



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:23 PM UTC

PDB ID : 1Q41 / pdb_00001q41
Title : GSK-3 Beta complexed with Indirubin-3'-monoxime
Authors : Bertrand, J.A.; Thieffine, S.; Vulpetti, A.; Cristiani, C.; Valsasina, B.; Knapp, S.; Kalisz, H.M.; Flocco, M.
Deposited on : 2003-08-01
Resolution : 2.10 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

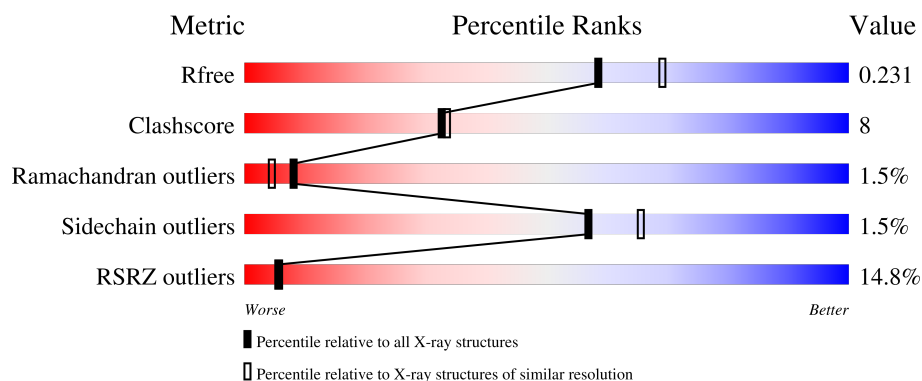
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	424	 11% 63% 15% • 20%
1	B	424	 12% 67% 13% • 19%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5617 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.

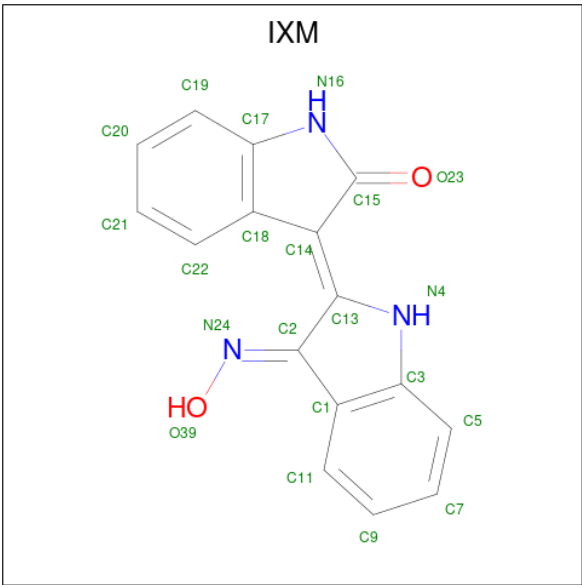
- Molecule 1 is a protein called GLYCOGEN SYNTHASE KINASE-3 BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	0	0
			2694	1736	463	484	11			
1	B	343	Total	C	N	O	S	0	0	0
			2734	1762	468	493	11			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	cloning artifact	UNP P49841
A	-2	PRO	-	cloning artifact	UNP P49841
A	-1	LEU	-	cloning artifact	UNP P49841
A	0	GLY	-	cloning artifact	UNP P49841
A	1	SER	-	cloning artifact	UNP P49841
B	-3	GLY	-	cloning artifact	UNP P49841
B	-2	PRO	-	cloning artifact	UNP P49841
B	-1	LEU	-	cloning artifact	UNP P49841
B	0	GLY	-	cloning artifact	UNP P49841
B	1	SER	-	cloning artifact	UNP P49841

- Molecule 2 is (Z)-1H,1'H-[2,3']BIINDOLYLIDENE-3,2'-DIONE-3-OXIME (CCD ID: IXM) (formula: C₁₆H₁₁N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	16	3	2		
2	B	1	Total	C	N	O	0	0
			21	16	3	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	69	Total	O	0	0
			69	69		
3	B	78	Total	O	0	0
			78	78		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.44Å 86.44Å 178.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.76 – 2.10 28.76 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.7 (28.76-2.10) 99.8 (28.76-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.83 (at 2.10Å)	Xtriage
Refinement program	CNX 2000	Depositor
R, R_{free}	0.229 , 0.245 (Not available) , 0.231	Depositor DCC
R_{free} test set	3827 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtriage
Anisotropy	0.709	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5617	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: IXM

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	0.67	1/2760 (0.0%)	1.01	17/3755 (0.5%)
1	B	0.69	0/2804	1.01	14/3818 (0.4%)
All	All	0.68	1/5564 (0.0%)	1.01	31/7573 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	162	MET	SD-CE	5.19	1.92	1.79

