



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2022 – 03:08 PM EDT

PDB ID : 7SO8  
Title : Crystal structure of Glutathione S-Transferase from Shrimp *Litopenaeus vannamei* in complex with silver ions and a molecules of Glutathione binding in G-site and H-site  
Authors : Escudero-Garcia, A.; Rudino-Pinera, E.; Miranda-Blancas, R.  
Deposited on : 2021-10-29  
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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.2
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)	:	2.31.2

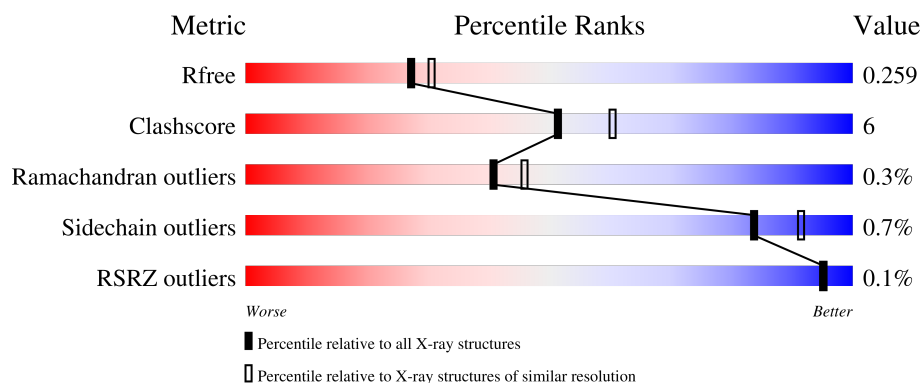
# 1 Overall quality at a glance

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))
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (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  $\geq 3$ , 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	219	 89% 11%
1	B	219	 87% 13%
1	C	219	 90% 9%
1	D	219	 87% 12% .

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Mol	Chain	Length	Quality of chain
1	E	219	 83% 16% .
1	F	219	 87% 13%
1	G	219	 86% 13%
1	H	219	 82% 18%

## 2 Entry composition

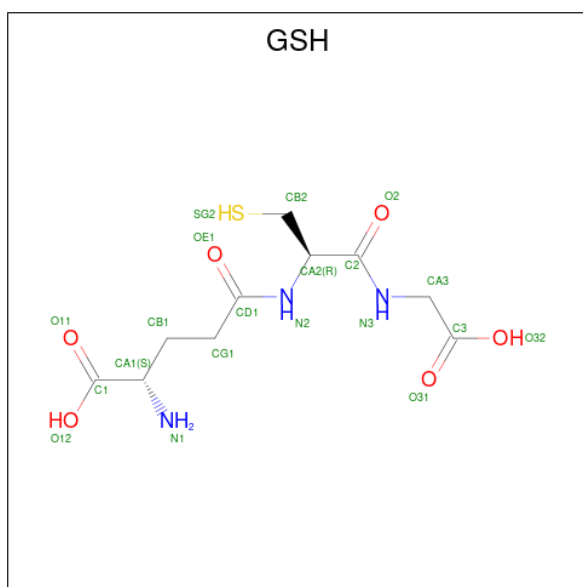
There are 5 unique types of molecules in this entry. The entry contains 15875 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 transferase.

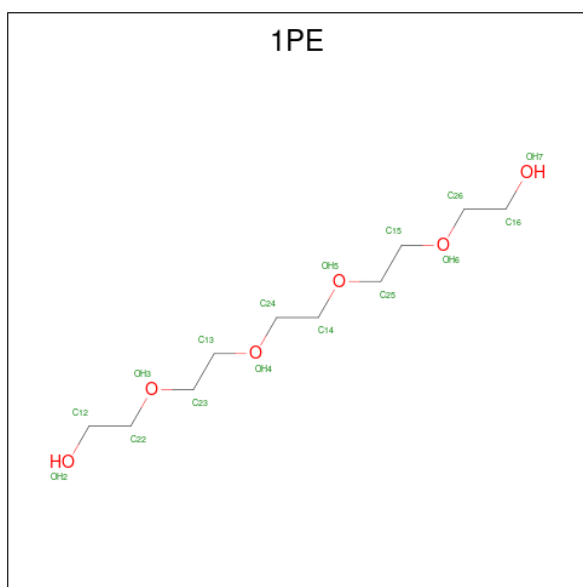
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	0	1	0
			1809	1178	292	328	11			
1	C	219	Total	C	N	O	S	0	2	0
			1818	1183	293	331	11			
1	E	219	Total	C	N	O	S	0	2	0
			1820	1184	296	329	11			
1	G	219	Total	C	N	O	S	0	1	0
			1809	1178	292	328	11			
1	H	219	Total	C	N	O	S	9	1	0
			1809	1178	292	328	11			
1	F	219	Total	C	N	O	S	0	1	0
			1809	1178	292	328	11			
1	B	219	Total	C	N	O	S	9	1	0
			1809	1178	292	328	11			
1	D	219	Total	C	N	O	S	0	1	0
			1811	1178	294	328	11			

- Molecule 2 is GLUTATHIONE (three-letter code: GSH) (formula: C<sub>10</sub>H<sub>17</sub>N<sub>3</sub>O<sub>6</sub>S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	C	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	E	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	G	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	H	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	F	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	F	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	B	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	B	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
2	D	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

- Molecule 3 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C<sub>10</sub>H<sub>22</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			16	10	6		

- Molecule 4 is SILVER ION (three-letter code: AG) (formula: Ag) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	1	Total	Ag	0	0
			1	1		
4	G	1	Total	Ag	0	0
			1	1		
4	D	1	Total	Ag	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	128	Total	O	0	0
			128	128		
5	C	178	Total	O	0	0
			178	178		
5	E	152	Total	O	0	0
			152	152		
5	G	145	Total	O	0	0
			145	145		
5	H	154	Total	O	0	0
			154	154		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	F	147	Total 147	O 147	0	0
5	B	127	Total 127	O 127	0	0
5	D	131	Total 131	O 131	0	0

