



## Full wwPDB X-ray Structure Validation Report ⓘ

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Title : Discovery of IRAK4 Inhibitor 16  
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Deposited on : 2022-08-22  
Resolution : 2.35 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](#) ①) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
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

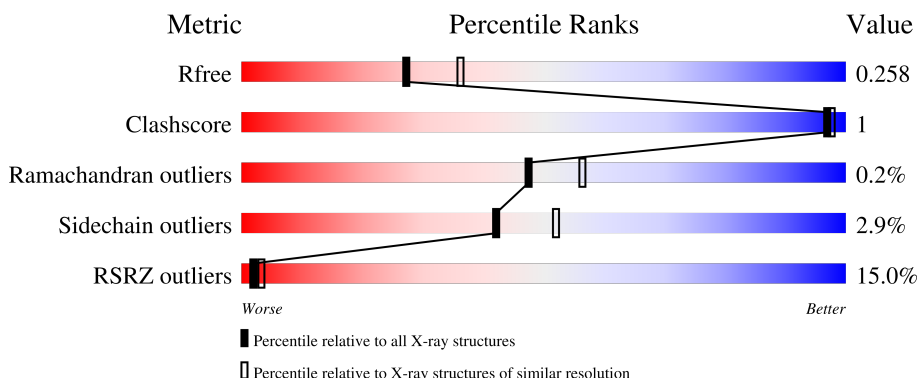
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

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	AAA	298	
1	BBB	298	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	SEP	BBB	346	-	-	-	X

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4594 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.

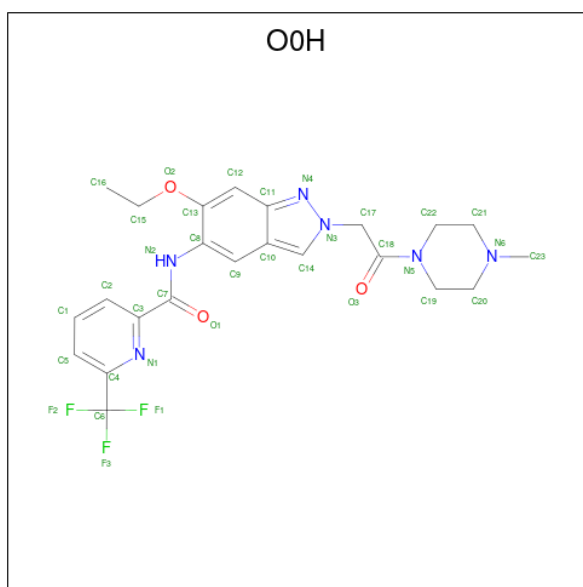
- Molecule 1 is a protein called Interleukin-1 receptor-associated kinase 4.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	AAA	280	2204	1384	370	434	2	14	0	0	0
1	BBB	282	2220	1394	373	436	2	15	0	1	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	163	GLY	-	expression tag	UNP Q9NWZ3
AAA	164	SER	-	expression tag	UNP Q9NWZ3
AAA	400	ALA	LYS	engineered mutation	UNP Q9NWZ3
AAA	401	ALA	GLU	engineered mutation	UNP Q9NWZ3
AAA	402	ALA	GLU	engineered mutation	UNP Q9NWZ3
BBB	163	GLY	-	expression tag	UNP Q9NWZ3
BBB	164	SER	-	expression tag	UNP Q9NWZ3
BBB	400	ALA	LYS	engineered mutation	UNP Q9NWZ3
BBB	401	ALA	GLU	engineered mutation	UNP Q9NWZ3
BBB	402	ALA	GLU	engineered mutation	UNP Q9NWZ3

- Molecule 2 is {N}-[6-ethoxy-2-[2-(4-methylpiperazin-1-yl)-2-oxidanylidene-ethyl]indazol-5-yl]-6-(trifluoromethyl)pyridine-2-carboxamide (three-letter code: OOH) (formula: C<sub>23</sub>H<sub>25</sub>F<sub>3</sub>N<sub>6</sub>O<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	F	N			O
2	AAA	1	35	23	3	6	3	0	0
2	BBB	1	35	23	3	6	3	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

