



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 10:22 PM UTC

PDB ID : 7ULC / pdb\_00007ulc  
Title : Crystal structure of queuine salvage enzyme DUF2419 mutant D231N, in complex with queuosine-5'-monophosphate  
Authors : Hung, S.-H.; Swairjo, M.A.  
Deposited on : 2022-04-04  
Resolution : 1.99 Å (reported)

This is a wwPDB X-ray Structure Validation Summary 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

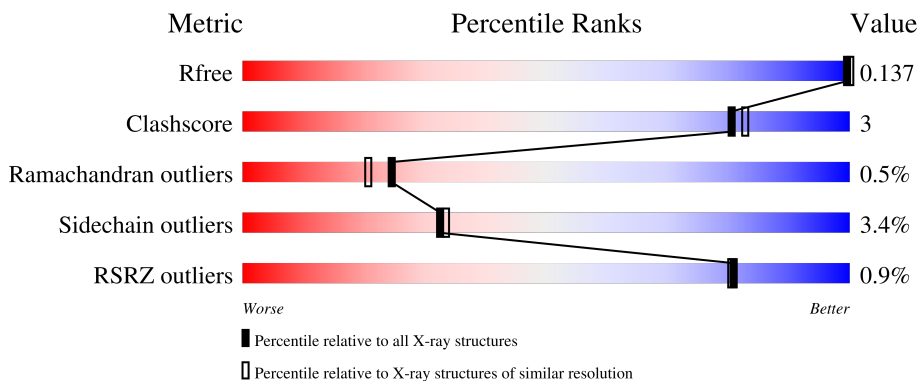
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


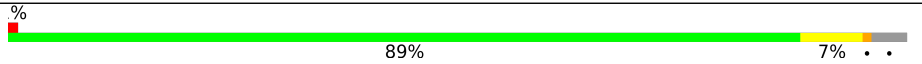
The reported resolution of this entry is 1.99 Å.

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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	331	
1	B	331	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5485 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 Queuosine salvage protein DUF2419.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	315	Total	C	N	O	S	0	4	0
			2534	1612	457	464	1			
1	B	319	Total	C	N	O	S	0	9	0
			2611	1664	469	477	1			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	231	ASN	ASP	engineered mutation	UNP D1C7A6
A	324	LEU	-	expression tag	UNP D1C7A6
A	325	GLU	-	expression tag	UNP D1C7A6
A	326	HIS	-	expression tag	UNP D1C7A6
A	327	HIS	-	expression tag	UNP D1C7A6
A	328	HIS	-	expression tag	UNP D1C7A6
A	329	HIS	-	expression tag	UNP D1C7A6
A	330	HIS	-	expression tag	UNP D1C7A6
A	331	HIS	-	expression tag	UNP D1C7A6
B	231	ASN	ASP	engineered mutation	UNP D1C7A6
B	324	LEU	-	expression tag	UNP D1C7A6
B	325	GLU	-	expression tag	UNP D1C7A6
B	326	HIS	-	expression tag	UNP D1C7A6
B	327	HIS	-	expression tag	UNP D1C7A6
B	328	HIS	-	expression tag	UNP D1C7A6
B	329	HIS	-	expression tag	UNP D1C7A6
B	330	HIS	-	expression tag	UNP D1C7A6
B	331	HIS	-	expression tag	UNP D1C7A6

- Molecule 2 is 2-amino-5-({[(1S,4S,5R)-4,5-dihydroxycyclopent-2-en-1-yl]amino}methyl)-7-(5-O-phosphono-beta-D-ribofuranosyl)-3,7-dihydro-4H-pyrrolo[2,3-d]pyrimidin-4-one (CCD ID: 56B) (formula: C<sub>17</sub>H<sub>24</sub>N<sub>5</sub>O<sub>10</sub>P) (labeled as "Ligand of Interest" by depositor).



