



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2024 – 11:08 PM EDT

PDB ID : 1LPP  
Title : ANALOGS OF REACTION INTERMEDIATES IDENTIFY A UNIQUE  
SUBSTRATE BINDING SITE IN CANDIDA RUGOSA LIPASE  
Authors : Grochulski, P.G.; Cygler, M.C.  
Deposited on : 1995-01-17  
Resolution : 2.18 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	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

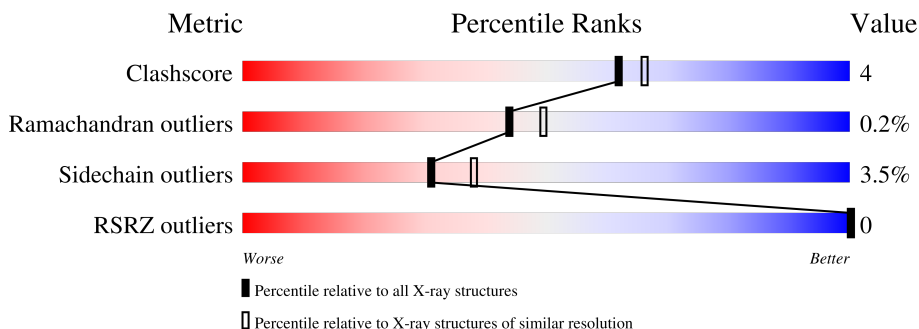
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.18 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))
Clashscore	180529	9404 (2.20-2.16)
Ramachandran outliers	177936	9297 (2.20-2.16)
Sidechain outliers	177891	9297 (2.20-2.16)
RSRZ outliers	164620	8337 (2.20-2.16)

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	549	
2	B	2	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4400 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 LIPASE.

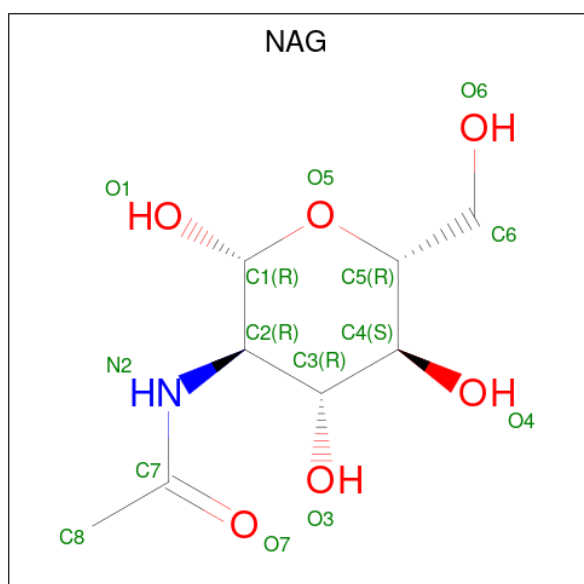
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4022	2556	659	788	19			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

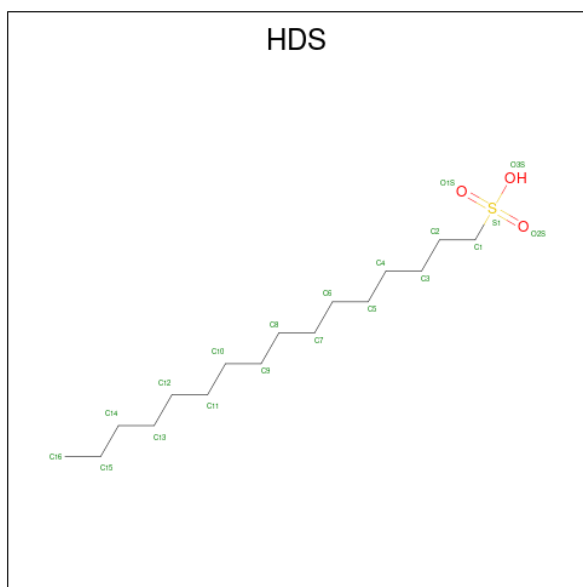


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Ca	0	0
			2	2		

- Molecule 5 is 1-HEXADECANOSULFONIC ACID (three-letter code: HDS) (formula: C<sub>16</sub>H<sub>34</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			19	16	2	1		
5	A	1	Total	C	O	S	0	0
			19	16	2	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	296	Total	O	0	0
			296	296		