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	136	PRO	N-CA-C	8.50	124.48	113.86
1	B	190	ASP	CA-C-N	7.76	127.31	119.24
1	B	190	ASP	C-N-CA	7.76	127.31	119.24
1	A	72	GLN	N-CA-C	-6.69	99.31	109.95
1	A	330	THR	N-CA-C	-6.33	102.85	110.13
1	A	190	ASP	CA-C-N	6.28	126.47	119.32
1	A	190	ASP	C-N-CA	6.28	126.47	119.32
1	B	149	ALA	N-CA-C	-6.16	104.16	113.89
1	A	69	VAL	N-CA-C	6.14	117.50	109.58
1	A	136	PRO	N-CA-C	6.12	122.69	114.18
1	B	78	SER	N-CA-C	6.00	120.21	113.01
1	B	377	LEU	N-CA-C	5.94	117.75	111.28
1	A	368	SER	N-CA-C	5.68	119.02	111.75
1	A	94	LYS	N-CA-C	-5.60	101.41	110.32
1	B	285	ASN	CA-C-N	5.57	125.93	119.47
1	B	285	ASN	C-N-CA	5.57	125.93	119.47
1	B	72	GLN	N-CA-C	-5.55	101.12	109.95
1	B	371	PRO	N-CA-C	5.41	117.30	110.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	48	PRO	N-CA-C	-5.36	103.43	111.57
1	A	371	PRO	N-CA-C	5.35	117.22	110.70
1	B	368	SER	N-CA-C	5.31	118.54	111.75
1	A	377	LEU	N-CA-C	5.30	117.05	111.28
1	B	192	ASP	N-CA-C	5.27	118.87	112.23
1	B	345	ASP	CA-C-N	5.26	125.57	119.47
1	B	345	ASP	C-N-CA	5.26	125.57	119.47
1	A	304	VAL	N-CA-C	5.21	115.43	110.53
1	A	252	LEU	N-CA-C	5.20	118.78	112.23
1	A	155	VAL	N-CA-C	5.19	117.53	111.05
1	A	192	ASP	N-CA-C	5.08	118.64	112.23
1	A	345	ASP	CA-C-N	5.04	125.32	119.47
1	A	345	ASP	C-N-CA	5.04	125.32	119.47

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	2694	0	2724	44	0
1	B	2734	0	2751	44	0
2	A	21	0	11	1	0
2	B	21	0	11	0	0
3	A	69	0	0	2	0
3	B	78	0	0	1	0
All	All	5617	0	5497	87	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 8.