- Molecule 4 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	A	115	Total 116	O 116	0	1
4	B	150	Total 151	O 151	0	1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.25Å 106.25Å 157.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.56 – 1.99 47.56 – 1.99	Depositor EDS
% Data completeness (in resolution range)	91.0 (47.56-1.99) 91.0 (47.56-1.99)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.20 (at 2.00Å)	Xtrriage
Refinement program	REFMAC 5.8.0222	Depositor
R, $R_{free}$	0.144 , 0.172 (Not available) , 0.137	Depositor DCC
$R_{free}$ test set	2819 reflections (4.51%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.1	Xtrriage
Anisotropy	0.068	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 49.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5485	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.08% 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: PEG, 56B

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	0.63	0/2596	0.90	0/3539
1	B	0.66	0/2678	0.87	0/3649
All	All	0.65	0/5274	0.88	0/7188

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	3
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	121	ARG	Sidechain
1	A	15	ARG	Sidechain
1	A	200	ARG	Sidechain
1	B	15[A]	ARG	Sidechain
1	B	58	ARG	Sidechain

## 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	2534	0	2466	15	0
1	B	2611	0	2541	14	0
2	A	33	0	0	0	0
2	B	33	0	0	0	0
3	B	7	0	10	0	0
4	A	116	0	0	1	0
4	B	151	0	0	3	0
All	All	5485	0	5017	27	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 3.

The worst 5 of 27 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:15:ARG:NH2	1:B:29[B]:ASP:OD2	2.22	0.66
1:A:233:LYS:NZ	4:A:501:HOH:O	2.27	0.66
1:A:209:TYR:CD1	1:A:223:LEU:HD11	2.40	0.56
1:A:233:LYS:HE3	1:A:236:GLN:OE1	2.05	0.56
1:B:209:TYR:CD1	1:B:223:LEU:HD11	2.41	0.55

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	317/331 (96%)	311 (98%)	4 (1%)	2 (1%)	21	17
1	B	326/331 (98%)	322 (99%)	3 (1%)	1 (0%)	36	35
All	All	643/662 (97%)	633 (98%)	7 (1%)	3 (0%)	24	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	56	ASP
1	A	74	CYS
1	A	319	ARG

### 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	249/259 (96%)	239 (96%)	10 (4%)	28	27
1	B	258/259 (100%)	251 (97%)	7 (3%)	39	42
All	All	507/518 (98%)	490 (97%)	17 (3%)	32	33

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	171	VAL
1	B	319	ARG
1	A	233	LYS
1	A	300	LEU
1	A	320	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	231	ASN
1	B	236	GLN
1	B	240	HIS

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Mol	Chain	Res	Type
1	A	251	HIS
1	B	22	GLN

### 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](#)

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$
2	56B	A	401	-	33,36,36	2.09	4 (12%)	41,55,55	2.03	10 (24%)
2	56B	B	402	-	33,36,36	1.86	2 (6%)	41,55,55	2.05	11 (26%)
3	PEG	B	401	-	6,6,6	0.61	0	5,5,5	0.45	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
2	56B	A	401	-	-	3/15/44/44	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	56B	B	402	-	-	4/15/44/44	0/4/4/4
3	PEG	B	401	-	-	0/4/4/4	-

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	56B	O6-C6	9.85	1.42	1.23
2	B	402	56B	O6-C6	8.79	1.40	1.23
2	A	401	56B	C6-N1	4.28	1.46	1.38
2	B	402	56B	C6-N1	3.87	1.46	1.38
2	A	401	56B	C8-N9	-2.02	1.34	1.38

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	56B	C5-C4-N3	-6.09	120.54	127.53
2	A	401	56B	C5-C4-N3	-5.80	120.87	127.53
2	A	401	56B	C7-C5-C4	-4.90	107.06	112.68
2	B	402	56B	C7-C5-C4	-4.56	107.45	112.68
2	B	402	56B	O4'-C1'-N9	-4.36	98.49	108.36