i

- Molecule 1: LIPASE

L486	H218	MET
S495	I219	GLU
	L220	ALA
M504	K231	LEU
I505		ALA
L508	R235	LEU
	A236	SER
L525	G237	LEU
	I238	ILE
		ALA
V534	S241	SER
	I253	VAL
	Y254	ALA
	D258	ALA
	D318	A1
	K327	TLI3
	Y328	N20
	I334	G25
	Q338	F38
	R359	K82
	K363	S55
	D371	S59
	G383	L73
	D384	P74
	S389	K85
	P390	E88
	L410	S93
	N424	N101
	Y432	TLI30
	K437	N155
	Q438	S174
	L439	Q182
	H449	R183
	S450	K202
	N451	I205
	S462	F206
	L465	S209
	I466	S214
	N469	

- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

NAG1  
NAG2

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	65.10Å 97.40Å 176.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.18 8.00 – 2.18	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.18) 87.3 (8.00-2.18)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.97 (at 2.05Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.137 , (Not available) 0.137 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.6	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 80.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4400	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% 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: CA, HDS, NAG

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	0.46	0/4118	0.71	1/5601 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	ASP	CB-CG-OD1	5.29	123.07	118.30

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	4022	0	3892	32	0
2	B	28	0	25	0	0
3	A	14	0	13	0	0
4	A	2	0	0	0	0
5	A	38	0	66	0	0
6	A	296	0	0	2	0
All	All	4400	0	3996	32	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 4.

All (32) 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:424:ASN:HD21	1:A:495:SER:H	1.23	0.84
1:A:338:GLN:HE21	1:A:451:ASN:HD21	1.38	0.70
1:A:183:ARG:HH11	1:A:218:HIS:HD1	1.39	0.66
1:A:182:GLN:HE22	1:A:214:SER:HB3	1.63	0.62
1:A:383:GLY:HA3	6:A:889:HOH:O	2.01	0.60
1:A:25:GLY:H	1:A:101:ASN:HD22	1.52	0.57
1:A:25:GLY:H	1:A:101:ASN:ND2	2.04	0.54
1:A:73:LEU:HB3	1:A:74:PRO:HD3	1.89	0.54
1:A:202:LYS:HA	1:A:235:ARG:HD2	1.91	0.52
1:A:130:THR:HG23	1:A:155:ASN:HD22	1.73	0.52
1:A:235:ARG:HD3	6:A:616:HOH:O	2.12	0.50
1:A:13:THR:HB	1:A:52:LYS:HG2	1.93	0.49
1:A:209:SER:HA	1:A:241:SER:O	2.13	0.49
1:A:338:GLN:NE2	1:A:451:ASN:HD21	2.07	0.48
1:A:449:HIS:O	1:A:450:SER:HB2	2.14	0.47
1:A:59:SER:O	1:A:93:SER:HA	2.16	0.46
1:A:206:PHE:HB2	1:A:238:ILE:HB	1.98	0.45
1:A:486:LEU:N	1:A:486:LEU:HD22	2.32	0.44
1:A:363:LYS:HE2	1:A:371:ASP:OD1	2.18	0.44
1:A:38:PHE:O	1:A:174:SER:HB3	2.18	0.43
1:A:465:LEU:O	1:A:469:ASN:HB2	2.18	0.43
1:A:205:ILE:O	1:A:237:GLY:HA2	2.18	0.43
1:A:220:LEU:HD23	1:A:328:TYR:HB3	2.01	0.42
1:A:253:ILE:HG23	1:A:254:TYR:N	2.34	0.42
1:A:389:SER:HA	1:A:390:PRO:C	2.40	0.42
1:A:130:THR:HG23	1:A:155:ASN:ND2	2.34	0.42
1:A:462:SER:O	1:A:465:LEU:HB2	2.19	0.41
1:A:437:LYS:HE3	1:A:504:MET:CE	2.51	0.41
1:A:334:ILE:HG12	1:A:432:TYR:HB2	2.02	0.41
1:A:465:LEU:HD12	1:A:465:LEU:HA	1.92	0.41
1:A:466:ILE:HG13	1:A:505:ILE:HD11	2.02	0.41
1:A:424:ASN:ND2	1:A:495:SER:H	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	532/549 (97%)	510 (96%)	21 (4%)	1 (0%)	44	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	384	ASP