All (87) 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:A:220:ARG:CD	1:A:221:TYR:H	1.61	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:ARG:HD2	1:A:221:TYR:H	1.15	1.01
1:A:220:ARG:HD2	1:A:221:TYR:N	1.78	0.97
1:B:220:ARG:CD	1:B:221:TYR:H	1.81	0.92
1:B:220:ARG:HD3	1:B:221:TYR:H	1.36	0.89
1:B:106:HIS:CD2	1:B:108:ASN:H	1.93	0.86
1:A:106:HIS:HD2	1:A:108:ASN:H	1.22	0.84
1:B:370:ASN:HD22	1:B:372:PRO:HD2	1.42	0.81
1:A:370:ASN:HD22	1:A:372:PRO:HD2	1.45	0.80
1:A:106:HIS:CD2	1:A:108:ASN:H	1.99	0.80
1:B:106:HIS:HD2	1:B:108:ASN:H	1.31	0.77
1:B:289:THR:O	1:B:290:GLU:HB2	1.84	0.76
1:B:220:ARG:HG2	1:B:220:ARG:HH11	1.49	0.76
1:B:349:LYS:HE2	1:B:355:ASP:OD1	1.88	0.74
1:A:94:LYS:HG3	1:A:99:GLN:NE2	2.04	0.73
1:A:220:ARG:HD3	1:A:221:TYR:H	1.50	0.73
1:A:349:LYS:HE2	1:A:355:ASP:OD1	1.90	0.71
1:B:370:ASN:ND2	1:B:372:PRO:HD2	2.07	0.70
1:A:220:ARG:CD	1:A:221:TYR:N	2.43	0.67
1:B:300:PRO:HG2	1:B:303:LYS:HB2	1.76	0.66
1:B:91:LYS:NZ	1:B:126:VAL:HG11	2.11	0.65
1:A:370:ASN:ND2	1:A:372:PRO:HD2	2.12	0.64
1:A:212:PRO:HB2	1:B:289:THR:HG23	1.81	0.61
1:A:98:LEU:O	1:A:102:ARG:HG3	2.01	0.61
1:B:220:ARG:HD2	1:B:221:TYR:H	1.63	0.61
1:B:205:LYS:NZ	1:B:213:ASN:OD1	2.26	0.59
1:B:67:PHE:CE2	1:B:202:GLY:HA3	2.38	0.59
1:A:205:LYS:NZ	1:A:213:ASN:OD1	2.25	0.58
1:B:67:PHE:HE2	1:B:202:GLY:HA3	1.68	0.58
1:B:193:THR:O	1:B:357:PRO:HG3	2.04	0.57
1:B:185:GLN:HG3	3:B:618:HOH:O	2.04	0.57
1:B:220:ARG:CD	1:B:221:TYR:N	2.61	0.56
1:B:300:PRO:HG2	1:B:303:LYS:CB	2.36	0.55
1:B:308:ARG:HH11	1:B:308:ARG:HG2	1.72	0.54
1:A:94:LYS:CG	1:A:99:GLN:NE2	2.69	0.54
1:A:307:PRO:O	1:A:308:ARG:HB2	2.08	0.54
1:A:225:PRO:HB2	1:A:284:MET:SD	2.48	0.54
1:A:276:PRO:HG3	1:A:323:TYR:CZ	2.42	0.54
1:A:106:HIS:HD2	1:A:108:ASN:N	2.01	0.54
1:A:94:LYS:HE3	1:A:99:GLN:HE22	1.73	0.53
1:A:220:ARG:HD3	1:A:221:TYR:CG	2.44	0.53
1:B:225:PRO:HB2	1:B:284:MET:SD	2.49	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:ARG:HD2	1:B:221:TYR:N	2.24	0.53
1:B:220:ARG:O	1:B:223:ARG:HG3	2.09	0.51
1:B:220:ARG:HD3	1:B:221:TYR:CG	2.46	0.50
1:B:98:LEU:O	1:B:102:ARG:HG3	2.12	0.50
1:A:38:THR:O	1:A:55:SER:HA	2.12	0.49
1:A:271:LYS:O	1:A:299:HIS:HB2	2.13	0.49
1:B:307:PRO:HB2	1:B:308:ARG:NH1	2.28	0.48
1:A:185:GLN:HG3	3:A:617:HOH:O	2.13	0.48
1:B:91:LYS:HZ3	1:B:126:VAL:HG11	1.77	0.48
1:A:144:ARG:O	1:A:148:ARG:HG2	2.14	0.47
1:B:106:HIS:HD2	1:B:108:ASN:N	2.05	0.47
1:B:153:LEU:HD12	1:B:154:PRO:HD2	1.96	0.46
1:A:119:SER:HB3	1:A:126:VAL:HA	1.98	0.46
1:B:115:PHE:HA	1:B:129:ASN:O	2.16	0.45
1:B:299:HIS:NE2	1:B:303:LYS:HB3	2.32	0.45
1:A:252:LEU:HG	1:A:305:PHE:CD1	2.51	0.45
1:B:370:ASN:HD22	1:B:370:ASN:C	2.24	0.44
1:B:196:LEU:C	1:B:196:LEU:HD23	2.43	0.44
1:B:101:MET:HA	1:B:101:MET:HE2	1.99	0.43
1:B:161:TYR:CE1	1:B:189:LEU:HD22	2.52	0.43
1:A:94:LYS:HG3	1:A:99:GLN:HE21	1.78	0.43
1:A:196:LEU:C	1:A:196:LEU:HD23	2.43	0.43
1:B:308:ARG:HG2	1:B:308:ARG:NH1	2.31	0.43
1:A:66:SER:HB2	1:B:267:VAL:HG11	2.01	0.43
1:B:299:HIS:O	1:B:300:PRO:C	2.61	0.43
1:B:369:SER:O	1:B:370:ASN:HB2	2.18	0.43
2:A:451:IXM:N24	2:A:451:IXM:H22	2.34	0.42
1:A:252:LEU:HG	1:A:305:PHE:CE1	2.54	0.42
1:A:161:TYR:CE1	1:A:189:LEU:HD22	2.54	0.42
1:A:299:HIS:CE1	1:A:303:LYS:HB3	2.54	0.42
1:A:370:ASN:HD22	1:A:370:ASN:C	2.26	0.42
1:B:220:ARG:HG2	1:B:220:ARG:NH1	2.24	0.42
1:B:275:THR:HA	1:B:276:PRO:HD3	1.90	0.42
1:A:300:PRO:O	1:A:302:THR:N	2.53	0.41
1:A:215:SER:HB3	1:A:231:ALA:O	2.20	0.41
1:A:193:THR:O	1:A:357:PRO:HG3	2.20	0.41
1:A:60:LYS:HE2	1:A:72:GLN:NE2	2.35	0.41
1:A:167:ARG:HD3	3:A:512:HOH:O	2.20	0.41
1:A:275:THR:HA	1:A:276:PRO:HD3	1.89	0.41
1:A:272:VAL:HA	1:A:299:HIS:HB3	2.03	0.41
1:A:68:GLY:HA3	1:A:87:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:GLN:OE1	1:A:89:GLN:HA	2.20	0.41
1:B:276:PRO:HG3	1:B:323:TYR:CZ	2.56	0.41
1:B:72:GLN:HG2	1:B:73:ALA:N	2.36	0.41
1:A:310:PRO:HA	1:A:311:PRO:HD3	1.86	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	333/424 (78%)	317 (95%)	11 (3%)	5 (2%)	8	4
1	B	339/424 (80%)	321 (95%)	13 (4%)	5 (2%)	8	4
All	All	672/848 (79%)	638 (95%)	24 (4%)	10 (2%)	8	4

