



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:06 PM UTC

PDB ID : 6F05 / pdb_00006f05
Title : ARABIDOPSIS THALIANA GSTF9, GSO3 BOUND
Authors : Tossounian, M.A.; Wahni, K.; VanMolle, I.; Vertommen, D.; Rosado, L.;
Messens, J.
Deposited on : 2017-11-17
Resolution : 2.20 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

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

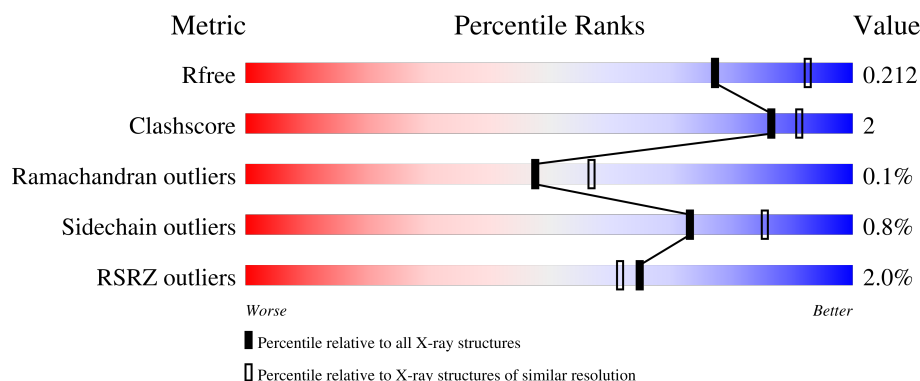
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 ≥ 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	215	 2% 87% 8% 5%
1	B	215	 % 87% 8% 5%
1	C	215	 2% 89% 5% 6%
1	D	215	 % 86% 7% 7%
1	E	215	 2% 88% 6% 7%

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Mol	Chain	Length	Quality of chain
1	F	215	2% 87% 7% 6%
1	G	215	87% 5% 8%
1	H	215	2% 87% 7% 7%
1	I	215	% 90% 6% •
1	J	215	5% 84% 6% 10%

2 Entry composition

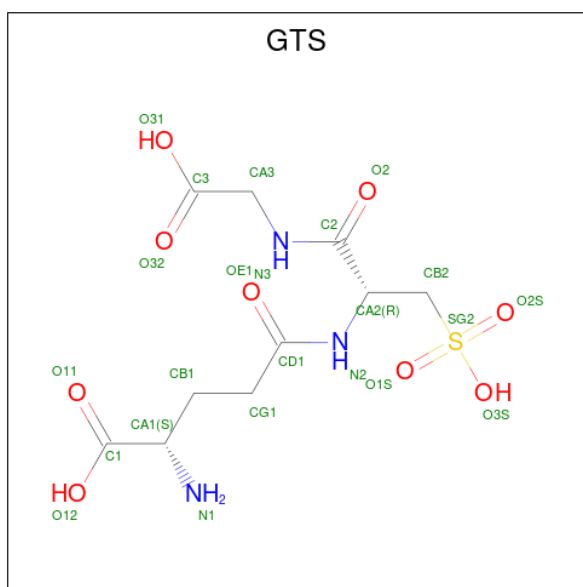
There are 5 unique types of molecules in this entry. The entry contains 16527 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glutathione S-transferase F9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	S	0	0	0
			1578	1024	260	291	3			
1	B	205	Total	C	N	O	S	0	2	0
			1617	1052	263	298	4			
1	C	202	Total	C	N	O	S	0	0	0
			1586	1030	260	293	3			
1	D	199	Total	C	N	O	S	0	0	0
			1524	989	250	282	3			
1	E	201	Total	C	N	O	S	0	0	0
			1557	1011	256	288	2			
1	F	203	Total	C	N	O	S	0	1	0
			1575	1020	259	294	2			
1	G	198	Total	C	N	O	S	0	0	0
			1515	983	248	282	2			
1	H	201	Total	C	N	O	S	0	0	0
			1547	1007	254	282	4			
1	I	206	Total	C	N	O	S	0	2	0
			1608	1040	261	303	4			
1	J	193	Total	C	N	O	S	0	0	0
			1422	918	235	267	2			

- Molecule 2 is GLUTATHIONE SULFONIC ACID (CCD ID: GTS) (formula: C₁₀H₁₇N₃O₉S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	B	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	C	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	D	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	E	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	F	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	G	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	H	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	I	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
2	J	1	Total	C	N	O	S	0	0
			23	10	3	9	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	E	1	Total	Cl	0	0
			1	1		
4	F	1	Total	Cl	0	0
			1	1		
4	G	1	Total	Cl	0	0
			1	1		
4	H	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	1	Total	Cl	0	0
			1	1		

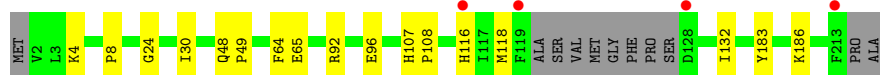
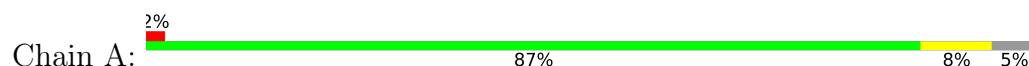
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	123	Total	O	0	0
			123	123		
5	B	127	Total	O	0	0
			127	127		
5	C	119	Total	O	0	0
			119	119		
5	D	47	Total	O	0	0
			47	47		
5	E	92	Total	O	0	0
			92	92		
5	F	52	Total	O	0	0
			52	52		
5	G	21	Total	O	0	0
			21	21		
5	H	44	Total	O	0	0
			44	44		
5	I	102	Total	O	0	0
			102	102		
5	J	8	Total	O	0	0
			8	8		

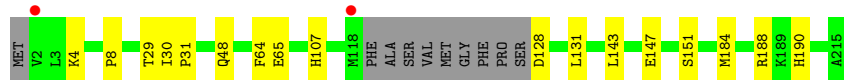
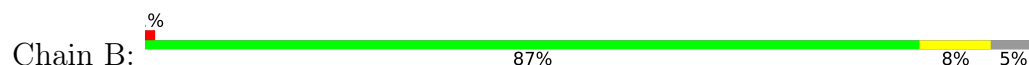
3 Residue-property plots [i](#)

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.

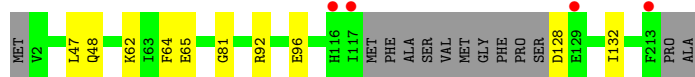
- Molecule 1: Glutathione S-transferase F9



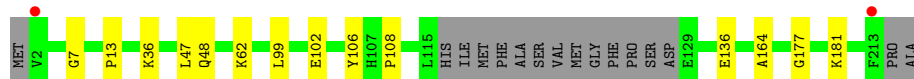
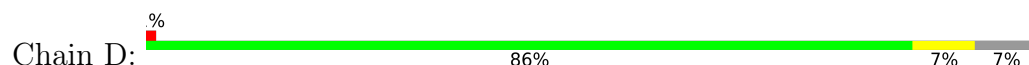
- Molecule 1: Glutathione S-transferase F9



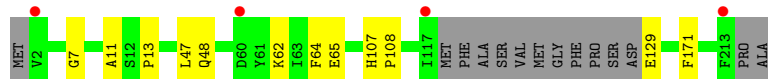
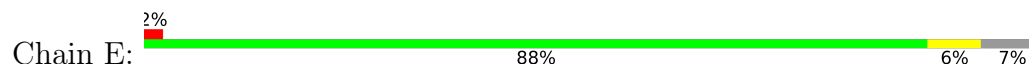
- Molecule 1: Glutathione S-transferase F9



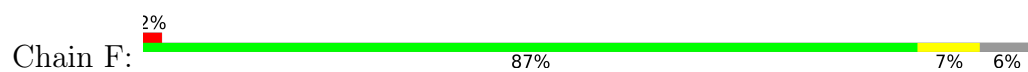
- Molecule 1: Glutathione S-transferase F9



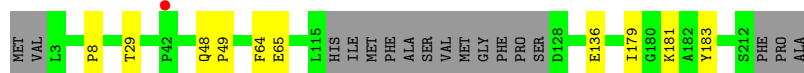
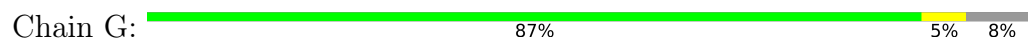
- Molecule 1: Glutathione S-transferase F9



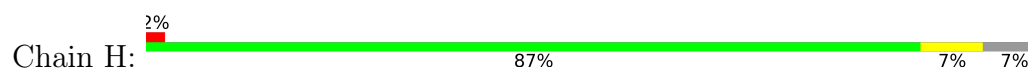
- Molecule 1: Glutathione S-transferase F9



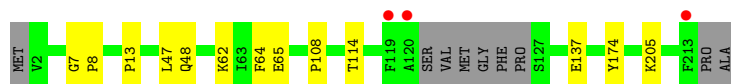
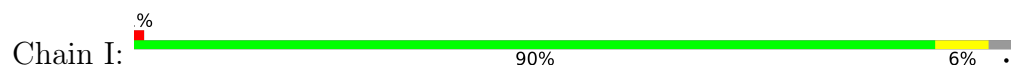
• Molecule 1: Glutathione S-transferase F9



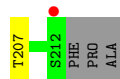
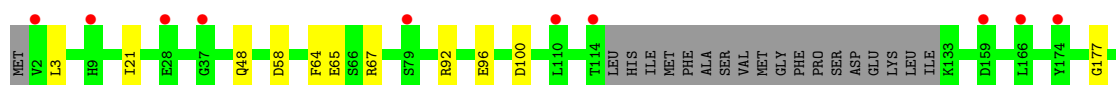
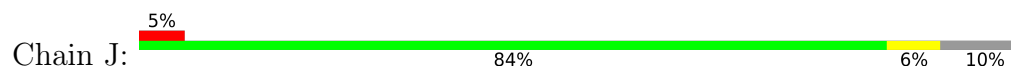
• Molecule 1: Glutathione S-transferase F9



• Molecule 1: Glutathione S-transferase F9



• Molecule 1: Glutathione S-transferase F9



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.43Å 93.97Å 107.71Å 93.19° 101.57° 101.97°	Depositor
Resolution (Å)	65.73 – 2.20 65.73 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.9 (65.73-2.20) 97.9 (65.73-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.50 (at 2.07Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.162 , 0.209 0.166 , 0.212	Depositor DCC
R_{free} test set	6862 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	34.1	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16527	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: GTS, CL, GOL

