	0.48
1:A:319:ARG:HB3	1:A:329:LEU:HD13	1.95	0.48
1:A:258:PRO:HD2	1:A:268:GLU:HG3	1.96	0.47
1:A:193:THR:OG1	1:A:195:VAL:HG13	2.14	0.47
1:A:139:VAL:CG2	1:A:189:LEU:HD23	2.45	0.47
1:B:258:PRO:HD2	1:B:268:GLU:HG3	1.97	0.47
2:B:501:IQ6:H3	2:B:501:IQ6:H2	1.98	0.46
1:B:319:ARG:HB3	1:B:329:LEU:HD13	1.97	0.45
1:B:305:PHE:CD2	1:B:314:ILE:HG12	2.52	0.45
1:B:271:LYS:O	1:B:299:HIS:HB2	2.17	0.44
2:B:501:IQ6:H3	2:B:501:IQ6:H17	1.38	0.43
1:B:88:LEU:HD12	1:B:127:TYR:CE1	2.54	0.43
1:B:278:ARG:HH12	1:B:282:ARG:HD2	1.82	0.43
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.88	0.43
2:A:501:IQ6:H3	2:A:501:IQ6:C2	2.48	0.43
1:A:173:HIS:C	1:A:175:PHE:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:ARG:HD3	1:B:203:SER:O	2.19	0.42
1:B:173:HIS:C	1:B:175:PHE:H	2.28	0.42
1:A:66:SER:HB2	1:B:267:VAL:HG11	2.01	0.41
1:A:305:PHE:CD1	1:A:314:ILE:HG12	2.55	0.41
1:B:224:ALA:HA	1:B:241:TRP:CD1	2.56	0.41
1:B:217:ILE:H	1:B:217:ILE:HG13	1.70	0.41
1:A:221:TYR:CD1	1:A:258:PRO:HA	2.56	0.41
1:A:224:ALA:HA	1:A:241:TRP:CD1	2.55	0.41
1:A:139:VAL:HG23	1:A:189:LEU:HD23	2.03	0.41
1:B:266:LEU:HD12	1:B:266:LEU:HA	1.88	0.41
1:B:337:HIS:CD2	1:B:339:PHE:H	2.24	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	318/426 (75%)	290 (91%)	22 (7%)	6 (2%)	6	25
1	B	326/426 (76%)	301 (92%)	22 (7%)	3 (1%)	14	42
All	All	644/852 (76%)	591 (92%)	44 (7%)	9 (1%)	9	31

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	ASP
1	B	49	ASP
1	A	221	TYR
1	B	221	TYR
1	A	301	TRP
1	A	174	SER
1	A	361	ASN

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Mol	Chain	Res	Type
1	A	381	HIS
1	B	174	SER

5.3.2 Protein sidechains ⓘ

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	273/367 (74%)	249 (91%)	24 (9%)	9	31
1	B	280/367 (76%)	249 (89%)	31 (11%)	6	22
All	All	553/734 (75%)	498 (90%)	55 (10%)	7	27

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	GLN
1	A	57	THR
1	A	59	THR
1	A	69	VAL
1	A	72	GLN
1	A	77	ASP
1	A	110	VAL
1	A	148	ARG
1	A	151	GLN
1	A	169	LEU
1	A	180	ARG
1	A	195	VAL
1	A	213	ASN
1	A	254	GLN
1	A	266	LEU
1	A	270	ILE
1	A	272	VAL
1	A	275	THR
1	A	279	GLU
1	A	324	THR
1	A	326	THR

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Mol	Chain	Res	Type
1	A	329	LEU
1	A	348	VAL
1	A	355	ASP
1	B	52	GLN
1	B	57	THR
1	B	59	THR
1	B	72	GLN
1	B	77	ASP
1	B	110	VAL
1	B	119	SER
1	B	148	ARG
1	B	151	GLN
1	B	152	THR
1	B	156	ILE
1	B	169	LEU
1	B	180	ARG
1	B	189	LEU
1	B	195	VAL
1	B	213	ASN
1	B	217	ILE
1	B	254	GLN
1	B	266	LEU
1	B	270	ILE
1	B	272	VAL
1	B	273	LEU
1	B	275	THR
1	B	279	GLU
1	B	297	LYS
1	B	318	SER
1	B	324	THR
1	B	326	THR
1	B	329	LEU
1	B	348	VAL
1	B	355	ASP