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	ARG
1	B	77	ASP
1	B	220	ARG
1	A	77	ASP
1	A	301	TRP
1	B	290	GLU
1	A	295	GLN
1	A	300	PRO
1	B	298	ALA
1	B	218	CYS

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	298/366 (81%)	294 (99%)	4 (1%)	61	69
1	B	303/366 (83%)	298 (98%)	5 (2%)	53	62
All	All	601/732 (82%)	592 (98%)	9 (2%)	57	65

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	77	ASP
1	A	217	ILE
1	A	220	ARG
1	B	77	ASP
1	B	99	GLN
1	B	101	MET
1	B	217	ILE
1	B	220	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	GLN
1	A	52	GLN
1	A	72	GLN
1	A	89	GLN
1	A	99	GLN
1	A	106	HIS
1	A	129	ASN
1	A	365	GLN
1	A	370	ASN
1	A	381	HIS
1	B	106	HIS
1	B	108	ASN
1	B	129	ASN

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Mol	Chain	Res	Type
1	B	365	GLN
1	B	370	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	IXM	B	452	-	24,24,24	1.80	8 (33%)	31,35,35	1.27	4 (12%)
2	IXM	A	451	-	24,24,24	1.64	6 (25%)	31,35,35	1.28	4 (12%)

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	IXM	B	452	-	-	0/6/30/30	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	IXM	A	451	-	-	0/6/30/30	0/4/4/4

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	452	IXM	C21-C22	3.16	1.44	1.38
2	B	452	IXM	C15-N16	2.97	1.39	1.36
2	A	451	IXM	C15-N16	2.96	1.39	1.36
2	A	451	IXM	C21-C22	2.94	1.43	1.38
2	A	451	IXM	C9-C11	2.88	1.43	1.38
2	B	452	IXM	C7-C5	2.81	1.43	1.38
2	B	452	IXM	C9-C11	2.72	1.43	1.38
2	B	452	IXM	C9-C7	2.47	1.43	1.38
2	A	451	IXM	C9-C7	2.38	1.43	1.38
2	B	452	IXM	C21-C20	2.37	1.43	1.38
2	A	451	IXM	C21-C20	2.34	1.43	1.38
2	B	452	IXM	C13-C2	2.21	1.54	1.46
2	B	452	IXM	C1-C2	2.16	1.50	1.45
2	A	451	IXM	C1-C2	2.10	1.50	1.45

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	451	IXM	C3-N4-C13	3.12	112.42	109.00
2	B	452	IXM	C3-N4-C13	3.10	112.40	109.00
2	B	452	IXM	C19-C17-N16	-2.60	125.57	130.83
2	A	451	IXM	C19-C17-N16	-2.60	125.58	130.83
2	B	452	IXM	C18-C17-N16	2.55	111.10	108.23
2	B	452	IXM	C1-C3-N4	2.49	111.04	108.23
2	A	451	IXM	C18-C17-N16	2.42	110.96	108.23
2	A	451	IXM	C1-C3-N4	2.41	110.95	108.23

There are no chirality outliers.

There are no torsion outliers.

