



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 12, 2024 – 02:10 PM EDT

PDB ID : 3F66  
Title : Human c-Met Kinase in complex with quinoxaline inhibitor  
Authors : Meier, C.; Ceska, T.  
Deposited on : 2008-11-05  
Resolution : 1.40 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

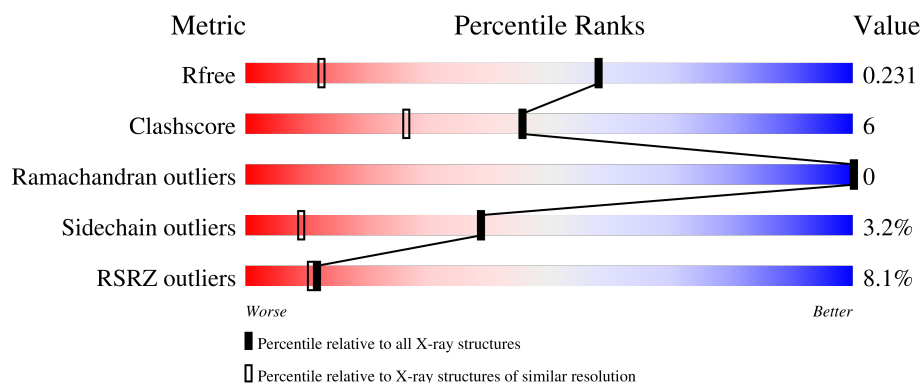
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1714 (1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

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  $\geq 3$ , 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	298	<div> <div>8%</div> <div> <div></div> <div>81%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	298	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>13%</div> <div>• 10%</div> </div> </div>

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5160 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.

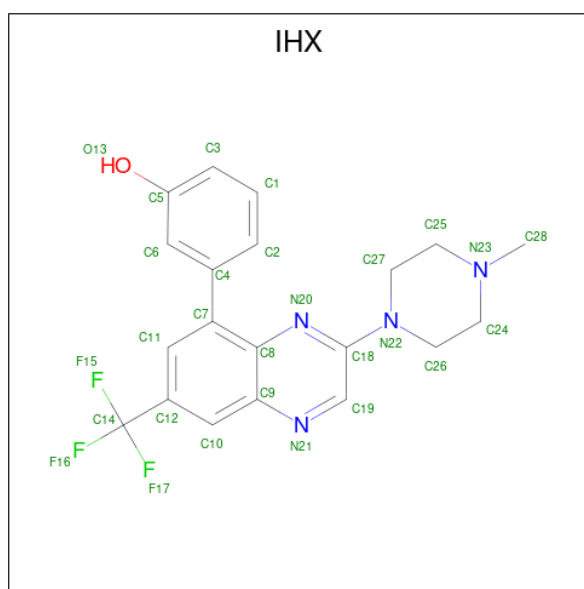
- Molecule 1 is a protein called Hepatocyte growth factor receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	8	7	0
			2352	1519	398	419	16			
1	B	269	Total	C	N	O	S	3	10	0
			2268	1465	393	394	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1272	LEU	VAL	SEE REMARK 999	UNP P08581
B	1272	LEU	VAL	SEE REMARK 999	UNP P08581

- Molecule 2 is 3-[3-(4-methylpiperazin-1-yl)-7-(trifluoromethyl)quinoxalin-5-yl]phenol (three-letter code: IHX) (formula: C<sub>20</sub>H<sub>19</sub>F<sub>3</sub>N<sub>4</sub>O).



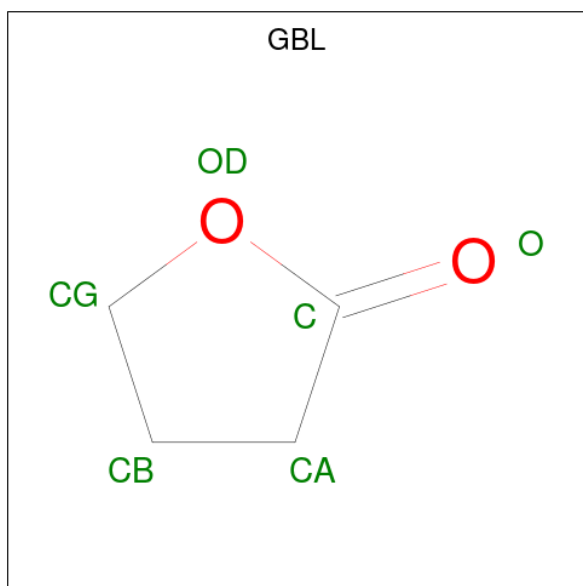
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	F	N	O	0	0
			28	20	3	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	F	N	O	0	0
			28	20	3	4	1		

- Molecule 3 is GAMMA-BUTYROLACTONE (three-letter code: GBL) (formula:  $C_4H_6O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	4	2		
3	A	1	Total	C	O	0	0
			6	4	2		
3	A	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		
3	B	1	Total	C	O	0	0
			6	4	2		

