



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 03:06 AM BST

PDB ID : 5G6U  
Title : Crystal structure of langerin carbohydrate recognition domain with GlcNS6S  
Authors : Porkolab, V.; Chabrol, E.; Varga, N.; Ordanini, S.; Sutkeviciute, I.; Thepaut, M.; Bernardi, A.; Fieschi, F.  
Deposited on : 2016-07-21  
Resolution : 1.84 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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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) : 2.13.1

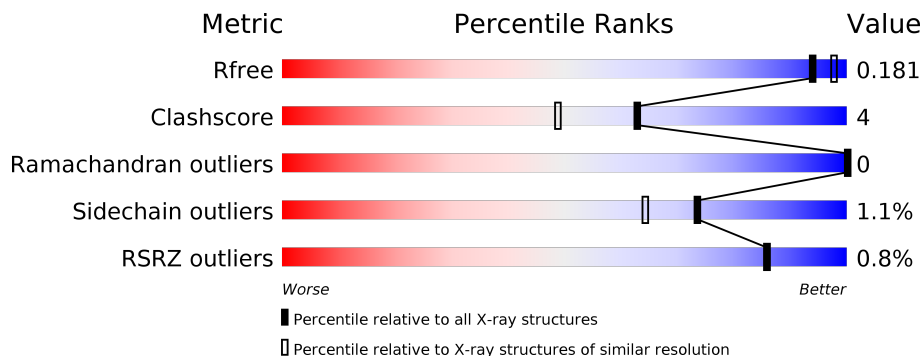
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.84 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	4003 (1.86-1.82)
Clashscore	141614	4233 (1.86-1.82)
Ramachandran outliers	138981	4185 (1.86-1.82)
Sidechain outliers	138945	4186 (1.86-1.82)
RSRZ outliers	127900	3957 (1.86-1.82)

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 on 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	261	
1	B	261	
1	C	261	
1	D	261	

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	CL	B	1330	-	-	X	-
6	CL	D	1327	-	-	X	-
6	CL	D	1328	-	-	X	-

## 2 Entry composition [i](#)

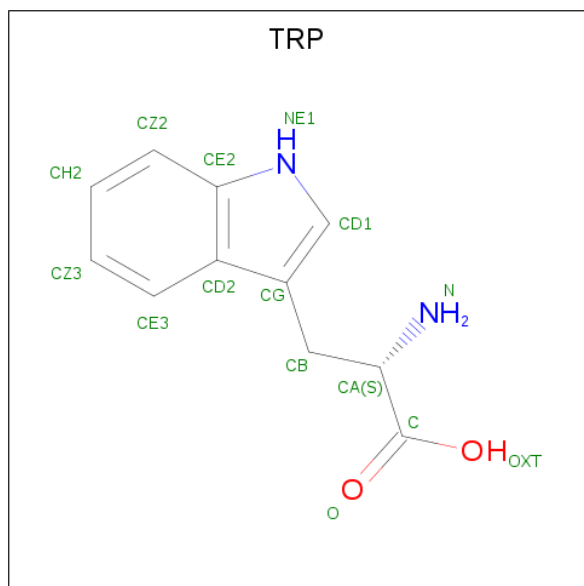
There are 7 unique types of molecules in this entry. The entry contains 8955 atoms, of which 4023 are hydrogens and 0 are deuteriums.

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 LANGERIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	131	Total 2098	C 702	H 1015	N 179	O 197	S 5	0	4	0
1	B	128	Total 2050	C 689	H 990	N 173	O 193	S 5	0	4	0
1	C	128	Total 2054	C 691	H 993	N 173	O 192	S 5	0	3	0
1	D	127	Total 2046	C 686	H 989	N 175	O 191	S 5	0	4	0

- Molecule 2 is TRYPTOPHAN (three-letter code: TRP) (formula: C<sub>11</sub>H<sub>12</sub>N<sub>2</sub>O<sub>2</sub>).



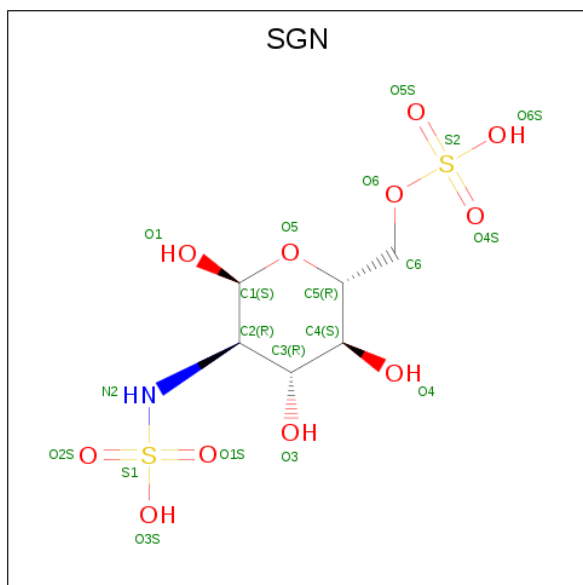
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	H	N	O		
2	A	1	Total 27	C 11	H 12	N 2	O 2	0	0
2	B	1	Total 27	C 11	H 12	N 2	O 2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	H	N			O
2	D	1	27	11	12	2	2	0	0