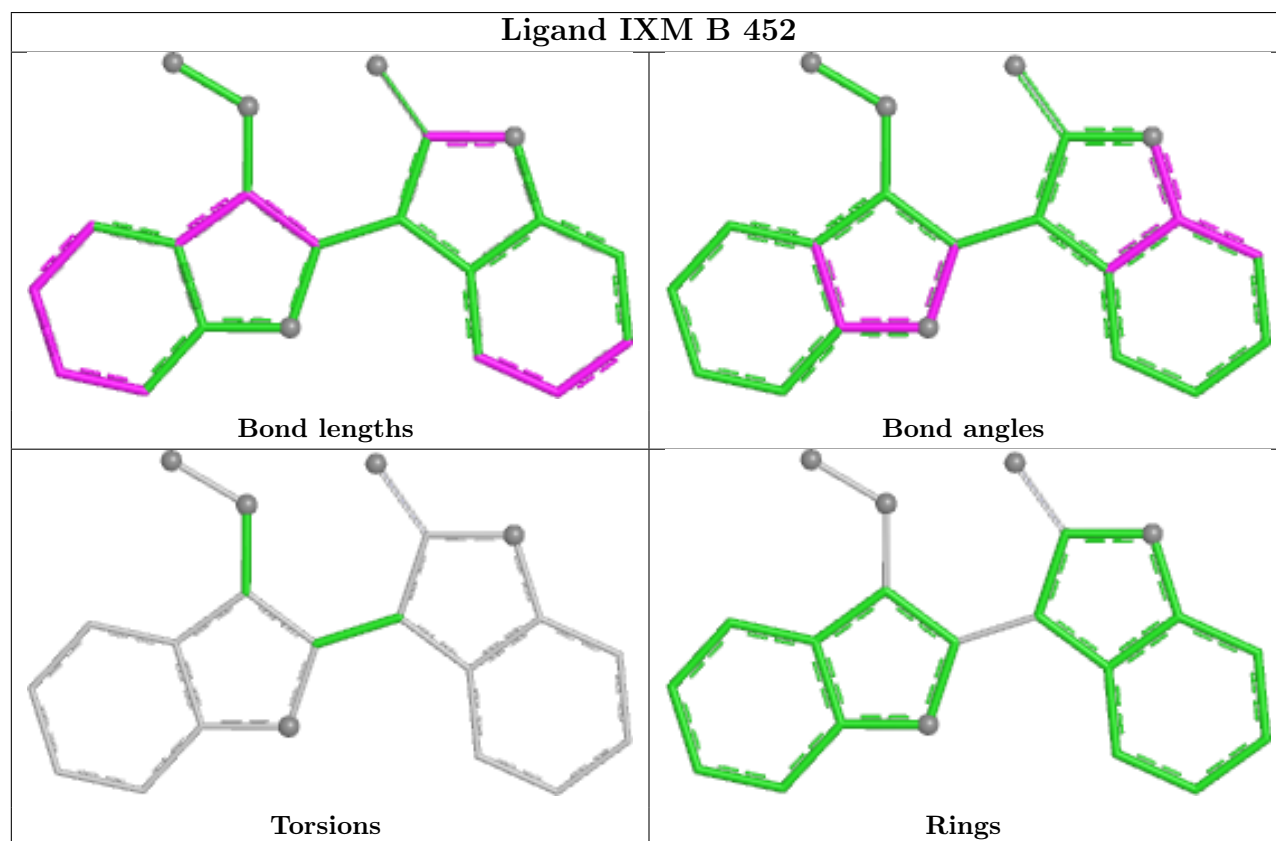
There are no ring outliers.