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Na	0	0
			1	1		

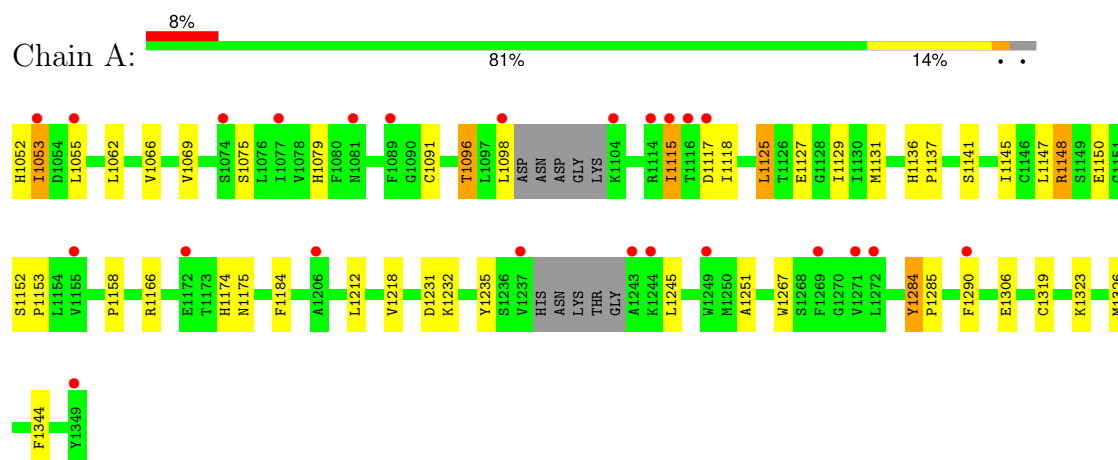
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	222	Total 222	O 222	0	0
5	B	231	Total 231	O 231	0	0

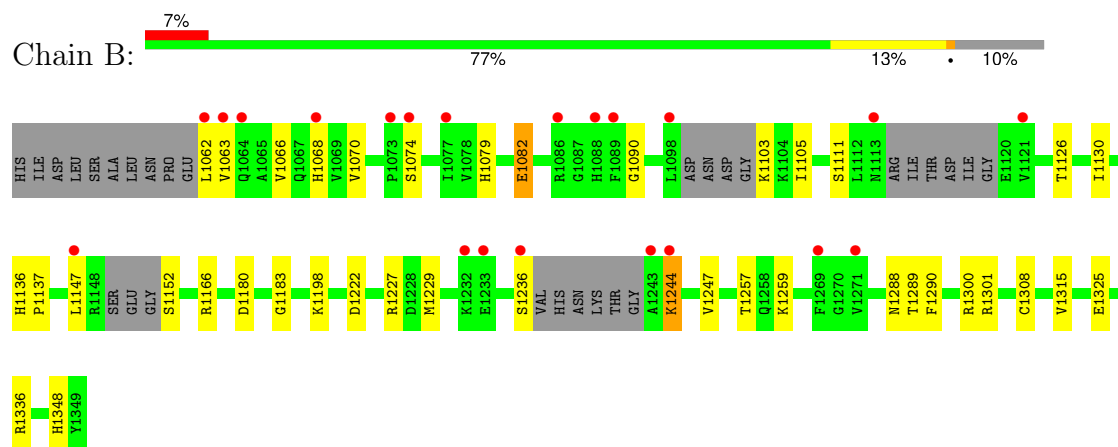
### 3 Residue-property plots

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: Hepatocyte growth factor receptor



- Molecule 1: Hepatocyte growth factor receptor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.38Å 47.59Å 75.60Å 100.70° 103.62° 98.29°	Depositor
Resolution (Å)	20.00 – 1.40 24.56 – 1.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.40) 88.6 (24.56-1.40)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.02 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.168 , 0.225 0.177 , 0.231	Depositor DCC
$R_{free}$ test set	3235 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	14.1	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 58.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5160	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.84% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GBL, NA, IHX

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	0.86	3/2410 (0.1%)	0.92	4/3263 (0.1%)
1	B	0.82	1/2323 (0.0%)	0.85	5/3134 (0.2%)
All	All	0.84	4/4733 (0.1%)	0.89	9/6397 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1284	TYR	CD2-CE2	-8.71	1.26	1.39
1	B	1308	CYS	CB-SG	-5.65	1.72	1.81
1	A	1184	PHE	CE1-CZ	5.11	1.47	1.37
1	A	1319	CYS	CB-SG	-5.09	1.73	1.81

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1284	TYR	CB-CG-CD2	-6.99	116.80	121.00
1	B	1300	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	A	1184	PHE	CB-CG-CD1	-6.87	115.99	120.80
1	B	1166	ARG	NE-CZ-NH1	6.54	123.57	120.30
1	A	1184	PHE	CB-CG-CD2	6.46	125.32	120.80
1	B	1166	ARG	NE-CZ-NH2	-6.13	117.24	120.30
1	A	1166	ARG	NE-CZ-NH2	-5.41	117.59	120.30
1	B	1227	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	B	1301	ARG	NE-CZ-NH1	5.04	122.82	120.30

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	2352	0	2357	34	0
1	B	2268	0	2281	28	0
2	A	28	0	18	0	0
2	B	28	0	18	0	0
3	A	18	0	18	0	0
3	B	12	0	12	2	0
4	B	1	0	0	0	0
5	A	222	0	0	8	0
5	B	231	0	0	4	0
All	All	5160	0	4704	60	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 6.