### 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	431/441 (98%)	416 (96%)	15 (4%)	31	38

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	55	SER
1	A	85	LYS
1	A	88	GLU
1	A	231	LYS
1	A	235	ARG
1	A	288	ASP
1	A	318	ASP
1	A	327	LYS
1	A	359	ARG

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Mol	Chain	Res	Type
1	A	410	LEU
1	A	439	LEU
1	A	465	LEU
1	A	508	LEU
1	A	525	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	101	ASN
1	A	137	GLN
1	A	155	ASN
1	A	182	GLN
1	A	187	GLN
1	A	240	GLN
1	A	330	ASN
1	A	338	GLN
1	A	339	ASN
1	A	424	ASN
1	A	456	GLN
1	A	500	ASN
1	A	528	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](#)

2 monosaccharides 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	NAG	B	1	2,1	14,14,15	0.92	0	17,19,21	1.04	0
2	NAG	B	2	2	14,14,15	0.86	0	17,19,21	1.31	3 (17%)

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	NAG	B	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	C4-C3-C2	2.51	114.70	111.02
2	B	2	NAG	C3-C4-C5	-2.14	106.35	110.23
2	B	2	NAG	O3-C3-C4	-2.11	105.40	110.38

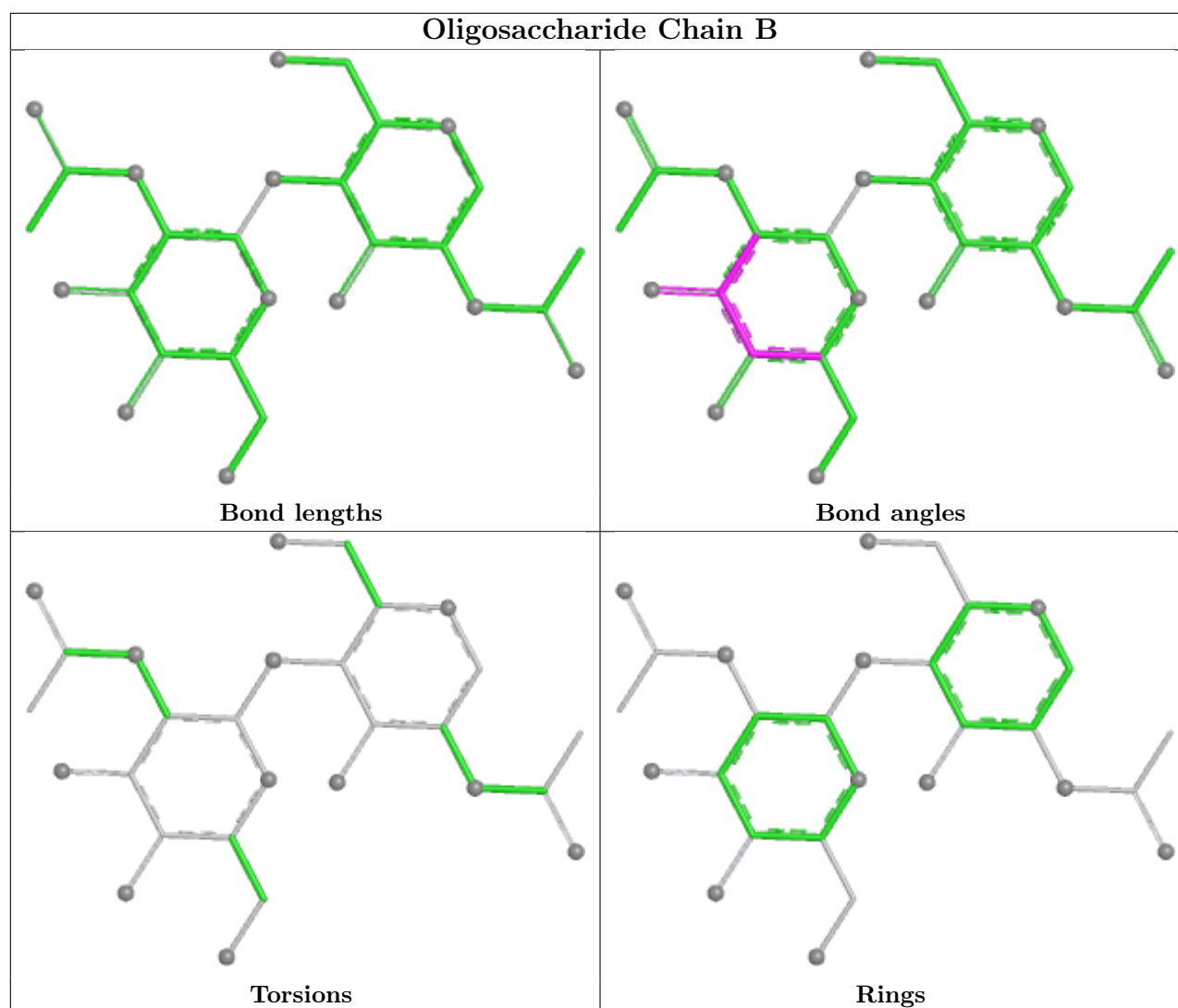
There are no chirality outliers.

