



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:35 PM UTC

PDB ID : 9BAS / pdb\_00009bas  
Title : Structure of S1\_15A, a lambda-carrageenan specific sulfatase, in complex with galactose-6-sulfate  
Authors : Hettle, J.A.; Vickers, C.; Boraston, A.B.  
Deposited on : 2024-04-04  
Resolution : 1.79 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>.

---

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

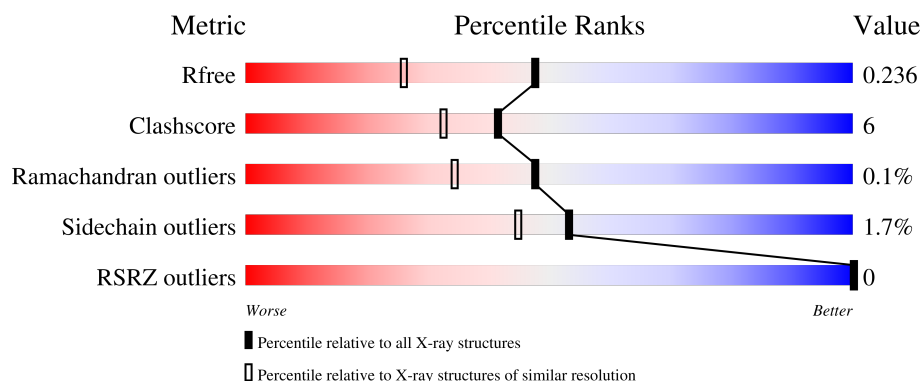
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



The reported resolution of this entry is 1.79 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	523	 80% 11% • 8%
1	B	523	 75% 16% • 8%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7697 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.

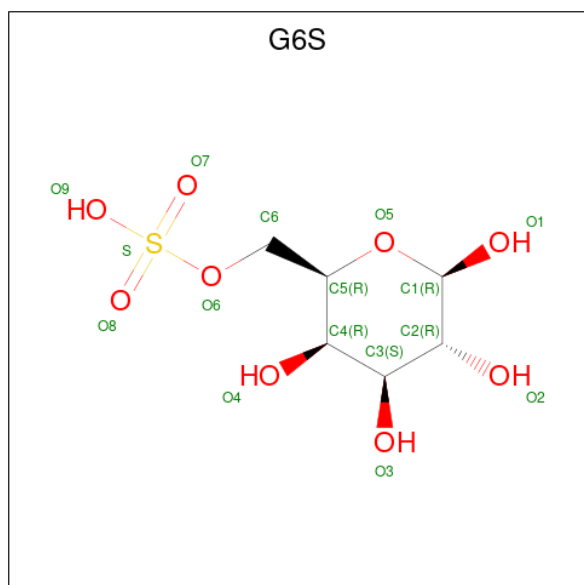
- Molecule 1 is a protein called S1\_15A sulfatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	479	Total	C	N	O	S	5	1	0
			3725	2370	629	713	13			
1	A	479	Total	C	N	O	S	7	3	0
			3719	2365	629	712	13			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is 6-O-sulfo-beta-D-galactopyranose (CCD ID: G6S) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>9</sub>S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	O	S	0	0
			16	6	9	1		
3	A	1	Total	C	O	S	0	0
			16	6	9	1		


- Molecule 4 is water.

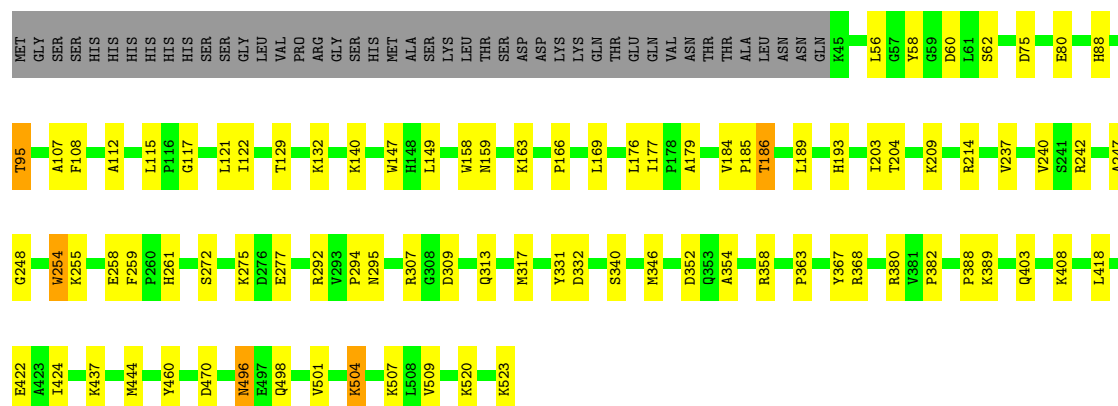
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	108	Total	O	0	0
			108	108		
4	A	111	Total	O	0	0
			111	111		

### 3 Residue-property plots


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.