All (60) 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:1284:TYR:OH	5:A:421:HOH:O	1.81	0.97
1:A:1066:VAL:HG22	1:A:1129:ILE:HD11	1.48	0.94
1:B:1247:VAL:HG21	1:B:1289:THR:HG23	1.53	0.90
1:A:1052:HIS:N	5:A:367:HOH:O	2.04	0.90
1:B:1236:SER:OG	1:B:1289:THR:HG21	1.75	0.87
1:A:1175:ASN:HB2	5:A:386:HOH:O	1.75	0.86
1:A:1174:HIS:ND1	5:A:349:HOH:O	2.08	0.86
1:A:1091:CYS:SG	5:A:408:HOH:O	2.35	0.83
1:B:1288[B]:ASN:HD21	1:B:1290:PHE:HB2	1.50	0.76
1:B:1062:LEU:O	1:B:1066:VAL:HG23	1.91	0.71
1:A:1231[B]:ASP:OD1	5:A:286:HOH:O	2.11	0.68
1:B:1198:LYS:O	5:B:426:HOH:O	2.13	0.65
1:A:1066:VAL:CG2	1:A:1129:ILE:HD11	2.26	0.64
1:B:1180:ASP:OD2	1:B:1348:HIS:HD2	1.82	0.62
1:A:1053:ILE:HD11	1:A:1118:ILE:HA	1.85	0.58
1:B:1257:THR:HB	1:B:1259:LYS:HE2	1.83	0.58
1:A:1285:PRO:O	1:B:1348:HIS:HE1	1.86	0.58
1:A:1212[A]:LEU:HD21	1:A:1218:VAL:HG22	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1247:VAL:CG2	1:B:1289:THR:HG23	2.33	0.57
1:B:1183:GLY:HA3	3:B:1350:GBL:HAC1	1.88	0.55
1:A:1125:LEU:HD11	1:A:1147:LEU:HD11	1.89	0.54
1:A:1069:VAL:HG21	1:A:1129:ILE:HA	1.88	0.54
1:B:1257:THR:HB	1:B:1259:LYS:CE	2.38	0.54
1:B:1079:HIS:HB2	1:B:1082:GLU:HG3	1.91	0.53
1:B:1222:ASP:HB2	3:B:1:GBL:HAC1	1.91	0.52
1:B:1229:MET:HE1	5:B:375:HOH:O	2.10	0.52
1:A:1212[A]:LEU:CD2	1:A:1218:VAL:HG22	2.39	0.51
1:B:1126:THR:HG22	1:B:1130:ILE:HD12	1.92	0.50
1:B:1068:HIS:NE2	5:B:393:HOH:O	2.35	0.50
1:B:1315:VAL:CG2	1:B:1336[D]:ARG:HG2	2.41	0.50
1:A:1136:HIS:CG	1:A:1137:PRO:HD2	2.47	0.49
1:A:1306:GLU:HG2	5:A:361:HOH:O	2.12	0.49
1:B:1325:GLU:H	1:B:1325:GLU:CD	2.15	0.49
1:A:1066:VAL:HG22	1:A:1129:ILE:CD1	2.32	0.48
1:B:1136:HIS:CG	1:B:1137:PRO:HD2	2.50	0.47
1:A:1127:GLU:O	1:A:1131:MET:HG2	2.14	0.47
1:A:1075:SER:HA	1:A:1098:LEU:HD12	1.95	0.47
1:A:1062:LEU:O	1:A:1066:VAL:HG23	2.15	0.46
1:B:1236:SER:HG	1:B:1289:THR:HG21	1.81	0.45
1:B:1066:VAL:HG12	1:B:1070:VAL:HG23	1.99	0.45
1:A:1079:HIS:NE2	1:A:1096:THR:HG23	2.31	0.45
1:B:1066:VAL:HG12	1:B:1066:VAL:O	2.16	0.45
1:A:1344:PHE:CE2	1:B:1103:LYS:HE3	2.52	0.44
1:B:1147:LEU:HD12	1:B:1147:LEU:N	2.32	0.44
1:A:1066:VAL:HG21	1:A:1125:LEU:HB3	1.99	0.44
1:A:1323:LYS:HE2	1:A:1326:MET:HG3	1.99	0.44
1:B:1244:LYS:HG3	1:B:1244:LYS:O	2.18	0.43
1:A:1148:ARG:HG3	1:A:1150:GLU:O	2.20	0.42
1:B:1180:ASP:OD2	1:B:1348:HIS:CD2	2.67	0.42
1:B:1090:GLY:HA3	1:B:1111:SER:O	2.20	0.42
1:A:1251:ALA:HA	1:A:1267:TRP:CD2	2.55	0.42
1:A:1066:VAL:HG11	1:A:1125:LEU:HG	2.01	0.41
1:A:1125:LEU:O	1:A:1129:ILE:HG13	2.19	0.41
1:A:1115:ILE:HG21	1:A:1153:PRO:HG2	2.03	0.41
1:A:1175:ASN:CB	5:A:386:HOH:O	2.50	0.41
1:B:1336[D]:ARG:CD	5:B:179:HOH:O	2.69	0.41
1:A:1069:VAL:CG2	1:A:1129:ILE:HA	2.51	0.40
1:A:1141[A]:SER:O	1:A:1158:PRO:HD3	2.20	0.40
1:A:1125:LEU:HD12	1:A:1145:ILE:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1235:TYR:HA	1:A:1245:LEU:O	2.21	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	289/298 (97%)	282 (98%)	7 (2%)	0	100	100
1	B	271/298 (91%)	266 (98%)	5 (2%)	0	100	100
All	All	560/596 (94%)	548 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	263/264 (100%)	253 (96%)	10 (4%)	33	6
1	B	252/264 (96%)	245 (97%)	7 (3%)	43	11
All	All	515/528 (98%)	498 (97%)	17 (3%)	39	9

