



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 01:18 AM UTC

PDB ID : 10DJ / pdb\_000010dj  
Title : Fyn Kinase Domain-Saracatinib Complex Structure  
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Deposited on : 2026-01-13  
Resolution : 2.22 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 : 4-5-2 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

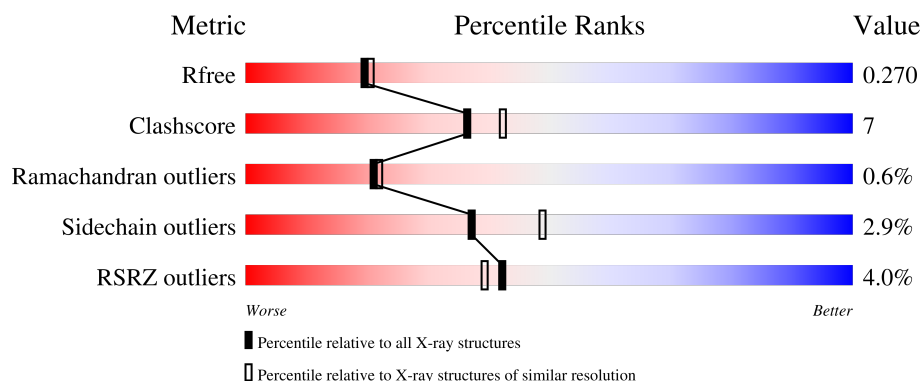
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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  $\geq 3$ , 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	282	
1	B	282	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9214 atoms, of which 4348 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.

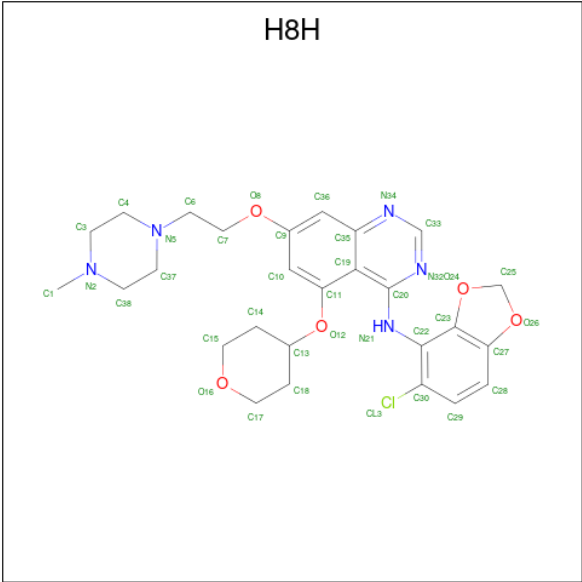
- Molecule 1 is a protein called Tyrosine-protein kinase Fyn.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	269	Total	C	H	N	O	S	0	0	0
			4273	1382	2131	355	391	14			
1	B	277	Total	C	H	N	O	S	0	0	0
			4343	1410	2153	367	399	14			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	538	VAL	-	expression tag	UNP P06241
A	539	ASP	-	expression tag	UNP P06241
A	540	GLU	-	expression tag	UNP P06241
A	541	ASN	-	expression tag	UNP P06241
A	542	LEU	-	expression tag	UNP P06241
A	543	TYR	-	expression tag	UNP P06241
A	544	PHE	-	expression tag	UNP P06241
B	538	VAL	-	expression tag	UNP P06241
B	539	ASP	-	expression tag	UNP P06241
B	540	GLU	-	expression tag	UNP P06241
B	541	ASN	-	expression tag	UNP P06241
B	542	LEU	-	expression tag	UNP P06241
B	543	TYR	-	expression tag	UNP P06241
B	544	PHE	-	expression tag	UNP P06241