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

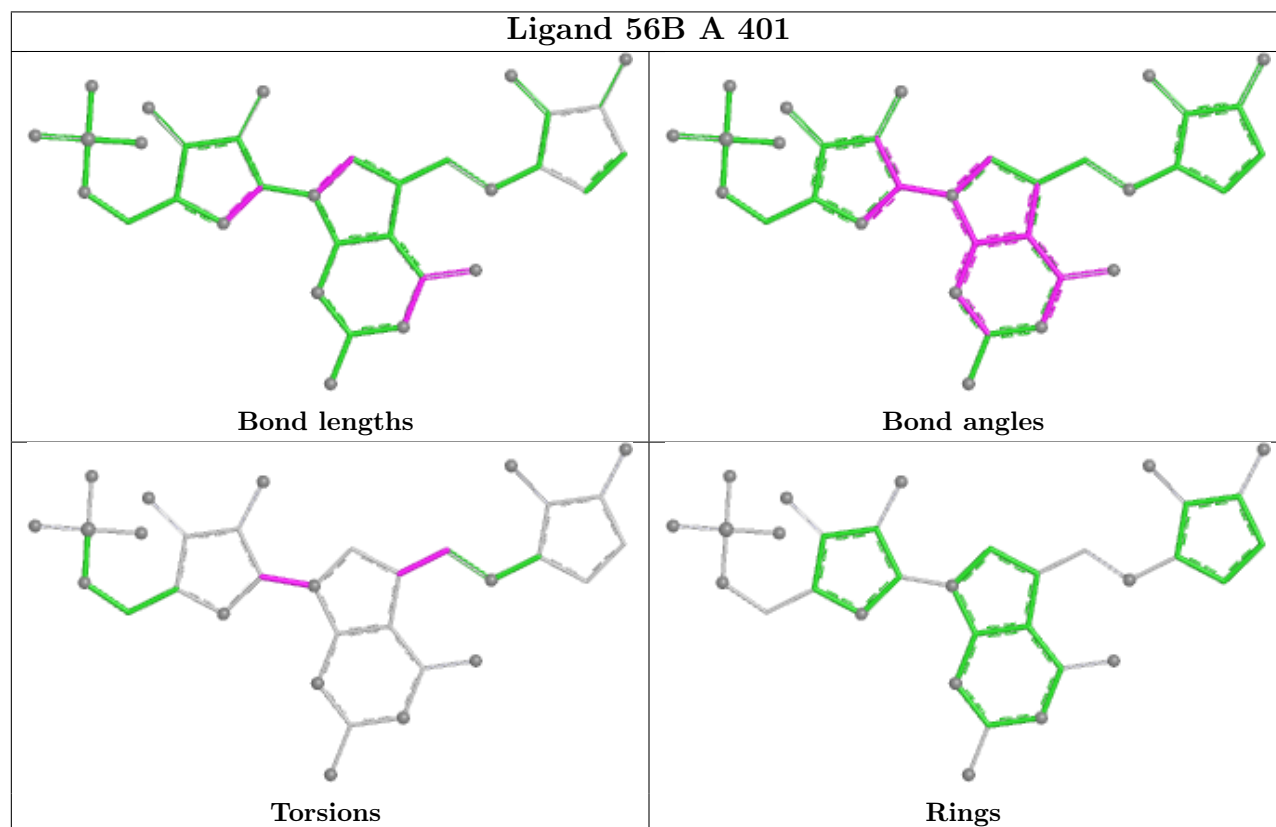
Mol	Chain	Res	Type	Atoms
2	A	401	56B	C5-C7-C9-N10
2	B	402	56B	C5'-O5'-P-OP1
2	A	401	56B	C2'-C1'-N9-C8
2	B	402	56B	C2'-C1'-N9-C8
2	A	401	56B	C8-C7-C9-N10