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	0.60	0/1619	0.85	4/2207 (0.2%)
1	B	0.66	2/1667 (0.1%)	0.93	4/2269 (0.2%)
1	C	0.59	0/1627	0.82	0/2214
1	D	0.49	0/1565	0.82	3/2140 (0.1%)
1	E	0.62	0/1598	0.83	4/2180 (0.2%)
1	F	0.59	2/1620 (0.1%)	0.89	7/2211 (0.3%)
1	G	0.42	0/1556	0.79	0/2129
1	H	0.48	0/1589	0.78	0/2169
1	I	0.55	0/1656	0.84	2/2259 (0.1%)
1	J	0.41	0/1462	0.80	2/2011 (0.1%)
All	All	0.55	4/15959 (0.0%)	0.84	26/21789 (0.1%)

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	F	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	184[A]	MET	CA-C	6.90	1.62	1.52
1	B	184[B]	MET	CA-C	6.90	1.62	1.52
1	F	41	GLN	C-N	5.45	1.40	1.34
1	F	42	PRO	N-CD	5.15	1.54	1.47

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	184[A]	MET	CA-C-O	9.13	130.43	120.20
1	B	184[B]	MET	CA-C-O	9.13	130.43	120.20
1	B	107	HIS	CA-C-N	-6.30	113.89	120.38
1	B	107	HIS	C-N-CA	-6.30	113.89	120.38
1	A	108	PRO	CA-C-N	-6.10	112.73	119.19
1	A	108	PRO	C-N-CA	-6.10	112.73	119.19
1	A	107	HIS	CA-C-N	-5.87	114.33	120.38
1	A	107	HIS	C-N-CA	-5.87	114.33	120.38
1	F	108	PRO	CA-C-N	-5.86	112.97	119.19
1	F	108	PRO	C-N-CA	-5.86	112.97	119.19
1	E	108	PRO	CA-C-N	-5.72	113.13	119.19
1	E	108	PRO	C-N-CA	-5.72	113.13	119.19
1	F	116	HIS	N-CA-C	5.68	120.05	113.18
1	F	78	ARG	N-CA-C	5.62	117.40	111.28
1	F	107	HIS	CB-CA-C	-5.50	102.53	112.76
1	E	107	HIS	CA-C-N	-5.48	114.73	120.38
1	E	107	HIS	C-N-CA	-5.48	114.73	120.38
1	F	41	GLN	CA-C-N	-5.46	113.40	119.19
1	F	41	GLN	C-N-CA	-5.46	113.40	119.19
1	J	177	GLY	CA-C-N	5.36	125.17	119.28
1	J	177	GLY	C-N-CA	5.36	125.17	119.28
1	D	177	GLY	CA-C-N	5.26	125.07	119.28
1	D	177	GLY	C-N-CA	5.26	125.07	119.28
1	D	108	PRO	N-CA-C	5.09	116.92	110.70
1	I	108	PRO	CA-C-N	-5.03	113.86	119.19
1	I	108	PRO	C-N-CA	-5.03	113.86	119.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	107	HIS	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	1578	0	1534	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1617	0	1589	8	0
1	C	1586	0	1567	5	0
1	D	1524	0	1456	7	0
1	E	1557	0	1512	4	0
1	F	1575	0	1520	9	0
1	G	1515	0	1443	5	0
1	H	1547	0	1492	8	0
1	I	1608	0	1553	8	0
1	J	1422	0	1283	6	0
2	A	23	0	15	0	0
2	B	23	0	15	0	0
2	C	23	0	15	0	0
2	D	23	0	15	0	0
2	E	23	0	15	0	0
2	F	23	0	15	0	0
2	G	23	0	15	0	0
2	H	23	0	15	0	0
2	I	23	0	15	0	0
2	J	23	0	15	0	0
3	A	12	0	16	2	0
3	B	12	0	16	1	0
4	A	1	0	0	1	0
4	B	1	0	0	1	0
4	C	1	0	0	0	0
4	D	1	0	0	1	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	1	0
4	I	1	0	0	1	0
5	A	123	0	0	0	0
5	B	127	0	0	0	0
5	C	119	0	0	0	0
5	D	47	0	0	0	0
5	E	92	0	0	0	0
5	F	52	0	0	0	0
5	G	21	0	0	0	0
5	H	44	0	0	1	0
5	I	102	0	0	0	0
5	J	8	0	0	0	0
All	All	16527	0	15131	64	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 2.

All (64) 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:4:LYS:HD3	1:B:30:ILE:HD11	1.73	0.69
1:I:114:THR:HG23	1:I:174:TYR:HE2	1.58	0.68
1:D:136:GLU:OE1	1:D:181:LYS:NZ	2.30	0.60
1:A:4:LYS:HD3	1:A:30:ILE:HD11	1.83	0.59
1:I:114:THR:HG23	1:I:174:TYR:CE2	2.36	0.59
1:F:15:ARG:HH22	1:F:107:HIS:CE1	2.23	0.56
1:D:47:LEU:HA	1:D:62:LYS:HD2	1.89	0.54
1:H:114:THR:HG22	1:H:179:ILE:HG13	1.90	0.53
1:F:107:HIS:CD2	1:F:111:LEU:HB2	2.43	0.53
1:D:36:LYS:HD3	1:I:137:GLU:CD	2.35	0.52
1:A:186:LYS:HG3	1:F:183:TYR:CZ	2.46	0.51
1:J:21:ILE:HD12	1:J:207:THR:HG23	1.91	0.51
1:A:24:GLY:HA3	3:A:302:GOL:H12	1.95	0.48
1:A:116:HIS:CB	1:A:132:ILE:HD11	2.43	0.48
1:E:47:LEU:HA	1:E:62:LYS:HD2	1.94	0.48
1:E:64:PHE:O	1:E:65:GLU:HB2	2.14	0.48
1:B:128:ASP:HB3	1:B:131:LEU:HB3	1.96	0.48
1:C:92:ARG:O	1:C:96:GLU:HG2	2.13	0.48
1:F:64:PHE:O	1:F:65:GLU:HB2	2.14	0.47
1:A:64:PHE:O	1:A:65:GLU:HB2	2.14	0.46
1:B:147:GLU:OE1	1:B:188:ARG:HB3	2.16	0.46
1:A:49:PRO:HG3	3:B:303:GOL:H32	1.97	0.46
1:F:112:ASN:O	1:F:116:HIS:HB2	2.16	0.46
1:G:64:PHE:O	1:G:65:GLU:HB2	2.14	0.46
1:A:183:TYR:HA	1:A:186:LYS:HB2	1.98	0.45
1:B:64:PHE:O	1:B:65:GLU:HB2	2.15	0.45
1:B:151[B]:SER:HG	1:B:190:HIS:CG	2.34	0.45
1:C:47:LEU:HA	1:C:62:LYS:HD2	1.99	0.45
1:F:100:ASP:OD2	1:J:67:ARG:HB2	2.17	0.45
1:I:47:LEU:HA	1:I:62:LYS:HD2	1.98	0.45
1:C:128:ASP:O	1:C:132:ILE:HG12	2.17	0.45
1:D:102:GLU:HA	1:D:106:TYR:HB3	1.99	0.44
3:A:303:GOL:H11	1:C:81:GLY:HA2	2.00	0.44
1:F:67:ARG:HB2	1:J:100:ASP:OD2	2.18	0.44
1:F:107:HIS:HD2	1:F:111:LEU:HB2	1.83	0.44
1:A:8:PRO:HD2	4:A:304:CL:CL	2.55	0.44
1:D:7:GLY:HA2	4:D:302:CL:CL	2.54	0.44
1:I:7:GLY:O	1:I:13:PRO:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:3:LEU:HA	1:J:58:ASP:HA	1.99	0.44
1:H:58:ASP:OD2	1:H:76:LYS:NZ	2.49	0.43
1:I:64:PHE:O	1:I:65:GLU:HB2	2.19	0.43
1:G:136:GLU:OE2	1:G:183:TYR:OH	2.30	0.43
1:C:64:PHE:O	1:C:65:GLU:HB2	2.19	0.43
1:E:7:GLY:O	1:E:13:PRO:HG2	2.19	0.43
1:H:2:VAL:HG13	1:H:4:LYS:HE3	2.00	0.43
1:H:7:GLY:HA2	4:H:302:CL:CL	2.55	0.43
1:J:92:ARG:O	1:J:96:GLU:HG2	2.20	0.42
1:B:31:PRO:HA	4:B:304:CL:CL	2.57	0.42
1:A:92:ARG:O	1:A:96:GLU:HG2	2.19	0.42
1:D:7:GLY:O	1:D:13:PRO:HG2	2.20	0.42
1:G:179:ILE:HG22	1:G:181:LYS:HG3	2.02	0.42
1:J:64:PHE:O	1:J:65:GLU:HB2	2.19	0.42
1:B:143:LEU:HD23	1:B:143:LEU:HA	1.91	0.41
1:H:110:LEU:HD11	1:H:175:LEU:HB2	2.01	0.41
1:F:15:ARG:NH2	1:F:107:HIS:ND1	2.68	0.41
1:I:8:PRO:HD2	4:I:302:CL:CL	2.58	0.41
1:G:49:PRO:HB2	1:H:145:VAL:HG13	2.02	0.41
1:G:8:PRO:HG3	1:G:29:THR:HB	2.01	0.41
1:H:40:LYS:NZ	5:H:405:HOH:O	2.53	0.41
1:D:99:LEU:HD13	1:D:164:ALA:HA	2.03	0.40
1:B:8:PRO:HG3	1:B:29:THR:HB	2.02	0.40
1:H:64:PHE:O	1:H:65:GLU:HB2	2.20	0.40
1:E:11:ALA:HB3	1:E:171:PHE:CZ	2.56	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	200/215 (93%)	196 (98%)	3 (2%)	1 (0%)	24 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	203/215 (94%)	200 (98%)	3 (2%)	0	100	100
1	C	198/215 (92%)	196 (99%)	2 (1%)	0	100	100
1	D	195/215 (91%)	193 (99%)	2 (1%)	0	100	100
1	E	197/215 (92%)	195 (99%)	2 (1%)	0	100	100
1	F	200/215 (93%)	198 (99%)	2 (1%)	0	100	100
1	G	194/215 (90%)	192 (99%)	2 (1%)	0	100	100
1	H	197/215 (92%)	195 (99%)	2 (1%)	0	100	100
1	I	204/215 (95%)	202 (99%)	2 (1%)	0	100	100
1	J	189/215 (88%)	187 (99%)	2 (1%)	0	100	100
All	All	1977/2150 (92%)	1954 (99%)	22 (1%)	1 (0%)	48	57

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	118	MET

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	161/182 (88%)	160 (99%)	1 (1%)	78	89
1	B	169/182 (93%)	168 (99%)	1 (1%)	78	89
1	C	166/182 (91%)	165 (99%)	1 (1%)	78	89
1	D	153/182 (84%)	152 (99%)	1 (1%)	76	87
1	E	159/182 (87%)	157 (99%)	2 (1%)	61	76
1	F	162/182 (89%)	160 (99%)	2 (1%)	63	78
1	G	152/182 (84%)	151 (99%)	1 (1%)	76	87
1	H	156/182 (86%)	155 (99%)	1 (1%)	78	89
1	I	167/182 (92%)	166 (99%)	1 (1%)	78	89