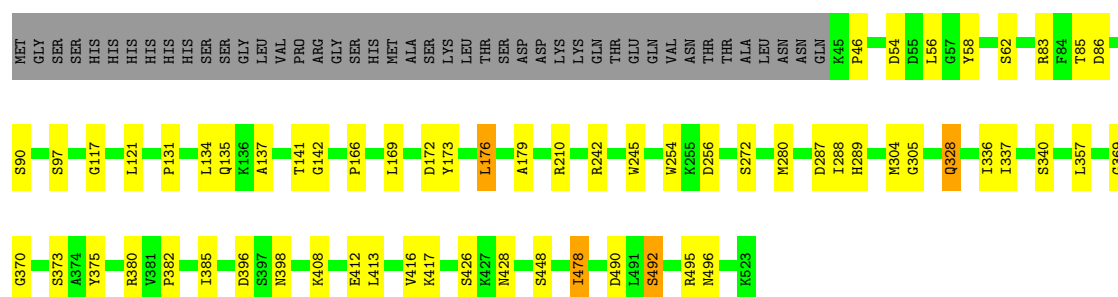
#### • Molecule 1: S1\_15A sulfatase

Chain B:  75% 16% 8%



#### • Molecule 1: S1\_15A sulfatase

Chain A:  80% 11% 8%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.87Å 110.01Å 71.79Å 90.00° 106.25° 90.00°	Depositor
Resolution (Å)	29.54 – 1.79 29.54 – 1.79	Depositor EDS
% Data completeness (in resolution range)	97.3 (29.54-1.79) 97.0 (29.54-1.79)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.16 (at 1.79Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.187 , 0.236 0.187 , 0.236	Depositor DCC
$R_{free}$ test set	4880 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.8	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 21.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.459 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7697	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.29% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, G6S

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	0.53	0/3819	0.85	3/5180 (0.1%)
1	B	0.55	0/3819	0.84	4/5177 (0.1%)
All	All	0.54	0/7638	0.84	7/10357 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	177	ILE	CA-C-N	5.86	127.17	119.84
1	B	177	ILE	C-N-CA	5.86	127.17	119.84
1	B	95	THR	CA-C-N	-5.34	113.53	119.19
1	B	95	THR	C-N-CA	-5.34	113.53	119.19
1	A	305	GLY	CA-C-N	5.09	125.12	119.32
1	A	305	GLY	C-N-CA	5.09	125.12	119.32
1	A	176	LEU	N-CA-C	5.01	117.42	109.50

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	3719	0	3625	34	0
1	B	3725	0	3646	60	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	16	0	11	0	0
3	B	16	0	12	0	0
4	A	111	0	0	1	0
4	B	108	0	0	2	0
All	All	7697	0	7294	94	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 6.