- Molecule 2 is N-(5-CHLORO-1,3-BENZODIOXOL-4-YL)-7-[2-(4-METHYLPYPERAZIN-1-YL)ETHOXY]-5-(TETRAHYDRO-2H-PYRAN-4-YLOXY)QUINAZOLIN-4-AMINE (CCD ID: H8H) (formula: C<sub>27</sub>H<sub>32</sub>ClN<sub>5</sub>O<sub>5</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	Cl	H	N	O	0	0
			70	27	1	32	5	5		
2	B	1	Total	C	Cl	H	N	O	0	0
			70	27	1	32	5	5		

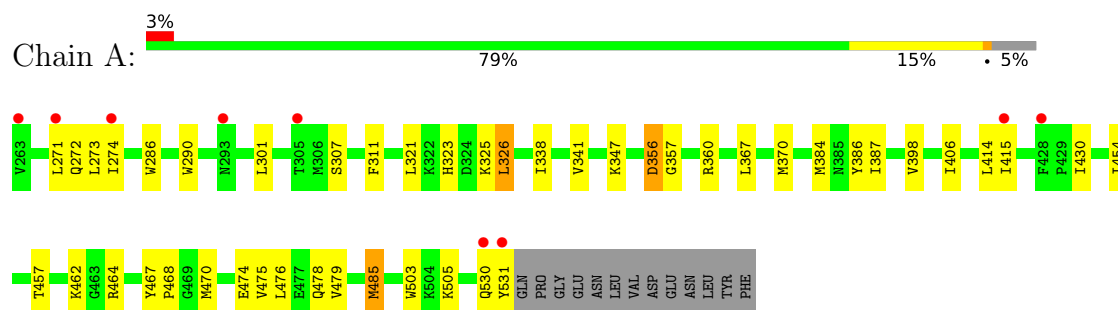
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	245	Total	O	0	0
			245	245		
3	B	213	Total	O	0	0
			213	213		

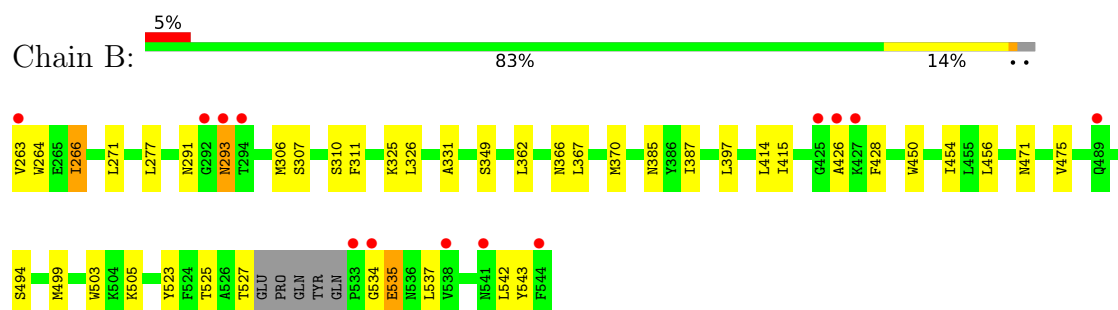
### 3 Residue-property plots [i](#)

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: Tyrosine-protein kinase Fyn



#### • Molecule 1: Tyrosine-protein kinase Fyn



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.69Å 89.60Å 92.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.85 – 2.22 33.85 – 2.22	Depositor EDS
% Data completeness (in resolution range)	99.7 (33.85-2.22) 99.6 (33.85-2.22)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 2.22Å)	Xtriage
Refinement program	PHENIX ("2.0_5885": ???)	Depositor
R, $R_{free}$	0.245 , 0.268 0.246 , 0.270	Depositor DCC
$R_{free}$ test set	1659 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.4	Xtriage
Anisotropy	0.674	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 66.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9214	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1616e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: H8H

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	0.17	0/2194	0.40	0/2973
1	B	0.17	0/2242	0.40	0/3038
All	All	0.17	0/4436	0.40	0/6011

There are no bond length outliers.

There are no bond angle outliers.