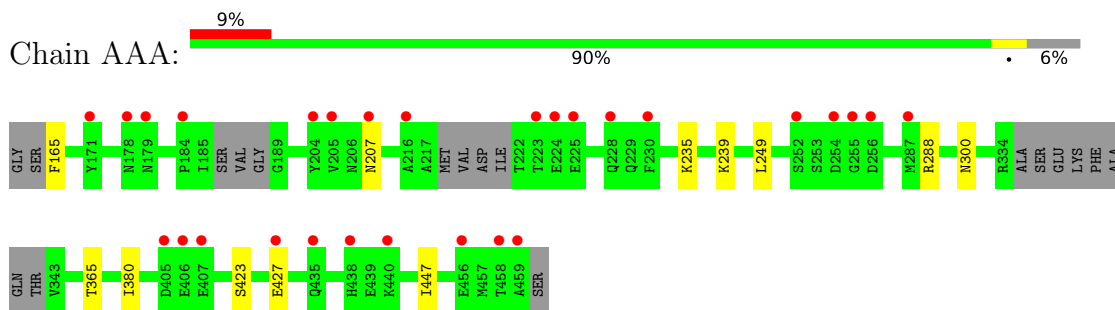
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AAA	55	Total 55	O 55	0	0
4	BBB	39	Total 39	O 39	0	0

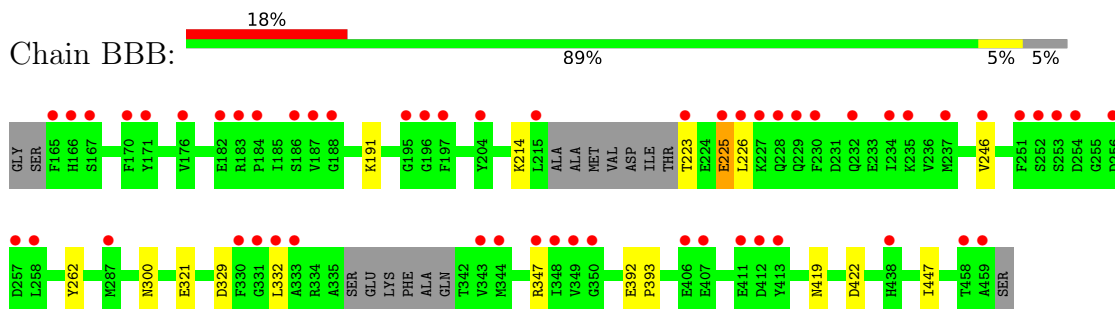
### 3 Residue-property plots [i](#)

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: Interleukin-1 receptor-associated kinase 4



- Molecule 1: Interleukin-1 receptor-associated kinase 4





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.20Å 117.50Å 136.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.89 – 2.35 48.89 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.89-2.35) 99.9 (48.89-2.35)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.58 (at 2.34Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.212 , 0.260 0.218 , 0.258	Depositor DCC
$R_{free}$ test set	888 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.4	Xtrriage
Anisotropy	0.112	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 47.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4594	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.48% 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: SEP, OOH, GOL, TPO, CSO

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	AAA	0.65	0/2202	0.70	0/2964
1	BBB	0.65	0/2219	0.70	0/2988
All	All	0.65	0/4421	0.70	0/5952

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 [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	AAA	2204	0	2166	4	0
1	BBB	2220	0	2182	4	0
2	AAA	35	0	0	0	0
2	BBB	35	0	0	0	0
3	AAA	6	0	8	0	0
4	AAA	55	0	0	0	0
4	BBB	39	0	0	0	0
All	All	4594	0	4356	8	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 1.