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1053	ILE
1	A	1055	LEU
1	A	1096	THR
1	A	1115	ILE
1	A	1117	ASP
1	A	1125	LEU
1	A	1148	ARG
1	A	1152	SER
1	A	1232	LYS
1	A	1290	PHE
1	B	1063	VAL
1	B	1074[A]	SER
1	B	1074[B]	SER
1	B	1082	GLU
1	B	1105	ILE
1	B	1152	SER
1	B	1244	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1068	HIS
1	A	1167	ASN
1	A	1256	GLN
1	A	1304	GLN
1	B	1167	ASN
1	B	1304	GLN
1	B	1348	HIS

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	GBL	A	3	-	6,6,6	0.70	0	7,7,7	0.55	0
3	GBL	A	5	-	6,6,6	1.33	1 (16%)	7,7,7	1.76	2 (28%)
2	IHX	A	1	-	30,31,31	1.43	4 (13%)	40,46,46	1.44	8 (20%)
3	GBL	A	4	-	6,6,6	0.48	0	7,7,7	0.95	1 (14%)
3	GBL	B	1	-	6,6,6	0.79	0	7,7,7	2.14	3 (42%)
2	IHX	B	2	-	30,31,31	1.44	5 (16%)	40,46,46	1.50	7 (17%)
3	GBL	B	1350	-	6,6,6	1.31	1 (16%)	7,7,7	2.76	4 (57%)

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	GBL	A	3	-	-	-	0/1/1/1
3	GBL	A	5	-	-	-	0/1/1/1
2	IHX	A	1	-	-	0/10/24/24	0/4/4/4
3	GBL	A	4	-	-	-	0/1/1/1
3	GBL	B	1	-	-	-	0/1/1/1
2	IHX	B	2	-	-	0/10/24/24	0/4/4/4
3	GBL	B	1350	-	-	-	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	IHX	C10-C9	-3.81	1.36	1.41
2	B	2	IHX	C18-N20	3.77	1.37	1.32
2	A	1	IHX	C18-N20	3.60	1.37	1.32
2	A	1	IHX	C7-C8	-3.15	1.39	1.45
2	A	1	IHX	C10-C9	-2.70	1.37	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	5	GBL	CA-C	-2.62	1.45	1.50
2	A	1	IHX	C19-N21	2.60	1.36	1.31
2	B	2	IHX	C19-N21	2.51	1.36	1.31
2	B	2	IHX	C7-C8	-2.37	1.40	1.45
3	B	1350	GBL	CA-C	-2.19	1.45	1.50
2	B	2	IHX	C26-N22	2.01	1.50	1.46