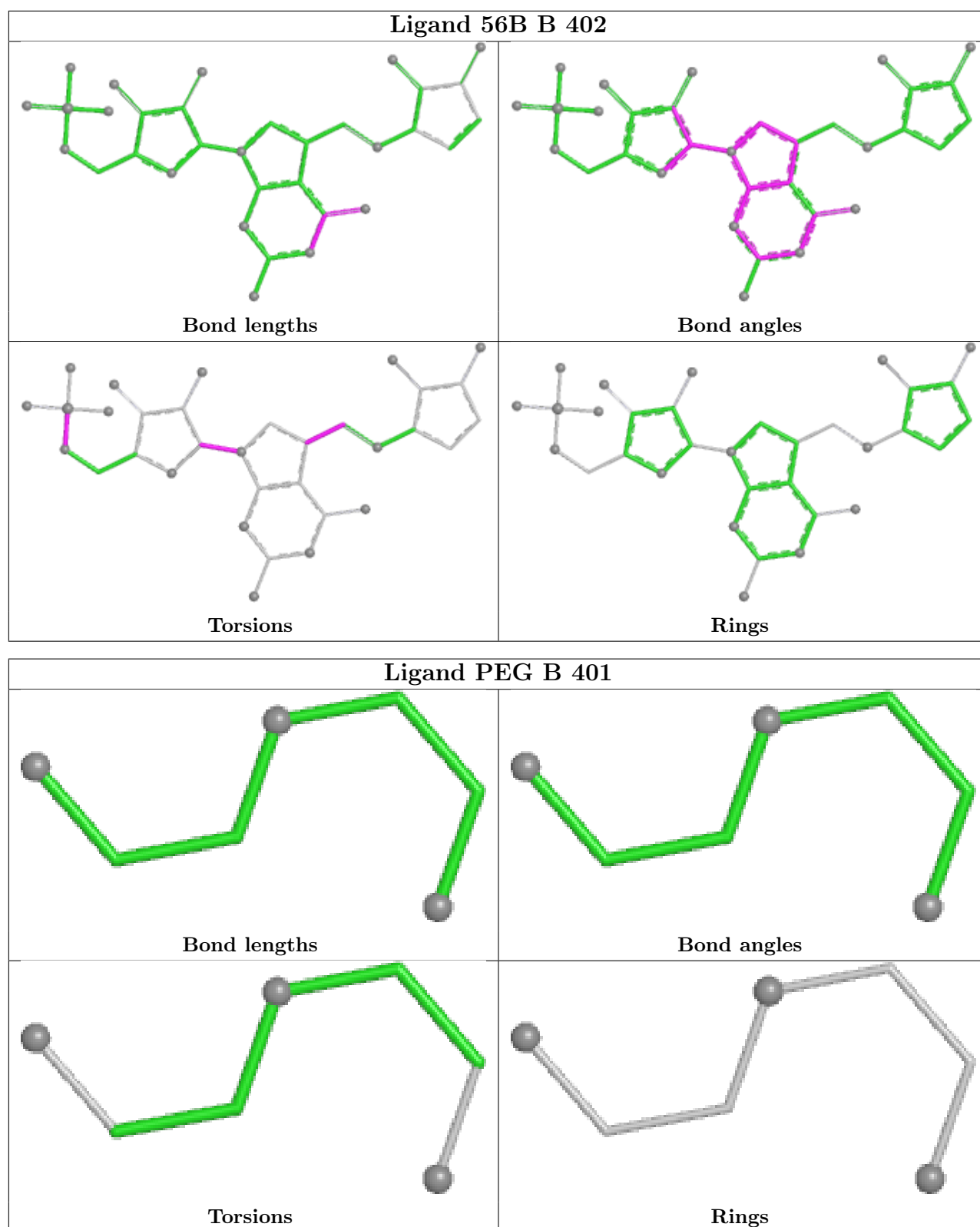
There are no ring outliers.

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

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, 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	315/331 (95%)	-0.06	2 (0%) 85 85	10, 34, 58, 68	4 (1%)
1	B	319/331 (96%)	-0.17	4 (1%) 75 74	9, 29, 50, 79	9 (2%)
All	All	634/662 (95%)	-0.12	6 (0%) 81 80	9, 31, 55, 79	13 (2%)

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	324	LEU	6.7
1	B	321[A]	ILE	5.5
1	B	323	TYR	3.8
1	A	320	THR	3.7
1	A	319	ARG	2.9