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	133/182 (73%)	132 (99%)	1 (1%)	73	85
All	All	1578/1820 (87%)	1566 (99%)	12 (1%)	73	85

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	GLN
1	B	48	GLN
1	C	48	GLN
1	D	48	GLN
1	E	48	GLN
1	E	129	GLU
1	F	48	GLN
1	F	117	ILE
1	G	48	GLN
1	H	48	GLN
1	I	48	GLN
1	J	48	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	HIS
1	C	107	HIS
1	E	107	HIS
1	F	80	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

Of 23 ligands modelled in this entry, 9 are monoatomic - leaving 14 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
2	GTS	J	301	-	20,22,22	2.06	5 (25%)	26,30,30	1.90	6 (23%)
2	GTS	D	301	-	20,22,22	1.99	5 (25%)	26,30,30	1.83	7 (26%)
3	GOL	B	303	-	5,5,5	0.40	0	5,5,5	0.77	0
2	GTS	E	301	-	20,22,22	2.00	6 (30%)	26,30,30	1.80	7 (26%)
2	GTS	B	301	-	20,22,22	1.84	4 (20%)	26,30,30	1.71	7 (26%)
3	GOL	A	302	-	5,5,5	0.47	0	5,5,5	0.31	0
2	GTS	A	301	-	20,22,22	1.83	4 (20%)	26,30,30	2.29	11 (42%)
2	GTS	H	301	-	20,22,22	2.05	7 (35%)	26,30,30	1.70	5 (19%)
3	GOL	A	303	-	5,5,5	0.38	0	5,5,5	0.36	0
3	GOL	B	302	-	5,5,5	0.33	0	5,5,5	0.32	0
2	GTS	G	301	-	20,22,22	2.04	6 (30%)	26,30,30	2.00	11 (42%)
2	GTS	I	301	-	20,22,22	1.98	4 (20%)	26,30,30	1.85	9 (34%)
2	GTS	F	301	-	20,22,22	2.02	6 (30%)	26,30,30	1.70	5 (19%)
2	GTS	C	301	-	20,22,22	2.01	7 (35%)	26,30,30	1.59	7 (26%)

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	GTS	J	301	-	-	4/27/27/27	-
2	GTS	D	301	-	-	3/27/27/27	-
3	GOL	B	303	-	-	0/4/4/4	-
2	GTS	E	301	-	-	3/27/27/27	-
2	GTS	B	301	-	-	5/27/27/27	-
3	GOL	A	302	-	-	4/4/4/4	-
2	GTS	A	301	-	-	6/27/27/27	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GTS	H	301	-	-	1/27/27/27	-
3	GOL	A	303	-	-	2/4/4/4	-
3	GOL	B	302	-	-	2/4/4/4	-
2	GTS	G	301	-	-	4/27/27/27	-
2	GTS	I	301	-	-	4/27/27/27	-
2	GTS	F	301	-	-	3/27/27/27	-
2	GTS	C	301	-	-	4/27/27/27	-

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	301	GTS	C2-N3	5.31	1.46	1.33
2	C	301	GTS	C2-N3	5.26	1.45	1.33
2	F	301	GTS	C2-N3	5.23	1.45	1.33
2	I	301	GTS	C2-N3	5.23	1.45	1.33
2	H	301	GTS	C2-N3	5.16	1.45	1.33
2	D	301	GTS	C2-N3	5.04	1.45	1.33
2	G	301	GTS	C2-N3	5.00	1.45	1.33
2	E	301	GTS	C2-N3	4.86	1.45	1.33
2	J	301	GTS	CD1-N2	4.56	1.43	1.34
2	B	301	GTS	C2-N3	4.52	1.44	1.33
2	E	301	GTS	CD1-N2	4.52	1.43	1.34
2	F	301	GTS	CD1-N2	4.49	1.43	1.34
2	G	301	GTS	CD1-N2	4.36	1.43	1.34
2	H	301	GTS	CD1-N2	4.21	1.42	1.34
2	B	301	GTS	CD1-N2	4.10	1.42	1.34
2	A	301	GTS	C2-N3	4.05	1.43	1.33
2	I	301	GTS	CD1-N2	3.98	1.42	1.34
2	D	301	GTS	CD1-N2	3.93	1.42	1.34
2	A	301	GTS	CD1-N2	3.87	1.42	1.34
2	C	301	GTS	CD1-N2	3.73	1.41	1.34
2	D	301	GTS	CG1-CD1	3.14	1.57	1.51
2	J	301	GTS	CG1-CD1	2.60	1.56	1.51
2	A	301	GTS	O1S-SG2	2.54	1.52	1.45
2	H	301	GTS	CB2-CA2	-2.54	1.51	1.53
2	C	301	GTS	CB2-CA2	-2.51	1.51	1.53
2	E	301	GTS	O1S-SG2	2.47	1.52	1.45
2	I	301	GTS	O11-C1	2.46	1.29	1.22
2	I	301	GTS	CG1-CD1	2.40	1.56	1.51
2	G	301	GTS	O2S-SG2	2.40	1.51	1.45
2	C	301	GTS	O11-C1	2.36	1.29	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	GTS	O1S-SG2	2.34	1.51	1.45
2	H	301	GTS	O11-C1	2.32	1.29	1.22
2	H	301	GTS	O1S-SG2	2.32	1.51	1.45
2	E	301	GTS	CB2-CA2	-2.30	1.51	1.53
2	D	301	GTS	O1S-SG2	2.28	1.51	1.45
2	E	301	GTS	CG1-CD1	2.26	1.55	1.51
2	A	301	GTS	O2S-SG2	2.21	1.51	1.45
2	J	301	GTS	O2S-SG2	2.18	1.51	1.45
2	C	301	GTS	O1S-SG2	2.15	1.51	1.45
2	F	301	GTS	O11-C1	2.15	1.28	1.22
2	G	301	GTS	O11-C1	2.13	1.28	1.22
2	F	301	GTS	O2S-SG2	2.13	1.51	1.45
2	J	301	GTS	O1S-SG2	2.12	1.51	1.45
2	G	301	GTS	O1S-SG2	2.11	1.51	1.45
2	E	301	GTS	O11-C1	2.11	1.28	1.22
2	C	301	GTS	CG1-CD1	2.10	1.55	1.51
2	G	301	GTS	CG1-CD1	2.10	1.55	1.51
2	D	301	GTS	O11-C1	2.09	1.28	1.22
2	F	301	GTS	CG1-CD1	2.08	1.55	1.51
2	B	301	GTS	CA2-N2	-2.08	1.41	1.45
2	C	301	GTS	O2S-SG2	2.06	1.50	1.45
2	H	301	GTS	O2S-SG2	2.05	1.50	1.45
2	B	301	GTS	CG1-CD1	2.02	1.55	1.51
2	H	301	GTS	CG1-CD1	2.01	1.55	1.51

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	GTS	C3-CA3-N3	-5.28	96.66	113.06
2	J	301	GTS	O1S-SG2-CB2	4.37	113.29	106.76
2	A	301	GTS	CA2-C2-N3	4.32	125.81	116.54
2	I	301	GTS	O3S-SG2-O2S	-4.31	100.63	111.40
2	A	301	GTS	O2-C2-N3	-4.25	113.99	122.98
2	G	301	GTS	O3S-SG2-O2S	-4.18	100.95	111.40
2	H	301	GTS	O2S-SG2-O1S	-4.12	100.42	113.82
2	H	301	GTS	O1S-SG2-CB2	4.12	112.90	106.76
2	D	301	GTS	O2S-SG2-O1S	-4.00	100.83	113.82
2	J	301	GTS	O3S-SG2-O2S	-3.97	101.48	111.40
2	I	301	GTS	O3S-SG2-CB2	3.81	113.32	105.97
2	B	301	GTS	O2S-SG2-CB2	3.80	112.44	106.76
2	D	301	GTS	OE1-CD1-N2	-3.59	116.86	122.95
2	F	301	GTS	O3S-SG2-CB2	3.59	112.90	105.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	301	GTS	CG1-CD1-N2	3.51	122.05	115.86
2	G	301	GTS	CG1-CD1-N2	3.39	121.83	115.86
2	G	301	GTS	O1S-SG2-CB2	3.37	111.79	106.76
2	E	301	GTS	O3S-SG2-O1S	-3.36	103.00	111.40
2	B	301	GTS	CG1-CB1-CA1	-3.31	106.27	113.86
2	E	301	GTS	O1S-SG2-CB2	3.31	111.69	106.76
2	E	301	GTS	CA2-CB2-SG2	-3.25	108.69	113.55
2	F	301	GTS	O2S-SG2-O1S	-3.24	103.27	113.82
2	A	301	GTS	CA3-N3-C2	3.20	129.47	121.38
2	E	301	GTS	CG1-CD1-N2	3.16	121.42	115.86
2	F	301	GTS	CG1-CD1-N2	3.08	121.30	115.86
2	C	301	GTS	O2S-SG2-O1S	-3.08	103.81	113.82
2	A	301	GTS	O3S-SG2-O1S	-3.06	103.75	111.40
2	C	301	GTS	CA2-CB2-SG2	-3.05	108.98	113.55
2	H	301	GTS	CA2-CB2-SG2	-3.04	109.01	113.55
2	D	301	GTS	CG1-CD1-N2	2.98	121.11	115.86
2	G	301	GTS	CA2-C2-N3	2.89	122.75	116.54
2	D	301	GTS	O1S-SG2-CB2	2.83	110.97	106.76
2	D	301	GTS	O3S-SG2-CB2	2.79	111.36	105.97
2	D	301	GTS	O2S-SG2-CB2	2.71	110.80	106.76
2	A	301	GTS	OE1-CD1-N2	-2.70	118.38	122.95
2	J	301	GTS	OE1-CD1-N2	-2.66	118.45	122.95
2	I	301	GTS	O2S-SG2-CB2	2.63	110.69	106.76
2	A	301	GTS	C2-CA2-N2	-2.62	104.02	111.11
2	J	301	GTS	O2S-SG2-CB2	2.62	110.67	106.76
2	I	301	GTS	CG1-CD1-N2	2.57	120.40	115.86
2	G	301	GTS	O2-C2-N3	-2.57	117.54	122.98
2	A	301	GTS	O12-C1-O11	-2.55	118.30	124.08
2	B	301	GTS	O12-C1-O11	-2.53	118.33	124.08
2	A	301	GTS	CG1-CD1-N2	2.53	120.32	115.86
2	G	301	GTS	O3S-SG2-CB2	2.51	110.81	105.97
2	G	301	GTS	O2S-SG2-CB2	2.51	110.50	106.76
2	H	301	GTS	CB2-CA2-C2	-2.50	103.89	109.56
2	C	301	GTS	O1S-SG2-CB2	2.50	110.48	106.76
2	I	301	GTS	OE1-CD1-N2	-2.49	118.73	122.95
2	F	301	GTS	CA2-CB2-SG2	-2.48	109.83	113.55
2	G	301	GTS	OE1-CD1-N2	-2.48	118.76	122.95
2	C	301	GTS	OE1-CD1-N2	-2.47	118.77	122.95
2	C	301	GTS	O2S-SG2-CB2	2.47	110.44	106.76
2	J	301	GTS	O3S-SG2-CB2	2.47	110.74	105.97
2	B	301	GTS	CA3-N3-C2	-2.42	115.28	121.38
2	G	301	GTS	O12-C1-O11	-2.39	118.67	124.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	GTS	OE1-CD1-N2	-2.37	118.93	122.95
2	G	301	GTS	C3-CA3-N3	-2.35	105.76	113.06
2	D	301	GTS	O12-C1-O11	-2.33	118.80	124.08
2	F	301	GTS	O2S-SG2-CB2	2.31	110.20	106.76
2	C	301	GTS	CB2-CA2-C2	-2.31	104.33	109.56
2	E	301	GTS	OE1-CD1-N2	-2.27	119.10	122.95
2	C	301	GTS	O3S-SG2-CB2	2.26	110.34	105.97
2	A	301	GTS	O3S-SG2-CB2	2.20	110.22	105.97
2	I	301	GTS	CG1-CB1-CA1	-2.16	108.90	113.86
2	A	301	GTS	O2S-SG2-CB2	2.14	109.96	106.76
2	I	301	GTS	O31-C3-CA3	2.14	120.94	112.81
2	I	301	GTS	CA2-CB2-SG2	-2.12	110.39	113.55
2	B	301	GTS	CG1-CD1-N2	2.09	119.55	115.86
2	I	301	GTS	CB2-CA2-C2	-2.08	104.83	109.56
2	G	301	GTS	CA2-CB2-SG2	-2.05	110.49	113.55
2	E	301	GTS	O31-C3-CA3	2.04	120.57	112.81
2	H	301	GTS	O3S-SG2-CB2	2.02	109.88	105.97
2	E	301	GTS	O3S-SG2-CB2	2.01	109.85	105.97
2	B	301	GTS	CA2-C2-N3	2.01	120.84	116.54