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1350	GBL	OD-C-O	5.06	133.64	120.43
3	B	1	GBL	OD-C-O	4.26	131.56	120.43
2	B	2	IHX	C10-C12-C11	3.87	121.26	117.75
2	A	1	IHX	C10-C12-C11	3.69	121.10	117.75
2	B	2	IHX	C28-N23-C25	3.33	116.97	110.63
3	B	1350	GBL	O-C-CA	-3.32	116.63	127.83
3	A	5	GBL	OD-C-O	3.15	128.66	120.43
2	B	2	IHX	C26-C24-N23	-3.14	105.81	110.86
2	A	1	IHX	F15-C14-C12	-3.09	106.28	112.90
3	B	1	GBL	CB-CA-C	2.95	109.87	104.50
3	A	5	GBL	CB-CA-C	2.87	109.72	104.50
3	B	1350	GBL	CB-CA-C	2.85	109.69	104.50
2	B	2	IHX	C26-N22-C27	2.83	117.94	111.57
3	B	1350	GBL	CA-CB-CG	-2.75	96.45	103.91
2	B	2	IHX	C11-C7-C8	-2.68	116.73	118.88
2	A	1	IHX	C26-N22-C27	2.56	117.33	111.57
2	A	1	IHX	C6-C4-C2	2.55	121.32	118.15
2	B	2	IHX	C24-C26-N22	2.44	115.92	110.78
2	A	1	IHX	C5-C6-C4	-2.34	118.00	120.40
2	A	1	IHX	C25-C27-N22	2.20	115.41	110.78
3	A	4	GBL	OD-C-O	2.15	126.05	120.43
3	B	1	GBL	O-C-CA	-2.13	120.67	127.83
2	A	1	IHX	C12-C11-C7	-2.10	120.30	122.48
2	A	1	IHX	C7-C8-N20	2.09	120.80	118.14
2	B	2	IHX	N20-C18-N22	-2.08	115.67	117.49

There are no chirality outliers.

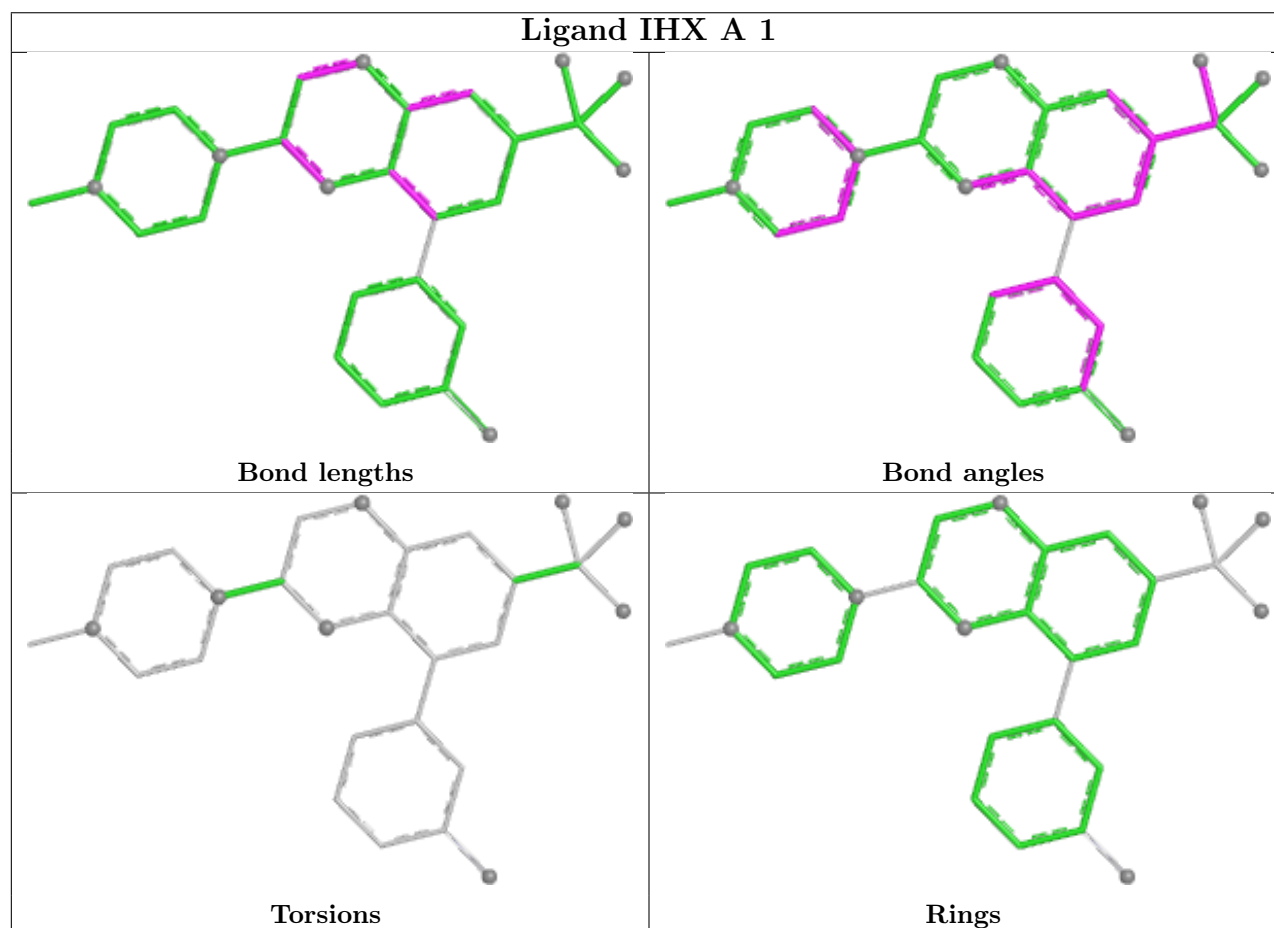
There are no torsion outliers.

