



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:25 PM UTC

PDB ID : 3TTI / pdb_00003tti
Title : Crystal Structure of JNK3 complexed with CC-930, an orally active anti-fibrotic JNK inhibitor
Authors : Plantevin-Krenitsky, V.; Nadolny, L.; Delgado, M.; Ayala, L.; Clareen, S.; Hilgraf, R.; Albers, R.; Hegde, S.; D'Sidocky, N.; Sapienza, J.; Wright, J.; McCarrick, M.; Bahmanyar, S.; Chamberlain, P.; Delker, S.L.; Muir, J.; Giegel, D.; Xu, L.; Celeridad, M.; Lachowitz, J.; Bennett, B.; Moghaddam, M.; Khatsenko, O.; Katz, J.; Fan, R.; Bai, A.; Tang, Y.; Shirley, M.A.; Benish, B.; Bodine, T.; Blease, K.; Raymon, H.; Cathers, B.E.; Satoh, Y.
Deposited on : 2011-09-14
Resolution : 2.20 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)

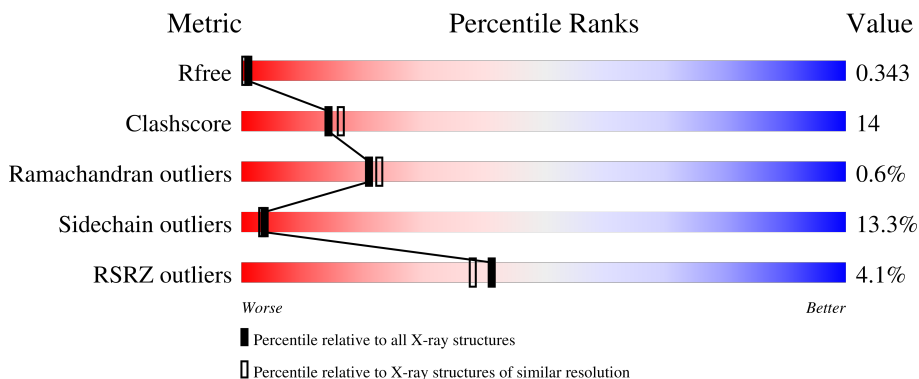
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	464	

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

2 Entry composition [i](#)

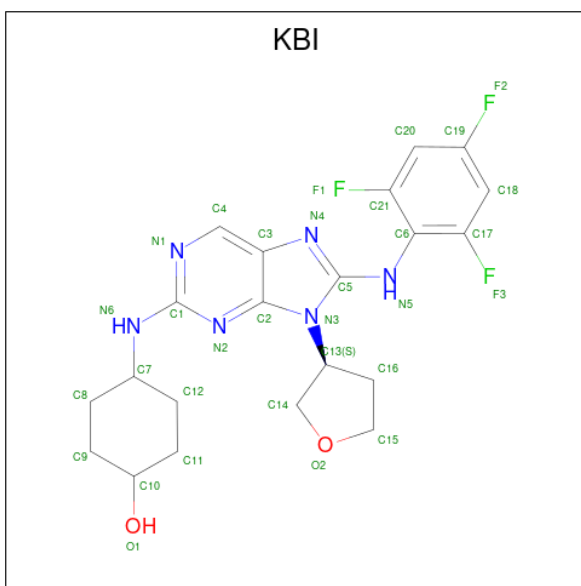
There are 4 unique types of molecules in this entry. The entry contains 3032 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Mitogen-activated protein kinase 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	341	2766	1779	469	497	21	0	0	0

- Molecule 2 is trans-4-({9-[(3S)-tetrahydrofuran-3-yl]-8-[(2,4,6-trifluorophenyl)amino]-9H-purin-2-yl}amino)cyclohexanol (CCD ID: KBI) (formula: C₂₁H₂₃F₃N₆O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	F	N	O		
2	A	1	32	21	3	6	2	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	228	Total O 228 228	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.19Å 71.20Å 106.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.06 – 2.20 33.06 – 2.20	Depositor EDS
% Data completeness (in resolution range)	92.4 (33.06-2.20) 92.3 (33.06-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.20Å)	Xtrriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.242 , 0.335 0.252 , 0.343	Depositor DCC
R_{free} test set	984 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	31.6	Xtrriage
Anisotropy	0.259	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3032	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: KBI, GOL