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	GTS	CA2-CB2-SG2-O3S
2	C	301	GTS	CA2-CB2-SG2-O1S
2	E	301	GTS	CA2-CB2-SG2-O1S
2	E	301	GTS	CA2-CB2-SG2-O2S
2	E	301	GTS	CA2-CB2-SG2-O3S
2	G	301	GTS	N1-CA1-CB1-CG1
2	G	301	GTS	C1-CA1-CB1-CG1
2	I	301	GTS	CA2-CB2-SG2-O3S
2	J	301	GTS	CA2-CB2-SG2-O1S
2	J	301	GTS	CA2-CB2-SG2-O2S
2	J	301	GTS	CA2-CB2-SG2-O3S
3	A	302	GOL	O1-C1-C2-C3
3	A	302	GOL	C1-C2-C3-O3
3	A	303	GOL	O1-C1-C2-O2
3	B	302	GOL	O1-C1-C2-C3
2	A	301	GTS	O2-C2-N3-CA3
2	G	301	GTS	O2-C2-N3-CA3
2	G	301	GTS	CA2-C2-N3-CA3

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Mol	Chain	Res	Type	Atoms
2	A	301	GTS	CA2-C2-N3-CA3
2	A	301	GTS	O32-C3-CA3-N3
2	A	301	GTS	O31-C3-CA3-N3
3	A	303	GOL	O1-C1-C2-C3
3	B	302	GOL	O1-C1-C2-O2
3	A	302	GOL	O1-C1-C2-O2
2	B	301	GTS	CA2-CB2-SG2-O1S
2	B	301	GTS	CA2-CB2-SG2-O2S
2	C	301	GTS	CA2-CB2-SG2-O2S
3	A	302	GOL	O2-C2-C3-O3
2	F	301	GTS	O31-C3-CA3-N3
2	D	301	GTS	CA2-CB2-SG2-O1S
2	C	301	GTS	CA2-CB2-SG2-O3S
2	F	301	GTS	O32-C3-CA3-N3
2	A	301	GTS	CA2-CB2-SG2-O1S
2	I	301	GTS	CA2-CB2-SG2-O1S
2	I	301	GTS	O12-C1-CA1-N1
2	J	301	GTS	O12-C1-CA1-N1
2	F	301	GTS	C3-CA3-N3-C2
2	I	301	GTS	C3-CA3-N3-C2
2	A	301	GTS	C3-CA3-N3-C2
2	B	301	GTS	O12-C1-CA1-N1
2	C	301	GTS	O12-C1-CA1-N1
2	H	301	GTS	C3-CA3-N3-C2
2	B	301	GTS	C3-CA3-N3-C2
2	D	301	GTS	O12-C1-CA1-N1
2	D	301	GTS	O11-C1-CA1-N1

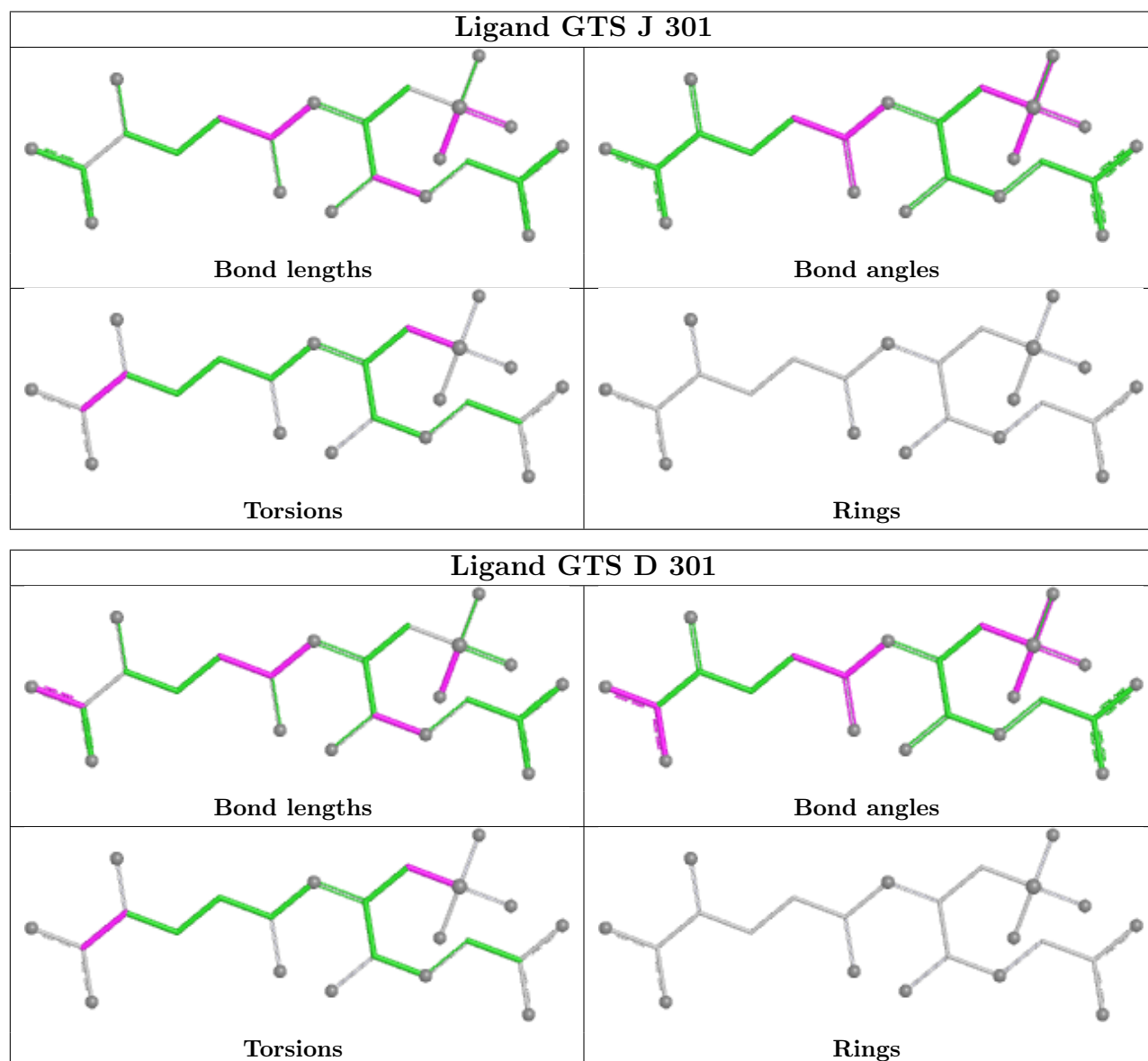
There are no ring outliers.

3 monomers are involved in 3 short contacts:

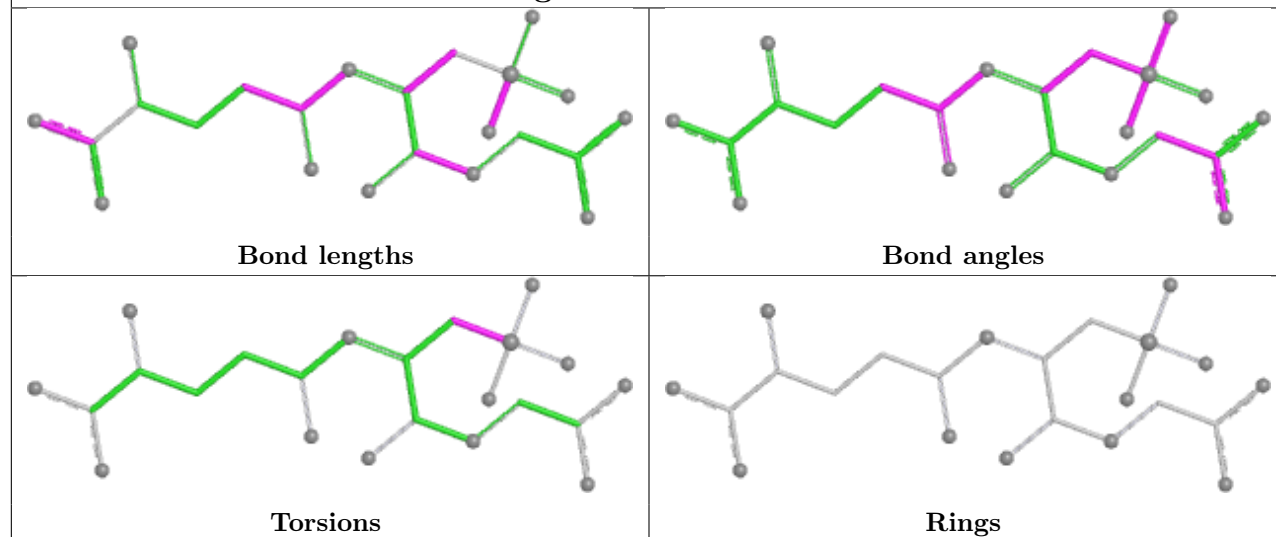
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	303	GOL	1	0
3	A	302	GOL	1	0
3	A	303	GOL	1	0