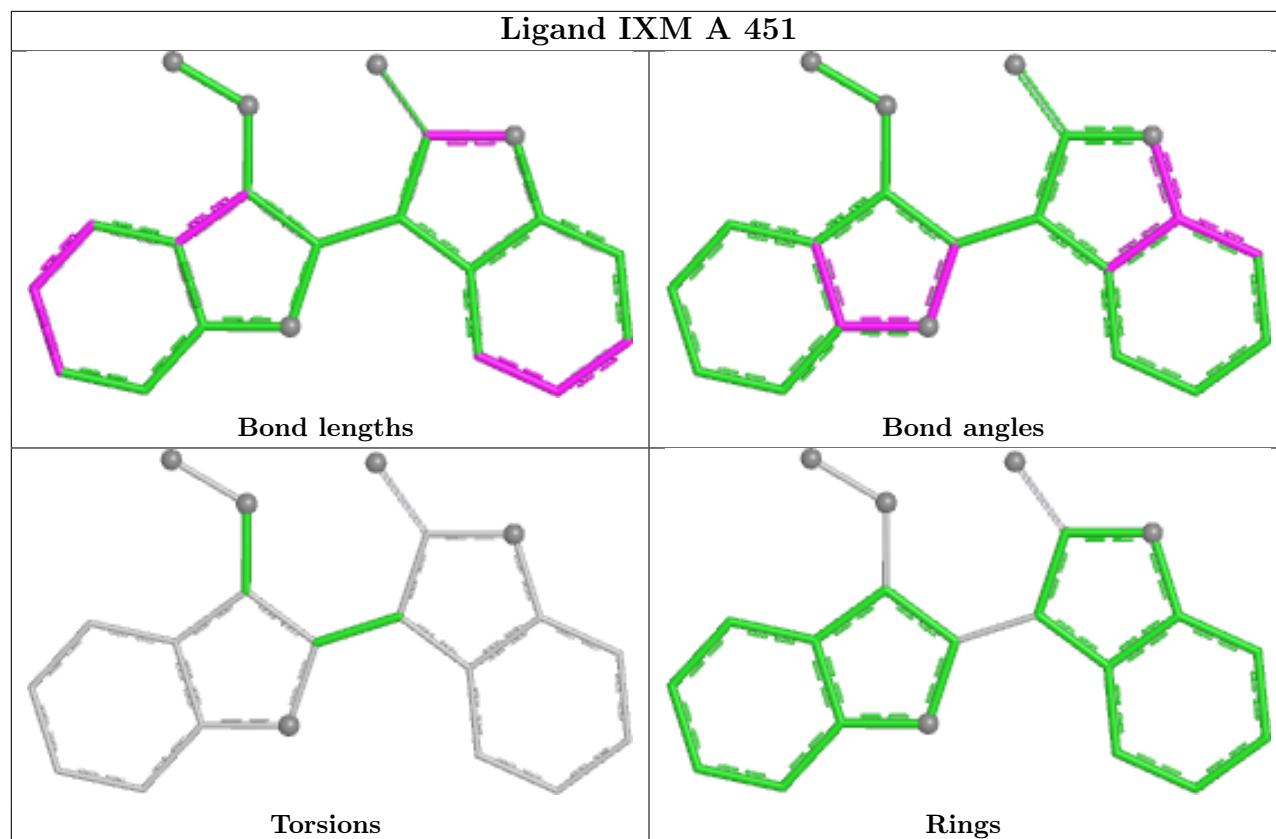
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	451	IXM	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	339/424 (79%)	0.90	48 (14%) 6 6	28, 45, 80, 110	0
1	B	343/424 (80%)	0.73	53 (15%) 5 5	28, 43, 80, 104	0
All	All	682/848 (80%)	0.82	101 (14%) 5 6	28, 44, 80, 110	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	ILE	9.0
1	A	298	ALA	7.5
1	A	67	PHE	7.3
1	B	288	TYR	6.5
1	A	93	PHE	6.3
1	B	67	PHE	6.2
1	A	300	PRO	6.1
1	B	291	PHE	6.1
1	A	297	LYS	6.0
1	A	294	PRO	5.4
1	A	292	LYS	5.2
1	B	294	PRO	5.1
1	B	296	ILE	5.1
1	A	386	ALA	4.8
1	B	293	PHE	4.6
1	A	66	SER	4.5
1	A	302	THR	4.5
1	A	68	GLY	4.4
1	A	293	PHE	4.4
1	B	278	ARG	4.2
1	B	287	ASN	3.8
1	A	299	HIS	3.7
1	B	383	ARG	3.7
1	B	65	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	308	ARG	3.7
1	A	91	LYS	3.7
1	B	35	SER	3.6
1	B	277	THR	3.6
1	B	220	ARG	3.5
1	B	382	ALA	3.5
1	B	37	VAL	3.4
1	B	290	GLU	3.4
1	B	149	ALA	3.4
1	B	302	THR	3.3
1	A	77	ASP	3.3
1	B	286	PRO	3.3
1	B	298	ALA	3.3
1	B	93	PHE	3.3
1	A	220	ARG	3.2
1	B	92	ARG	3.2
1	B	90	ASP	3.1
1	A	353	GLY	3.1
1	B	91	LYS	3.0
1	A	295	GLN	3.0
1	B	295	GLN	3.0
1	B	66	SER	2.9
1	A	301	TRP	2.9
1	B	282	ARG	2.9
1	A	92	ARG	2.8
1	B	77	ASP	2.8
1	A	145	HIS	2.7
1	B	292	LYS	2.7
1	A	35	SER	2.7
1	B	210	GLY	2.7
1	A	52	GLN	2.6
1	B	308	ARG	2.6
1	B	299	HIS	2.6
1	B	71	TYR	2.6
1	B	61	VAL	2.6
1	A	148	ARG	2.6
1	A	76	CYS	2.6
1	A	149	ALA	2.6
1	B	209	ARG	2.6
1	A	63	GLY	2.5
1	B	285	ASN	2.5
1	A	59	THR	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	279	GLU	2.5
1	B	88	LEU	2.5
1	A	151	GLN	2.4
1	B	289	THR	2.4
1	A	119	SER	2.4
1	A	127	TYR	2.4
1	A	252	LEU	2.4
1	A	69	VAL	2.4
1	A	61	VAL	2.3
1	B	306	ARG	2.3
1	A	99	GLN	2.3
1	B	148	ARG	2.3
1	B	297	LYS	2.3
1	A	50	ARG	2.2
1	A	371	PRO	2.2
1	B	36	LYS	2.2
1	A	88	LEU	2.2
1	A	281	ILE	2.2
1	B	48	PRO	2.2
1	B	126	VAL	2.1
1	B	49	ASP	2.1
1	B	301	TRP	2.1
1	A	363	THR	2.1
1	B	276	PRO	2.1
1	B	145	HIS	2.1
1	B	68	GLY	2.1
1	A	87	VAL	2.1
1	B	147	SER	2.1
1	A	48	PRO	2.1
1	A	37	VAL	2.1
1	A	126	VAL	2.1
1	B	57	THR	2.0
1	B	59	THR	2.0
1	A	283	GLU	2.0
1	B	358	ALA	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 [i](#)