- Molecule 3 is 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose (three-letter code: SGN) (formula: C<sub>6</sub>H<sub>13</sub>NO<sub>11</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	20	6	1	11	2	0	0
3	B	1	20	6	1	11	2	0	0
3	C	1	20	6	1	11	2	0	0
3	D	1	20	6	1	11	2	0	0

- Molecule 4 is EUROPIUM ION (three-letter code: EU) (formula: Eu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Eu 2 2	0	1
4	A	2	Total Eu 2 2	0	1
4	C	2	Total Eu 2 2	0	1

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total Ca 1 1	0	1
5	A	1	Total Ca 1 1	0	1
5	D	1	Total Ca 1 1	0	0
5	C	1	Total Ca 1 1	0	1

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	5	Total Cl 5 5	0	0
6	A	1	Total Cl 1 1	0	0
6	D	3	Total Cl 3 3	0	0
6	C	3	Total Cl 3 3	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	155	Total O 155 155	0	0
7	B	138	Total O 138 138	0	0
7	C	126	Total O 126 126	0	0
7	D	105	Total O 105 105	0	0







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	79.48Å 79.48Å 90.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.40 – 1.84 47.78 – 1.84	Depositor EDS
% Data completeness (in resolution range)	99.6 (36.40-1.84) 99.6 (47.78-1.84)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.46 (at 1.84Å)	Xtrriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, $R_{free}$	0.152 , 0.180 0.153 , 0.181	Depositor DCC
$R_{free}$ test set	2442 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	24.1	Xtrriage
Anisotropy	0.075	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 51.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.043 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8955	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.35% 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: EU, CA, CL, SGN

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	0.32	0/1142	0.52	0/1553
1	B	0.31	0/1116	0.52	0/1520
1	C	0.31	0/1105	0.52	0/1506
1	D	0.32	0/1112	0.50	0/1514
All	All	0.32	0/4475	0.51	0/6093

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	A	1083	1015	987	9	0
1	B	1060	990	974	6	0
1	C	1061	993	987	5	0
1	D	1057	989	973	7	0
2	A	15	12	9	0	0
2	B	15	12	9	2	0
2	D	15	12	9	0	0
3	A	20	0	7	1	0
3	B	20	0	7	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	20	0	7	1	0
3	D	20	0	7	1	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	1	0	0	0	0
6	B	5	0	0	2	0
6	C	3	0	0	2	0
6	D	3	0	0	4	0
7	A	155	0	0	7	0
7	B	138	0	0	7	1
7	C	126	0	0	5	0
7	D	105	0	0	2	1
All	All	4932	4023	3976	36	1

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 (36) 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)
6:C:1330:CL:CL	7:C:2009:HOH:O	2.10	1.06
7:B:2051:HOH:O	1:D:221[A]:GLN:NE2	1.91	1.04
1:B:236:GLU:OE1	7:B:2039:HOH:O	1.86	0.91
1:A:247:GLY:O	7:A:2052:HOH:O	1.92	0.88
1:C:198:GLY:N	7:C:2001:HOH:O	2.14	0.80
1:C:304[A]:GLN:NE2	7:C:2109:HOH:O	2.15	0.78
2:B:1334:TRP:N	7:B:2106:HOH:O	2.19	0.74
1:A:236:GLU:OE1	7:A:2050:HOH:O	2.10	0.70
1:B:263:ASP:OD2	7:B:2058:HOH:O	2.12	0.68
1:A:226:ARG:NH2	7:A:2004:HOH:O	2.26	0.64
1:D:278:VAL:HG13	6:D:1327:CL:CL	2.36	0.62
1:A:226:ARG:NE	7:A:2004:HOH:O	2.31	0.59
3:C:1326:SGN:N2	7:C:2120:HOH:O	2.32	0.55
1:A:200:LYS:HD2	7:A:2079:HOH:O	2.06	0.54
1:B:198:GLY:N	7:B:2001:HOH:O	2.41	0.54
6:B:1330:CL:CL	6:B:1332:CL:CL	3.02	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:278:VAL:HG12	6:D:1328:CL:CL	2.50	0.49
6:D:1327:CL:CL	6:D:1328:CL:CL	3.05	0.48
1:A:326:SER:HB2	7:A:2149:HOH:O	2.13	0.48
1:D:285:GLU:OE2	3:D:1325:SGN:O3	2.33	0.46
1:D:263:ASP:OD2	7:D:2052:HOH:O	2.21	0.45
1:A:274:LYS:O	1:A:278[B]:VAL:HG13	2.17	0.44
1:B:314[B]:THR:HG22	7:B:2128:HOH:O	2.19	0.42
6:D:1328:CL:CL	6:D:1329:CL:CL	3.11	0.42
1:B:299:LYS:NZ	3:B:1326:SGN:O1	2.53	0.42
1:D:240:GLU:CD	7:D:2036:HOH:O	2.58	0.41
2:B:1334:TRP:HB3	7:B:2106:HOH:O	2.21	0.41
1:A:302:SER:O	1:A:305:ALA:HB2	2.20	0.41
1:C:212[A]:ILE:HG12	6:C:1331:CL:CL	2.58	0.40
1:C:322:PRO:HB3	7:C:2012:HOH:O	2.20	0.40
1:D:275:VAL:O	1:D:278:VAL:HG12	2.20	0.40
1:A:226:ARG:CZ	7:A:2004:HOH:O	2.62	0.40
6:B:1330:CL:CL	6:B:1333:CL:CL	3.13	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:2005:HOH:O	7:D:2095:HOH:O[3_655]	2.12	0.08