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

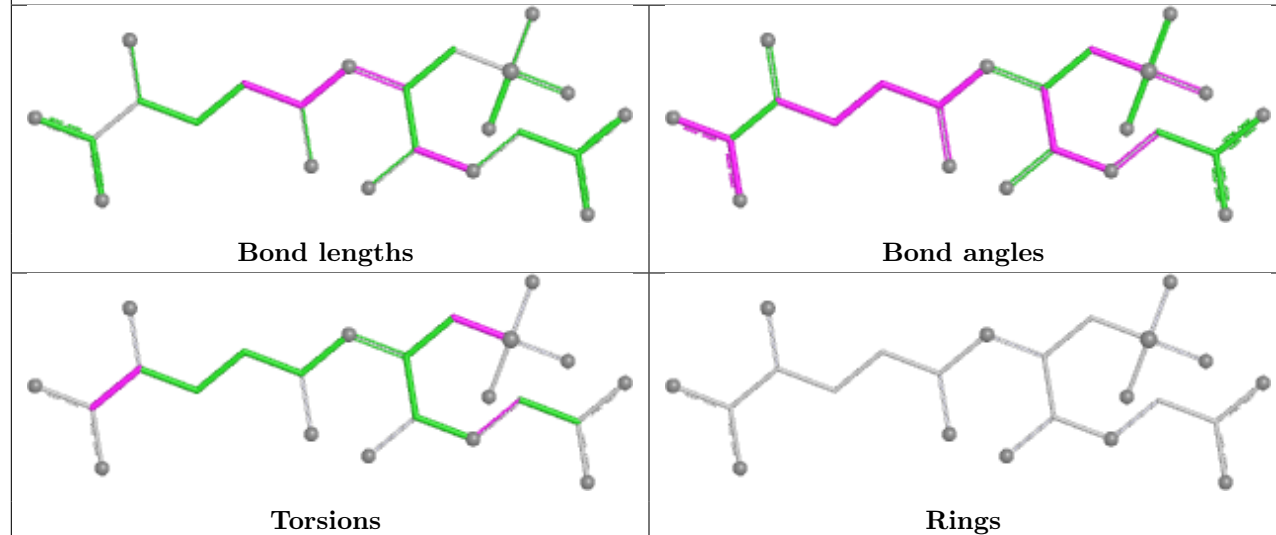
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.