### 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.

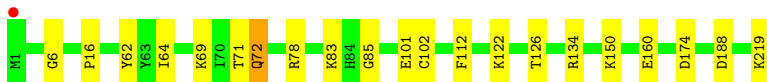
- Molecule 1: Glutathione transferase

Chain A:  89% 11%




- Molecule 1: Glutathione transferase

Chain C:  90% 9%



- Molecule 1: Glutathione transferase

Chain E:  83% 16%




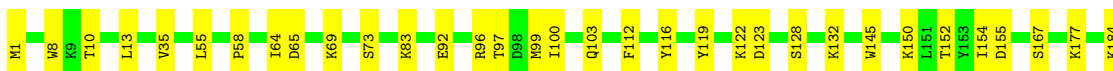
- Molecule 1: Glutathione transferase

Chain G:  86% 13%



- Molecule 1: Glutathione transferase

Chain H:  82% 18%







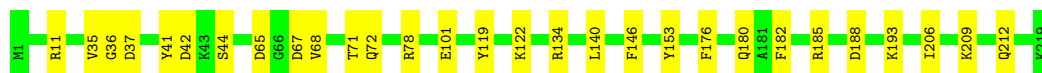
- Molecule 1: Glutathione transferase

Chain F: 87% 13%



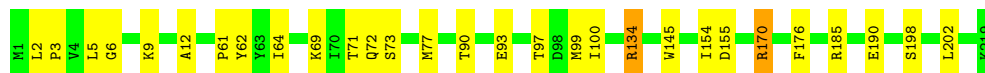
- Molecule 1: Glutathione transferase

Chain B: 87% 13%



- Molecule 1: Glutathione transferase

Chain D: 87% 12%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.39Å 93.02Å 169.11Å 90.00° 90.75° 90.00°	Depositor
Resolution (Å)	56.37 – 2.20 56.37 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.2 (56.37-2.20) 98.2 (56.37-2.20)	Depositor EDS
$R_{merge}$	0.62	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.68 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.191 , 0.260 0.191 , 0.259	Depositor DCC
$R_{free}$ test set	4472 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	13.8	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 44.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.039 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	15875	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 28.09 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.9581e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: GSH, 1PE, AG

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	0.44	0/1858	0.62	0/2507
1	B	0.41	0/1858	0.57	0/2507
1	C	0.46	0/1867	0.58	0/2519
1	D	0.43	0/1860	0.61	0/2510
1	E	0.46	0/1869	0.61	0/2521
1	F	0.43	0/1858	0.62	0/2507
1	G	0.45	0/1858	0.59	0/2507
1	H	0.44	0/1858	0.62	0/2507
All	All	0.44	0/14886	0.60	0/20085