## 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	133/261 (51%)	128 (96%)	5 (4%)	0	100	100
1	B	130/261 (50%)	126 (97%)	4 (3%)	0	100	100
1	C	129/261 (49%)	124 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	129/261 (49%)	125 (97%)	4 (3%)	0	100	100
All	All	521/1044 (50%)	503 (96%)	18 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	118/230 (51%)	117 (99%)	1 (1%)	81	75
1	B	114/230 (50%)	113 (99%)	1 (1%)	78	71
1	C	113/230 (49%)	112 (99%)	1 (1%)	78	71
1	D	113/230 (49%)	111 (98%)	2 (2%)	59	44
All	All	458/920 (50%)	453 (99%)	5 (1%)	73	64

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	243	TYR
1	B	243	TYR
1	C	243	TYR
1	D	243	TYR
1	D	304	GLN

Some 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](#)

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 22 are monoatomic - leaving 7 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	SGN	D	1325	5	19,20,20	2.31	10 (52%)	24,31,31	1.56	4 (16%)
3	SGN	A	1327	5	19,20,20	2.17	10 (52%)	24,31,31	1.46	2 (8%)
3	SGN	B	1326	5	19,20,20	2.12	9 (47%)	24,31,31	1.58	2 (8%)
3	SGN	C	1326	5	19,20,20	2.30	9 (47%)	24,31,31	1.59	3 (12%)

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	SGN	D	1325	5	-	2/11/31/31	0/1/1/1
3	SGN	A	1327	5	-	0/11/31/31	0/1/1/1
3	SGN	B	1326	5	-	1/11/31/31	0/1/1/1
3	SGN	C	1326	5	-	1/11/31/31	0/1/1/1

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1325	SGN	O2S-S1	4.69	1.47	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1326	SGN	O2S-S1	4.64	1.47	1.42
3	A	1327	SGN	O6-S2	-3.90	1.46	1.56
3	D	1325	SGN	O6-S2	-3.70	1.46	1.56
3	C	1326	SGN	O6-S2	-3.68	1.46	1.56
3	B	1326	SGN	O6-S2	-3.63	1.47	1.56
3	A	1327	SGN	S1-N2	3.44	1.64	1.59
3	D	1325	SGN	S1-N2	3.15	1.63	1.59
3	C	1326	SGN	S1-N2	3.13	1.63	1.59
3	B	1326	SGN	S1-N2	3.10	1.63	1.59
3	A	1327	SGN	C1-C2	-2.96	1.49	1.52
3	B	1326	SGN	C1-C2	-2.93	1.49	1.52
3	A	1327	SGN	O4-C4	-2.86	1.36	1.43
3	B	1326	SGN	O4-C4	-2.78	1.36	1.43
3	C	1326	SGN	C1-C2	-2.76	1.49	1.52
3	D	1325	SGN	O1S-S1	2.70	1.45	1.42
3	D	1325	SGN	C1-C2	-2.60	1.49	1.52
3	C	1326	SGN	O4-C4	-2.60	1.36	1.43
3	C	1326	SGN	C6-C5	2.54	1.59	1.51
3	A	1327	SGN	O2S-S1	2.50	1.45	1.42
3	D	1325	SGN	C4-C5	2.45	1.58	1.53
3	B	1326	SGN	O2S-S1	2.44	1.44	1.42
3	B	1326	SGN	C6-C5	2.41	1.59	1.51
3	D	1325	SGN	C6-C5	2.37	1.59	1.51
3	C	1326	SGN	O1S-S1	2.37	1.44	1.42
3	B	1326	SGN	O1S-S1	2.35	1.44	1.42
3	A	1327	SGN	C6-C5	2.33	1.58	1.51
3	A	1327	SGN	O1S-S1	2.32	1.44	1.42
3	C	1326	SGN	C4-C5	2.31	1.57	1.53
3	D	1325	SGN	O4-C4	-2.27	1.37	1.43
3	A	1327	SGN	O6-C6	-2.19	1.37	1.46
3	B	1326	SGN	C4-C5	2.17	1.57	1.53
3	D	1325	SGN	O6-C6	-2.14	1.37	1.46
3	B	1326	SGN	O6-C6	-2.12	1.38	1.46
3	C	1326	SGN	O6-C6	-2.12	1.38	1.46
3	D	1325	SGN	O3-C3	-2.03	1.38	1.43
3	A	1327	SGN	O3-C3	-2.02	1.38	1.43
3	A	1327	SGN	C4-C5	2.01	1.57	1.53

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1326	SGN	O2S-S1-O1S	-4.73	108.98	120.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1325	SGN	O2S-S1-O1S	-4.52	109.48	120.16
3	C	1326	SGN	O5-C5-C6	4.47	115.69	106.67
3	A	1327	SGN	O5-C5-C6	4.44	115.62	106.67
3	C	1326	SGN	O2S-S1-O1S	-4.36	109.84	120.16
3	B	1326	SGN	O5-C5-C6	4.25	115.24	106.67
3	A	1327	SGN	O2S-S1-O1S	-4.08	110.52	120.16
3	D	1325	SGN	C1-O5-C5	3.30	119.89	113.66
3	D	1325	SGN	O5-C5-C6	3.27	113.26	106.67
3	D	1325	SGN	O5-C5-C4	2.36	113.98	109.69
3	C	1326	SGN	C1-O5-C5	2.28	117.96	113.66