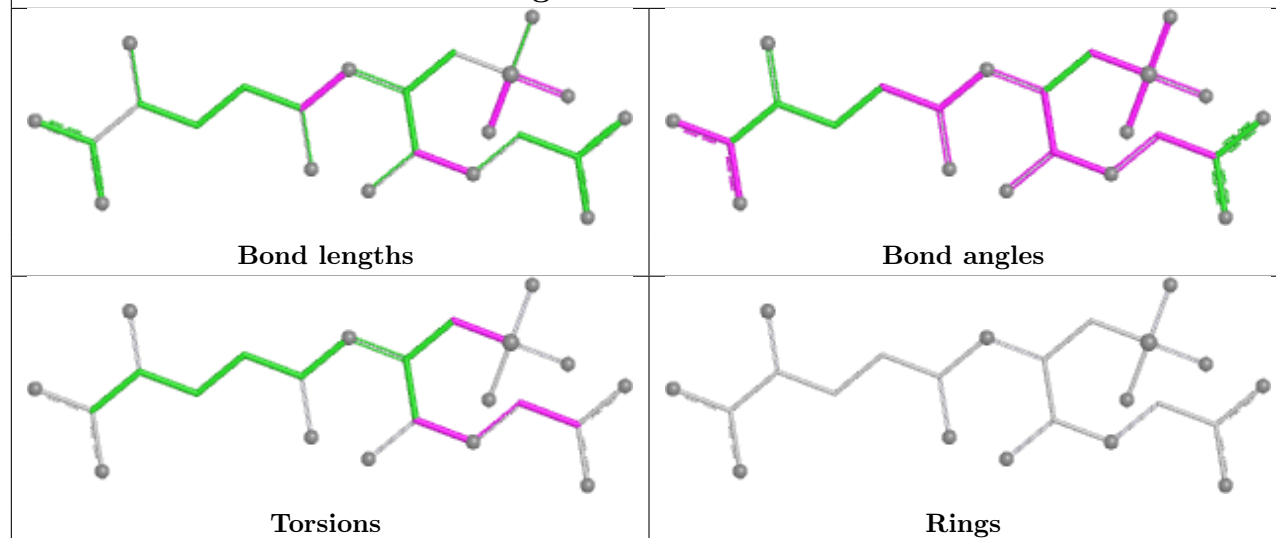
Ligand GTS E 301



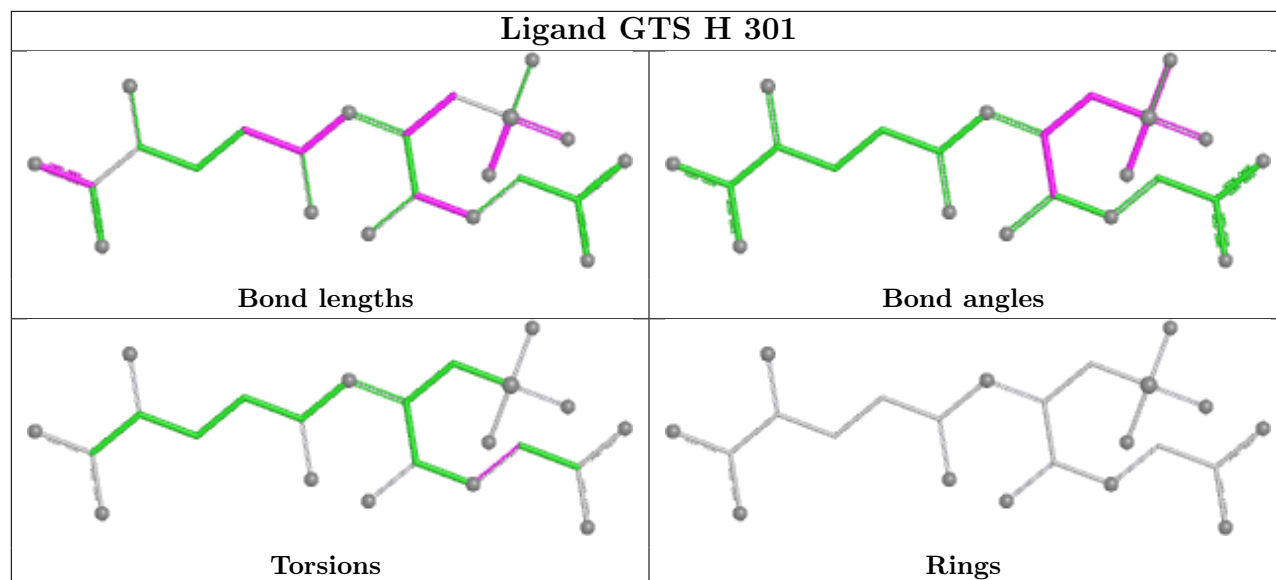
Ligand GTS B 301



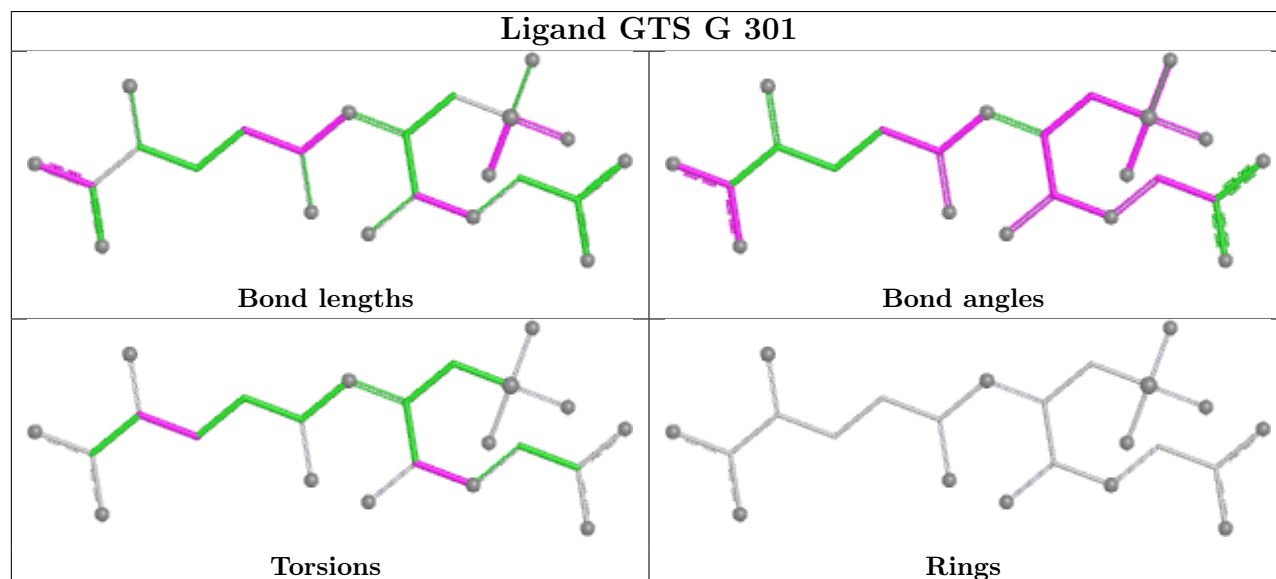
Ligand GTS A 301



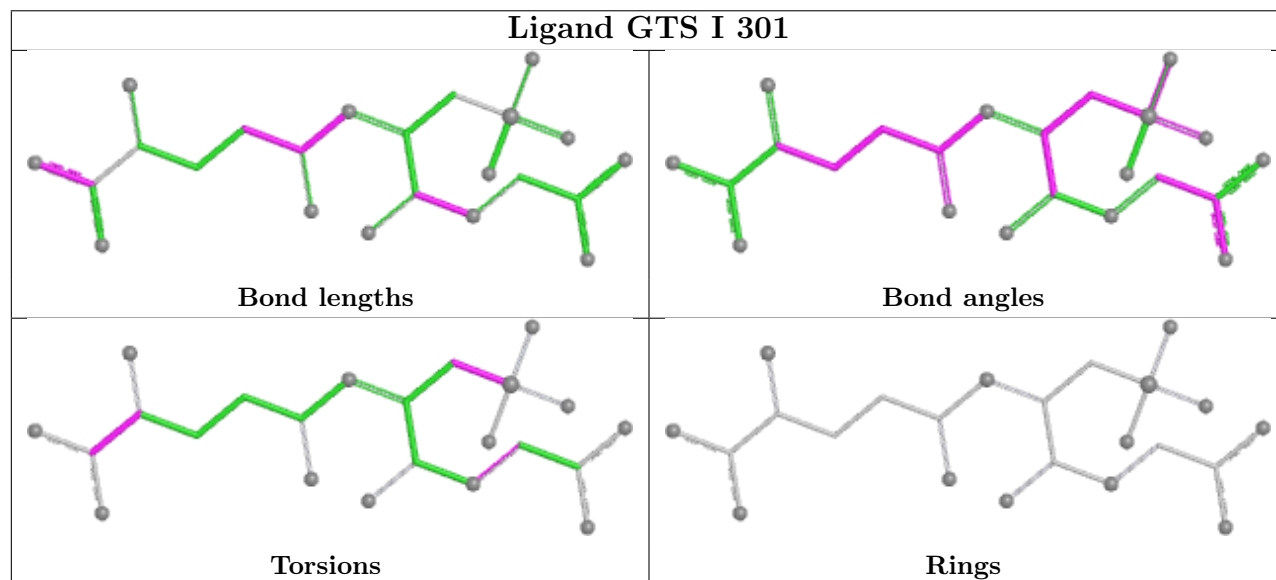
Ligand GTS H 301

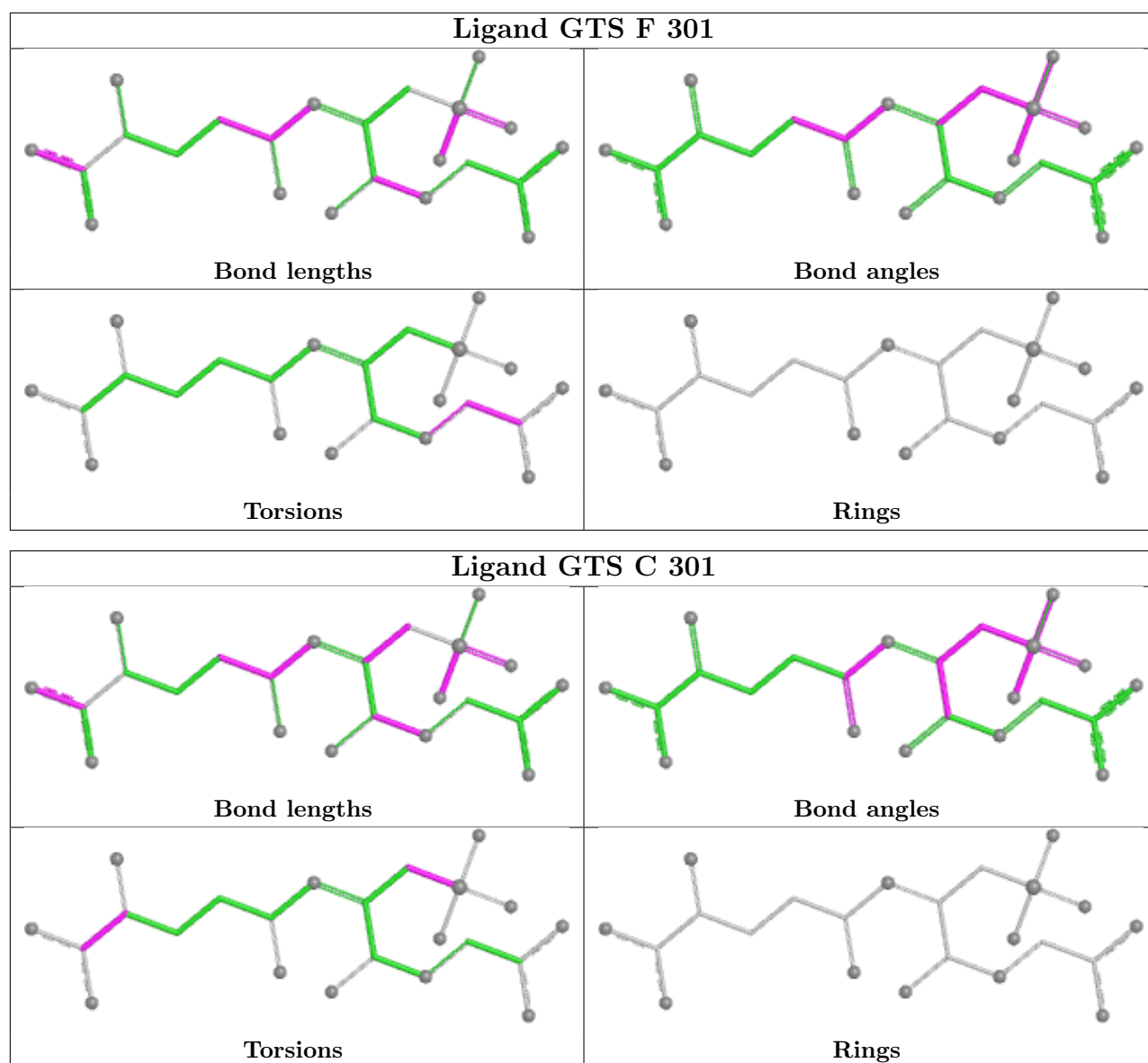


Ligand GTS G 301



Ligand GTS I 301





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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	204/215 (94%)	-0.57	4 (1%) 65 62	22, 32, 54, 91	0
1	B	205/215 (95%)	-0.60	2 (0%) 79 77	21, 30, 62, 97	3 (1%)
1	C	202/215 (93%)	-0.44	4 (1%) 65 62	23, 35, 62, 89	0
1	D	199/215 (92%)	-0.07	2 (1%) 79 77	29, 55, 70, 78	0
1	E	201/215 (93%)	-0.35	4 (1%) 65 62	23, 37, 74, 99	0
1	F	203/215 (94%)	-0.09	4 (1%) 65 62	27, 49, 72, 102	1 (0%)
1	G	198/215 (92%)	0.08	1 (0%) 87 85	38, 57, 76, 87	1 (0%)
1	H	201/215 (93%)	0.09	5 (2%) 58 55	34, 52, 71, 86	0
1	I	206/215 (95%)	-0.40	3 (1%) 72 69	24, 39, 57, 86	2 (0%)
1	J	193/215 (89%)	0.68	11 (5%) 29 26	52, 77, 97, 102	0
All	All	2012/2150 (93%)	-0.17	40 (1%) 65 62	21, 46, 81, 102	7 (0%)

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	213	PHE	6.1
1	A	213	PHE	5.4
1	D	213	PHE	5.1
1	H	174	TYR	4.7
1	E	213	PHE	4.6
1	H	213	PHE	4.3
1	H	131	LEU	4.1
1	F	213	PHE	3.9
1	A	128	ASP	3.6
1	E	117	ILE	3.4
1	I	120	ALA	3.2
1	I	213	PHE	3.2
1	C	129	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	116	HIS	2.8
1	A	119	PHE	2.8
1	I	119	PHE	2.8
1	C	117	ILE	2.7
1	E	2	VAL	2.7
1	F	2	VAL	2.7
1	E	60	ASP	2.7
1	J	2	VAL	2.7
1	F	117	ILE	2.6
1	J	79	SER	2.6
1	H	132	ILE	2.6
1	F	41	GLN	2.6
1	J	159	ASP	2.6
1	J	174	TYR	2.5
1	A	116	HIS	2.4
1	J	114	THR	2.3
1	J	212	SER	2.3
1	B	118	MET	2.3
1	D	2	VAL	2.3
1	G	42	PRO	2.2
1	J	37	GLY	2.2
1	J	28	GLU	2.2
1	J	9	HIS	2.2
1	H	2	VAL	2.1
1	J	110	LEU	2.1
1	B	2	VAL	2.1
1	J	166	LEU	2.0

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,

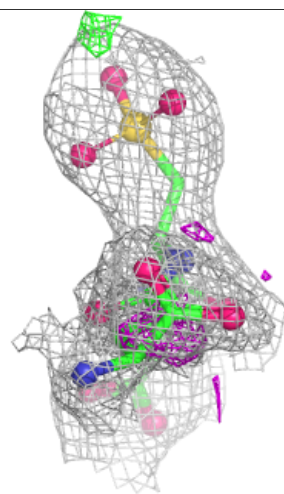
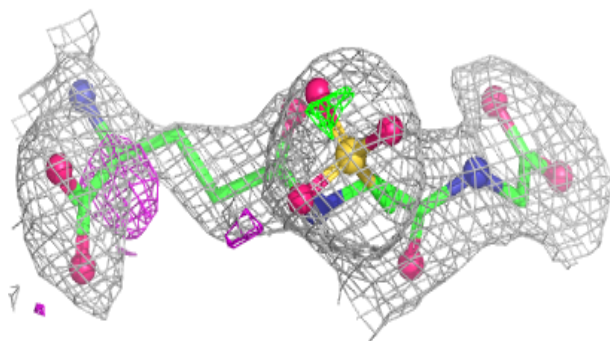
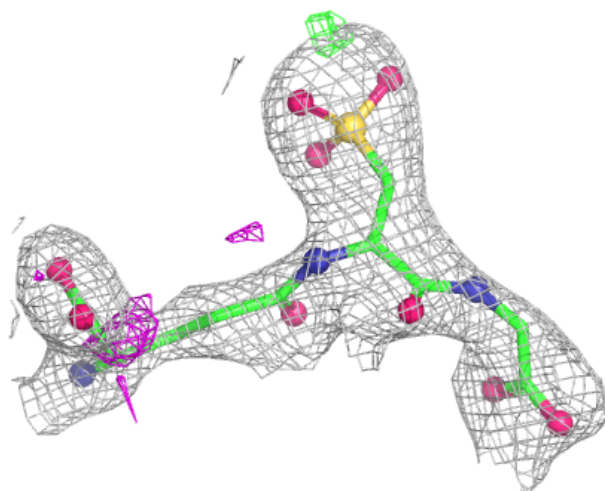
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	CL	G	302	1/1	0.82	0.24	130,130,130,130	0
3	GOL	A	303	6/6	0.84	0.15	50,66,68,75	0
3	GOL	B	303	6/6	0.86	0.13	56,60,61,61	0
2	GTS	J	301	23/23	0.90	0.11	42,71,76,80	0
2	GTS	G	301	23/23	0.90	0.11	50,65,101,104	0
3	GOL	B	302	6/6	0.91	0.12	68,74,78,83	0
4	CL	D	302	1/1	0.93	0.15	48,48,48,48	0
3	GOL	A	302	6/6	0.93	0.10	55,63,65,68	0
2	GTS	E	301	23/23	0.95	0.09	30,47,80,85	0
2	GTS	F	301	23/23	0.95	0.08	45,53,62,67	0
2	GTS	H	301	23/23	0.96	0.07	40,43,53,55	0
2	GTS	A	301	23/23	0.96	0.07	28,34,47,58	0
4	CL	H	302	1/1	0.96	0.14	48,48,48,48	0
2	GTS	B	301	23/23	0.97	0.06	23,28,43,48	0
2	GTS	C	301	23/23	0.97	0.06	30,37,59,65	0
2	GTS	D	301	23/23	0.97	0.06	25,36,47,51	0
4	CL	C	302	1/1	0.98	0.05	44,44,44,44	0
2	GTS	I	301	23/23	0.98	0.06	22,30,49,54	0
4	CL	E	302	1/1	0.98	0.05	52,52,52,52	0
4	CL	F	302	1/1	0.98	0.05	50,50,50,50	0
4	CL	A	304	1/1	0.98	0.05	35,35,35,35	0
4	CL	B	304	1/1	0.98	0.07	34,34,34,34	0
4	CL	I	302	1/1	0.99	0.07	36,36,36,36	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.