### 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](#)

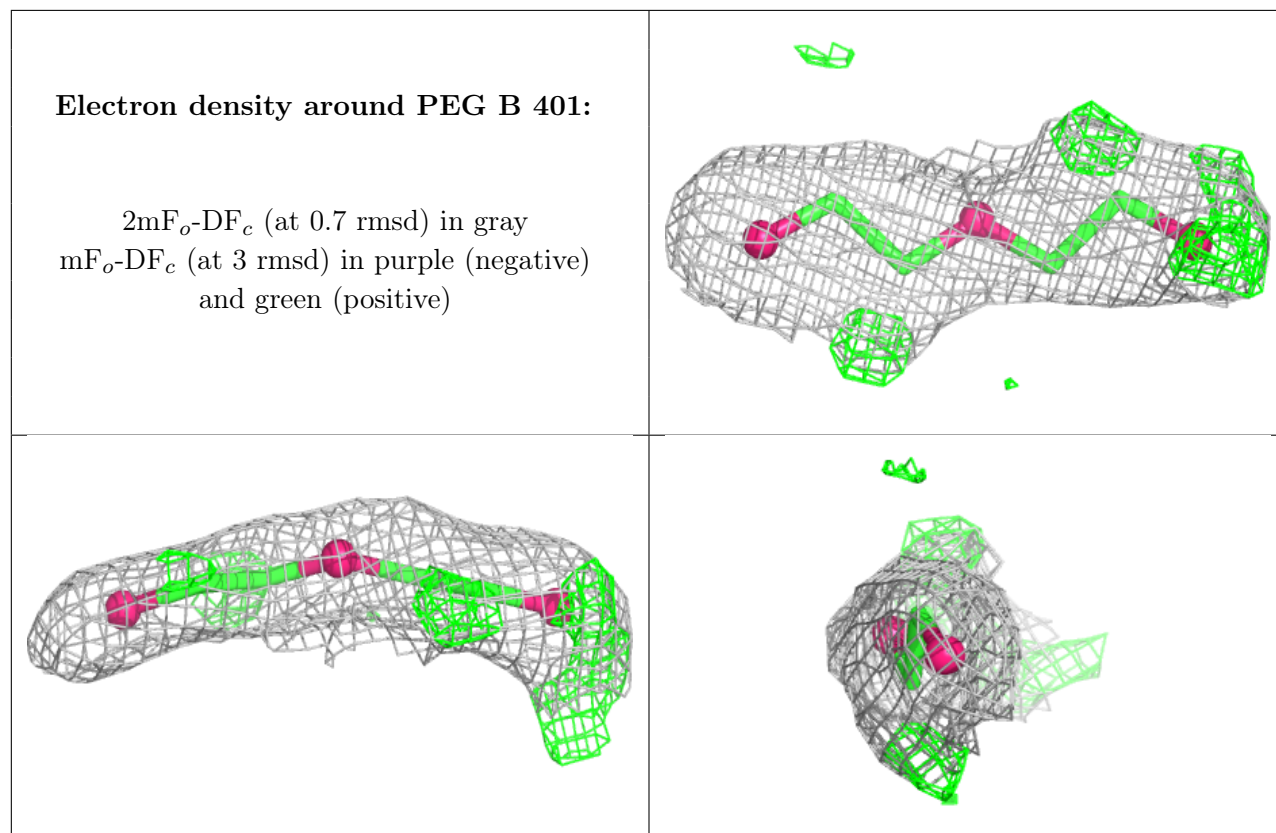
There are no oligosaccharides in this entry.

### 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.

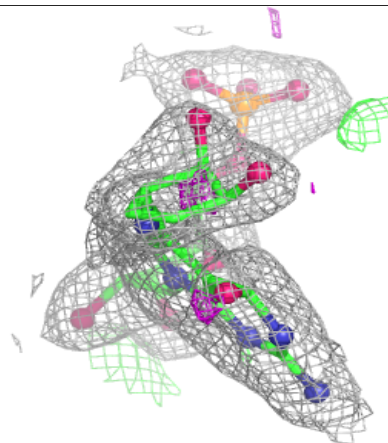
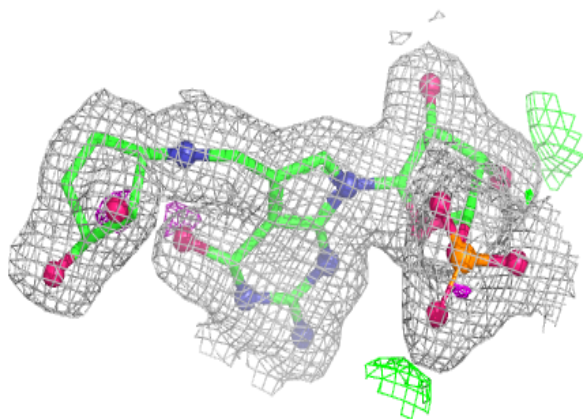
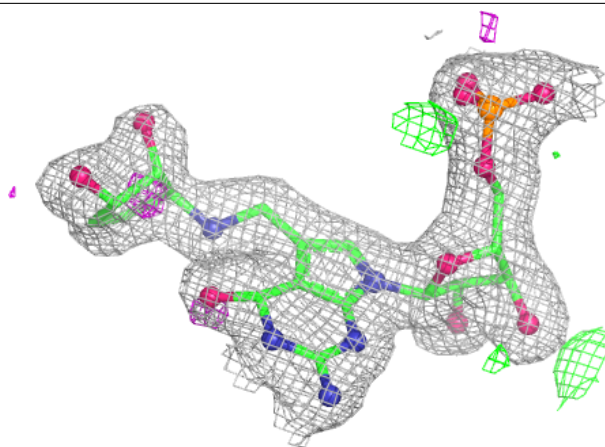
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	PEG	B	401	7/7	0.88	0.15	58,66,72,78	0
2	56B	A	401	33/33	0.97	0.06	24,28,48,54	0
2	56B	B	402	33/33	0.98	0.05	18,23,40,42	0