All (8) 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:AAA:423:SER:O	1:AAA:427:GLU:HG3	2.15	0.46
1:BBB:321:GLU:H	1:BBB:321:GLU:CD	2.22	0.43
1:BBB:300:ASN:HA	1:BBB:447:ILE:HG21	2.01	0.42
1:AAA:300:ASN:HA	1:AAA:447:ILE:HG21	2.02	0.42
1:BBB:246:VAL:CG1	1:BBB:262:TYR:CD2	3.04	0.41
1:AAA:165:PHE:HA	1:AAA:249:LEU:O	2.20	0.41
1:AAA:288:ARG:HB3	1:AAA:380:ILE:HG23	2.03	0.40
1:BBB:392:GLU:HA	1:BBB:393:PRO:HA	1.93	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	AAA	268/298 (90%)	259 (97%)	9 (3%)	0	100	100
1	BBB	273/298 (92%)	260 (95%)	12 (4%)	1 (0%)	34	38
All	All	541/596 (91%)	519 (96%)	21 (4%)	1 (0%)	47	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	225	GLU

### 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	AAA	237/251 (94%)	233 (98%)	4 (2%)	60	72
1	BBB	239/251 (95%)	229 (96%)	10 (4%)	30	36
All	All	476/502 (95%)	462 (97%)	14 (3%)	42	52

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	207	ASN
1	AAA	235	LYS
1	AAA	239	LYS
1	AAA	365	THR
1	BBB	191	LYS
1	BBB	214	LYS
1	BBB	223	THR
1	BBB	225	GLU
1	BBB	226	LEU
1	BBB	329	ASP
1	BBB	332	LEU
1	BBB	347	ARG
1	BBB	419	ASN
1	BBB	422	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

8 non-standard protein/DNA/RNA residues 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
1	CSO	AAA	240	1	3,6,7	0.77	0	0,6,8	-	-
1	SEP	BBB	346	1	8,9,10	0.59	0	8,12,14	0.64	0
1	CSO	AAA	289	1	3,6,7	0.67	0	0,6,8	-	-
1	TPO	BBB	345	1	8,10,11	0.81	0	10,14,16	0.78	0
1	SEP	AAA	346	1	8,9,10	0.61	0	8,12,14	0.65	0
1	CSO	BBB	240	1	3,6,7	0.74	0	0,6,8	-	-
1	CSO	BBB	289	1	3,6,7	0.67	0	0,6,8	-	-
1	TPO	AAA	345	1	8,10,11	0.68	0	10,14,16	0.84	0

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
1	CSO	AAA	240	1	-	1/1/5/7	-
1	SEP	BBB	346	1	-	0/5/8/10	-
1	CSO	AAA	289	1	-	0/1/5/7	-
1	TPO	BBB	345	1	-	3/9/11/13	-
1	SEP	AAA	346	1	-	1/5/8/10	-
1	CSO	BBB	240	1	-	0/1/5/7	-
1	CSO	BBB	289	1	-	0/1/5/7	-
1	TPO	AAA	345	1	-	2/9/11/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	AAA	345	TPO	N-CA-CB-OG1
1	AAA	346	SEP	N-CA-CB-OG
1	BBB	345	TPO	N-CA-CB-OG1
1	BBB	345	TPO	O-C-CA-CB
1	BBB	345	TPO	CB-OG1-P-O1P
1	AAA	240	CSO	N-CA-CB-SG
1	AAA	345	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

3 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	AAA	502	-	5,5,5	0.10	0	5,5,5	0.29	0
2	O0H	AAA	501	-	35,38,38	0.44	0	50,55,55	1.12	4 (8%)
2	O0H	BBB	501	-	35,38,38	0.37	0	50,55,55	0.97	3 (6%)

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	GOL	AAA	502	-	-	2/4/4/4	-
2	O0H	AAA	501	-	-	7/24/35/35	1/4/4/4
2	O0H	BBB	501	-	-	6/24/35/35	0/4/4/4

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	501	O0H	C18-C17-N3	3.70	115.17	110.76
2	BBB	501	O0H	C14-N3-N4	3.65	114.31	111.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	501	O0H	C14-N3-N4	3.59	114.26	111.45
2	AAA	501	O0H	C17-N3-C14	-3.21	125.10	129.19
2	BBB	501	O0H	C18-C17-N3	2.84	114.14	110.76
2	AAA	501	O0H	O2-C13-C12	-2.11	121.62	125.19
2	BBB	501	O0H	C12-C11-N4	-2.01	127.74	130.28