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	64	ASN
1	A	72	GLN
1	A	254	GLN
1	A	381	HIS

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Mol	Chain	Res	Type
1	B	72	GLN
1	B	254	GLN
1	B	299	HIS
1	B	337	HIS
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	IQ6	B	501	-	26,26,26	0.96	3 (11%)	33,36,36	1.80	12 (36%)
2	IQ6	A	501	-	26,26,26	1.13	4 (15%)	33,36,36	2.01	10 (30%)

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	IQ6	B	501	-	-	5/8/16/16	0/4/4/4
2	IQ6	A	501	-	-	3/8/16/16	0/4/4/4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	IQ6	C8-C9	-3.36	1.44	1.49
2	B	501	IQ6	C8-C9	-2.66	1.45	1.49
2	A	501	IQ6	C11-N22	2.49	1.42	1.37
2	A	501	IQ6	C7-C10	-2.28	1.39	1.41
2	B	501	IQ6	C7-C9	-2.08	1.42	1.47
2	A	501	IQ6	C7-C9	-2.07	1.42	1.47
2	B	501	IQ6	C11-N22	2.00	1.41	1.37

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	IQ6	C9-C6-N21	-4.58	107.50	110.46
2	A	501	IQ6	C10-N21-C6	4.20	110.46	108.48
2	A	501	IQ6	N22-C11-N20	4.18	123.08	117.03
2	B	501	IQ6	C16-C18-N22	3.04	115.68	110.77
2	B	501	IQ6	C5-N19-C10	-3.03	110.84	116.74
2	A	501	IQ6	C4-C8-C9	2.75	124.31	120.74
2	B	501	IQ6	C9-C6-N21	-2.73	108.70	110.46
2	B	501	IQ6	C7-C9-C8	2.69	131.49	126.62
2	A	501	IQ6	C7-C9-C6	2.63	107.63	105.98
2	A	501	IQ6	C5-N19-C10	-2.60	111.69	116.74
2	A	501	IQ6	C3-C11-N22	-2.56	116.53	122.07
2	A	501	IQ6	C8-C3-C11	2.55	122.45	118.45
2	A	501	IQ6	C3-C8-C9	-2.52	117.47	120.74
2	B	501	IQ6	C4-C12-CL1	-2.49	115.72	118.92
2	B	501	IQ6	C10-N21-C6	2.40	109.61	108.48
2	B	501	IQ6	C1-C2-C7	-2.40	115.46	119.80
2	A	501	IQ6	C7-C10-N21	-2.39	107.55	108.42
2	B	501	IQ6	C10-C7-C9	2.28	107.01	106.22
2	B	501	IQ6	C17-C18-N22	2.26	114.41	110.77
2	B	501	IQ6	C1-C5-N19	2.15	126.82	123.42
2	B	501	IQ6	C4-C8-C3	-2.13	117.16	119.65
2	B	501	IQ6	C14-C16-C18	-2.04	107.44	111.09

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	IQ6	C3-C11-N22-C18
2	A	501	IQ6	N20-C11-N22-C18
2	B	501	IQ6	C3-C8-C9-C7
2	B	501	IQ6	C4-C8-C9-C7
2	A	501	IQ6	C4-C8-C9-C7
2	B	501	IQ6	C3-C8-C9-C6
2	B	501	IQ6	C4-C8-C9-C6
2	B	501	IQ6	N20-C11-N22-C18