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1325	SGN	C4-C5-C6-O6
3	B	1326	SGN	C4-C5-C6-O6
3	C	1326	SGN	C4-C5-C6-O6
3	D	1325	SGN	C2-N2-S1-O1S

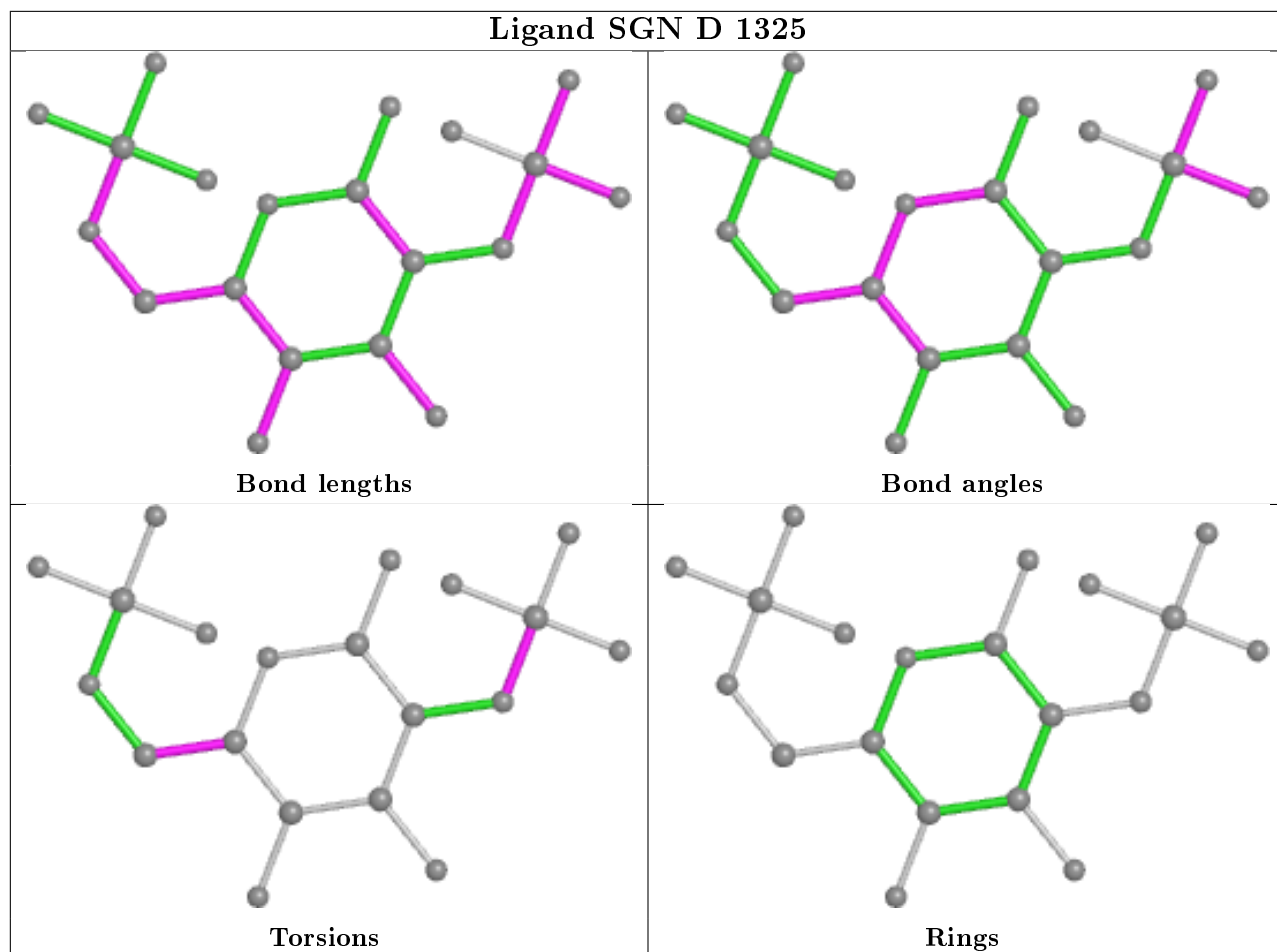
There are no ring outliers.