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	501	O0H	C17-C18-N5-C22
2	AAA	501	O0H	O3-C18-N5-C22
2	BBB	501	O0H	C17-C18-N5-C19
2	BBB	501	O0H	O3-C18-N5-C19
3	AAA	502	GOL	O1-C1-C2-C3
2	AAA	501	O0H	C8-C13-O2-C15
2	AAA	501	O0H	C12-C13-O2-C15
2	AAA	501	O0H	C16-C15-O2-C13
2	BBB	501	O0H	C8-C13-O2-C15
2	AAA	501	O0H	O3-C18-N5-C19
2	BBB	501	O0H	C12-C13-O2-C15
2	AAA	501	O0H	C17-C18-N5-C19
2	BBB	501	O0H	C2-C3-C7-N2
3	AAA	502	GOL	O1-C1-C2-O2
2	BBB	501	O0H	N1-C3-C7-N2

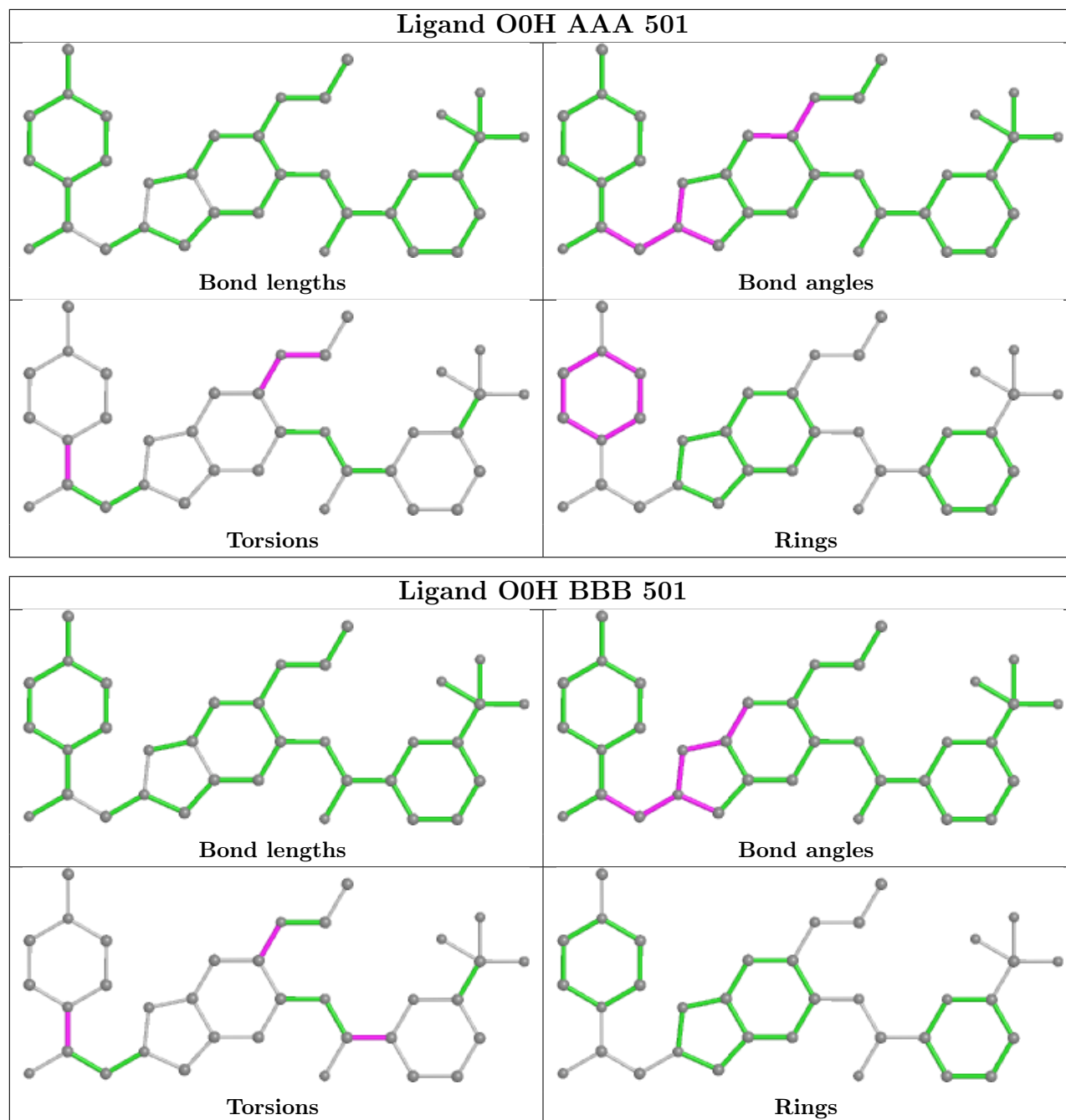
All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	501	O0H	C19-C20-C21-C22-N5-N6