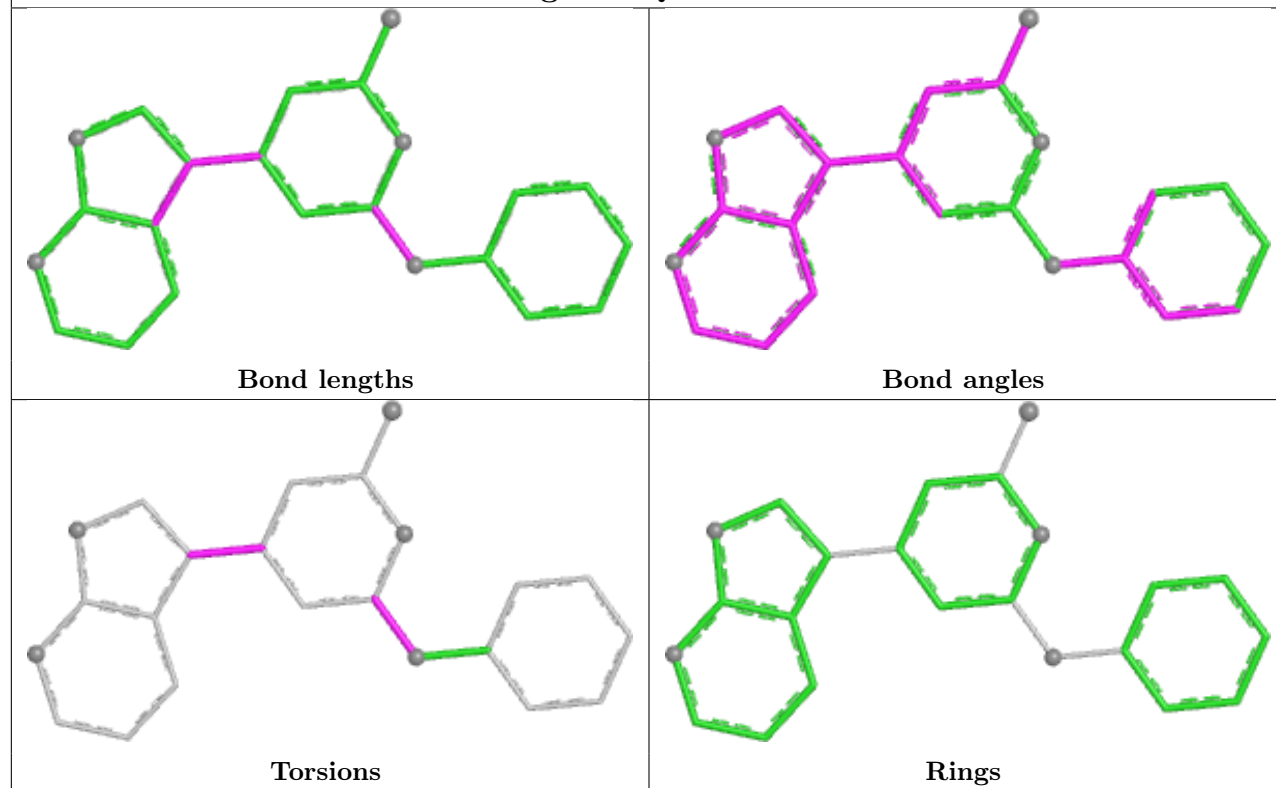
There are no ring outliers.

2 monomers are involved in 6 short contacts:

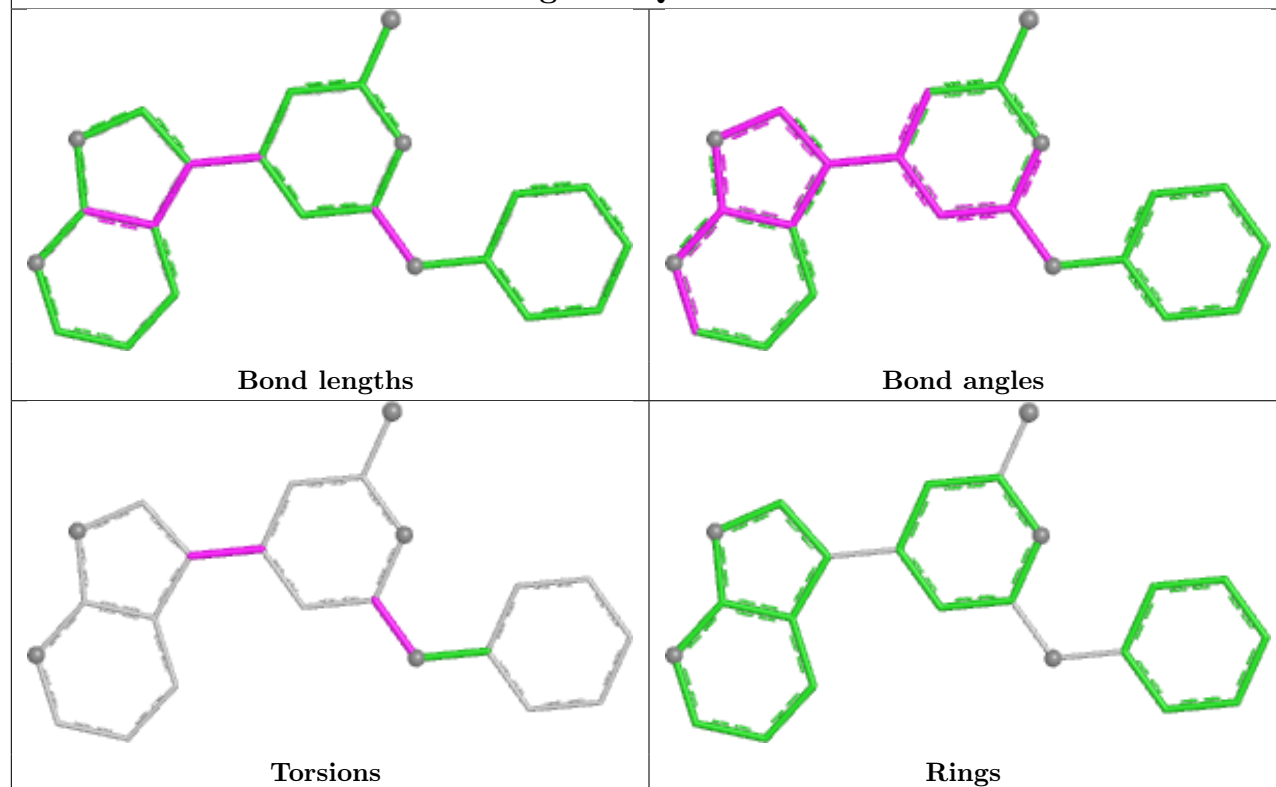
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	IQ6	2	0
2	A	501	IQ6	4	0

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.

Ligand IQ6 B 501



Ligand IQ6 A 501



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 [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	324/426 (76%)	0.05	15 (4%) 37 20	36, 55, 92, 103	0
1	B	332/426 (77%)	-0.06	6 (1%) 67 46	32, 51, 93, 111	0
All	All	656/852 (76%)	-0.01	21 (3%) 50 30	32, 53, 93, 111	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	300	PRO	4.6
1	A	285	ASN	4.4
1	B	293	PHE	4.2
1	B	90	ASP	4.2
1	A	90	ASP	3.8
1	A	127	TYR	3.5
1	A	118	SER	3.2
1	B	294	PRO	3.1
1	A	65	GLY	3.0
1	A	66	SER	2.9
1	B	119	SER	2.8
1	A	93	PHE	2.6
1	B	283	GLU	2.6
1	A	382	ALA	2.5
1	A	67	PHE	2.5
1	A	91	LYS	2.5
1	A	46	GLN	2.2
1	A	64	ASN	2.2
1	A	145	HIS	2.2
1	A	94	LYS	2.0
1	B	150	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](#)