All (94) 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:B:186:THR:HB	1:B:254:TRP:CZ3	2.21	0.76
1:B:166:PRO:HG3	1:B:176:LEU:HD11	1.74	0.69
1:B:254:TRP:HD1	1:B:259:PHE:CZ	2.10	0.68
1:B:254:TRP:CD1	1:B:259:PHE:CZ	2.81	0.68
1:B:107:ALA:HB1	1:B:112:ALA:HB3	1.74	0.67
1:B:62:SER:OG	1:B:75:ASP:OD2	2.13	0.65
1:A:496:ASN:O	1:A:496:ASN:OD1	2.15	0.65
1:A:490:ASP:OD1	1:A:492:SER:HB3	1.98	0.64
1:A:412:GLU:OE2	1:A:428:ASN:ND2	2.25	0.63
1:B:254:TRP:H	1:B:254:TRP:HE3	1.46	0.62
1:B:346:MET:HE3	1:B:354:ALA:HB1	1.81	0.61
1:B:254:TRP:HD1	1:B:259:PHE:CE1	2.17	0.61
1:B:496:ASN:HD22	1:B:498:GLN:HG2	1.65	0.61
1:A:166:PRO:HG3	1:A:176:LEU:HD11	1.82	0.61
1:B:240:VAL:HB	1:B:242:ARG:NH2	2.17	0.60
1:B:80:GLU:OE2	4:B:701:HOH:O	2.15	0.60
1:B:242:ARG:CZ	1:B:254:TRP:HZ2	2.15	0.59
1:B:58:TYR:HA	1:B:382:PRO:HD3	1.83	0.59
1:B:332:ASP:HA	1:B:388:PRO:HG2	1.84	0.59
1:B:272:SER:O	1:B:275:LYS:HG2	2.04	0.58
1:A:117:GLY:HA2	1:A:179:ALA:HB2	1.85	0.57
1:B:159:ASN:OD1	1:B:203:ILE:N	2.36	0.56
1:B:367:TYR:OH	1:B:380:ARG:NH1	2.40	0.55
1:A:272:SER:HA	1:A:328:GLN:HE22	1.72	0.55
1:B:80:GLU:OE1	1:B:331:TYR:OH	2.23	0.54
1:B:294:PRO:HG3	1:B:309:ASP:HB2	1.90	0.54
1:B:460:TYR:HE1	1:B:509:VAL:HG13	1.73	0.54

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:242:ARG:CZ	1:B:254:TRP:CZ2	2.91	0.53
1:B:368:ARG:HB2	1:B:496:ASN:O	2.08	0.53
1:B:255:LYS:HG2	1:B:258:GLU:HG2	1.90	0.53
1:B:60:ASP:OD2	1:B:307:ARG:NH2	2.42	0.53
1:B:408:LYS:HG2	1:B:418:LEU:HD12	1.91	0.53
1:A:495:ARG:O	4:A:701:HOH:O	2.18	0.52
1:B:108:PHE:CE1	1:B:520:LYS:HG2	2.44	0.52
1:A:121:LEU:HD11	1:A:169:LEU:HD11	1.91	0.52
1:B:122:ILE:HG12	1:B:149:LEU:O	2.11	0.51
1:B:88:HIS:CD2	1:B:444:MET:HE1	2.46	0.51
1:A:56:LEU:HB3	1:A:340:SER:HB2	1.93	0.50
1:A:370:GLY:N	1:A:478:ILE:HD11	2.27	0.49
1:B:117:GLY:HA2	1:B:179:ALA:HB2	1.94	0.49
1:B:147:TRP:HB2	1:B:176:LEU:HD13	1.93	0.49
1:B:88:HIS:O	1:B:403:GLN:HG3	2.13	0.48
1:B:121:LEU:HD11	1:B:169:LEU:HD11	1.94	0.48
1:B:496:ASN:HD22	1:B:498:GLN:CG	2.25	0.48
1:B:214:ARG:NH2	1:B:248:GLY:O	2.39	0.48
1:B:209:LYS:HE2	1:B:209:LYS:HB2	1.63	0.47
1:A:370:GLY:H	1:A:373:SER:HB3	1.79	0.47
1:B:460:TYR:CE1	1:B:509:VAL:HG13	2.49	0.47
1:B:129:THR:HG23	1:B:132:LYS:HD2	1.95	0.47
1:A:86:ASP:HB2	1:A:398:ASN:OD1	2.14	0.47
1:A:134:LEU:HD12	1:A:280:MET:HE3	1.96	0.47
1:A:142:GLY:HA2	1:A:173:TYR:O	2.15	0.47
1:B:292:ARG:HD3	1:B:307:ARG:HA	1.98	0.46
1:B:261:HIS:HE1	1:B:295:ASN:ND2	2.13	0.46
1:A:242:ARG:NH1	1:A:254:TRP:HH2	2.13	0.46
1:B:158:TRP:CZ2	1:B:185:PRO:HG3	2.52	0.45
1:A:85:THR:O	1:A:380:ARG:NH1	2.48	0.45
1:A:254:TRP:HZ3	1:A:256:ASP:OD1	2.00	0.45
1:B:189:LEU:HD12	1:B:193:HIS:C	2.42	0.44
1:B:418:LEU:HD22	1:B:422:GLU:CD	2.43	0.44
1:B:292:ARG:NH1	1:B:352:ASP:OD2	2.50	0.44
1:A:369:GLY:HA3	1:A:375:TYR:HB2	1.99	0.44
1:B:237:VAL:HB	1:B:254:TRP:CZ3	2.53	0.44
1:A:134:LEU:HD23	1:A:134:LEU:HA	1.85	0.44
1:B:470:ASP:N	1:B:470:ASP:OD1	2.49	0.44
1:A:58:TYR:HA	1:A:382:PRO:HD3	2.00	0.44
1:B:204:THR:OG1	1:B:247:ALA:HB3	2.18	0.43
1:B:358:ARG:HH21	1:B:363:PRO:HD2	1.84	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:ILE:O	1:A:385:ILE:HA	2.18	0.43
1:B:140:LYS:HA	1:B:140:LYS:HD2	1.85	0.43
1:B:389:LYS:HB3	1:B:389:LYS:HE3	1.64	0.43
1:B:437:LYS:HB3	1:B:437:LYS:HE2	1.56	0.43
1:A:141:THR:O	1:A:172:ASP:HB2	2.19	0.43
1:B:129:THR:H	1:B:132:LYS:HB2	1.83	0.43
1:B:307:ARG:HD3	4:B:778:HOH:O	2.18	0.42
1:B:88:HIS:CG	1:B:444:MET:HE1	2.55	0.42
1:A:90:SER:OG	1:A:97:SER:HA	2.19	0.42
1:A:54:ASP:CG	1:A:289:HIS:HB2	2.45	0.42
1:B:95:THR:HG21	1:B:115:LEU:HG	2.02	0.42
1:A:337:ILE:HG12	1:A:385:ILE:HG12	2.02	0.42
1:A:131:PRO:O	1:A:135:GLN:HG3	2.18	0.41
1:A:137:ALA:HB2	1:A:416:VAL:CG2	2.50	0.41
1:A:408:LYS:HD2	1:A:426:SER:O	2.20	0.41
1:A:287:ASP:HA	1:A:288:ILE:HA	1.65	0.41
1:A:304:MET:HA	1:A:357:LEU:O	2.20	0.41
1:B:501:VAL:HG12	1:B:504:LYS:HD2	2.02	0.41
1:A:46:PRO:HG2	1:A:413:LEU:HD21	2.03	0.41
1:A:83:ARG:HB3	1:A:396:ASP:OD1	2.20	0.41
1:B:163:LYS:HB2	1:B:163:LYS:HE3	1.80	0.41
1:B:313:GLN:O	1:B:317:MET:HG3	2.20	0.41
1:B:184:VAL:HA	1:B:186:THR:HG22	2.03	0.41
1:B:56:LEU:HB3	1:B:340:SER:HB2	2.03	0.40
1:A:210:ARG:HA	1:A:245:TRP:CG	2.57	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	480/523 (92%)	455 (95%)	24 (5%)	1 (0%)	43	31