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	0.83	1/2830 (0.0%)	1.07	8/3830 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	LYS	CA-C	5.61	1.57	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	221	THR	N-CA-C	6.62	113.96	108.07
1	A	313	PHE	CA-C-N	6.27	126.63	119.92
1	A	313	PHE	C-N-CA	6.27	126.63	119.92
1	A	231	ALA	CA-C-N	5.63	126.88	119.84
1	A	231	ALA	C-N-CA	5.63	126.88	119.84
1	A	343	ASP	CA-C-N	5.29	124.95	119.56
1	A	343	ASP	C-N-CA	5.29	124.95	119.56
1	A	328	LYS	N-CA-C	5.12	116.86	111.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2766	0	2789	78	0
2	A	32	0	23	2	0
3	A	6	0	8	0	0
4	A	228	0	0	12	0
All	All	3032	0	2820	79	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 14.

All (79) 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:102:GLN:HE22	1:A:385:HIS:H	1.16	0.94
1:A:295:ARG:HG3	1:A:299:GLU:OE1	1.79	0.81
1:A:111:GLU:O	1:A:115:MET:HB2	1.83	0.79
1:A:290:LEU:CD1	1:A:298:VAL:HG21	2.13	0.78
1:A:188:ARG:HD2	1:A:240:TYR:CE1	2.19	0.78
1:A:399:MET:O	1:A:400:ASN:HB2	1.82	0.78
1:A:103:THR:HG21	1:A:219:MET:HE2	1.67	0.76
1:A:102:GLN:NE2	1:A:385:HIS:H	1.84	0.73
1:A:155:GLN:HA	1:A:158:GLN:HE21	1.53	0.72
1:A:227:ARG:HH12	1:A:230:ARG:NH1	1.92	0.68
1:A:67:LEU:HD22	1:A:81:ALA:HB2	1.76	0.67
2:A:465:KBI:H16	2:A:465:KBI:N2	2.10	0.65
1:A:283:CYS:HB3	4:A:586:HOH:O	1.96	0.64
1:A:173:MET:O	1:A:177:ILE:HD12	1.97	0.64
1:A:120:HIS:HE1	1:A:122:ASN:OD1	1.84	0.60
1:A:319:PRO:O	1:A:325:ASN:HB3	2.01	0.60
1:A:88:ARG:HH21	1:A:147:GLU:HG3	1.67	0.60
1:A:291:GLN:NE2	4:A:667:HOH:O	2.35	0.59
1:A:120:HIS:HB2	1:A:179:HIS:CD2	2.38	0.58
1:A:118:VAL:HG13	1:A:179:HIS:CD2	2.38	0.58
1:A:156:VAL:HG21	1:A:197:VAL:HG21	1.85	0.57
1:A:55:ASP:N	1:A:55:ASP:OD1	2.37	0.57
1:A:398:VAL:HG12	1:A:399:MET:HE2	1.87	0.57
1:A:284:PRO:HD2	4:A:515:HOH:O	2.04	0.57
1:A:116:LYS:NZ	4:A:572:HOH:O	2.40	0.55
1:A:192:PRO:HG3	1:A:252:ILE:HG12	1.88	0.55
1:A:227:ARG:NH1	1:A:230:ARG:NH1	2.54	0.54
1:A:250:GLY:C	1:A:340:LEU:HD21	2.32	0.54
1:A:152:ASN:O	1:A:156:VAL:HG23	2.07	0.54
1:A:99:PHE:O	1:A:387:ILE:HG22	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:CYS:HA	1:A:262:LEU:HD12	1.92	0.52