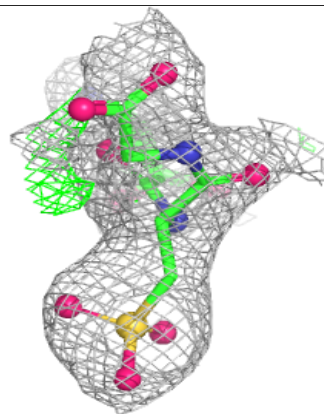
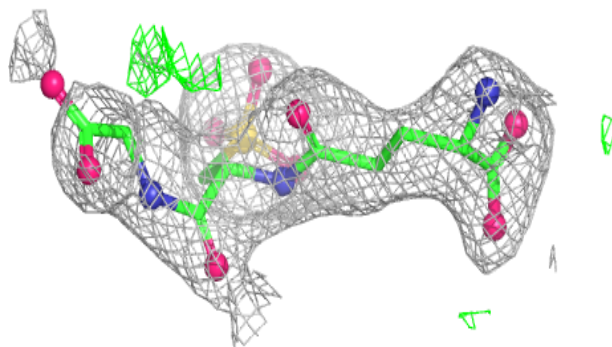
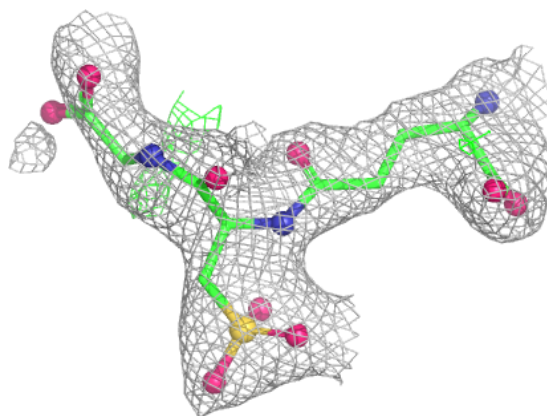
Electron density around GTS J 301:

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



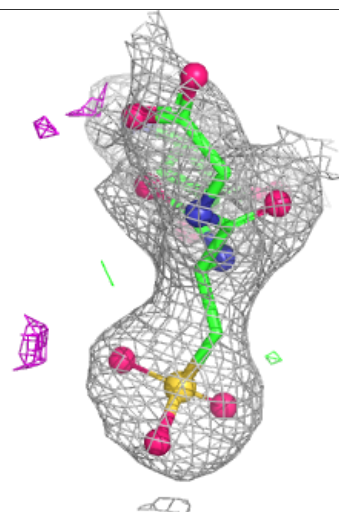
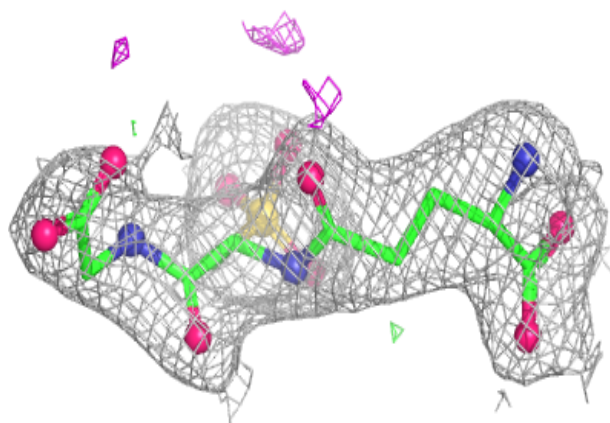
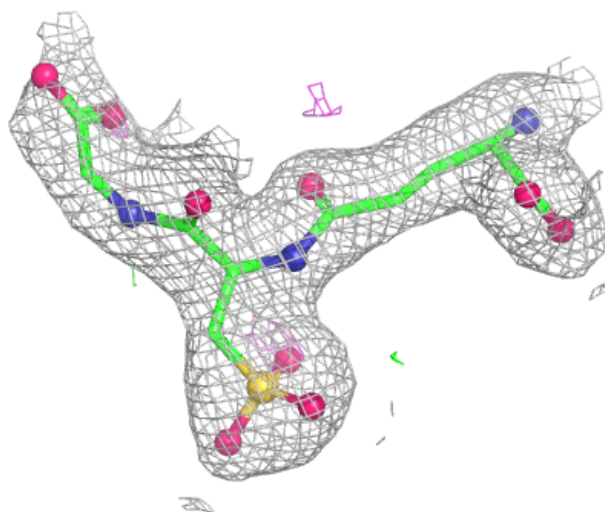
Electron density around GTS G 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



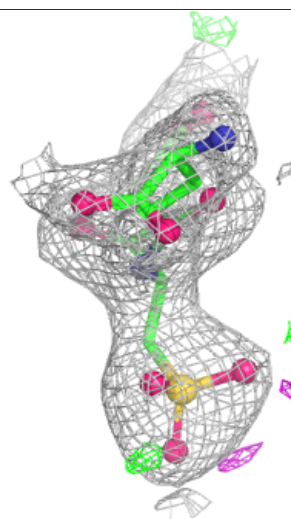
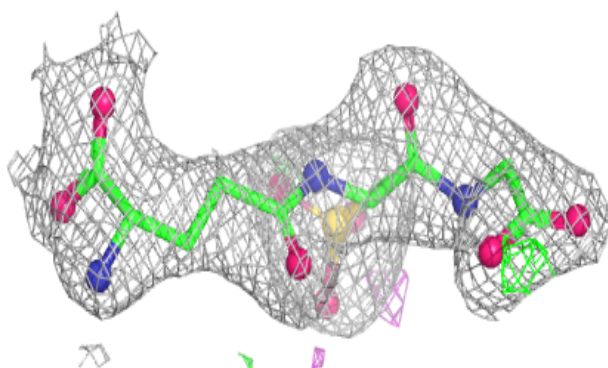
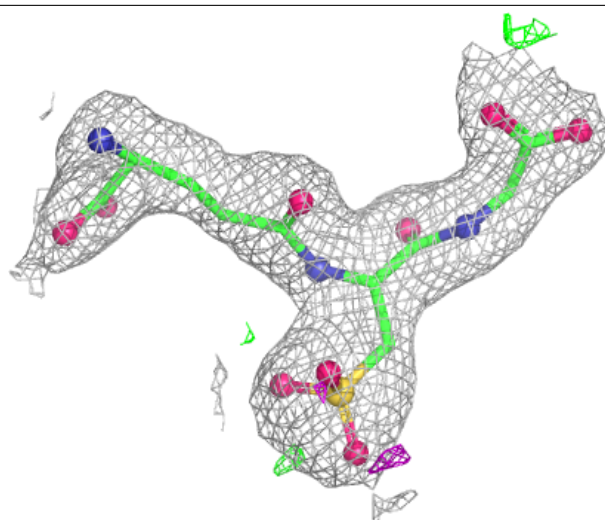
Electron density around GTS E 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



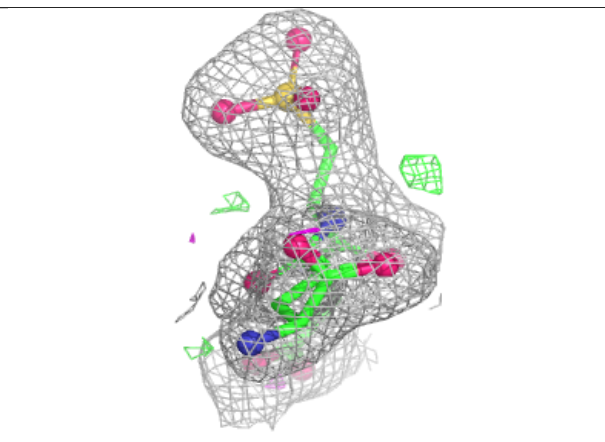
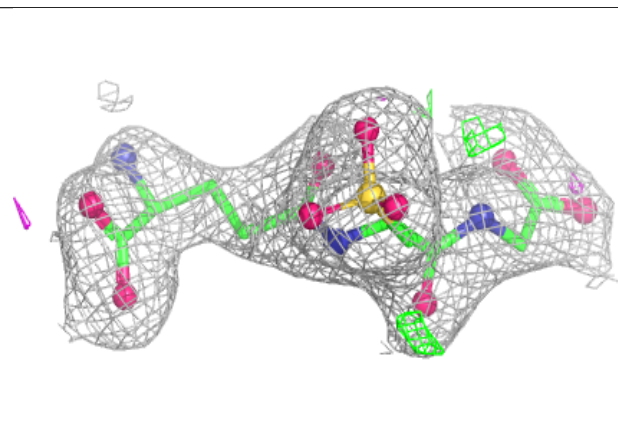
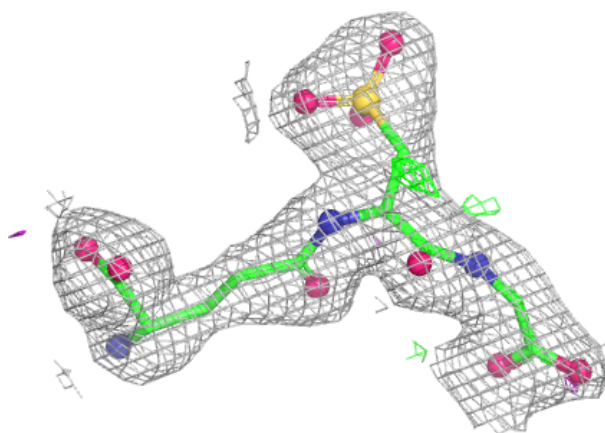
Electron density around GTS F 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