Continued on next page...

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	478/523 (91%)	454 (95%)	24 (5%)	0	100	100
All	All	958/1046 (92%)	909 (95%)	48 (5%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	448	SER

### 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	394/438 (90%)	389 (99%)	5 (1%)	61	54
1	B	397/438 (91%)	389 (98%)	8 (2%)	48	38
All	All	791/876 (90%)	778 (98%)	13 (2%)	53	47

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

Mol	Chain	Res	Type
1	B	186	THR
1	B	254	TRP
1	B	277	GLU
1	B	424	ILE
1	B	496	ASN
1	B	504	LYS
1	B	507	LYS
1	B	523	LYS
1	A	62	SER
1	A	328	GLN
1	A	417	LYS
1	A	478	ILE
1	A	492	SER

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

Mol	Chain	Res	Type
1	B	70	GLN
1	B	261	HIS
1	B	328	GLN
1	B	496	ASN
1	B	513	GLN
1	A	105	GLN
1	A	320	GLN
1	A	328	GLN
1	A	465	ASN
1	A	487	GLN
1	A	513	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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	G6S	A	602	2	16,16,16	1.20	0	23,24,24	1.27	2 (8%)
3	G6S	B	602	2	16,16,16	1.24	0	23,24,24	1.08	2 (8%)

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	G6S	A	602	2	-	3/6/26/26	0/1/1/1
3	G6S	B	602	2	-	2/6/26/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	602	G6S	C1-C2-C3	2.63	115.72	110.36
3	B	602	G6S	C6-C5-C4	-2.50	106.83	112.07
3	A	602	G6S	O6-C6-C5	-2.16	103.71	107.57
3	B	602	G6S	O9-S-O6	2.12	111.24	106.37

There are no chirality outliers.