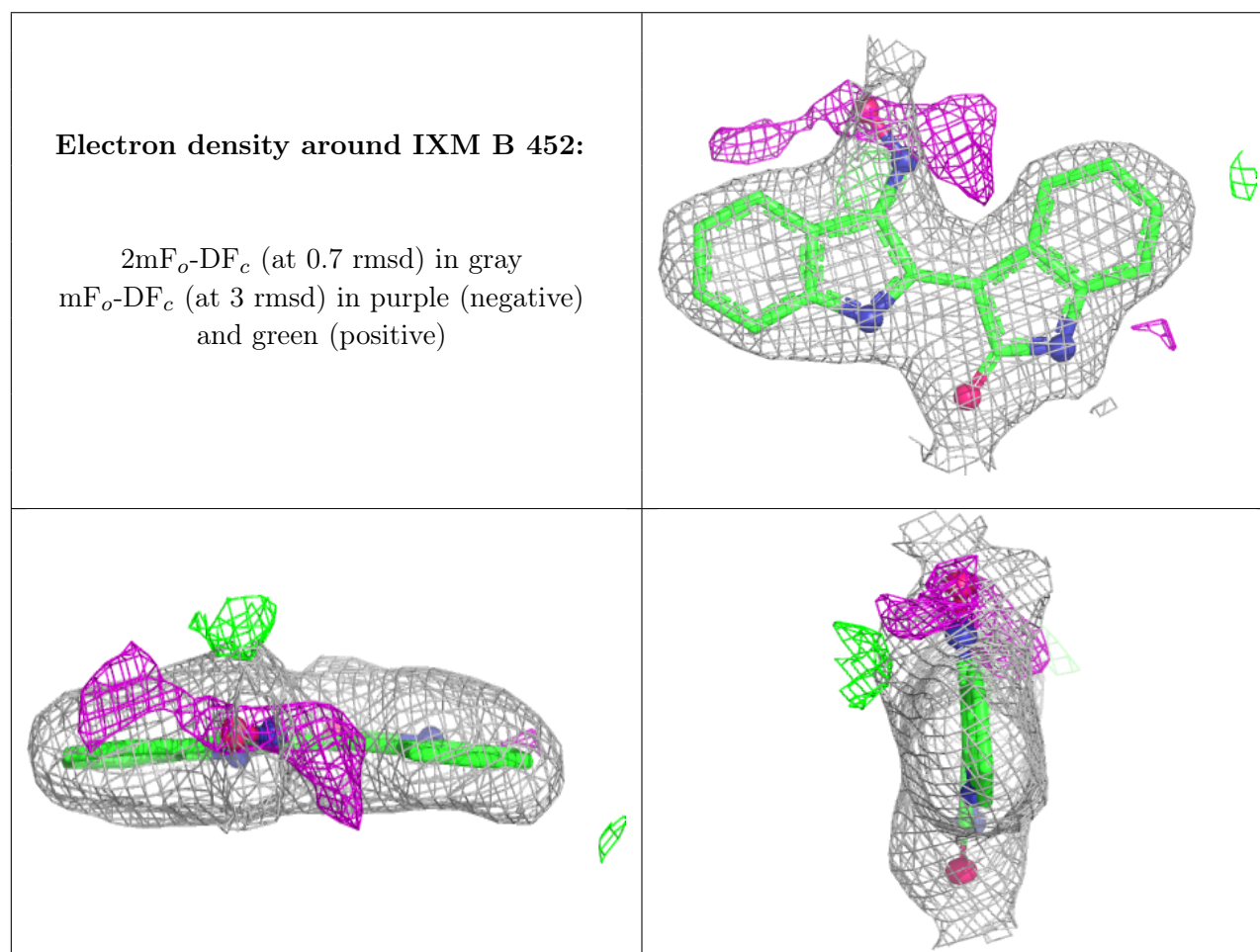
There are no oligosaccharides in this entry.