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	2142	2131	2125	38	0
1	B	2190	2153	2151	25	0
2	A	38	32	31	0	0
2	B	38	32	31	0	0
3	A	245	0	0	0	0
3	B	213	0	0	0	0
All	All	4866	4348	4338	63	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 7.

All (63) 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:457:THR:HG21	1:A:485:MET:HE3	1.69	0.73
1:B:426:ALA:HB3	1:B:428:PHE:CZ	2.33	0.64
1:A:386:TYR:CZ	1:A:414:LEU:HD23	2.35	0.61
1:A:387:ILE:HG12	1:A:415:ILE:HD11	1.83	0.60
1:A:387:ILE:HD13	1:A:415:ILE:HD11	1.83	0.60
1:A:387:ILE:CD1	1:A:415:ILE:HD11	2.32	0.59
1:B:367:LEU:HD23	1:B:370:MET:CE	2.32	0.59
1:A:387:ILE:CG1	1:A:415:ILE:HD11	2.32	0.59
1:A:414:LEU:HD13	1:A:415:ILE:N	2.18	0.59
1:B:471:ASN:O	1:B:475:VAL:HG23	2.03	0.58
1:A:271:LEU:HD23	1:A:290:TRP:HA	1.85	0.58
1:A:301:LEU:HD22	1:A:311:PHE:CB	2.34	0.57
1:A:301:LEU:HD23	1:A:338:ILE:HD12	1.87	0.57
1:A:387:ILE:HG23	1:A:415:ILE:HD12	1.88	0.55
1:B:266:ILE:HG12	1:B:331:ALA:HB1	1.86	0.55
1:B:385:ASN:O	1:B:414:LEU:HD12	2.06	0.55
1:B:362:LEU:HD22	1:B:366:ASN:CB	2.36	0.54
1:A:367:LEU:HD23	1:A:370:MET:CE	2.37	0.54
1:A:387:ILE:HG12	1:A:415:ILE:CD1	2.38	0.54
1:A:273:LEU:HD23	1:A:286:TRP:CG	2.45	0.52
1:A:384:MET:HE2	1:A:384:MET:HA	1.91	0.52
1:B:325:LYS:C	1:B:326:LEU:HD12	2.34	0.52
1:A:474:GLU:O	1:A:478:GLN:HG2	2.11	0.51
1:A:467:TYR:HB3	1:A:470:MET:HE3	1.93	0.51
1:B:454:ILE:HD13	1:B:503:TRP:CZ2	2.45	0.50
1:A:301:LEU:HD22	1:A:311:PHE:HB2	1.93	0.50
1:A:454:ILE:HD13	1:A:503:TRP:CZ2	2.47	0.49
1:A:464:ARG:NH1	1:A:468:PRO:HD3	2.26	0.49
1:A:321:LEU:HD21	1:A:386:TYR:CD2	2.48	0.49
1:A:387:ILE:HG23	1:A:415:ILE:CD1	2.42	0.48