All (5) torsion outliers are listed below:

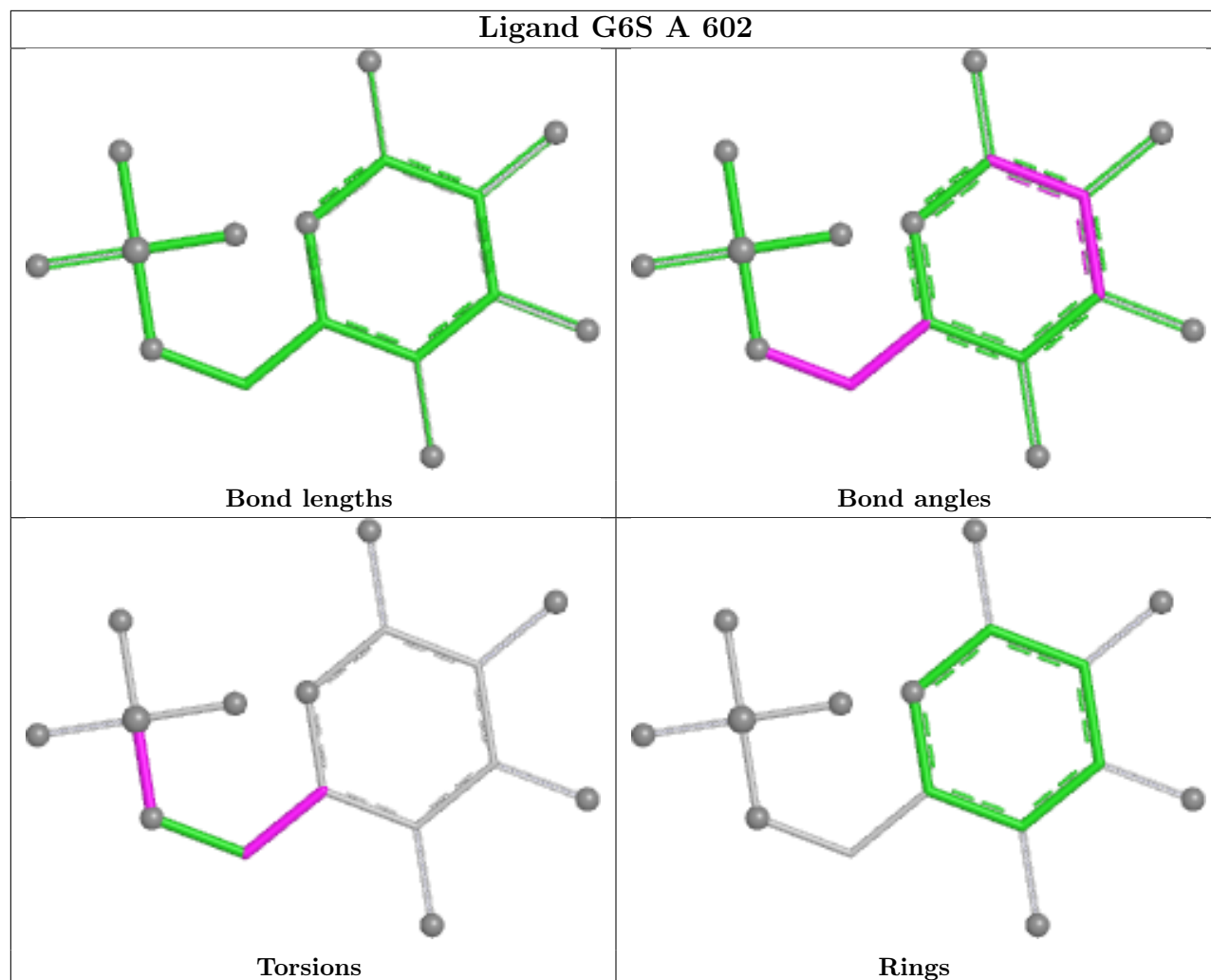
Mol	Chain	Res	Type	Atoms
3	B	602	G6S	C4-C5-C6-O6
3	B	602	G6S	O5-C5-C6-O6
3	A	602	G6S	C6-O6-S-O7
3	A	602	G6S	O5-C5-C6-O6
3	A	602	G6S	C4-C5-C6-O6