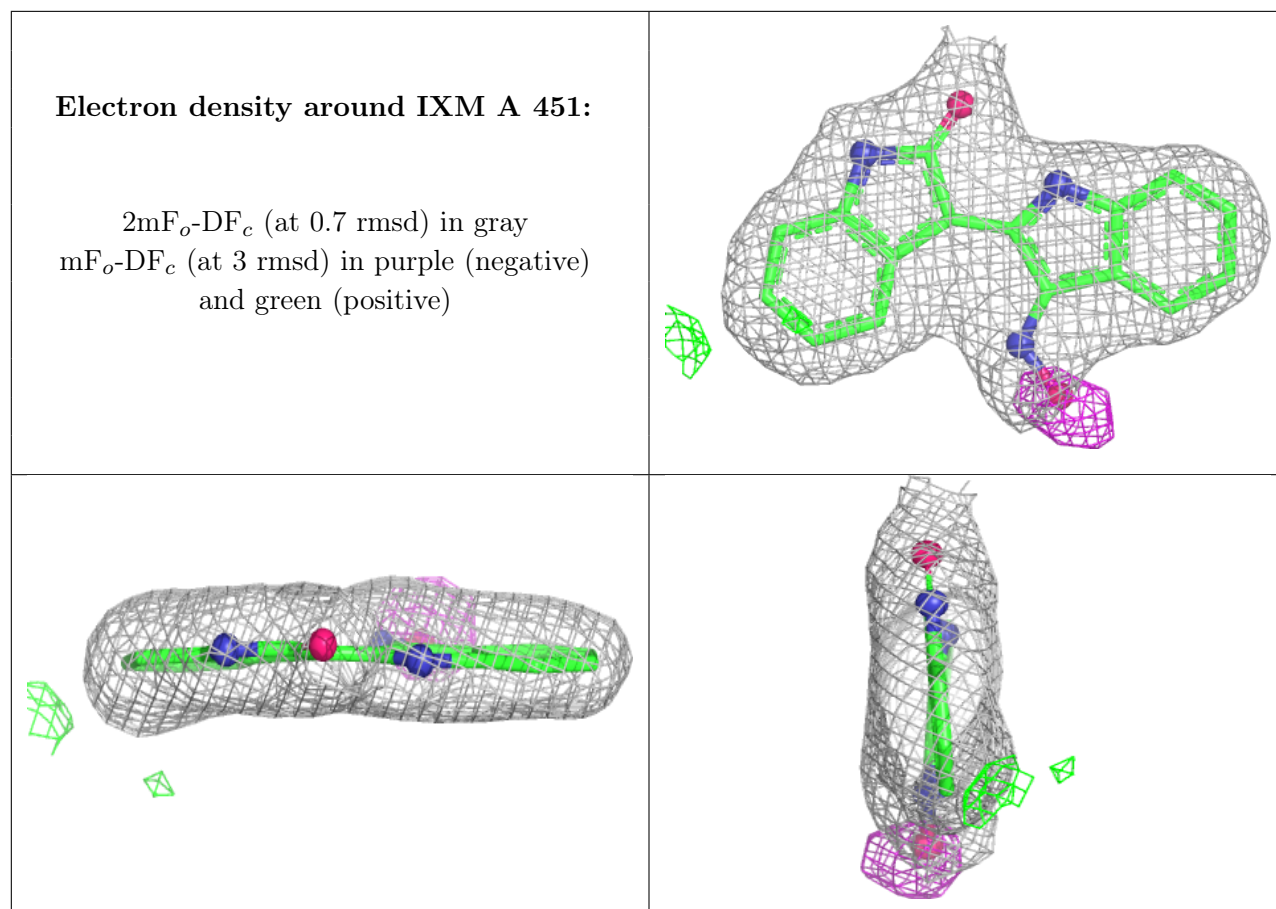
6.4 Ligands [i](#)

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	IXM	B	452	21/21	0.91	0.11	44,50,55,63	0
2	IXM	A	451	21/21	0.94	0.09	43,46,48,58	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.





6.5 Other polymers ⓘ

There are no such residues in this entry.