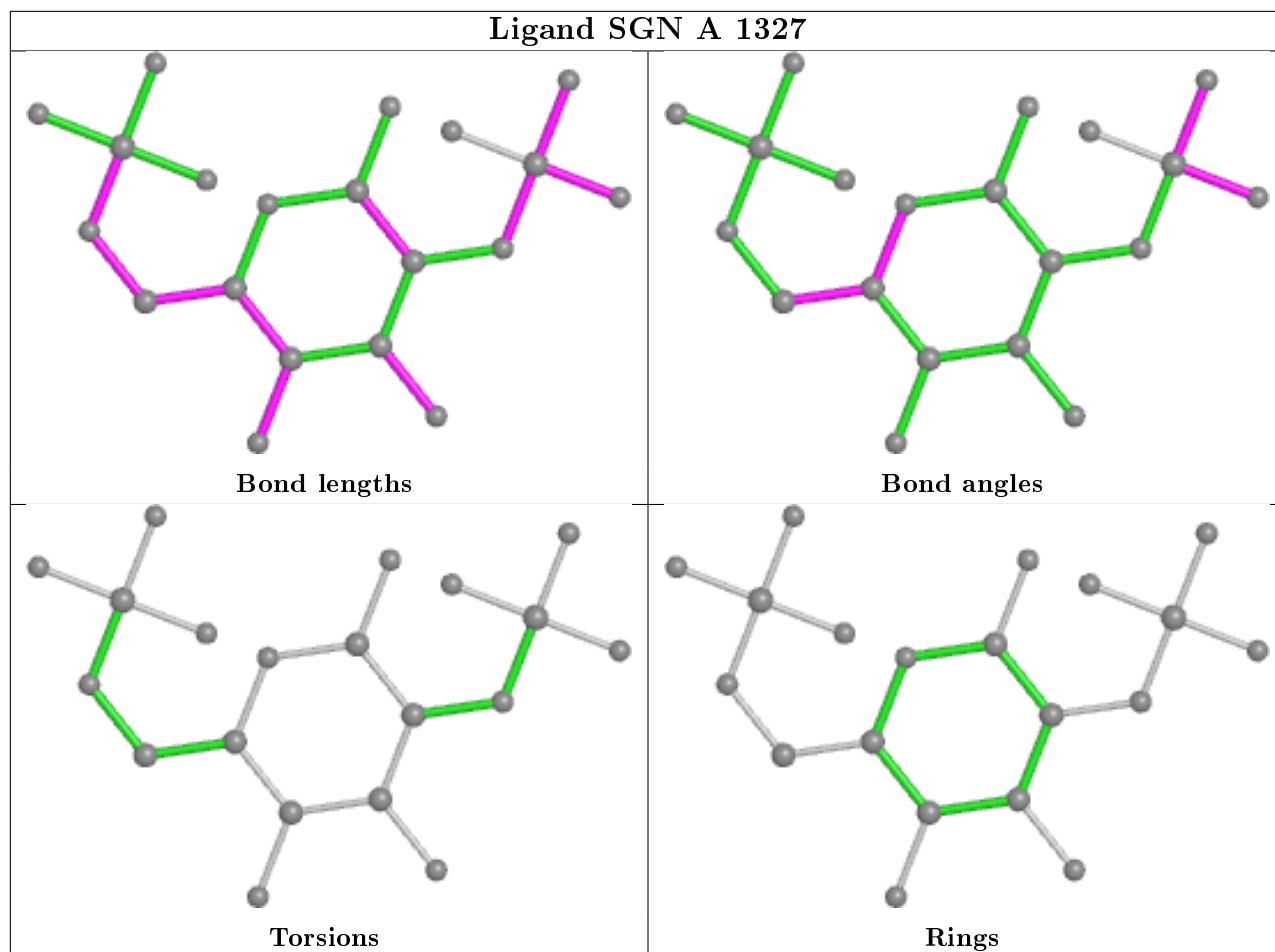
4 monomers are involved in 4 short contacts:

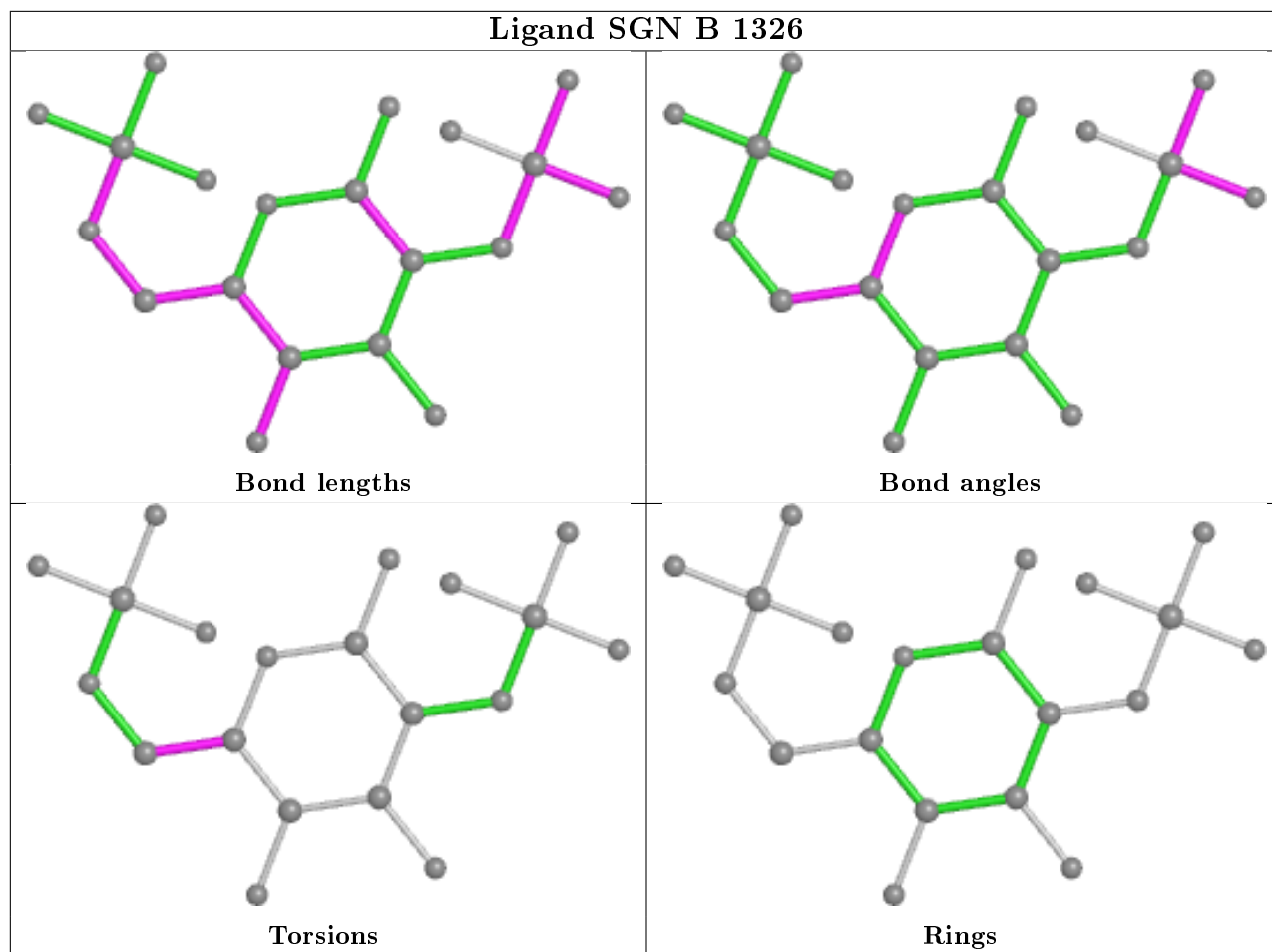
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1325	SGN	1	0
3	A	1327	SGN	1	0
3	B	1326	SGN	1	0
3	C	1326	SGN	1	0