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	1809	0	1794	15	0
1	B	1809	0	1794	17	0
1	C	1818	0	1799	17	0
1	D	1811	0	1794	20	0
1	E	1820	0	1806	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1809	0	1794	21	0
1	G	1809	0	1794	21	0
1	H	1809	0	1794	31	0
2	A	20	0	15	2	0
2	B	40	0	30	1	0
2	C	20	0	15	1	0
2	D	20	0	14	0	0
2	E	20	0	15	2	0
2	F	40	0	30	3	0
2	G	20	0	15	0	0
2	H	20	0	15	2	0
3	C	16	0	22	4	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
5	A	128	0	0	5	0
5	B	127	0	0	2	0
5	C	178	0	0	8	0
5	D	131	0	0	4	0
5	E	152	0	0	4	0
5	F	147	0	0	5	0
5	G	145	0	0	4	0
5	H	154	0	0	9	0
All	All	15875	0	14540	164	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 (164) 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:G:5:LEU:HD12	5:G:416:HOH:O	1.48	1.12
1:E:132:LYS:HG2	5:E:546:HOH:O	1.78	0.84
1:G:121:GLN:HG2	5:G:528:HOH:O	1.79	0.83
1:G:70:ILE:HG23	1:H:99:MET:HE2	1.61	0.82
1:C:174:ASP:HB3	3:C:302:1PE:H261	1.64	0.78
3:C:302:1PE:H121	1:H:92:GLU:HG3	1.66	0.75
1:F:145:TRP:HB2	5:F:409:HOH:O	1.87	0.74
1:H:1:MET:N	5:H:403:HOH:O	2.21	0.73
1:C:134:ARG:NH2	5:C:404:HOH:O	2.26	0.69
1:C:134:ARG:NH2	5:C:403:HOH:O	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:ARG:NH2	5:C:405:HOH:O	2.27	0.68
1:A:97:THR:HA	1:A:154:ILE:HD13	1.77	0.66
1:G:11:ARG:HB2	1:G:206:ILE:HA	1.77	0.66
1:H:97:THR:HA	1:H:154:ILE:HD13	1.78	0.65
1:B:134:ARG:HG3	1:B:176:PHE:HE1	1.61	0.64
1:D:2:LEU:HD12	1:D:3:PRO:HD2	1.80	0.63
1:D:202:LEU:HB2	5:D:506:HOH:O	1.99	0.63
1:A:55:LEU:HD22	1:A:71:THR:HG23	1.82	0.61
2:B:302:GSH:HA31	2:B:302:GSH:HB23	1.81	0.61
1:E:203:LYS:HE3	5:E:466:HOH:O	2.01	0.61
1:B:78:ARG:NH2	5:B:405:HOH:O	2.32	0.60
1:E:170[A]:ARG:HD3	5:E:474:HOH:O	2.00	0.60
1:H:132:LYS:HE2	5:H:435:HOH:O	2.02	0.60
1:F:145:TRP:CH2	1:F:185:ARG:HG2	2.37	0.59
1:E:118:HIS:HB3	1:E:121:GLN:HE21	1.68	0.59
1:G:145:TRP:CH2	1:G:185:ARG:HG2	2.38	0.59
1:D:134[A]:ARG:NH2	5:D:404:HOH:O	2.37	0.58
1:C:64:ILE:HD12	1:C:69:LYS:HG2	1.86	0.58
1:C:174:ASP:HB3	3:C:302:1PE:C26	2.32	0.58
1:H:119:TYR:CZ	1:H:122:LYS:HD2	2.38	0.58
1:A:140:LEU:HB2	1:A:146:PHE:CE1	2.39	0.58
1:B:65:ASP:HB3	1:B:67:ASP:H	1.69	0.58
1:C:83:LYS:NZ	5:C:411:HOH:O	2.36	0.57
1:H:150:LYS:HE2	5:F:477:HOH:O	2.03	0.57
1:B:188:ASP:HA	1:B:193:LYS:HE2	1.86	0.56
1:F:184[B]:LYS:NZ	5:F:407:HOH:O	2.36	0.56
1:E:13:LEU:HD11	1:E:61:PRO:HD3	1.88	0.56
1:G:194:LYS:NZ	5:G:403:HOH:O	2.35	0.56
1:H:55:LEU:HD21	1:H:64:ILE:HD11	1.87	0.56
1:A:78:ARG:NH2	5:A:411:HOH:O	2.39	0.55
1:G:134:ARG:HG3	1:G:134:ARG:HH11	1.70	0.55
1:H:64:ILE:HG23	1:H:69:LYS:HG2	1.88	0.55
1:F:208:ASN:HB2	2:F:301:GSH:HA31	1.87	0.55
1:G:35:VAL:HB	1:G:209:LYS:HD3	1.87	0.55
1:D:2:LEU:HD12	1:D:3:PRO:CD	2.36	0.55
1:D:190:GLU:HG2	5:D:439:HOH:O	2.06	0.55
1:C:134:ARG:HH11	1:C:134:ARG:HG2	1.72	0.54
1:D:97:THR:HA	1:D:154:ILE:HD13	1.90	0.54
1:E:60:LEU:HB2	2:E:302:GSH:HG12	1.90	0.53
1:D:145:TRP:CH2	1:D:185:ARG:HG2	2.44	0.53
1:G:7:TYR:O	1:G:32:ASN:HA	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:8:TRP:HE1	2:H:301:GSH:HA2	1.75	0.52
1:H:58:PRO:HG2	5:H:514:HOH:O	2.09	0.52
1:A:134:ARG:NH1	5:A:406:HOH:O	2.32	0.52
1:B:182:PHE:HA	1:B:185:ARG:NH1	2.25	0.51
1:E:29:GLU:HB3	1:F:199:PRO:HG2	1.93	0.51
1:E:194:LYS:HD2	1:E:195:TYR:N	2.26	0.51
1:D:100:ILE:HG21	1:D:154:ILE:HD12	1.92	0.51
1:E:97:THR:HA	1:E:154:ILE:HD13	1.93	0.51
1:E:65:ASP:HB3	1:E:67:ASP:H	1.75	0.51
1:H:35:VAL:HB	1:H:209:LYS:HD3	1.94	0.50
1:A:188:ASP:HA	1:A:193:LYS:HE2	1.93	0.50
1:D:64:ILE:HG23	1:D:69:LYS:HG2	1.93	0.50
1:C:85:GLY:O	1:C:150:LYS:NZ	2.45	0.50
1:B:11:ARG:HB2	1:B:206:ILE:HA	1.93	0.50
1:C:16:PRO:HD3	1:C:160:GLU:HG2	1.94	0.50
1:G:97:THR:HA	1:G:154:ILE:HD13	1.93	0.49
1:C:219:LYS:O	5:C:401:HOH:O	2.20	0.49
1:H:167:SER:O	1:H:219:LYS:NZ	2.36	0.49
1:H:123:ASP:OD1	5:H:401:HOH:O	2.20	0.49
1:E:7:TYR:O	1:E:32:ASN:HA	2.13	0.49
1:H:177:LYS:HG2	5:H:552:HOH:O	2.12	0.48
1:B:36:GLY:O	1:B:41:TYR:HA	2.12	0.48
1:G:89:THR:OG1	1:G:150:LYS:NZ	2.45	0.48
3:C:302:1PE:H121	1:H:92:GLU:CG	2.40	0.48
1:E:31:LYS:NZ	5:E:409:HOH:O	2.40	0.48
1:E:64:ILE:HG23	1:E:69:LYS:HG2	1.96	0.47
1:H:96:ARG:NH2	5:H:417:HOH:O	2.46	0.47
1:H:145:TRP:HB3	1:H:155:ASP:OD2	2.14	0.47