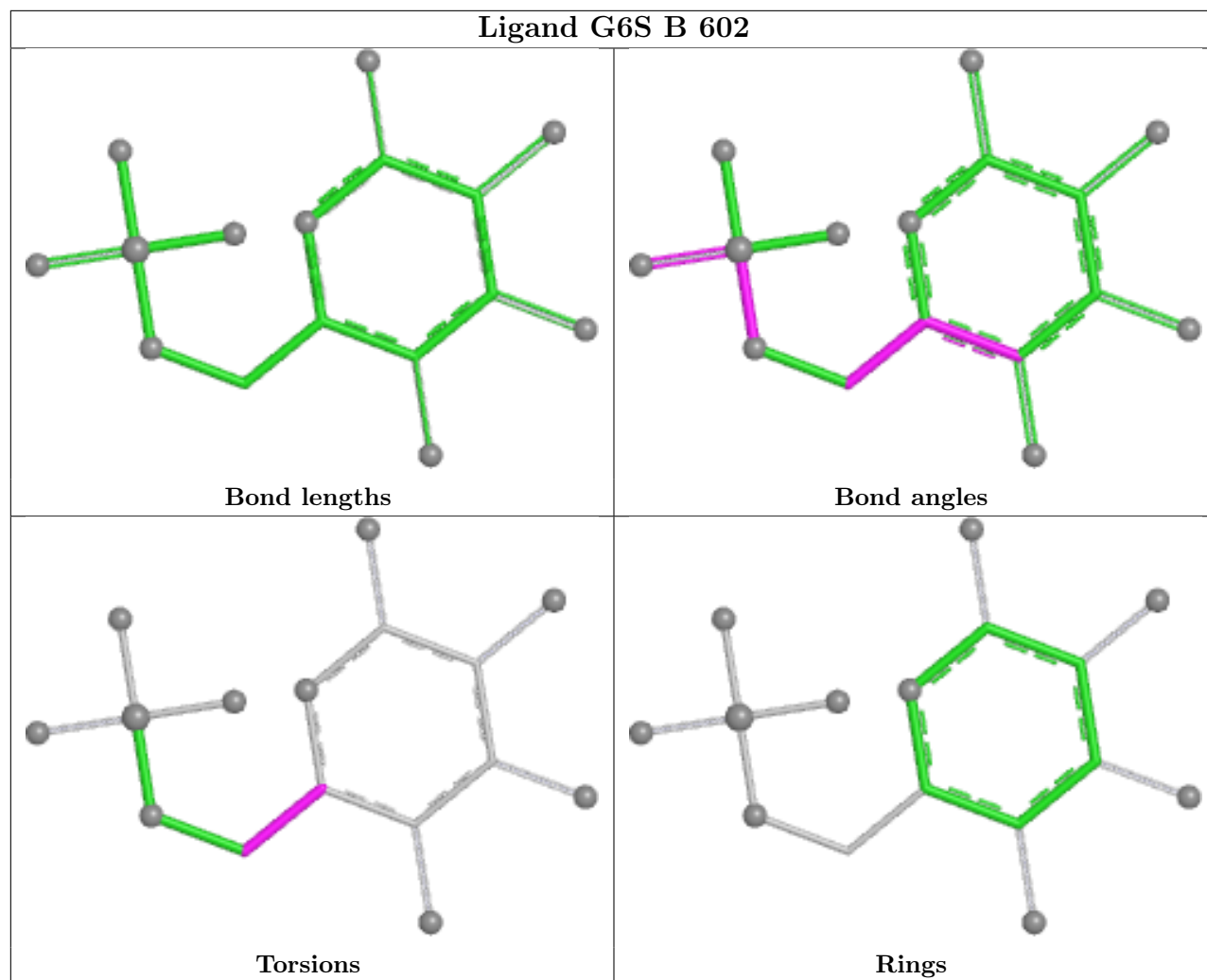
There are no ring outliers.

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.