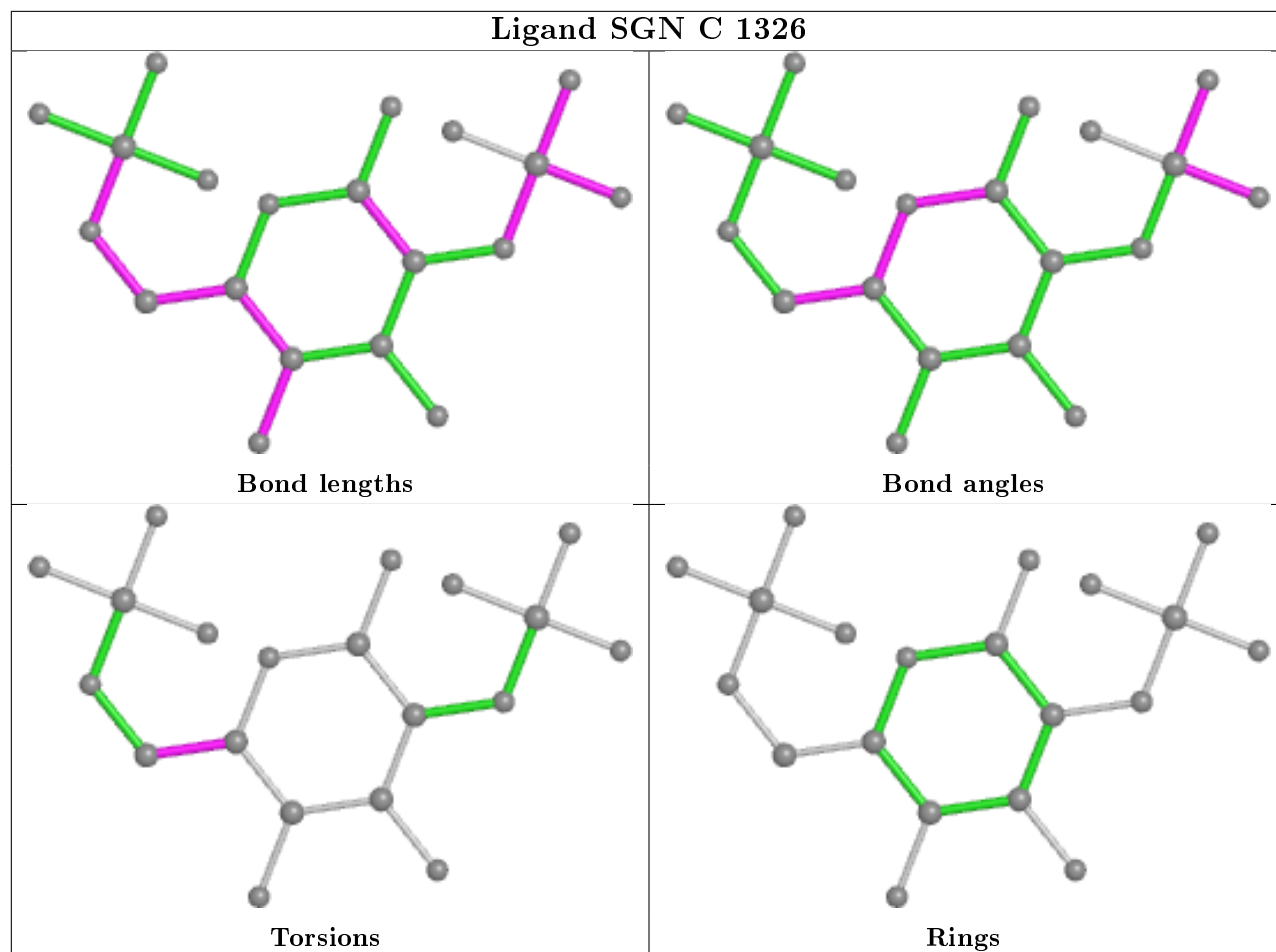
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	131/261 (50%)	-0.19	1 (0%) 86 86	17, 27, 48, 68	0
1	B	128/261 (49%)	-0.37	0 100 100	18, 26, 39, 58	0
1	C	128/261 (49%)	-0.25	1 (0%) 86 86	17, 25, 50, 71	0
1	D	127/261 (48%)	-0.27	2 (1%) 72 71	19, 29, 45, 71	0
All	All	514/1044 (49%)	-0.27	4 (0%) 86 86	17, 27, 48, 71	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	278[A]	VAL	3.8
1	C	291	ASN	3.3
1	D	324	VAL	2.3
1	D	278	VAL	2.3