1:A:301:ARG:HG3	1:A:302:PRO:HD2	1.91	0.51
1:A:394:ILE:O	1:A:398:VAL:HG23	2.10	0.51
1:A:46:ASN:N	4:A:665:HOH:O	2.43	0.51
1:A:102:GLN:NE2	1:A:385:HIS:O	2.45	0.50
1:A:291:GLN:HG3	1:A:292:PRO:HD2	1.92	0.50
1:A:78:VAL:HG21	2:A:465:KBI:H13	1.93	0.49
1:A:97:ARG:HD3	1:A:140:GLN:HE21	1.78	0.49
1:A:308:THR:HB	1:A:310:PRO:HD2	1.94	0.49
1:A:287:MET:O	1:A:295:ARG:HD3	2.13	0.49
1:A:365:PRO:O	1:A:369:GLU:HG2	2.13	0.49
1:A:97:ARG:HD3	1:A:140:GLN:NE2	2.28	0.49
1:A:290:LEU:HD12	1:A:298:VAL:HG21	1.94	0.49
1:A:53:VAL:HG11	1:A:79:CYS:SG	2.53	0.48
1:A:383:ARG:N	1:A:383:ARG:HD3	2.29	0.48
1:A:369:GLU:HB3	4:A:508:HOH:O	2.15	0.47
1:A:356:HIS:O	1:A:360:ASN:N	2.45	0.47
1:A:290:LEU:HD13	1:A:298:VAL:HG21	1.96	0.46
1:A:312:LEU:C	1:A:314:PRO:HD3	2.40	0.46
1:A:196:VAL:CG2	1:A:206:LEU:HD21	2.45	0.46
1:A:102:GLN:HE22	1:A:385:HIS:N	1.98	0.46
1:A:227:ARG:CD	4:A:617:HOH:O	2.63	0.46
1:A:227:ARG:NE	4:A:617:HOH:O	2.47	0.45
1:A:254:GLY:HA3	1:A:262:LEU:HD11	1.98	0.45
1:A:177:ILE:O	1:A:178:LYS:C	2.57	0.45
1:A:251:CYS:HB3	4:A:677:HOH:O	2.17	0.44
1:A:83:ASP:O	1:A:87:ASP:N	2.50	0.44
1:A:97:ARG:NH1	1:A:99:PHE:HE1	2.16	0.44
1:A:243:ASN:HB3	4:A:537:HOH:O	2.17	0.44
1:A:117:CYS:SG	1:A:118:VAL:HG23	2.57	0.43
1:A:133:GLN:HG2	1:A:138:GLU:O	2.18	0.43
1:A:324:HIS:HD2	1:A:325:ASN:OD1	2.01	0.43
1:A:153:LEU:HB3	1:A:192:PRO:HB3	2.01	0.43
1:A:121:LYS:HE2	1:A:367:GLU:HG2	2.01	0.42
1:A:295:ARG:CG	1:A:299:GLU:OE1	2.61	0.42
1:A:311:LYS:HA	1:A:311:LYS:HD3	1.86	0.42
1:A:103:THR:O	1:A:104:HIS:C	2.63	0.41
1:A:271:GLN:O	1:A:275:VAL:HG23	2.19	0.41
1:A:396:LYS:HE3	1:A:396:LYS:HB2	1.96	0.41
1:A:107:ARG:O	1:A:111:GLU:HB2	2.20	0.41
1:A:229:TYR:CD1	1:A:229:TYR:N	2.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:LEU:O	1:A:336:LEU:C	2.63	0.41
1:A:53:VAL:HG21	1:A:79:CYS:SG	2.61	0.41
1:A:77:ILE:N	1:A:77:ILE:HD13	2.36	0.41
1:A:233:GLU:CG	1:A:244:VAL:HG23	2.50	0.41
1:A:106:LYS:HA	1:A:106:LYS:HD2	1.81	0.41
1:A:383:ARG:N	4:A:661:HOH:O	2.54	0.40
1:A:284:PRO:CD	4:A:473:HOH:O	2.69	0.40
1:A:199:SER:C	1:A:201:CYS:H	2.29	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	335/464 (72%)	309 (92%)	24 (7%)	2 (1%)	21 23