1:B:306:MET:HE3	1:B:310:SER:OG	2.14	0.48
1:B:494:SER:OG	1:B:523:TYR:OH	2.31	0.48
1:A:325:LYS:C	1:A:326:LEU:HD23	2.38	0.48
1:A:530:GLN:O	1:A:531:TYR:CB	2.63	0.47
1:A:325:LYS:O	1:A:326:LEU:HD23	2.14	0.47
1:A:530:GLN:O	1:A:531:TYR:HB2	2.15	0.47
1:A:430:ILE:HG23	1:A:476:LEU:HD11	1.98	0.46
1:B:263:VAL:HG22	1:B:264:TRP:CD1	2.50	0.46
1:A:386:TYR:CE2	1:A:414:LEU:HD23	2.50	0.46
1:B:542:LEU:HD21	1:B:543:TYR:CD2	2.52	0.45
1:A:475:VAL:O	1:A:479:VAL:HG22	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:ASP:HA	1:A:360:ARG:HD3	1.99	0.45
1:B:263:VAL:HG13	1:B:264:TRP:N	2.32	0.44
1:A:414:LEU:HD13	1:A:414:LEU:C	2.43	0.44
1:B:291:ASN:C	1:B:293:ASN:H	2.25	0.44
1:A:323:HIS:HB3	1:A:326:LEU:HG	1.99	0.44
1:A:272:GLN:HG3	1:A:274:ILE:CD1	2.48	0.44
1:B:349:SER:HA	1:B:397:LEU:HD23	2.00	0.44
1:A:457:THR:HG21	1:A:485:MET:CE	2.43	0.44
1:B:456:LEU:HB3	1:B:499:MET:HE2	2.00	0.43
1:B:325:LYS:O	1:B:326:LEU:HD12	2.18	0.42
1:A:387:ILE:CG2	1:A:415:ILE:CD1	2.97	0.42
1:B:266:ILE:HD11	1:B:271:LEU:HD11	2.00	0.42
1:B:362:LEU:HD22	1:B:366:ASN:HB3	2.01	0.42
1:B:263:VAL:HG13	1:B:264:TRP:H	1.84	0.42
1:B:450:TRP:CE3	1:B:503:TRP:HA	2.55	0.41
1:B:306:MET:HE2	1:B:311:PHE:HB2	2.03	0.41
1:A:271:LEU:HD11	1:A:341:VAL:HG21	2.02	0.41
1:A:326:LEU:HD22	1:A:406:ILE:HB	2.02	0.41
1:B:387:ILE:HG23	1:B:415:ILE:CD1	2.51	0.41
1:B:534:GLY:O	1:B:535:GLU:CB	2.69	0.40
1:B:456:LEU:CB	1:B:499:MET:HE2	2.52	0.40
1:A:347:LYS:HB2	1:A:398:VAL:HB	2.03	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	267/282 (95%)	260 (97%)	6 (2%)	1 (0%)	30	33
1	B	273/282 (97%)	264 (97%)	7 (3%)	2 (1%)	18	18
All	All	540/564 (96%)	524 (97%)	13 (2%)	3 (1%)	21	22