There are no torsion outliers.

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 oligosaccharide.



## 5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 2 are monoatomic - leaving 3 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	NAG	A	991	1	14,14,15	1.27	2 (14%)	17,19,21	1.83	5 (29%)
5	HDS	A	560	-	17,18,19	1.88	1 (5%)	15,18,21	0.59	0
5	HDS	A	561	-	17,18,19	2.00	1 (5%)	15,18,21	0.52	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
3	NAG	A	991	1	-	0/6/23/26	0/1/1/1
5	HDS	A	560	-	-	2/15/16/17	-
5	HDS	A	561	-	-	5/15/16/17	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	561	HDS	O1S-S1	-8.11	1.40	1.47
5	A	560	HDS	O1S-S1	-7.67	1.40	1.47
3	A	991	NAG	C3-C2	2.76	1.58	1.52
3	A	991	NAG	C6-C5	2.23	1.59	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	991	NAG	C1-O5-C5	4.75	118.55	112.19
3	A	991	NAG	O7-C7-C8	-2.36	117.86	122.05
3	A	991	NAG	O3-C3-C2	2.33	114.25	109.40
3	A	991	NAG	C1-C2-N2	2.28	114.03	110.43
3	A	991	NAG	O3-C3-C4	-2.27	105.02	110.38

There are no chirality outliers.

All (7) torsion outliers are listed below:

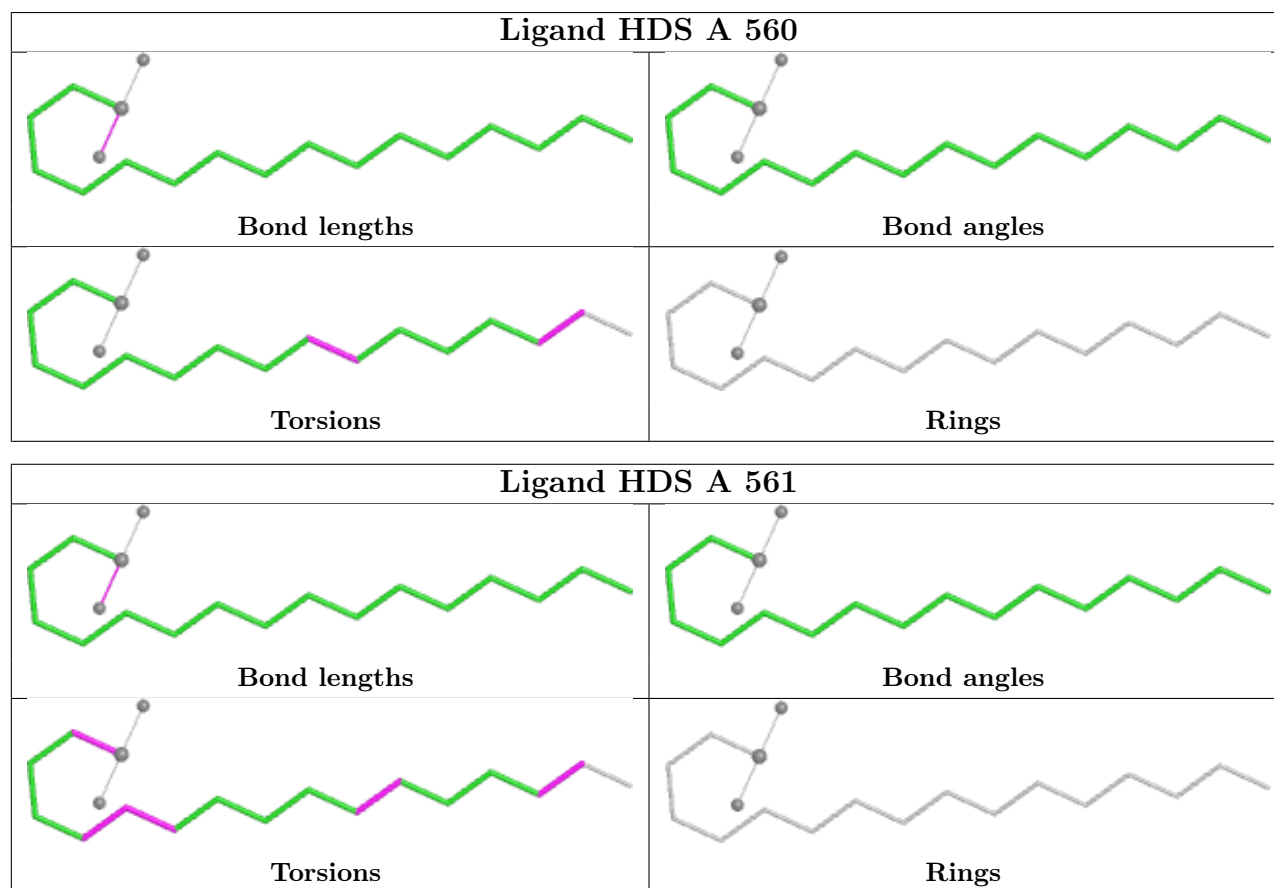
Mol	Chain	Res	Type	Atoms
5	A	561	HDS	C2-C1-S1-O1S
5	A	561	HDS	C13-C14-C15-C16
5	A	561	HDS	C4-C5-C6-C7
5	A	561	HDS	C3-C4-C5-C6
5	A	561	HDS	C9-C10-C11-C12
5	A	560	HDS	C13-C14-C15-C16
5	A	560	HDS	C11-C10-C9-C8