1:F:166:LEU:HD21	1:F:173:LEU:HG	1.96	0.47
1:A:13:LEU:HD13	1:A:73:SER:HB3	1.96	0.47
1:H:65:ASP:OD2	5:H:402:HOH:O	2.21	0.47
1:G:57:PHE:CD2	1:H:103:GLN:HG2	2.50	0.47
1:F:148:GLY:N	5:F:409:HOH:O	2.46	0.46
1:G:6:GLY:O	1:G:61:PRO:HA	2.15	0.46
1:E:104:LEU:HB3	1:E:161:ILE:HD12	1.97	0.46
1:A:196:MET:HG2	1:A:201:PHE:CE1	2.50	0.46
1:F:7:TYR:O	1:F:32:ASN:HA	2.16	0.46
1:F:64:ILE:HG12	1:F:69:LYS:HG2	1.98	0.46
1:D:6:GLY:O	1:D:61:PRO:HA	2.15	0.46
1:D:154:ILE:HG13	1:D:155:ASP:N	2.31	0.46
1:D:134[A]:ARG:HG2	1:D:176:PHE:HE1	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:134:ARG:HG3	1:G:134:ARG:NH1	2.30	0.45
1:B:140:LEU:HB2	1:B:146:PHE:CE1	2.51	0.45
1:C:6:GLY:HA3	1:C:62:TYR:CZ	2.51	0.45
1:C:112:PHE:HB3	2:C:301:GSH:HA32	1.98	0.45
1:A:11:ARG:HG2	1:A:15:GLN:OE1	2.16	0.45
1:A:95:VAL:HG21	1:B:68:VAL:HG21	1.97	0.45
1:E:33:TYR:HD2	1:E:45:GLU:HG2	1.82	0.45
2:A:301:GSH:HA31	2:A:301:GSH:HA2	1.61	0.45
1:C:101[B]:GLU:HG3	1:C:102:CYS:N	2.31	0.45
1:A:9:LYS:HE2	1:A:202:LEU:HD11	1.99	0.44
1:D:90:THR:OG1	1:D:93:GLU:HG3	2.17	0.44
1:G:184[A]:LYS:HG3	1:G:188:ASP:OD2	2.18	0.44
1:G:34:PRO:HG2	1:G:45:GLU:OE1	2.17	0.44
1:E:93:GLU:O	1:E:97:THR:HG23	2.16	0.44
1:D:5:LEU:HD12	1:D:62:TYR:O	2.17	0.44
1:G:56:ALA:HB3	5:G:406:HOH:O	2.17	0.44
1:G:204:LYS:HB2	1:G:205:PRO:HA	2.00	0.44
1:A:122:LYS:NZ	5:A:420:HOH:O	2.50	0.44
2:E:302:GSH:HA2	2:E:302:GSH:HG13	1.56	0.44
1:F:4:VAL:HA	1:F:29:GLU:O	2.18	0.44
1:B:101:GLU:OE1	1:B:153:TYR:OH	2.25	0.44
1:G:71:THR:O	1:G:72:GLN:HB2	2.18	0.43
1:F:71:THR:O	1:F:72:GLN:HB2	2.17	0.43
1:D:9:LYS:O	1:D:9:LYS:HG3	2.18	0.43
1:G:150:LYS:HE3	1:G:150:LYS:HB2	1.69	0.43
1:E:81:ALA:HB2	1:E:153:TYR:HB3	1.99	0.43
1:C:188:ASP:HB2	5:C:500:HOH:O	2.18	0.43
1:H:10:THR:HG22	1:H:205:PRO:HG2	2.00	0.43
1:A:150:LYS:HE2	5:C:436:HOH:O	2.18	0.43
1:H:100:ILE:CG2	1:H:154:ILE:HD12	2.49	0.43
1:H:112:PHE:O	1:H:116:TYR:HB2	2.18	0.43
1:F:140:LEU:HB2	1:F:146:PHE:CE1	2.54	0.43
5:A:411:HOH:O	1:B:78:ARG:HD3	2.18	0.43
1:E:108:HIS:HB2	1:E:161:ILE:HD13	2.01	0.43
1:H:128:SER:OG	1:H:132:LYS:NZ	2.49	0.43
1:B:71:THR:O	1:B:72:GLN:HB2	2.19	0.42
2:A:301:GSH:HB23	5:A:501:HOH:O	2.19	0.42
1:E:67:ASP:OD1	1:G:1:MET:HB3	2.20	0.42
1:H:119:TYR:OH	1:H:122:LYS:HD2	2.18	0.42
1:D:99:MET:HE3	1:D:99:MET:HB3	1.92	0.42
1:H:116:TYR:CE1	1:H:207:CYS:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:122:LYS:O	1:C:126:THR:HG23	2.19	0.42
1:B:42:ASP:OD1	1:B:44:SER:OG	2.36	0.42
1:A:71:THR:O	1:A:72:GLN:HB2	2.20	0.42
1:D:71:THR:O	1:D:72:GLN:HB2	2.19	0.42
1:D:73:SER:O	1:D:77:MET:HG3	2.19	0.42
1:B:37:ASP:OD1	1:B:209:LYS:NZ	2.51	0.42
1:H:152:THR:OG1	1:H:154:ILE:HG12	2.19	0.42
5:C:494:HOH:O	1:F:193:LYS:HE3	2.20	0.42
1:F:97:THR:HA	1:F:154:ILE:HD13	2.02	0.42
1:B:35:VAL:HB	1:B:209:LYS:HD2	2.01	0.42
1:E:71:THR:O	1:E:72:GLN:HB2	2.20	0.41
1:E:191:ALA:O	1:E:194:LYS:HE3	2.19	0.41
1:E:6:GLY:HA3	1:E:62:TYR:CZ	2.56	0.41
1:B:180:GLN:HG2	5:B:479:HOH:O	2.19	0.41
1:D:202:LEU:HD13	5:D:506:HOH:O	2.21	0.41
1:H:202:LEU:HB2	5:H:428:HOH:O	2.20	0.41
1:F:40:ASP:OD1	1:F:40:ASP:N	2.53	0.41
1:F:193:LYS:HE2	1:F:193:LYS:HB2	1.73	0.41
1:H:13:LEU:HD13	1:H:73:SER:HB3	2.02	0.41
1:F:9:LYS:NZ	5:F:403:HOH:O	2.28	0.41
5:H:459:HOH:O	1:F:150:LYS:HE2	2.20	0.41
1:F:5:LEU:O	1:F:30:GLU:HA	2.21	0.41
1:H:116:TYR:CZ	2:H:301:GSH:HG13	2.55	0.40
1:H:145:TRP:CH2	1:H:185:ARG:HG2	2.56	0.40
1:C:71:THR:O	1:C:72:GLN:HB2	2.20	0.40
1:E:16:PRO:O	1:E:20:MET:HG3	2.20	0.40
1:F:208:ASN:HD22	2:F:301:GSH:CA3	2.34	0.40
1:D:134[A]:ARG:HG2	1:D:176:PHE:CE1	2.56	0.40
1:F:73:SER:H	2:F:302:GSH:HN11	1.69	0.40
1:A:36:GLY:O	1:A:41:TYR:HA	2.21	0.40
1:B:119:TYR:HB2	1:B:212:GLN:HB2	2.02	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	218/219 (100%)	214 (98%)	3 (1%)	1 (0%)	29	31
1	B	218/219 (100%)	213 (98%)	5 (2%)	0	100	100
1	C	219/219 (100%)	211 (96%)	7 (3%)	1 (0%)	29	31
1	D	218/219 (100%)	211 (97%)	5 (2%)	2 (1%)	17	16
1	E	219/219 (100%)	212 (97%)	6 (3%)	1 (0%)	29	31
1	F	218/219 (100%)	212 (97%)	5 (2%)	1 (0%)	29	31
1	G	218/219 (100%)	212 (97%)	6 (3%)	0	100	100
1	H	218/219 (100%)	210 (96%)	8 (4%)	0	100	100
All	All	1746/1752 (100%)	1695 (97%)	45 (3%)	6 (0%)	41	46