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	THR
1	A	222	PRO

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	308/416 (74%)	267 (87%)	41 (13%)	4 3

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	VAL
1	A	52	GLU
1	A	55	ASP
1	A	65	GLN
1	A	67	LEU
1	A	68	LYS
1	A	88	ARG
1	A	111	GLU
1	A	117	CYS
1	A	121	LYS
1	A	125	SER
1	A	127	LEU
1	A	157	ILE
1	A	164	GLU
1	A	170	LEU
1	A	177	ILE
1	A	178	LYS
1	A	190	LEU
1	A	196	VAL
1	A	201	CYS
1	A	220	MET
1	A	221	THR
1	A	224	VAL
1	A	234	VAL
1	A	242	GLU
1	A	271	GLN
1	A	285	GLU
1	A	291	GLN
1	A	295	ARG
1	A	296	ASN
1	A	299	GLU
1	A	308	THR
1	A	330	SER
1	A	338	LYS
1	A	360	ASN
1	A	369	GLU
1	A	383	ARG
1	A	386	THR

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Mol	Chain	Res	Type
1	A	387	ILE
1	A	392	GLU
1	A	400	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	102	GLN
1	A	120	HIS
1	A	140	GLN
1	A	158	GLN
1	A	243	ASN
1	A	271	GLN
1	A	291	GLN
1	A	324	HIS
1	A	355	GLN
1	A	385	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
3	GOL	A	466	-	5,5,5	0.45	0	5,5,5	0.28	0
2	KBI	A	465	-	36,36,36	1.35	4 (11%)	45,52,52	2.34	16 (35%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	466	-	-	2/4/4/4	-
2	KBI	A	465	-	-	3/12/29/29	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	465	KBI	C6-N5	-4.61	1.35	1.43
2	A	465	KBI	C2-N3	-3.42	1.33	1.38
2	A	465	KBI	C3-N4	-2.62	1.34	1.39
2	A	465	KBI	C5-N3	-2.24	1.34	1.37

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	465	KBI	N3-C5-N4	-7.45	108.92	114.42
2	A	465	KBI	N1-C1-N2	-6.04	120.58	126.42
2	A	465	KBI	C3-C2-N2	-5.11	119.65	126.55
2	A	465	KBI	C17-C18-C19	3.72	120.64	116.67
2	A	465	KBI	C4-N1-C1	3.68	120.85	115.81
2	A	465	KBI	C20-C19-C18	-3.37	119.41	123.50
2	A	465	KBI	C14-C13-N3	-3.22	108.11	112.87
2	A	465	KBI	C21-C20-C19	3.09	119.96	116.67
2	A	465	KBI	C21-C6-C17	2.83	118.76	115.70
2	A	465	KBI	F3-C17-C6	2.76	120.51	117.64
2	A	465	KBI	F1-C21-C6	2.27	119.99	117.64
2	A	465	KBI	C2-C3-N4	-2.27	108.01	110.34
2	A	465	KBI	C18-C17-C6	-2.25	120.25	123.39
2	A	465	KBI	C3-N4-C5	2.22	107.53	104.06
2	A	465	KBI	C1-N2-C2	2.07	120.62	113.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	465	KBI	C21-C6-N5	-2.03	119.56	122.12

There are no chirality outliers.