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 [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	534/549 (97%)	-1.11	0 100 100	9, 17, 35, 52	0

There are no RSRZ outliers to report.

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

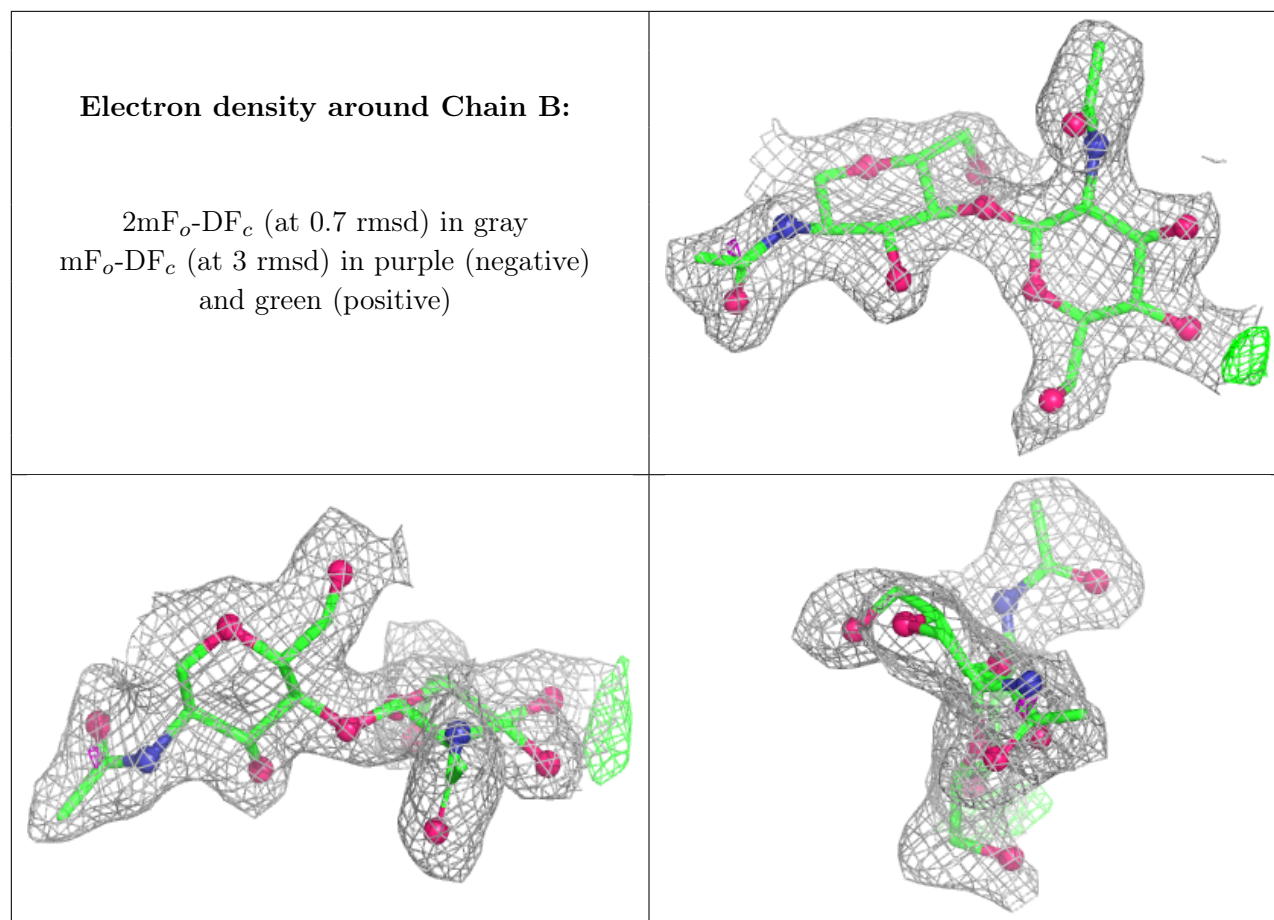
There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