### 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 monosaccharides in this entry.

### 6.4 Ligands [i](#)

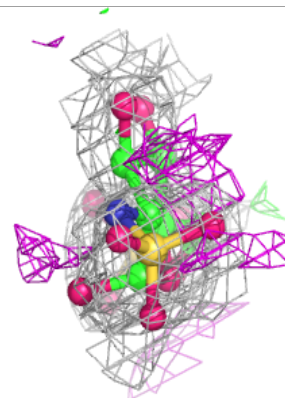
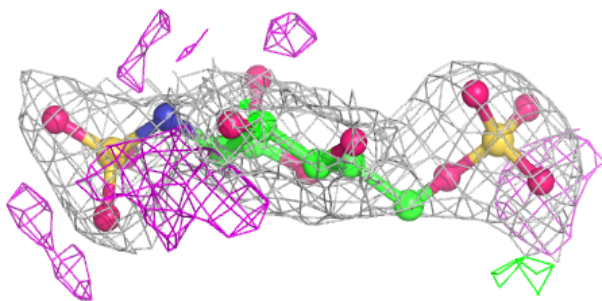
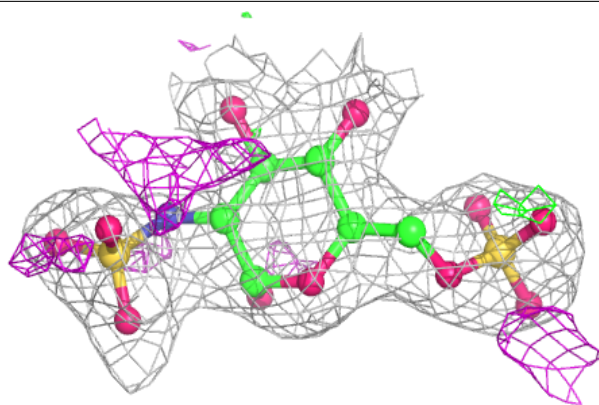
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	CL	B	1333	1/1	0.56	0.21	77,77,77,77	0
2	TRP	B	1334	15/15	0.58	0.38	69,80,95,96	0
6	CL	C	1330	1/1	0.73	0.36	67,67,67,67	1
2	TRP	A	1331	15/15	0.81	0.22	41,52,63,65	0
6	CL	B	1331	1/1	0.82	0.19	66,66,66,66	0
6	CL	D	1327	1/1	0.86	0.22	44,44,44,44	1
3	SGN	D	1325	20/20	0.89	0.16	49,58,72,74	0
2	TRP	D	1330	15/15	0.90	0.14	36,45,57,57	0
3	SGN	C	1326	20/20	0.90	0.17	51,62,85,86	0
3	SGN	B	1326	20/20	0.90	0.17	41,58,89,90	0
6	CL	D	1328	1/1	0.91	0.15	37,37,37,37	1
6	CL	B	1332	1/1	0.91	0.18	38,38,38,38	1
3	SGN	A	1327	20/20	0.92	0.15	34,52,69,69	0
6	CL	D	1329	1/1	0.93	0.15	38,38,38,38	1
4	EU	A	2000	1/1	0.93	0.08	78,78,78,78	1
4	EU	C	1329	1/1	0.95	0.15	28,28,28,28	1
6	CL	C	1332	1/1	0.96	0.11	59,59,59,59	0
5	CA	A	1329[A]	1/1	0.98	0.06	32,32,32,32	1
4	EU	A	1328[B]	1/1	0.98	0.06	32,32,32,32	1
4	EU	B	1327[B]	1/1	0.98	0.11	28,28,28,28	1
6	CL	B	1329	1/1	0.98	0.05	38,38,38,38	0
6	CL	A	1330	1/1	0.98	0.11	30,30,30,30	0
5	CA	B	1328[A]	1/1	0.98	0.11	24,24,24,24	1
6	CL	C	1331	1/1	0.98	0.07	33,33,33,33	0
4	EU	C	1328[B]	1/1	0.99	0.09	34,34,34,34	1
5	CA	C	1327[A]	1/1	0.99	0.09	34,34,34,34	1
5	CA	D	1326	1/1	0.99	0.10	22,22,22,22	0
6	CL	B	1330	1/1	0.99	0.12	32,32,32,32	1
4	EU	B	2000	1/1	0.99	0.05	30,30,30,30	1