## Ligand G6S A 602





## 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	479/523 (91%)	-0.75	0 100 100	9, 17, 26, 34	4 (0%)
1	B	479/523 (91%)	-0.78	0 100 100	11, 17, 24, 34	3 (0%)
All	All	958/1046 (91%)	-0.77	0 100 100	9, 17, 25, 34	7 (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](#)

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(Å <sup>2</sup> )	Q<0.9
3	G6S	B	602	16/16	0.98	0.05	13,15,18,20	0
2	CA	A	601	1/1	0.99	0.15	27,27,27,27	0
2	CA	B	601	1/1	0.99	0.11	27,27,27,27	0
3	G6S	A	602	16/16	0.99	0.03	12,14,17,19	0

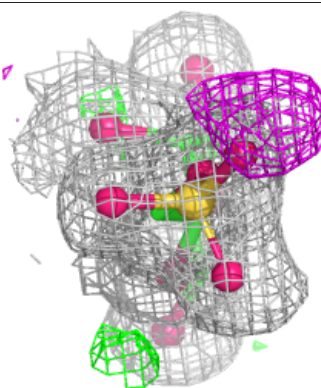
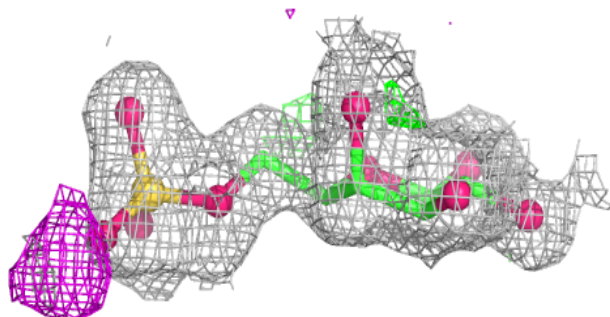
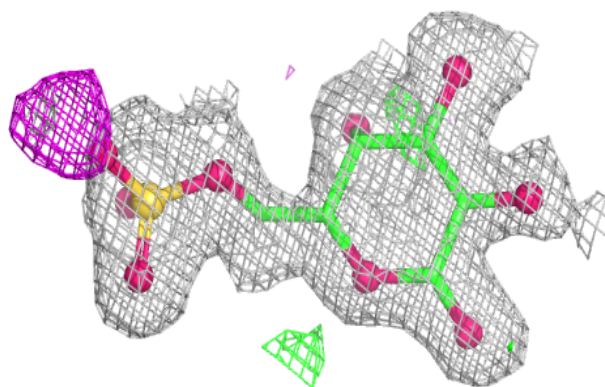
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers



as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

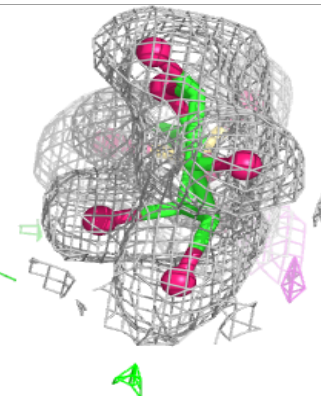
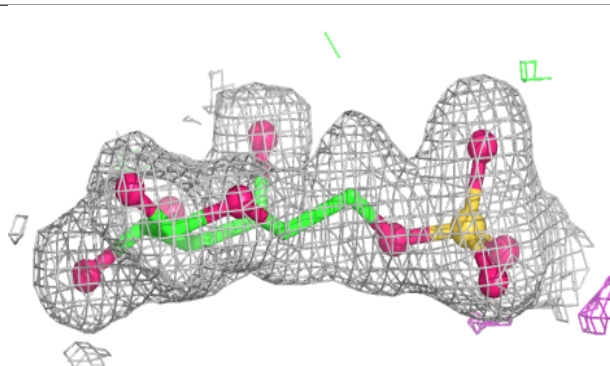
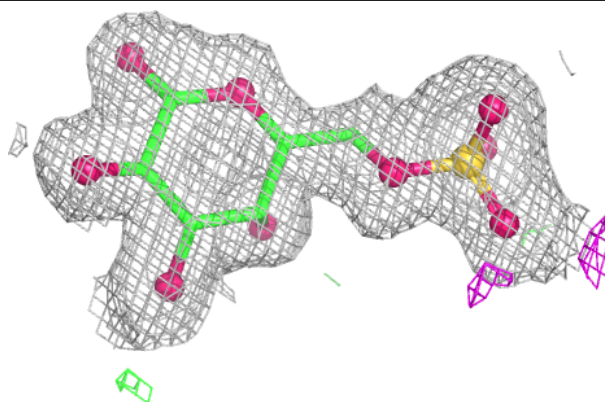
**Electron density around G6S B 602:**

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



**Electron density around G6S A 602:**

$2mF_o-DF_c$  (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.