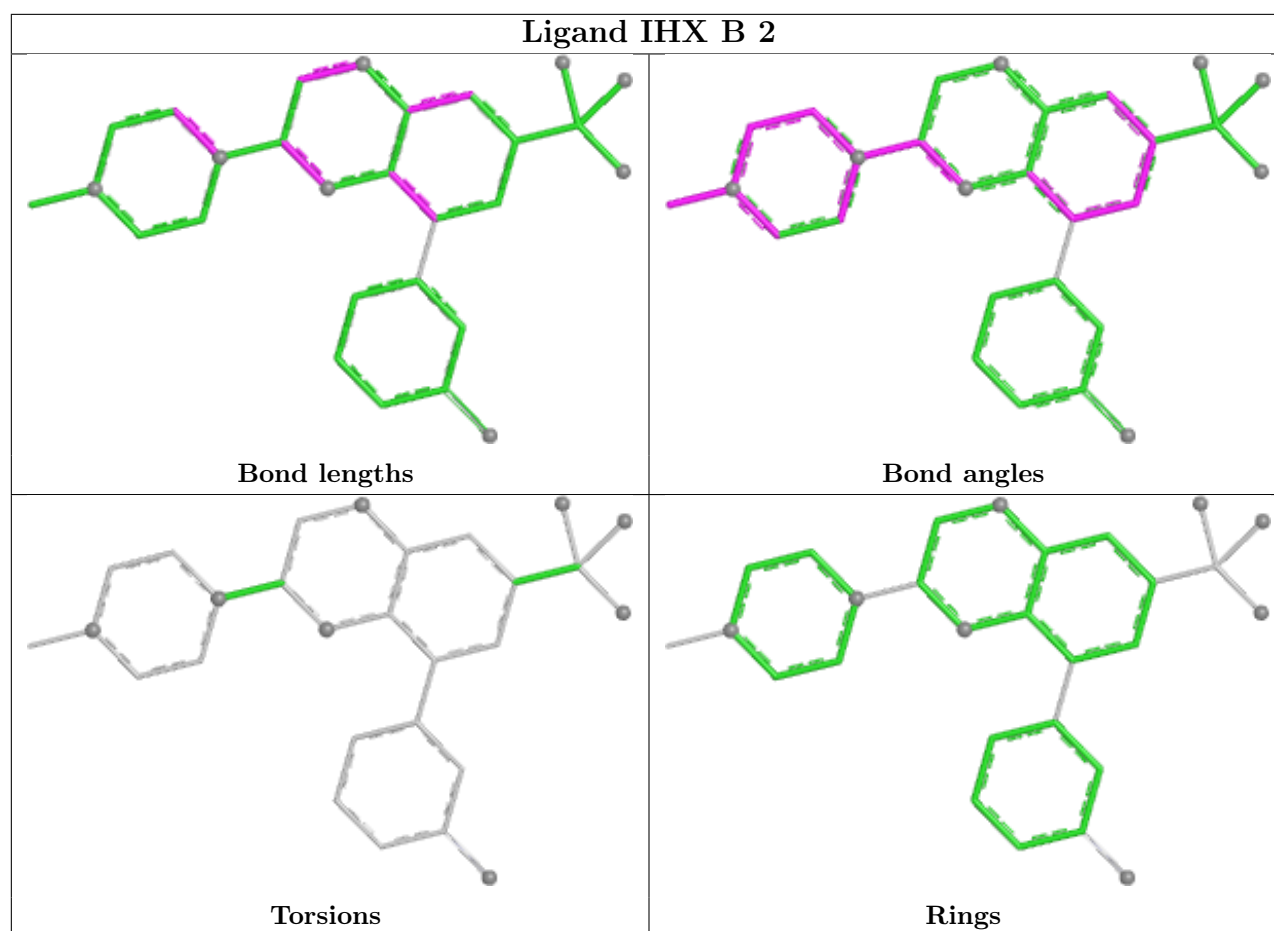
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1	GBL	1	0
3	B	1350	GBL	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q < 0.9
1	A	288/298 (96%)	0.23	24 (8%)	11 10	11, 24, 50, 63	6 (2%)
1	B	269/298 (90%)	0.21	21 (7%)	13 12	12, 23, 47, 58	4 (1%)
All	All	557/596 (93%)	0.22	45 (8%)	12 11	11, 24, 48, 63	10 (1%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1290	PHE	6.6
1	A	1098	LEU	5.8
1	A	1237	VAL	5.7
1	B	1089	PHE	5.5
1	B	1243	ALA	5.2
1	A	1114	ARG	5.1
1	B	1147	LEU	4.4
1	B	1236	SER	4.1
1	A	1053	ILE	3.9
1	B	1113	ASN	3.9
1	B	1088[A]	HIS	3.8
1	A	1243	ALA	3.6
1	A	1172	GLU	3.4
1	B	1074[A]	SER	3.3
1	B	1098	LEU	3.2
1	A	1272	LEU	3.1
1	B	1121	VAL	3.1
1	B	1232	LYS	3.0
1	B	1244	LYS	3.0
1	A	1244	LYS	2.8
1	B	1062	LEU	2.7
1	A	1116	THR	2.5
1	A	1271	VAL	2.5
1	B	1063	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	1064	GLN	2.4
1	A	1115	ILE	2.4
1	B	1233	GLU	2.4
1	A	1089	PHE	2.3
1	A	1104	LYS	2.3
1	B	1073	PRO	2.3
1	A	1055	LEU	2.3
1	B	1269	PHE	2.2
1	A	1081[A]	ASN	2.2
1	A	1349	TYR	2.2
1	A	1155	VAL	2.2
1	A	1206	ALA	2.2
1	B	1271	VAL	2.2
1	A	1249	TRP	2.2
1	A	1077	ILE	2.1
1	A	1074[A]	SER	2.1
1	B	1077	ILE	2.1
1	B	1086	ARG	2.1
1	A	1269	PHE	2.1
1	A	1117	ASP	2.0
1	B	1068	HIS	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> 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.