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

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	AAA	276/298 (92%)	0.58	28 (10%) <b>7</b> <b>11</b>	27, 52, 90, 112	0
1	BBB	278/298 (93%)	1.17	55 (19%) <b>1</b> <b>2</b>	26, 54, 116, 145	0
All	All	554/596 (92%)	0.87	83 (14%) <b>2</b> <b>3</b>	26, 53, 108, 145	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	350	GLY	8.5
1	BBB	331	GLY	8.2
1	BBB	196	GLY	7.7
1	BBB	223	THR	6.9
1	BBB	187	VAL	6.4
1	BBB	230	PHE	6.2
1	BBB	225	GLU	5.8
1	BBB	226	LEU	5.8
1	BBB	258	LEU	5.8
1	BBB	256	ASP	5.4
1	BBB	165	PHE	5.4
1	BBB	215	LEU	5.3
1	BBB	254	ASP	4.9
1	AAA	256	ASP	4.8
1	AAA	459	ALA	4.8
1	BBB	253	SER	4.6
1	BBB	182	GLU	4.6
1	BBB	347	ARG	4.5
1	BBB	348	ILE	4.4
1	AAA	438	HIS	4.2
1	BBB	343	VAL	4.2
1	BBB	171	TYR	3.8
1	BBB	229	GLN	3.8
1	BBB	459	ALA	3.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	AAA	255	GLY	3.7
1	BBB	234	ILE	3.7
1	BBB	166	HIS	3.6
1	BBB	411	GLU	3.6
1	BBB	333	ALA	3.6
1	BBB	252	SER	3.6
1	AAA	178	ASN	3.4
1	BBB	332	LEU	3.3
1	BBB	197	PHE	3.3
1	BBB	406	GLU	3.2
1	BBB	407	GLU	3.2
1	BBB	183	ARG	3.0
1	BBB	349	VAL	3.0
1	BBB	235	LYS	3.0
1	AAA	224	GLU	2.9
1	BBB	170	PHE	2.9
1	BBB	167	SER	2.9
1	AAA	254	ASP	2.8
1	AAA	225	GLU	2.8
1	AAA	405	ASP	2.8
1	BBB	195	GLY	2.8
1	AAA	223	THR	2.8
1	AAA	207	ASN	2.8
1	BBB	186	SER	2.7
1	BBB	287[A]	MET	2.7
1	AAA	230	PHE	2.7
1	AAA	216	ALA	2.7
1	BBB	257	ASP	2.7
1	BBB	204	TYR	2.6
1	BBB	330	PHE	2.6
1	AAA	228	GLN	2.5
1	BBB	438	HIS	2.5
1	AAA	440	LYS	2.5
1	BBB	251	PHE	2.5
1	BBB	458	THR	2.5
1	AAA	171	TYR	2.4
1	BBB	412	ASP	2.4
1	BBB	227	LYS	2.4
1	AAA	179	ASN	2.4
1	AAA	252	SER	2.3
1	BBB	228	GLN	2.3
1	AAA	205	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	BBB	188	GLY	2.2
1	BBB	232	GLN	2.2
1	AAA	287	MET	2.2
1	AAA	406	GLU	2.2
1	BBB	237	MET	2.2
1	BBB	344	MET	2.2
1	BBB	176	VAL	2.1
1	BBB	413	TYR	2.1
1	AAA	435	GLN	2.1
1	AAA	456	GLU	2.1
1	AAA	204	TYR	2.1
1	BBB	246	VAL	2.1
1	AAA	184	PRO	2.1
1	AAA	427	GLU	2.1
1	AAA	407	GLU	2.0
1	AAA	458	THR	2.0
1	BBB	184	PRO	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains [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(Å <sup>2</sup> )	Q<0.9
1	SEP	BBB	346	10/11	0.37	0.49	125,147,174,174	0
1	SEP	AAA	346	10/11	0.59	0.30	100,118,139,140	0
1	TPO	BBB	345	11/12	0.70	0.28	103,112,119,120	0
1	CSO	AAA	240	7/8	0.83	0.18	56,62,67,69	0
1	TPO	AAA	345	11/12	0.87	0.17	84,89,94,94	0
1	CSO	BBB	240	7/8	0.89	0.16	55,59,65,66	0
1	CSO	BBB	289	7/8	0.91	0.18	35,37,42,44	0
1	CSO	AAA	289	7/8	0.92	0.15	34,36,40,42	0