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	170	ARG
1	A	12	ALA
1	E	72	GLN
1	F	12	ALA
1	D	12	ALA
1	C	72	GLN

### 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	192/191 (100%)	192 (100%)	0	100	100
1	B	192/191 (100%)	191 (100%)	1 (0%)	88	94
1	C	193/191 (101%)	193 (100%)	0	100	100
1	D	192/191 (100%)	188 (98%)	4 (2%)	53	67
1	E	193/191 (101%)	190 (98%)	3 (2%)	62	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	192/191 (100%)	192 (100%)	0	100	100
1	G	192/191 (100%)	191 (100%)	1 (0%)	88	94
1	H	192/191 (100%)	189 (98%)	3 (2%)	62	76
All	All	1538/1528 (101%)	1526 (99%)	12 (1%)	84	90

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

Mol	Chain	Res	Type
1	E	2	LEU
1	E	194	LYS
1	E	200	LYS
1	G	204	LYS
1	H	83	LYS
1	H	184[A]	LYS
1	H	184[B]	LYS
1	B	122	LYS
1	D	134[A]	ARG
1	D	134[B]	ARG
1	D	170	ARG
1	D	198	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	121	GLN
1	G	59	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry

Of 14 ligands modelled in this entry, 3 are monoatomic - leaving 11 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	GSH	F	302	-	18,19,19	0.71	1 (5%)	23,24,24	0.99	1 (4%)
3	1PE	C	302	-	15,15,15	1.07	1 (6%)	14,14,14	0.92	0
2	GSH	G	302	-	18,19,19	0.76	1 (5%)	23,24,24	0.78	0
2	GSH	B	301	-	18,19,19	0.84	1 (5%)	23,24,24	0.65	0
2	GSH	B	302	-	18,19,19	2.33	3 (16%)	23,24,24	2.87	10 (43%)
2	GSH	H	301	-	18,19,19	0.74	0	23,24,24	0.72	0
2	GSH	C	301	-	18,19,19	0.69	0	23,24,24	1.18	1 (4%)
2	GSH	E	302	-	18,19,19	0.78	0	23,24,24	0.74	0
2	GSH	F	301	-	18,19,19	2.19	4 (22%)	23,24,24	2.26	7 (30%)
2	GSH	D	302	-	18,19,19	0.79	1 (5%)	23,24,24	0.83	0
2	GSH	A	301	-	18,19,19	0.71	0	23,24,24	0.82	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GSH	F	302	-	-	6/24/24/24	-
3	1PE	C	302	-	-	8/13/13/13	-
2	GSH	G	302	-	-	15/24/24/24	-
2	GSH	B	301	-	-	3/24/24/24	-
2	GSH	B	302	-	-	10/24/24/24	-
2	GSH	H	301	-	-	8/24/24/24	-
2	GSH	C	301	-	-	13/24/24/24	-
2	GSH	E	302	-	-	12/24/24/24	-
2	GSH	F	301	-	-	6/24/24/24	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GSH	D	302	-	-	14/24/24/24	-
2	GSH	A	301	-	-	14/24/24/24	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	302	GSH	C2-N3	6.30	1.47	1.33
2	B	302	GSH	CD1-N2	6.05	1.47	1.34
2	F	301	GSH	CD1-N2	5.99	1.46	1.34
2	F	301	GSH	C2-N3	5.33	1.45	1.33
2	B	301	GSH	O12-C1	-2.68	1.21	1.30
3	C	302	1PE	C22-C12	2.30	1.61	1.49
2	D	302	GSH	O32-C3	-2.16	1.23	1.30
2	F	301	GSH	CA3-C3	2.15	1.55	1.51
2	F	302	GSH	O32-C3	-2.10	1.23	1.30
2	G	302	GSH	O32-C3	-2.06	1.23	1.30
2	B	302	GSH	O2-C2	-2.04	1.19	1.23
2	F	301	GSH	O2-C2	-2.03	1.19	1.23

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	302	GSH	CB2-CA2-N2	-8.14	99.67	111.28
2	F	301	GSH	CA2-CB2-SG2	-6.45	106.94	114.19
2	B	302	GSH	CA2-CB2-SG2	-6.02	107.43	114.19
2	B	302	GSH	CA2-N2-CD1	5.30	135.27	121.65
2	B	302	GSH	CA2-C2-N3	3.34	123.28	116.54
2	F	301	GSH	CA3-N3-C2	3.16	129.21	121.37
2	F	301	GSH	O32-C3-O31	-3.11	115.55	123.30
2	C	301	GSH	CB1-CG1-CD1	-2.99	106.35	113.04
2	F	301	GSH	O32-C3-CA3	2.88	122.74	112.74
2	B	302	GSH	CA3-N3-C2	2.83	128.38	121.37
2	B	302	GSH	O2-C2-N3	-2.80	116.99	122.99
2	F	301	GSH	C3-CA3-N3	-2.73	104.65	113.06
2	F	301	GSH	C2-CA2-N2	2.44	117.81	111.16
2	F	302	GSH	CB1-CG1-CD1	2.43	118.47	113.04
2	F	301	GSH	O2-C2-CA2	2.37	125.45	120.45
2	B	302	GSH	O12-C1-CA1	2.34	121.34	113.38
2	B	302	GSH	CB1-CG1-CD1	-2.26	108.00	113.04
2	B	302	GSH	O12-C1-O11	-2.18	119.15	124.09
2	B	302	GSH	C2-CA2-N2	2.09	116.85	111.16

There are no chirality outliers.