All (5) torsion outliers are listed below:

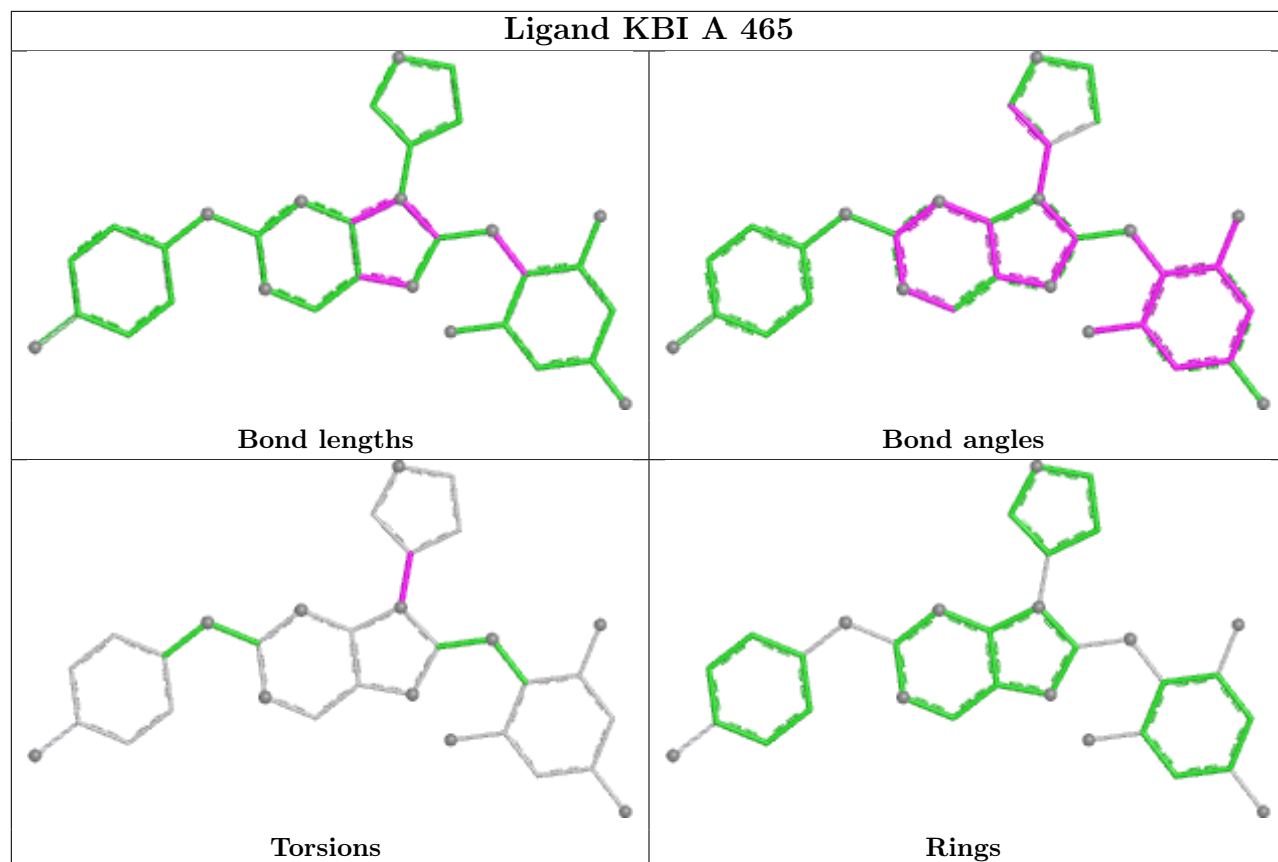
Mol	Chain	Res	Type	Atoms
2	A	465	KBI	C16-C13-N3-C2
2	A	465	KBI	C16-C13-N3-C5
3	A	466	GOL	C1-C2-C3-O3
3	A	466	GOL	O2-C2-C3-O3
2	A	465	KBI	C14-C13-N3-C5

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	465	KBI	2	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 [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	341/464 (73%)	0.47	14 (4%) 41 38	19, 36, 59, 78	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	224	VAL	4.3
1	A	74	ALA	3.7
1	A	211	ALA	3.1
1	A	54	GLY	3.0
1	A	223	TYR	2.7
1	A	294	VAL	2.7
1	A	75	GLN	2.6
1	A	293	THR	2.6
1	A	218	PHE	2.4
1	A	371	PRO	2.4
1	A	304	TYR	2.2
1	A	305	ALA	2.1
1	A	117	CYS	2.1
1	A	373	PRO	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](#)