## 6.3 Carbohydrates [i](#)

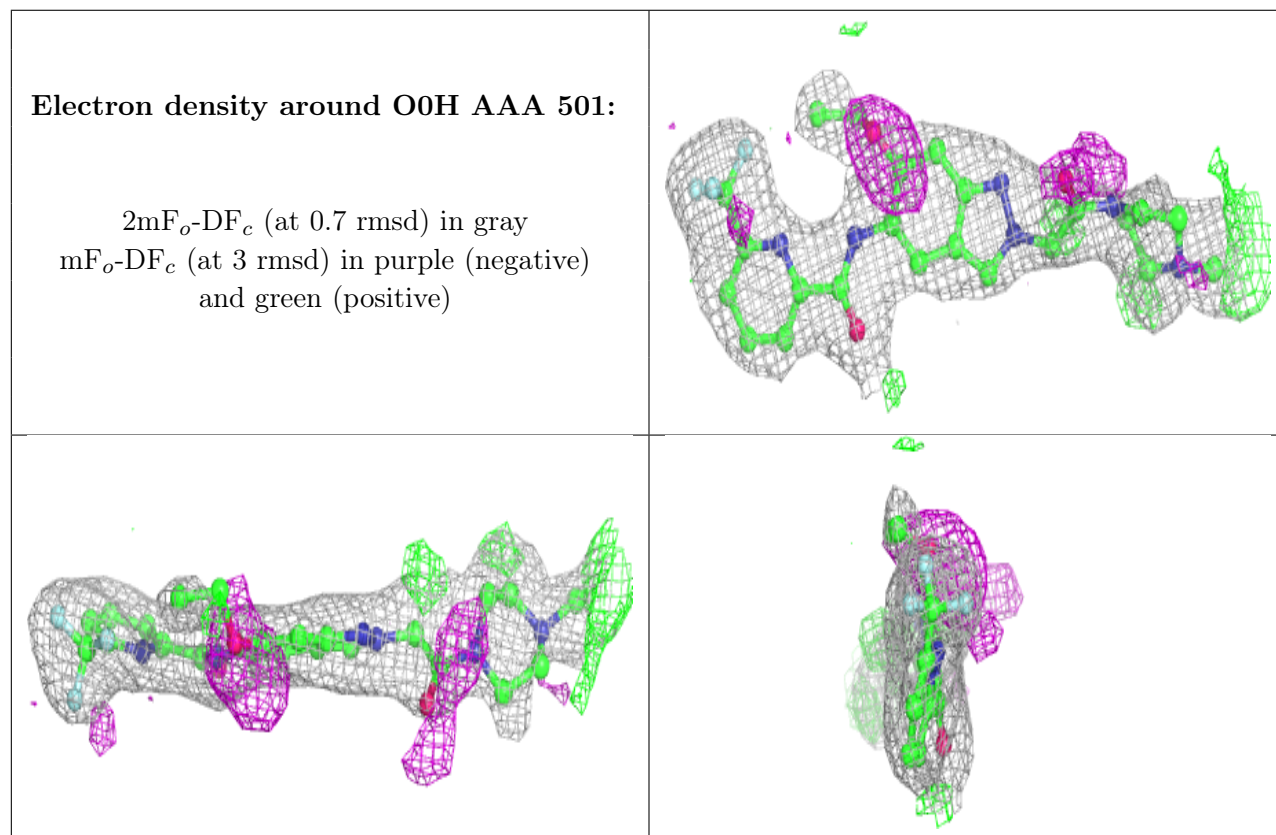
There are no monosaccharides in this entry.