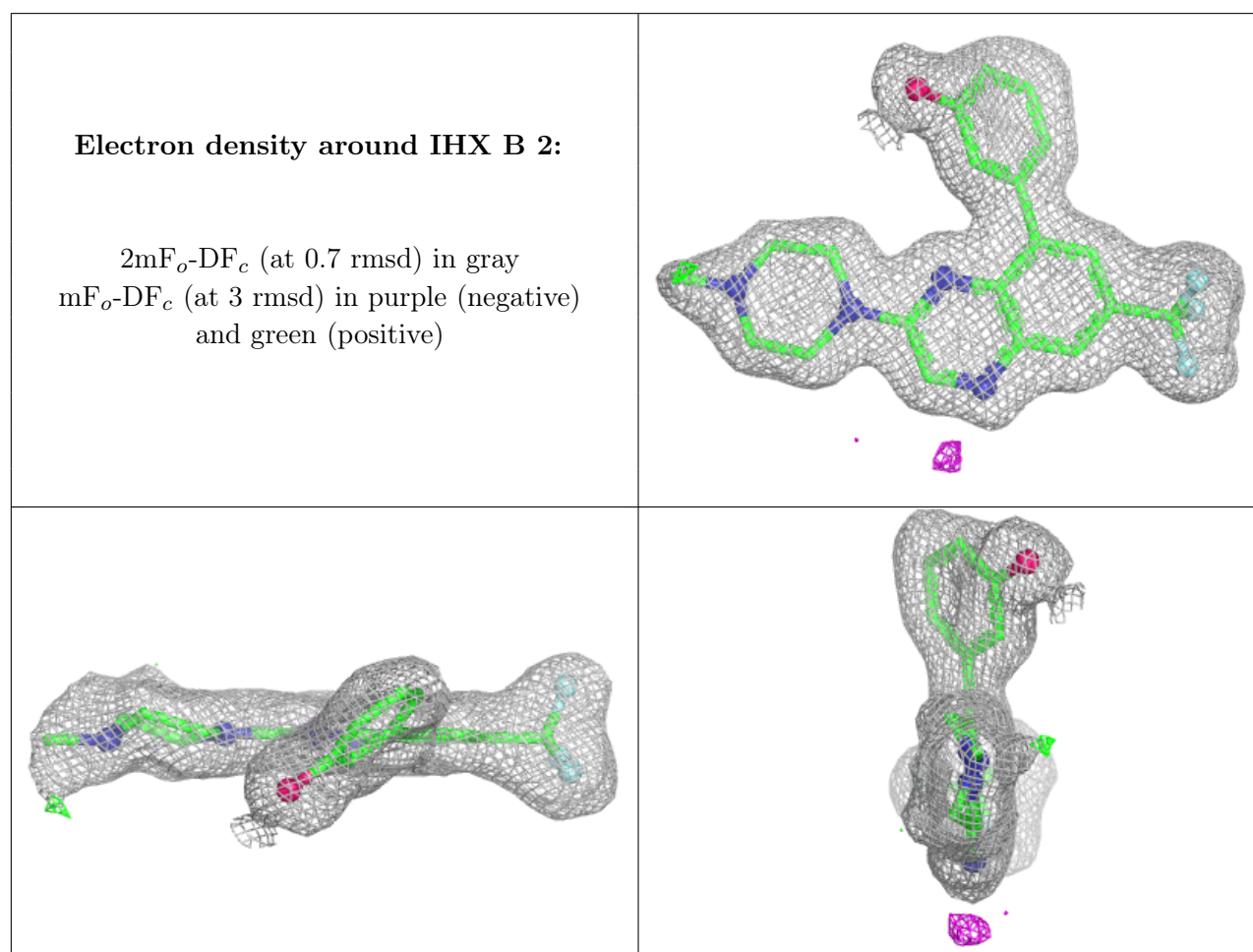
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	GBL	A	5	6/6	0.95	0.10	18,27,32,35	0
2	IHX	B	2	28/28	0.96	0.08	18,26,41,46	0