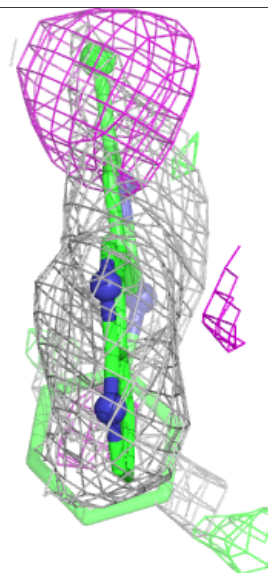
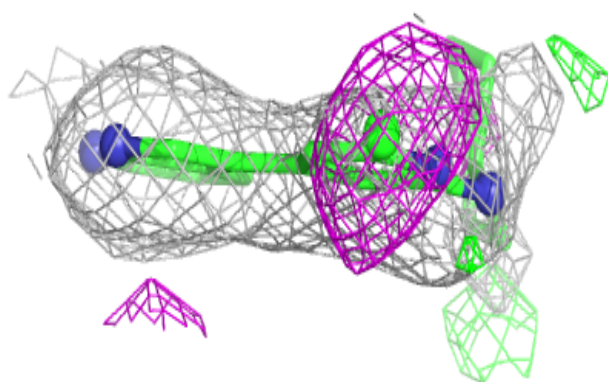
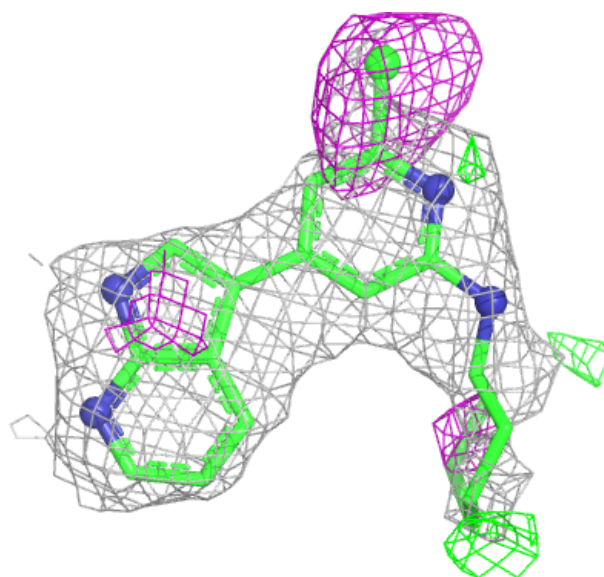
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	IQ6	B	501	23/23	0.82	0.22	69,82,87,88	0
2	IQ6	A	501	23/23	0.83	0.20	73,79,96,97	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.

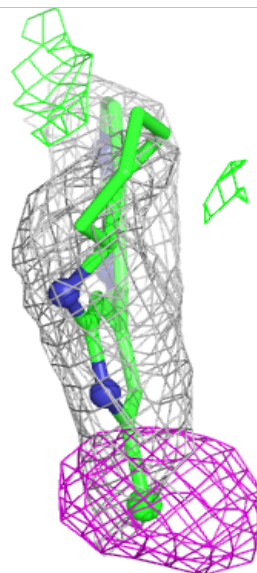
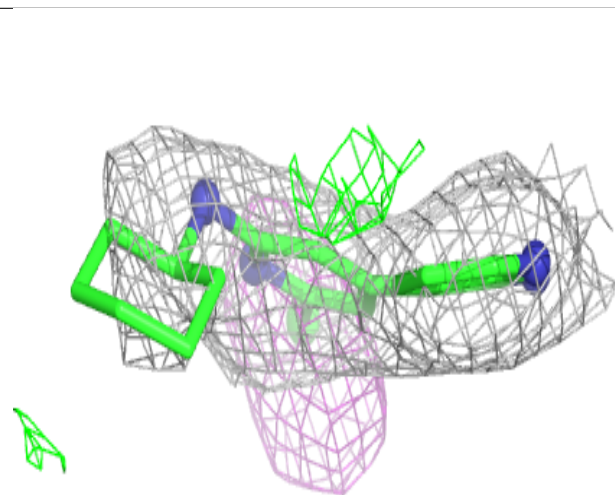
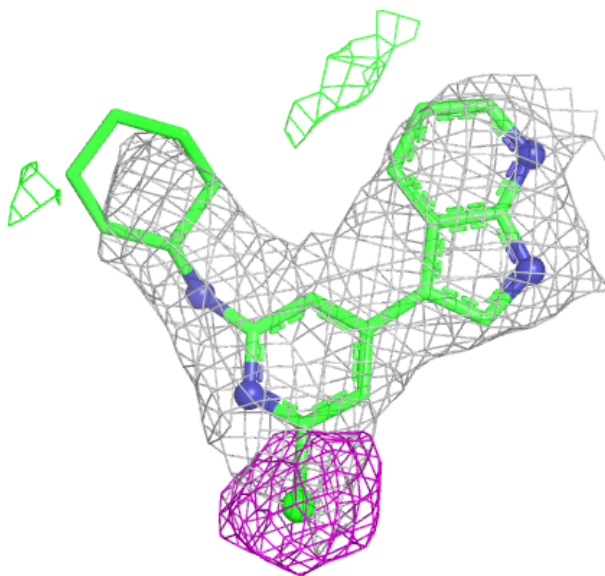
Electron density around IQ6 B 501:

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



Electron density around IQ6 A 501:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.