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	293	ASN
1	B	535	GLU
1	A	357	GLY

### 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	226/244 (93%)	220 (97%)	6 (3%)	39	52
1	B	228/244 (93%)	221 (97%)	7 (3%)	35	46
All	All	454/488 (93%)	441 (97%)	13 (3%)	37	49

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	307	SER
1	A	326	LEU
1	A	356	ASP
1	A	462	LYS
1	A	485	MET
1	A	505	LYS
1	B	266	ILE
1	B	277	LEU
1	B	307	SER
1	B	505	LYS
1	B	525	THR
1	B	527	THR
1	B	537	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	366	ASN

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Mol	Chain	Res	Type
1	B	272	GLN
1	B	316	GLN
1	B	418	ASN

### 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	H8H	B	601	-	43,43,43	3.16	17 (39%)	54,60,60	2.93	15 (27%)
2	H8H	A	601	-	43,43,43	3.12	17 (39%)	54,60,60	3.04	14 (25%)

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	H8H	B	601	-	-	3/14/38/38	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	H8H	A	601	-	-	5/14/38/38	0/6/6/6

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	H8H	O24-C23	-10.61	1.23	1.38
2	B	601	H8H	O24-C23	-10.59	1.23	1.38
2	B	601	H8H	O26-C27	-8.92	1.23	1.38
2	A	601	H8H	O26-C27	-8.80	1.23	1.38
2	B	601	H8H	C35-N34	-6.15	1.27	1.37
2	A	601	H8H	C35-N34	-6.13	1.27	1.37
2	B	601	H8H	C20-N21	-5.22	1.28	1.36
2	A	601	H8H	C20-N21	-4.99	1.28	1.36
2	A	601	H8H	C36-C35	-4.80	1.34	1.41
2	B	601	H8H	C36-C35	-4.77	1.34	1.41
2	B	601	H8H	C20-N32	-4.67	1.27	1.34
2	A	601	H8H	C20-N32	-4.42	1.28	1.34
2	B	601	H8H	C22-N21	-4.31	1.28	1.41
2	A	601	H8H	C22-N21	-4.26	1.28	1.41
2	B	601	H8H	C33-N32	-4.22	1.26	1.33
2	A	601	H8H	C33-N32	-4.12	1.26	1.33
2	A	601	H8H	C33-N34	-3.76	1.26	1.32
2	B	601	H8H	C19-C35	-3.76	1.36	1.42
2	B	601	H8H	C33-N34	-3.61	1.26	1.32
2	A	601	H8H	C19-C35	-3.61	1.36	1.42
2	A	601	H8H	C11-C19	-2.86	1.36	1.42
2	B	601	H8H	C11-C19	-2.81	1.36	1.42
2	B	601	H8H	C22-C30	-2.80	1.36	1.40
2	A	601	H8H	C28-C27	-2.69	1.34	1.39
2	A	601	H8H	C29-C28	-2.69	1.34	1.38
2	B	601	H8H	C28-C27	-2.61	1.34	1.39
2	B	601	H8H	C10-C9	-2.57	1.34	1.39
2	B	601	H8H	C29-C28	-2.56	1.34	1.38
2	A	601	H8H	C20-C19	-2.54	1.36	1.43
2	B	601	H8H	C20-C19	-2.48	1.37	1.43
2	A	601	H8H	C10-C9	-2.35	1.35	1.39
2	A	601	H8H	C22-C30	-2.33	1.36	1.40
2	A	601	H8H	C27-C23	-2.11	1.34	1.40
2	B	601	H8H	C27-C23	-2.06	1.34	1.40

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	H8H	C19-C20-N21	14.07	133.95	119.75
2	B	601	H8H	C19-C20-N21	13.39	133.26	119.75
2	A	601	H8H	N21-C20-N32	-9.13	108.45	118.71
2	B	601	H8H	N21-C20-N32	-9.13	108.46	118.71
2	B	601	H8H	C22-N21-C20	6.42	139.03	123.53
2	A	601	H8H	C22-N21-C20	5.91	137.80	123.53
2	A	601	H8H	C33-N32-C20	5.59	120.94	116.60
2	A	601	H8H	C36-C9-C10	-4.82	115.52	121.14
2	B	601	H8H	C36-C9-C10	-4.76	115.59	121.14
2	B	601	H8H	C33-N32-C20	4.62	120.18	116.60
2	A	601	H8H	C11-C10-C9	4.34	125.13	120.33
2	B	601	H8H	C11-C10-C9	4.14	124.91	120.33
2	A	601	H8H	C22-C30-CL3	4.12	123.97	119.28
2	A	601	H8H	O26-C25-O24	-3.85	102.06	108.09
2	B	601	H8H	O26-C25-O24	-3.81	102.11	108.09
2	A	601	H8H	C25-O24-C23	3.01	108.63	105.05
2	A	601	H8H	C36-C35-N34	-2.82	114.77	118.01
2	B	601	H8H	C25-O24-C23	2.81	108.39	105.05
2	B	601	H8H	C30-C22-N21	-2.64	118.49	121.97
2	B	601	H8H	C22-C30-CL3	2.49	122.11	119.28
2	A	601	H8H	O26-C27-C23	2.42	112.07	109.62
2	A	601	H8H	C11-O12-C13	2.40	122.89	119.11
2	B	601	H8H	O26-C27-C23	2.32	111.97	109.62
2	A	601	H8H	O12-C13-C14	2.22	113.50	108.31
2	B	601	H8H	C36-C35-N34	-2.14	115.56	118.01
2	A	601	H8H	C36-C35-C19	2.13	122.63	120.40
2	B	601	H8H	C11-O12-C13	2.09	122.39	119.11
2	B	601	H8H	O8-C7-C6	2.05	112.97	107.79
2	B	601	H8H	O12-C13-C14	2.03	113.05	108.31

There are no chirality outliers.