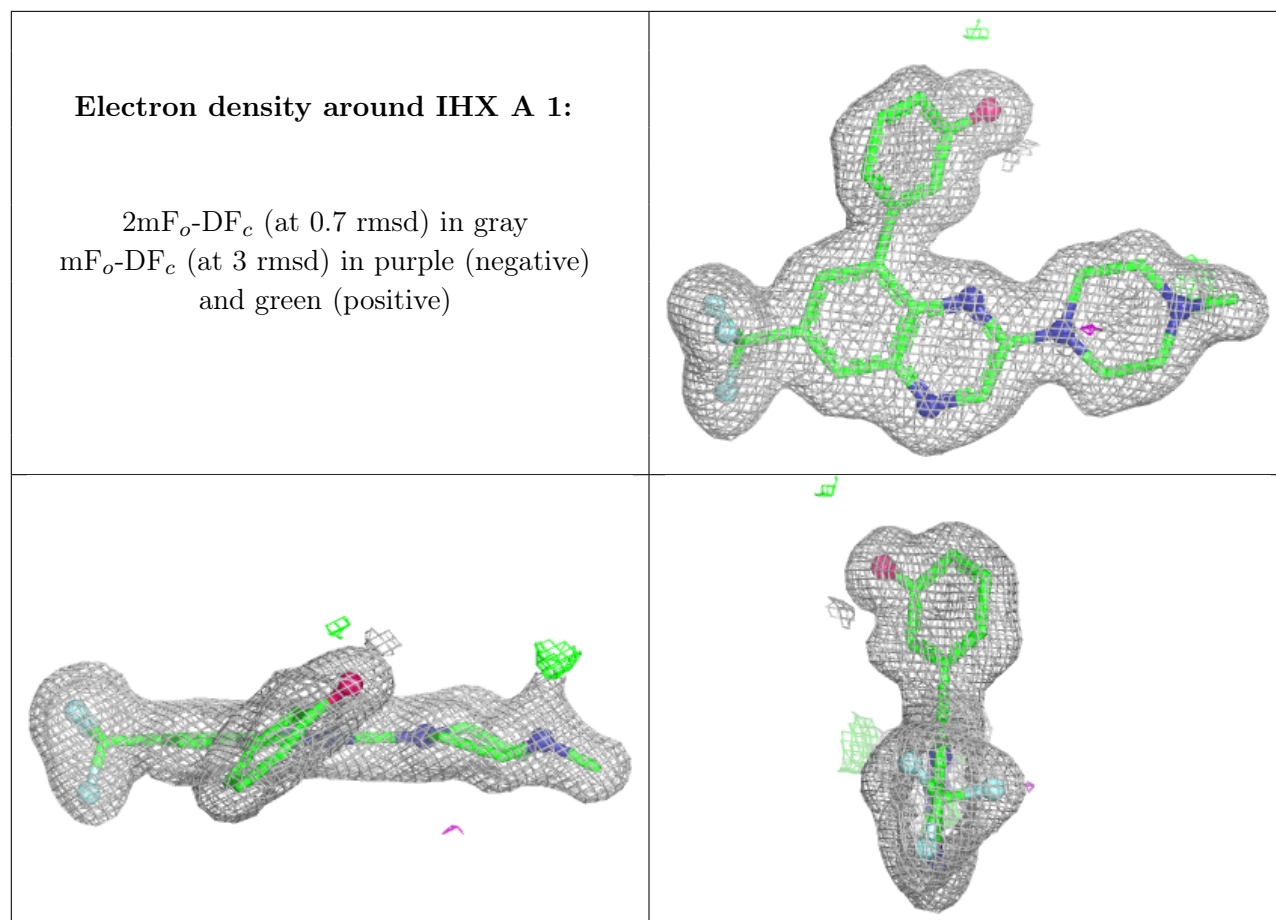
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	IHX	A	1	28/28	0.96	0.07	18,26,34,42	0
3	GBL	B	1	6/6	0.97	0.13	19,21,21,29	0
3	GBL	A	4	6/6	0.98	0.11	29,31,35,37	0
3	GBL	B	1350	6/6	0.98	0.05	12,19,24,30	0
3	GBL	A	3	6/6	0.99	0.09	20,22,22,25	0
4	NA	B	6	1/1	0.99	0.09	34,34,34,34	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 [i](#)

There are no such residues in this entry.