All (109) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	GSH	O11-C1-CA1-N1
2	A	301	GSH	CG1-CD1-N2-CA2
2	A	301	GSH	OE1-CD1-N2-CA2
2	A	301	GSH	C2-CA2-CB2-SG2
2	C	301	GSH	CG1-CD1-N2-CA2
2	C	301	GSH	OE1-CD1-N2-CA2
2	C	301	GSH	N2-CA2-CB2-SG2
2	C	301	GSH	C2-CA2-CB2-SG2
2	C	301	GSH	O31-C3-CA3-N3
2	C	301	GSH	O32-C3-CA3-N3
2	E	302	GSH	N2-CA2-CB2-SG2
2	E	302	GSH	C2-CA2-CB2-SG2
2	G	302	GSH	O11-C1-CA1-N1
2	G	302	GSH	N1-CA1-CB1-CG1
2	G	302	GSH	CG1-CD1-N2-CA2
2	G	302	GSH	OE1-CD1-N2-CA2
2	G	302	GSH	N2-CA2-CB2-SG2
2	G	302	GSH	C2-CA2-CB2-SG2
2	H	301	GSH	CG1-CD1-N2-CA2
2	H	301	GSH	OE1-CD1-N2-CA2
2	F	301	GSH	N1-CA1-CB1-CG1
2	F	301	GSH	C2-CA2-CB2-SG2
2	D	302	GSH	O11-C1-CA1-N1
2	D	302	GSH	CG1-CD1-N2-CA2
2	D	302	GSH	OE1-CD1-N2-CA2
2	D	302	GSH	O2-C2-CA2-CB2
2	D	302	GSH	N3-C2-CA2-CB2
2	D	302	GSH	N2-CA2-CB2-SG2
2	D	302	GSH	C2-CA2-CB2-SG2
2	E	302	GSH	OE1-CD1-N2-CA2
2	A	301	GSH	O2-C2-N3-CA3
2	B	302	GSH	O2-C2-N3-CA3
2	E	302	GSH	CG1-CD1-N2-CA2
3	C	302	1PE	OH4-C13-C23-OH3
2	C	301	GSH	CA2-C2-N3-CA3
2	B	302	GSH	CA2-C2-N3-CA3
2	G	302	GSH	O12-C1-CA1-N1
2	F	302	GSH	O12-C1-CA1-N1
2	D	302	GSH	O12-C1-CA1-N1
2	A	301	GSH	CA2-C2-N3-CA3

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Mol	Chain	Res	Type	Atoms
2	C	301	GSH	O2-C2-N3-CA3
3	C	302	1PE	OH6-C15-C25-OH5
2	B	301	GSH	CG1-CD1-N2-CA2
2	E	302	GSH	CA1-CB1-CG1-CD1
2	G	302	GSH	CA1-CB1-CG1-CD1
2	B	301	GSH	OE1-CD1-N2-CA2
2	A	301	GSH	O12-C1-CA1-N1
2	B	302	GSH	N2-CD1-CG1-CB1
2	B	302	GSH	OE1-CD1-CG1-CB1
2	H	301	GSH	N2-CD1-CG1-CB1
2	H	301	GSH	OE1-CD1-CG1-CB1
2	A	301	GSH	O31-C3-CA3-N3
2	A	301	GSH	O32-C3-CA3-N3
2	D	302	GSH	O31-C3-CA3-N3
2	D	302	GSH	O32-C3-CA3-N3
3	C	302	1PE	OH2-C12-C22-OH3
2	A	301	GSH	CA1-CB1-CG1-CD1
2	H	301	GSH	CA1-CB1-CG1-CD1
2	F	302	GSH	CA1-CB1-CG1-CD1
2	F	302	GSH	O11-C1-CA1-CB1
2	F	302	GSH	O12-C1-CA1-CB1
2	C	301	GSH	O2-C2-CA2-CB2
2	C	301	GSH	N3-C2-CA2-CB2
2	H	301	GSH	O2-C2-CA2-CB2
2	H	301	GSH	N3-C2-CA2-CB2
2	F	301	GSH	CB2-CA2-N2-CD1
3	C	302	1PE	OH7-C16-C26-OH6
2	G	302	GSH	N3-C2-CA2-N2
2	G	302	GSH	O2-C2-CA2-N2
2	A	301	GSH	N2-CA2-CB2-SG2
2	C	301	GSH	O2-C2-CA2-N2
2	E	302	GSH	O12-C1-CA1-N1
2	F	301	GSH	O12-C1-CA1-N1
2	C	301	GSH	N3-C2-CA2-N2
2	E	302	GSH	O11-C1-CA1-N1
2	F	301	GSH	O11-C1-CA1-N1
2	F	302	GSH	O11-C1-CA1-N1
2	G	302	GSH	C1-CA1-CB1-CG1
2	F	301	GSH	C1-CA1-CB1-CG1
3	C	302	1PE	C25-C15-OH6-C26
2	A	301	GSH	O12-C1-CA1-CB1
2	B	302	GSH	O11-C1-CA1-CB1

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Mol	Chain	Res	Type	Atoms
2	C	301	GSH	N1-CA1-CB1-CG1
2	A	301	GSH	O11-C1-CA1-CB1
2	D	302	GSH	O12-C1-CA1-CB1
3	C	302	1PE	C15-C25-OH5-C14
3	C	302	1PE	C23-C13-OH4-C24
2	G	302	GSH	O11-C1-CA1-CB1
2	G	302	GSH	O12-C1-CA1-CB1
2	B	302	GSH	O12-C1-CA1-CB1
2	D	302	GSH	O11-C1-CA1-CB1
3	C	302	1PE	C12-C22-OH3-C23
2	B	302	GSH	O2-C2-CA2-N2
2	H	301	GSH	O11-C1-CA1-N1
2	E	302	GSH	O2-C2-CA2-N2
2	E	302	GSH	O11-C1-CA1-CB1
2	B	302	GSH	N3-C2-CA2-N2
2	D	302	GSH	C1-CA1-CB1-CG1
2	F	302	GSH	N2-CA2-CB2-SG2
2	E	302	GSH	O12-C1-CA1-CB1
2	E	302	GSH	N3-C2-CA2-N2
2	B	302	GSH	C3-CA3-N3-C2
2	D	302	GSH	C3-CA3-N3-C2
2	G	302	GSH	O2-C2-CA2-CB2
2	G	302	GSH	N3-C2-CA2-CB2
2	B	302	GSH	O31-C3-CA3-N3
2	A	301	GSH	O2-C2-CA2-N2
2	E	302	GSH	N1-CA1-CB1-CG1
2	B	301	GSH	O32-C3-CA3-N3