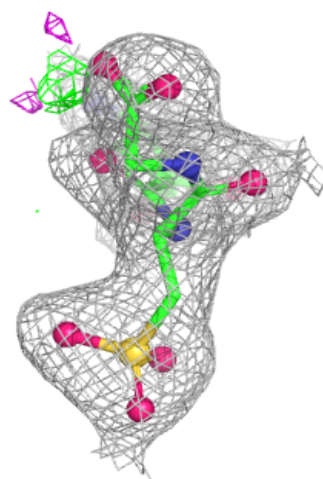
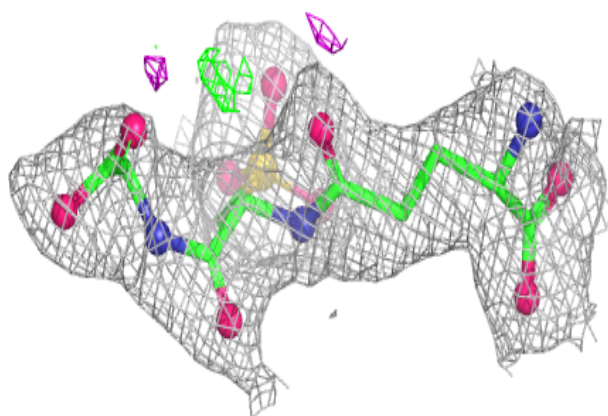
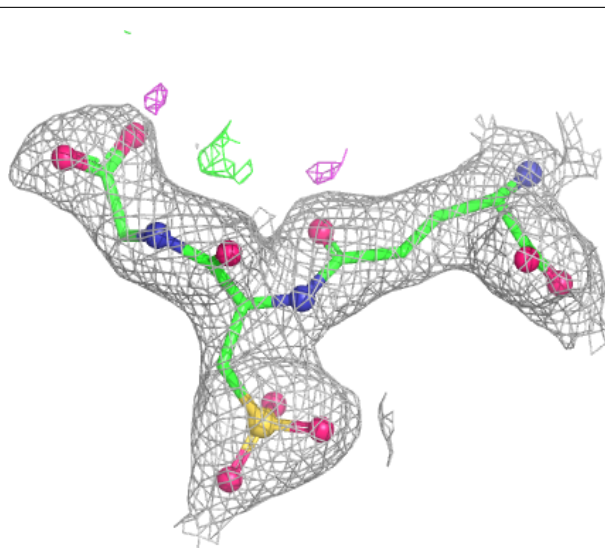
Electron density around GTS H 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



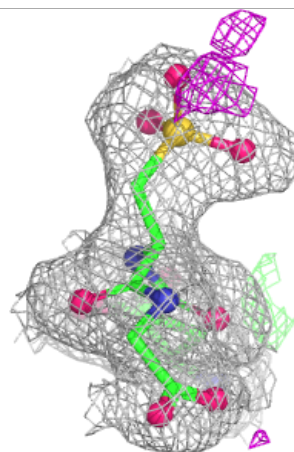
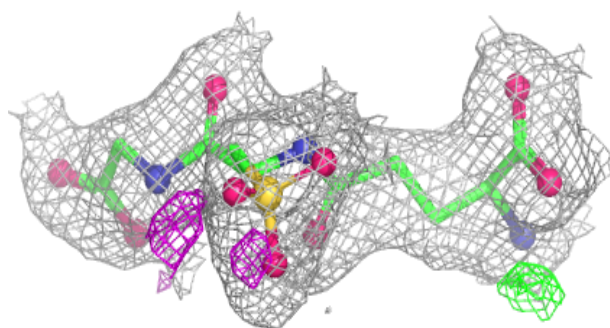
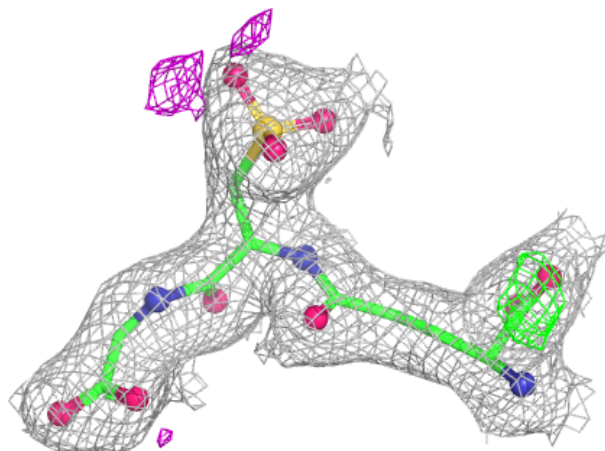
Electron density around GTS A 301:

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and green (positive)



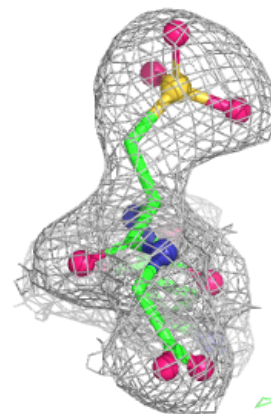
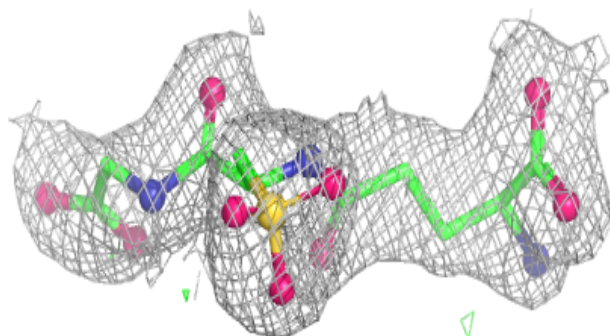
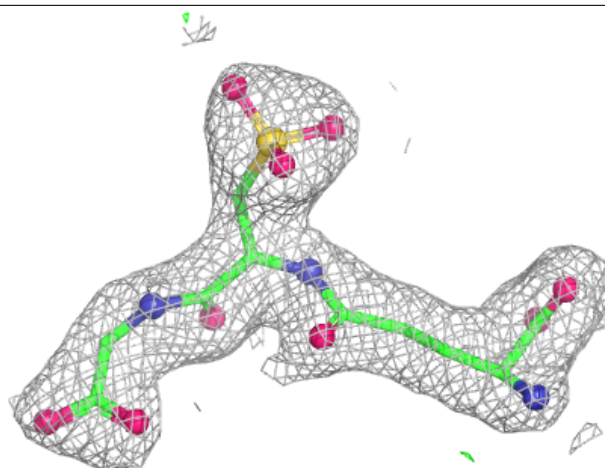
Electron density around GTS B 301:

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



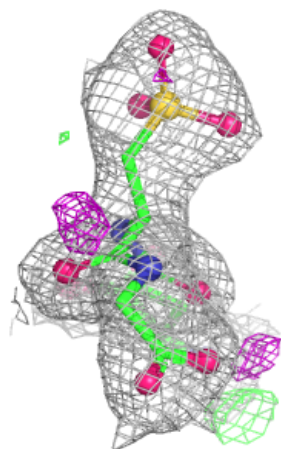
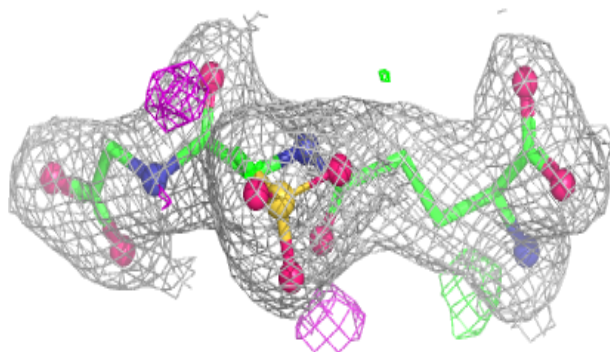
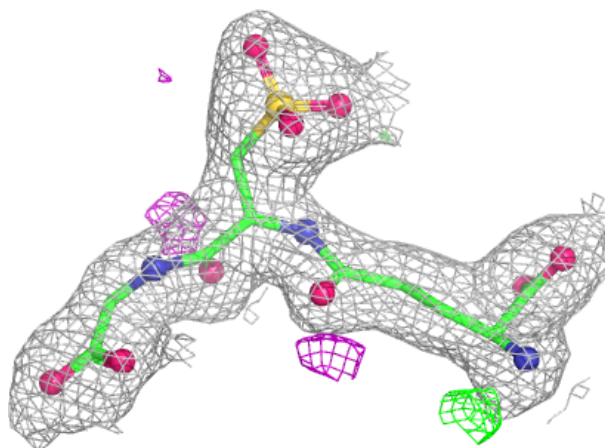
Electron density around GTS C 301:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



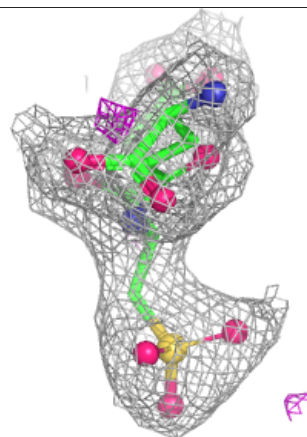
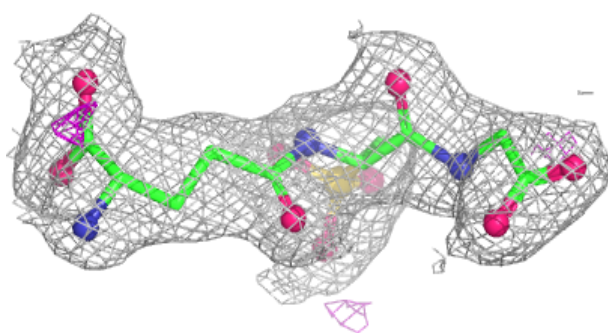
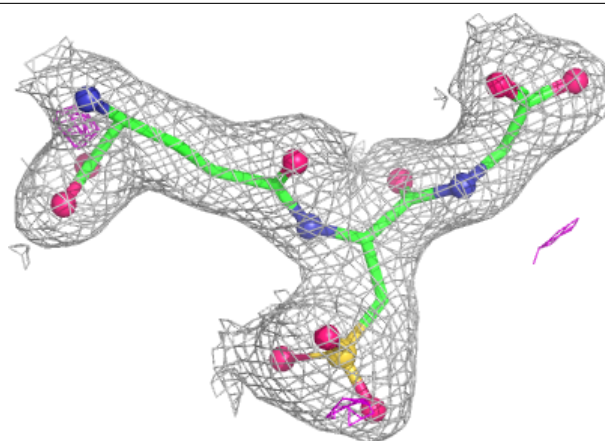
Electron density around GTS D 301:

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



Electron density around GTS I 301:

$2mF_o - DF_c$ (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.