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
2	NAG	B	2	14/15	0.92	0.06	16,23,32,38	14
2	NAG	B	1	14/15	0.94	0.05	21,24,33,36	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



## 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.

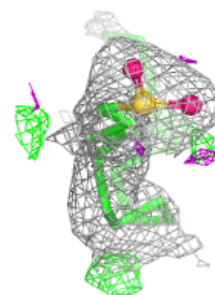
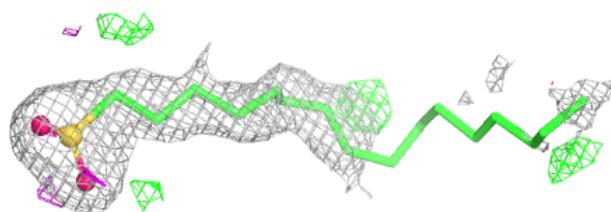
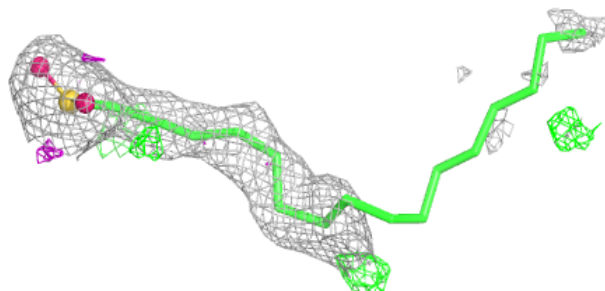
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	A	991	14/15	0.92	0.07	26,31,38,44	0
4	CA	A	589	1/1	0.95	0.04	38,38,38,38	0
4	CA	A	590	1/1	0.97	0.03	12,12,12,12	1
5	HDS	A	561	19/20	0.98	0.08	17,43,48,48	10
5	HDS	A	560	19/20	0.99	0.04	13,19,26,28	0

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.

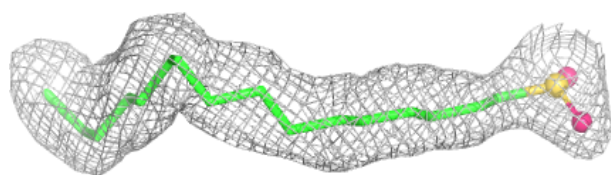
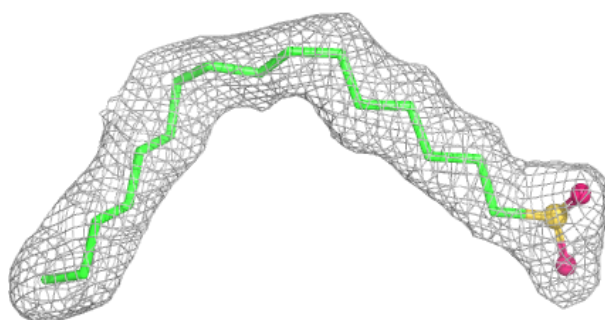


**Electron density around HDS A 561:**

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

**Electron density around HDS A 560:**

$2mF_o-DF_c$  (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.