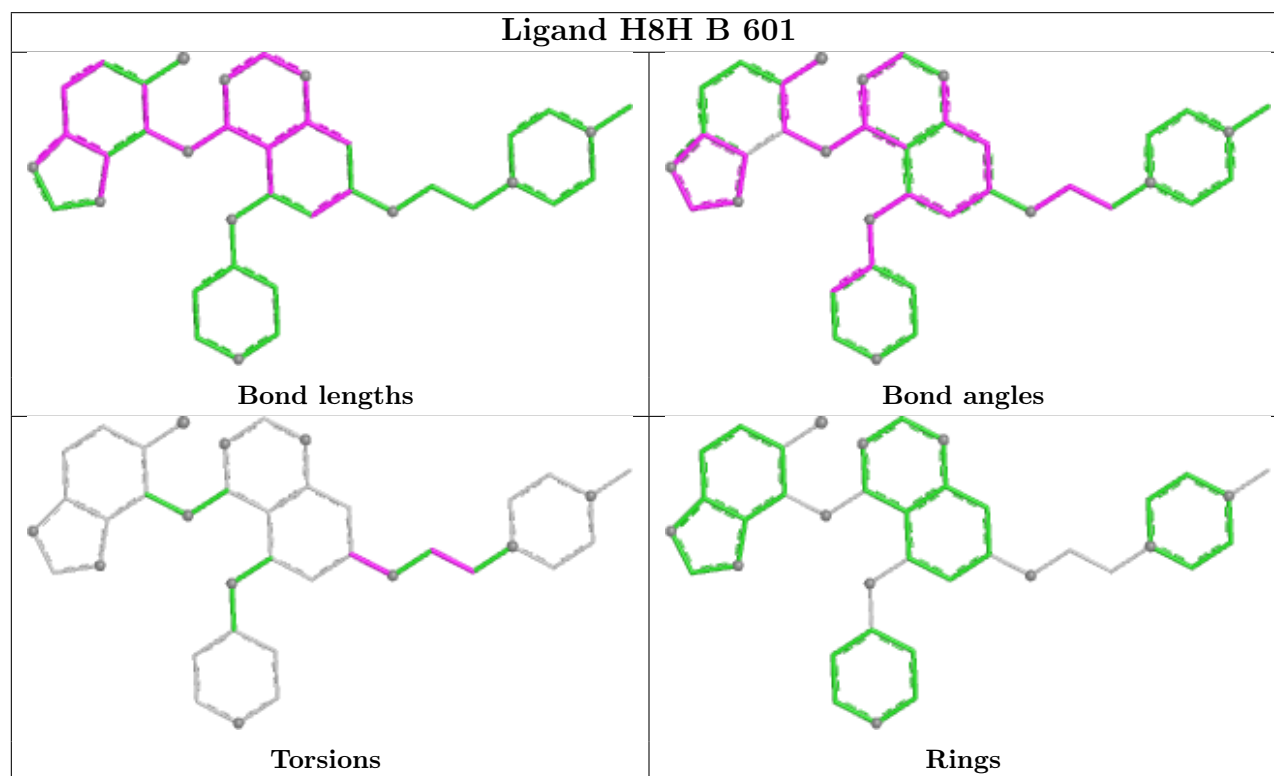
All (8) torsion outliers are listed below:

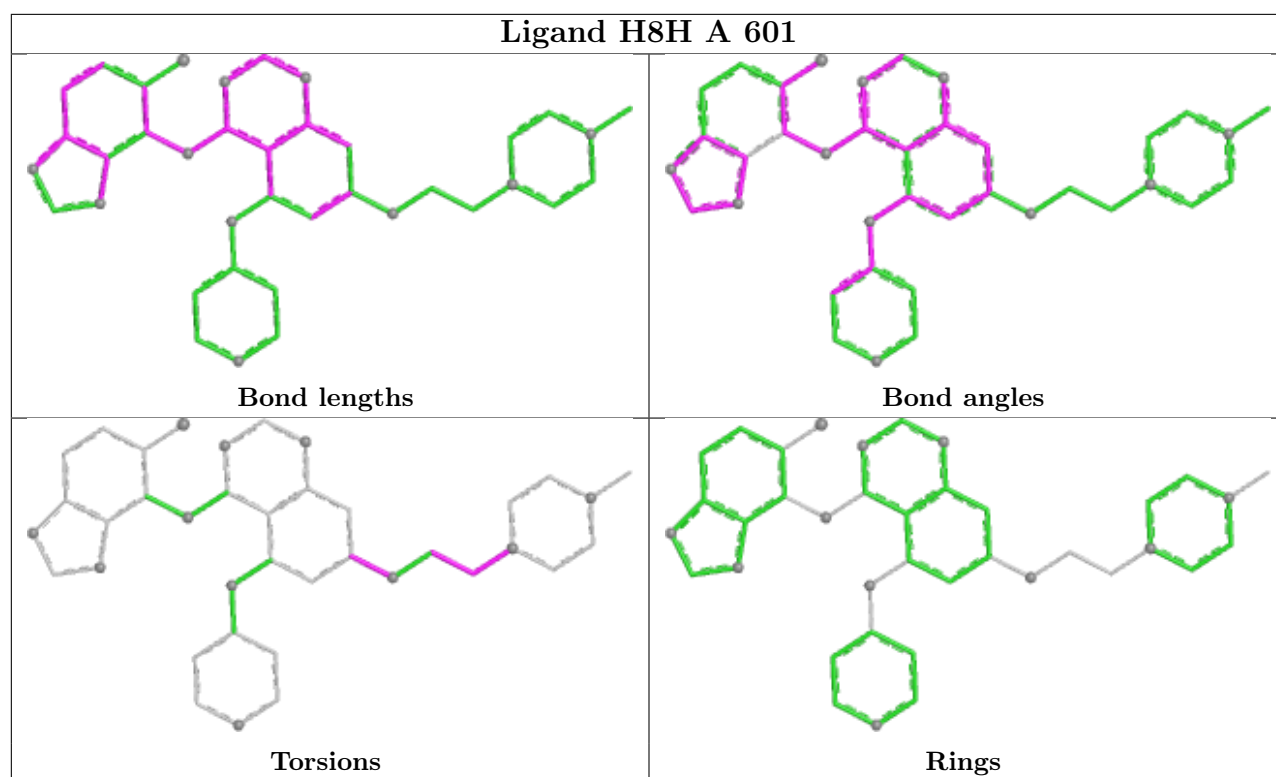
Mol	Chain	Res	Type	Atoms
2	A	601	H8H	N5-C6-C7-O8
2	B	601	H8H	N5-C6-C7-O8
2	B	601	H8H	C10-C9-O8-C7
2	A	601	H8H	C10-C9-O8-C7
2	B	601	H8H	C36-C9-O8-C7
2	A	601	H8H	C36-C9-O8-C7
2	A	601	H8H	C7-C6-N5-C37
2	A	601	H8H	C7-C6-N5-C4

There are no ring outliers.

No monomer is involved in short contacts.

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	269/282 (95%)	0.45	9 (3%) 49 46	32, 48, 90, 124	0
1	B	277/282 (98%)	0.44	13 (4%) 36 33	28, 44, 83, 120	0
All	All	546/564 (96%)	0.44	22 (4%) 42 39	28, 46, 85, 124	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	VAL	6.2
1	A	428	PHE	5.1
1	B	263	VAL	4.0
1	B	533	PRO	3.9
1	B	293	ASN	3.5
1	A	293	ASN	3.2
1	A	531	TYR	3.2
1	B	294	THR	3.2
1	A	415	ILE	3.2
1	B	427	LYS	2.9
1	B	425	GLY	2.8
1	A	271	LEU	2.6
1	B	541	ASN	2.6
1	B	538	VAL	2.6
1	B	534	GLY	2.5
1	B	292	GLY	2.3
1	A	274	ILE	2.3
1	B	544	PHE	2.3
1	A	305	THR	2.2
1	B	426	ALA	2.2
1	A	530	GLN	2.1
1	B	489	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](#)