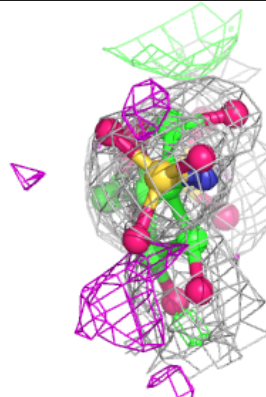
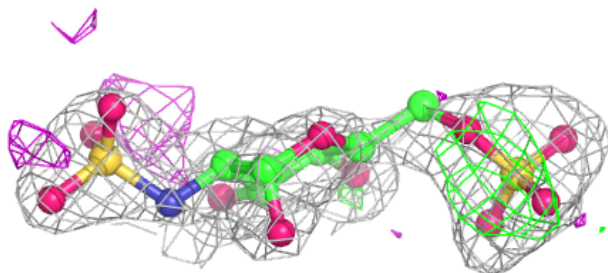
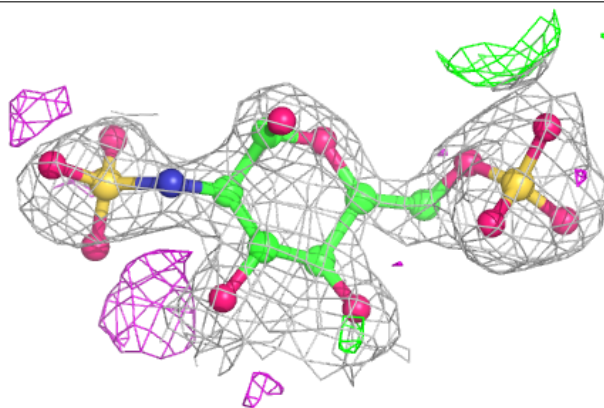
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around SGN D 1325:**

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

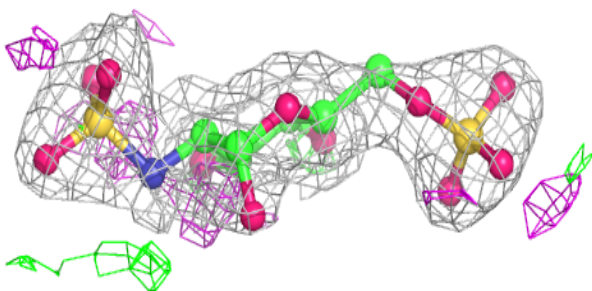
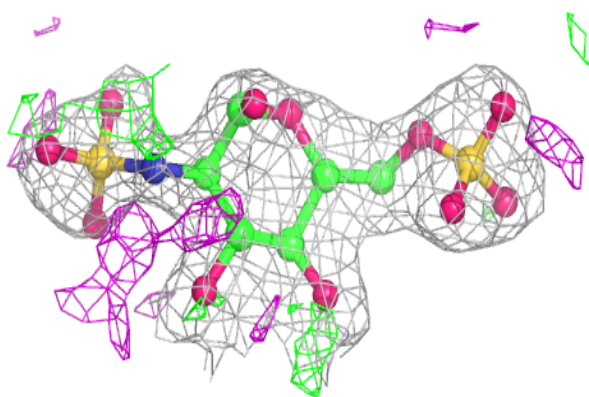
**Electron density around SGN C 1326:**

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

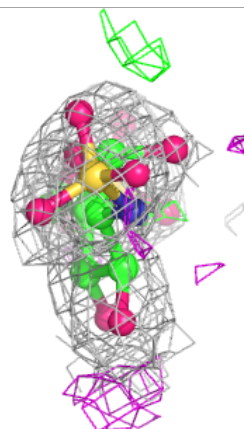
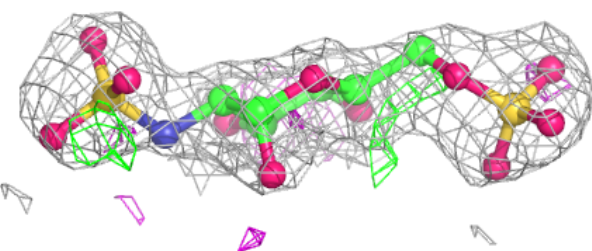
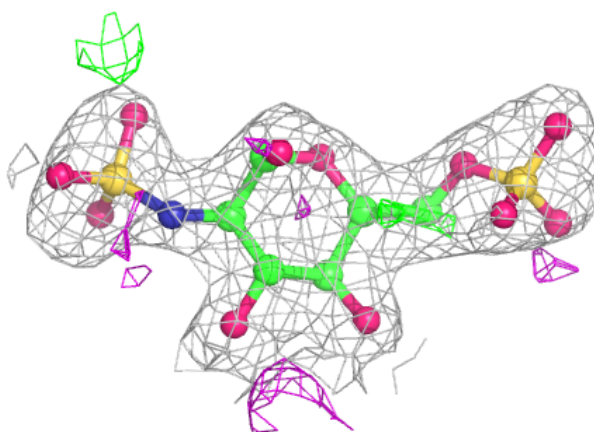


**Electron density around SGN B 1326:**

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

**Electron density around SGN A 1327:**

$2mF_o-DF_c$  (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.