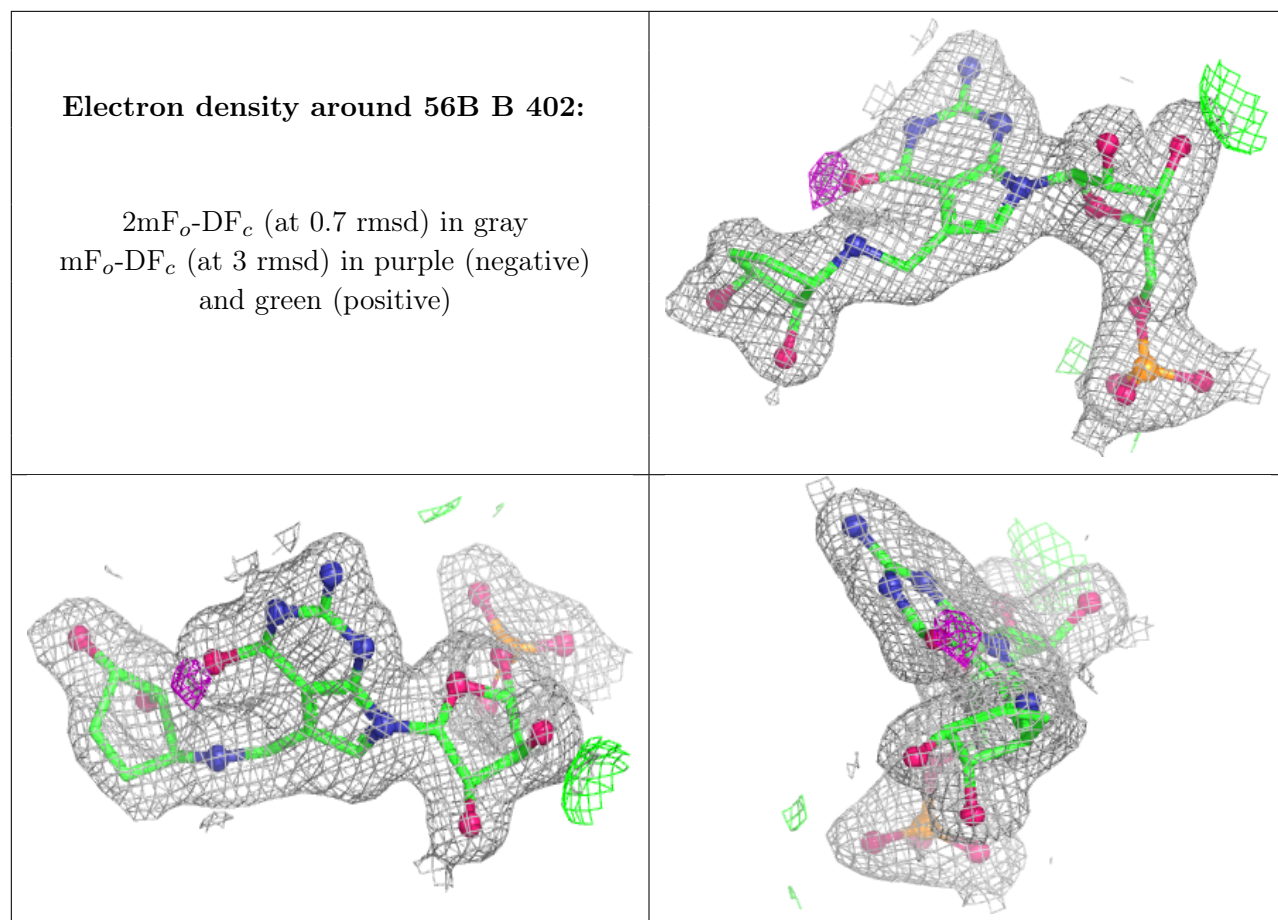
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



**Electron density around 56B A 401:**

$2mF_o-DF_c$  (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.