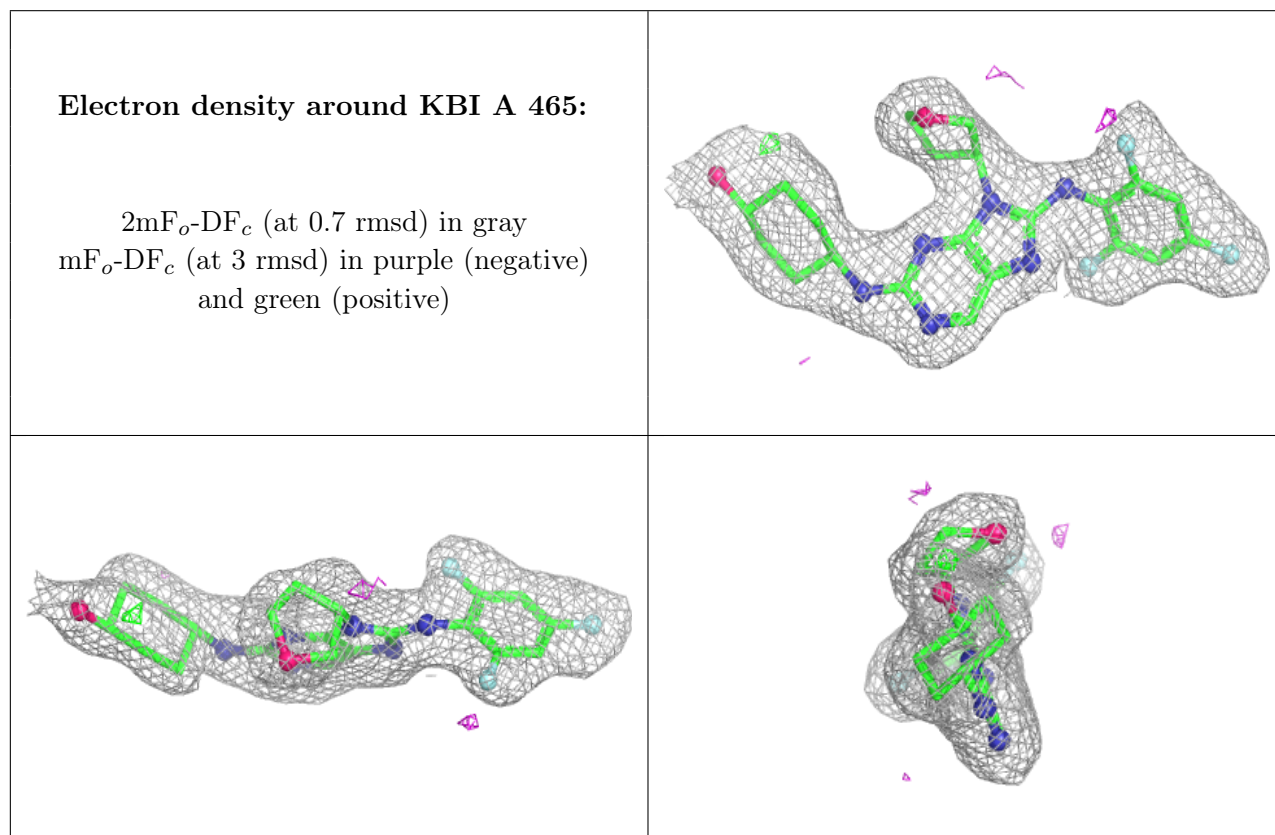
There are no oligosaccharides in this entry.

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(Å ²)	Q<0.9
3	GOL	A	466	6/6	0.66	0.19	61,63,64,64	0
2	KBI	A	465	32/32	0.93	0.08	24,29,32,34	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.