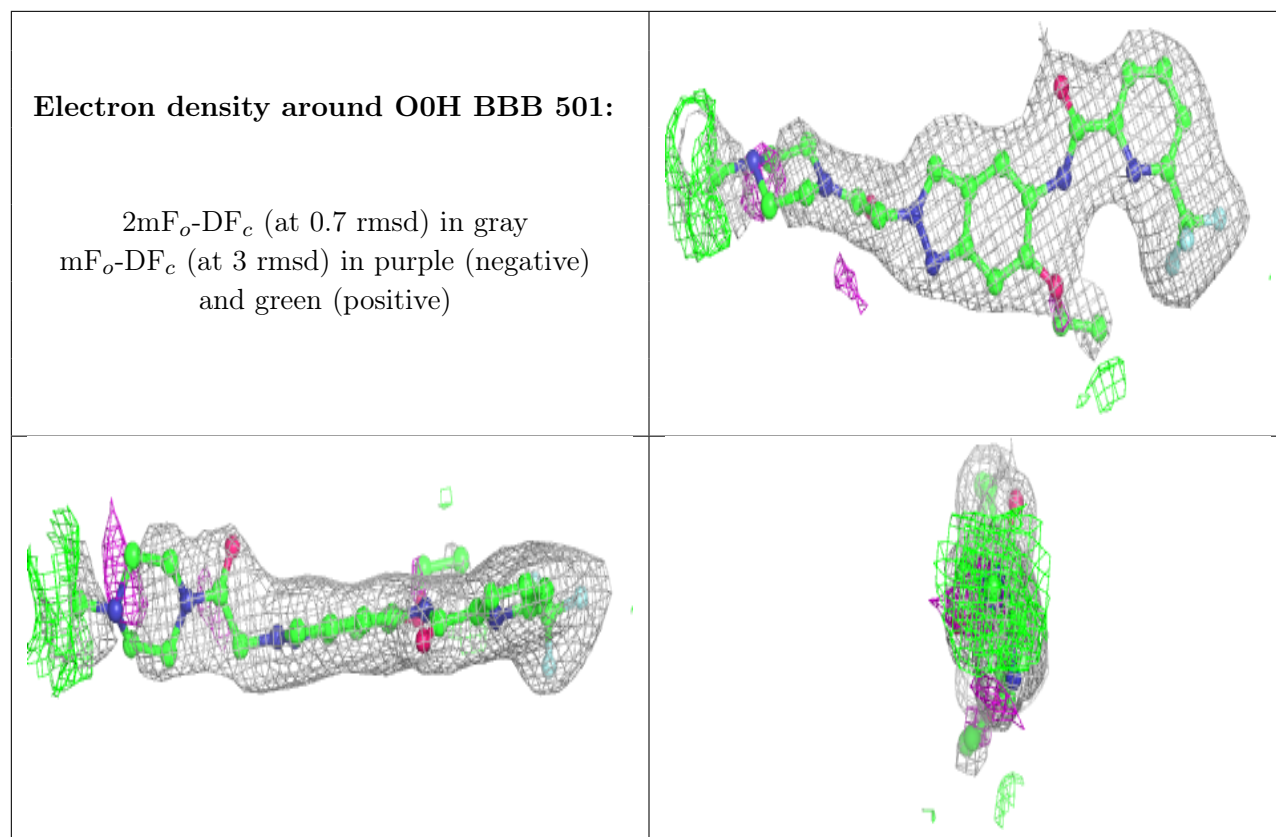
## 6.4 Ligands [i](#)

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(Å <sup>2</sup> )	Q<0.9
3	GOL	AAA	502	6/6	0.83	0.13	77,81,83,84	0
2	O0H	AAA	501	35/35	0.85	0.22	39,47,76,78	0
2	O0H	BBB	501	35/35	0.90	0.19	44,56,81,84	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.