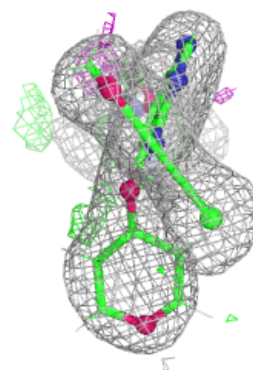
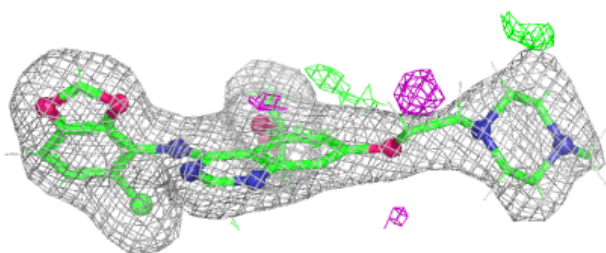
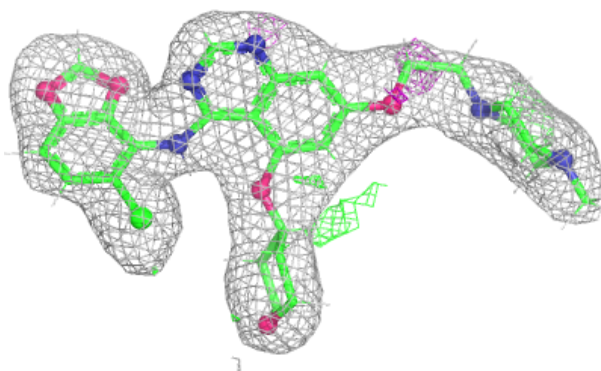
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
2	H8H	A	601	38/38	0.92	0.10	35,47,66,67	0
2	H8H	B	601	38/38	0.93	0.08	28,42,56,64	0

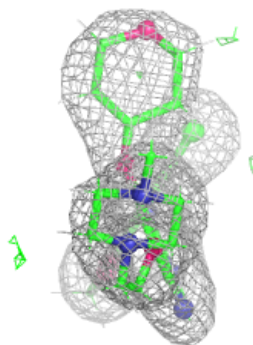
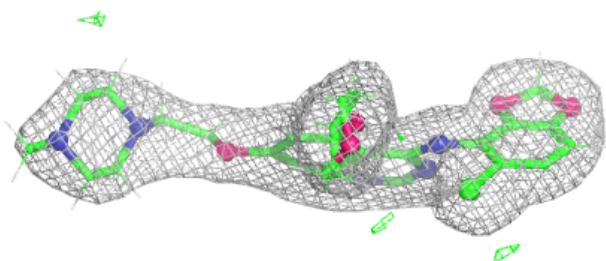
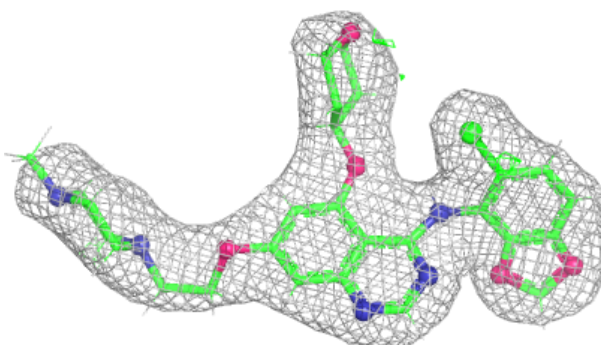
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around H8H A 601:**

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

**Electron density around H8H B 601:**

$2mF_o-DF_c$  (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.