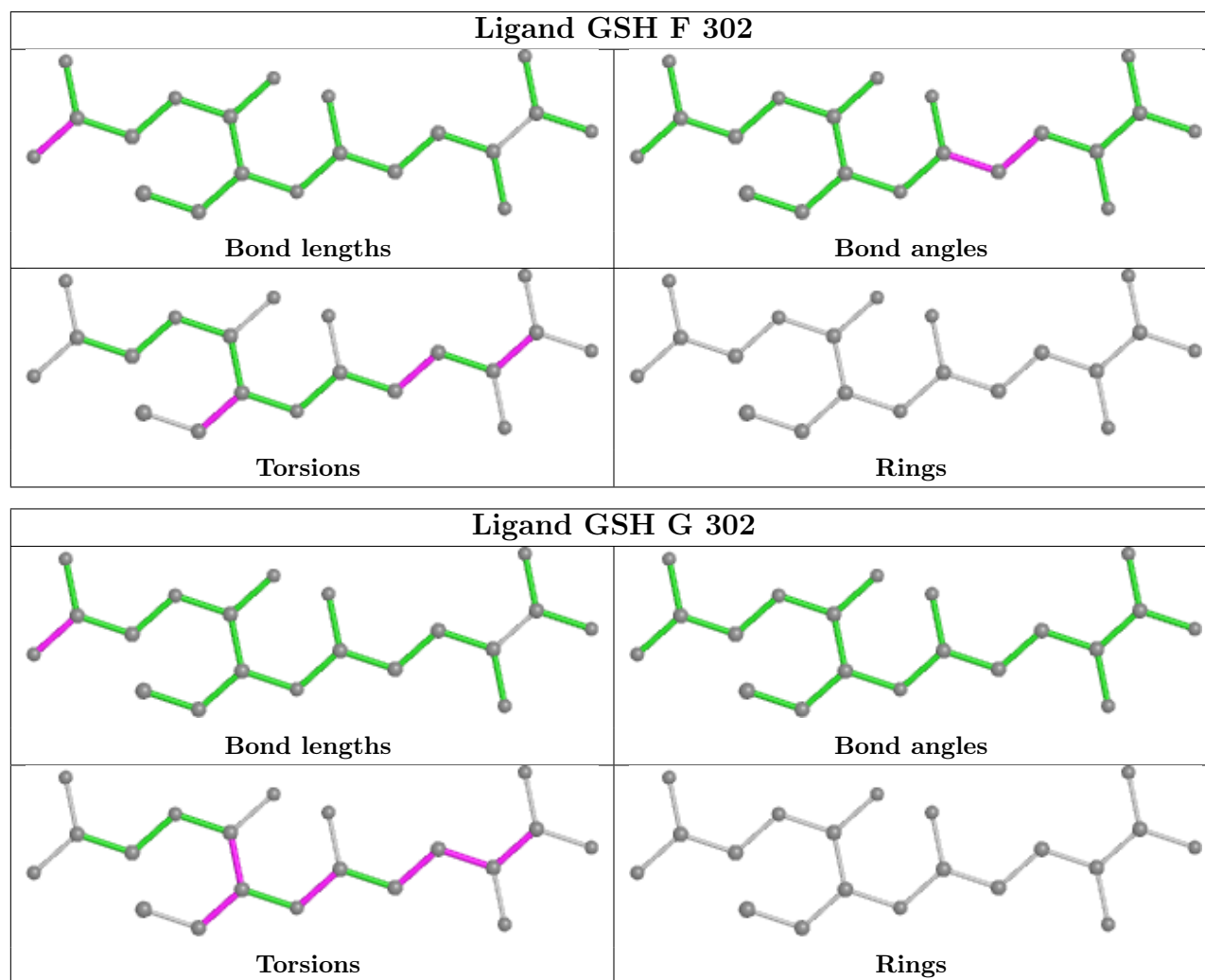
There are no ring outliers.

8 monomers are involved in 15 short contacts:

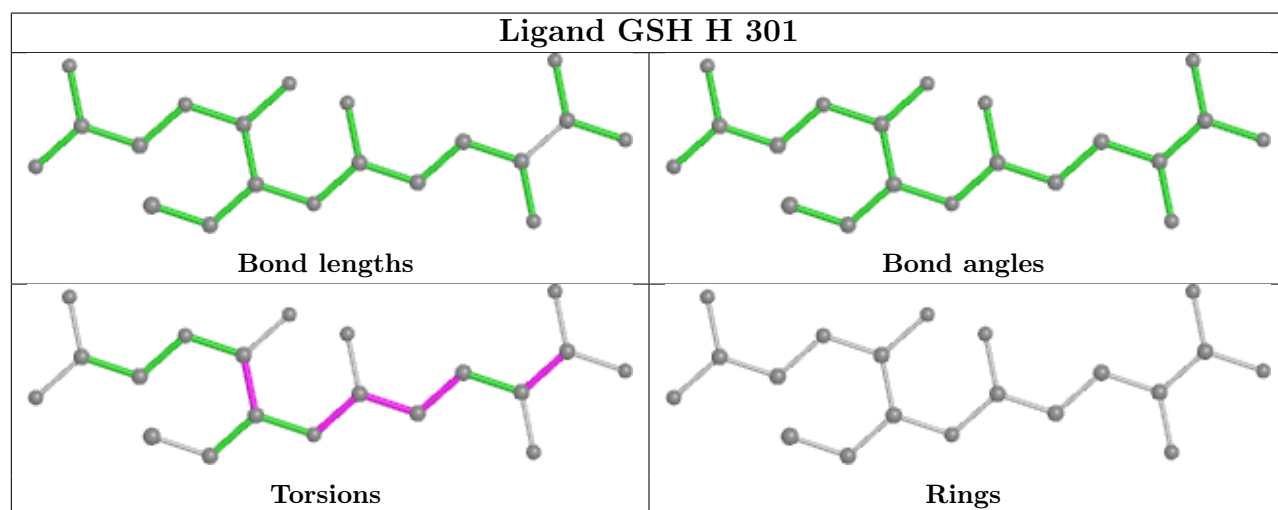
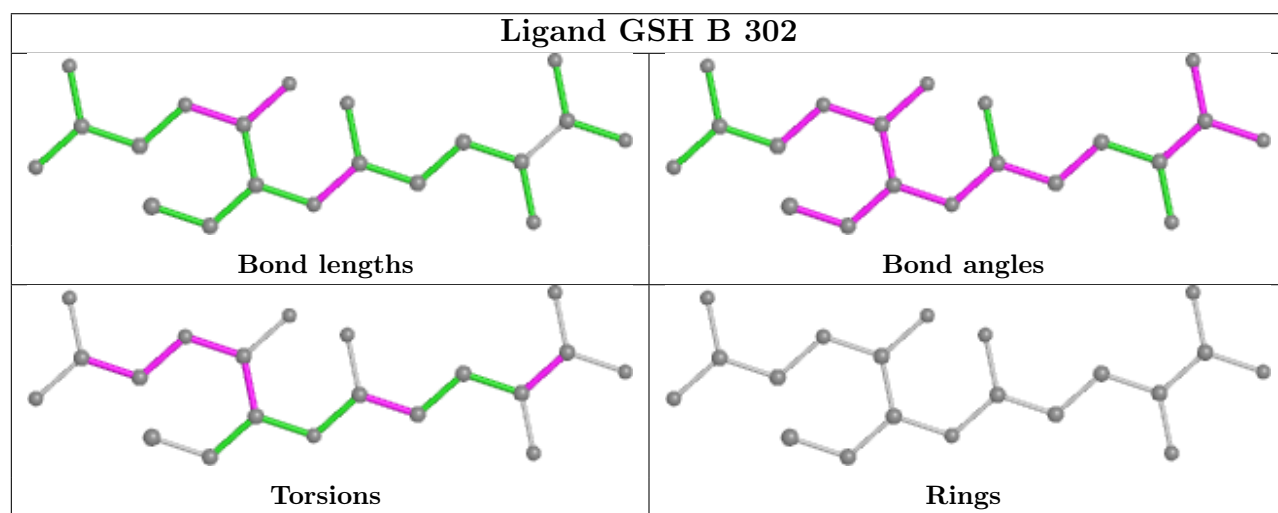
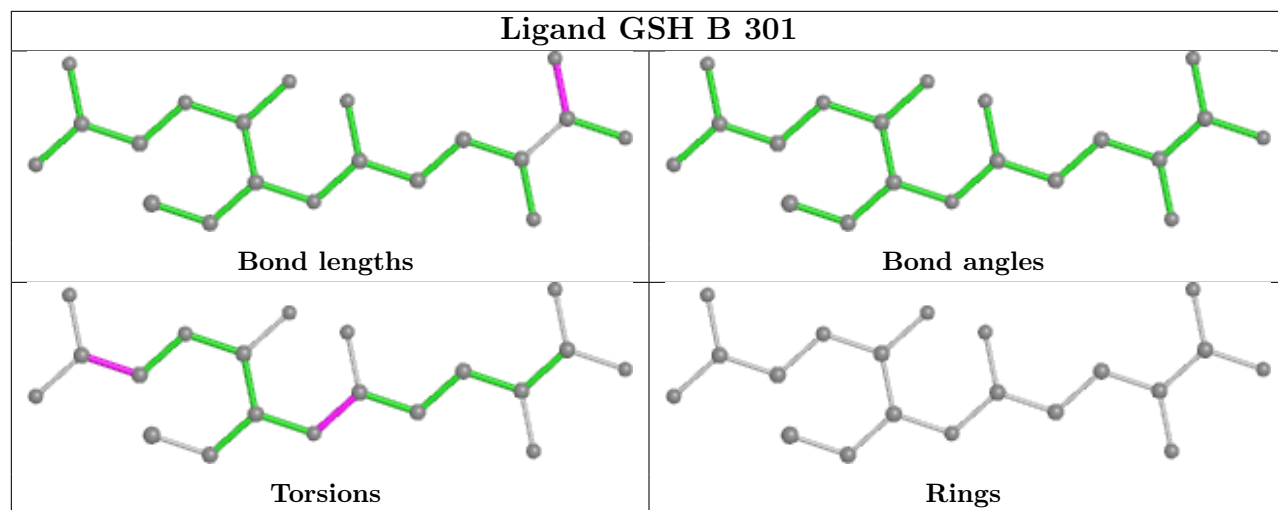
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	302	GSH	1	0
3	C	302	1PE	4	0
2	B	302	GSH	1	0
2	H	301	GSH	2	0
2	C	301	GSH	1	0
2	E	302	GSH	2	0
2	F	301	GSH	2	0
2	A	301	GSH	2	0

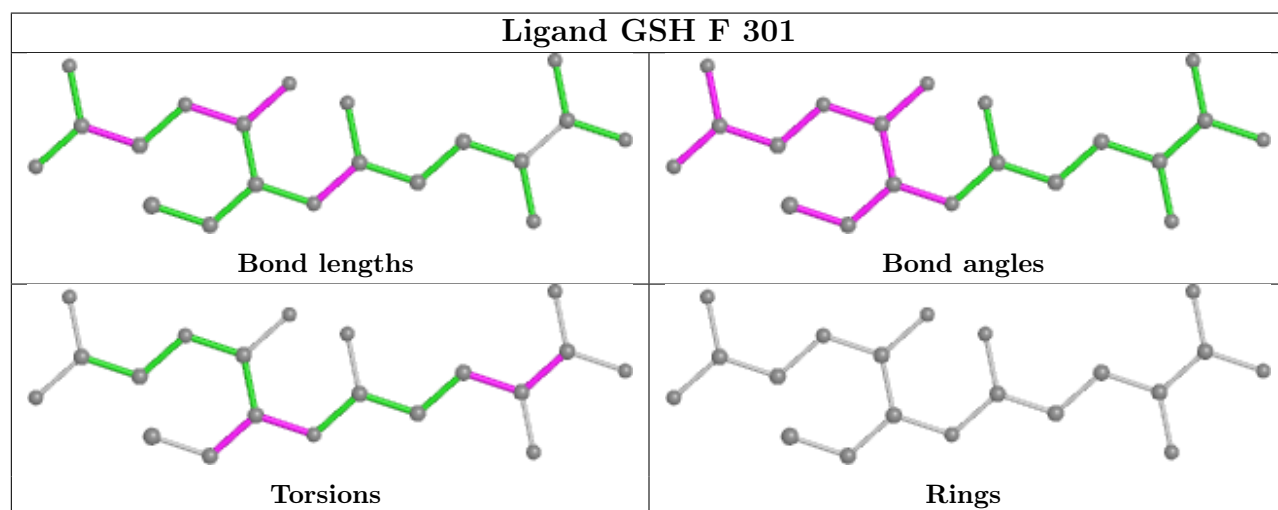
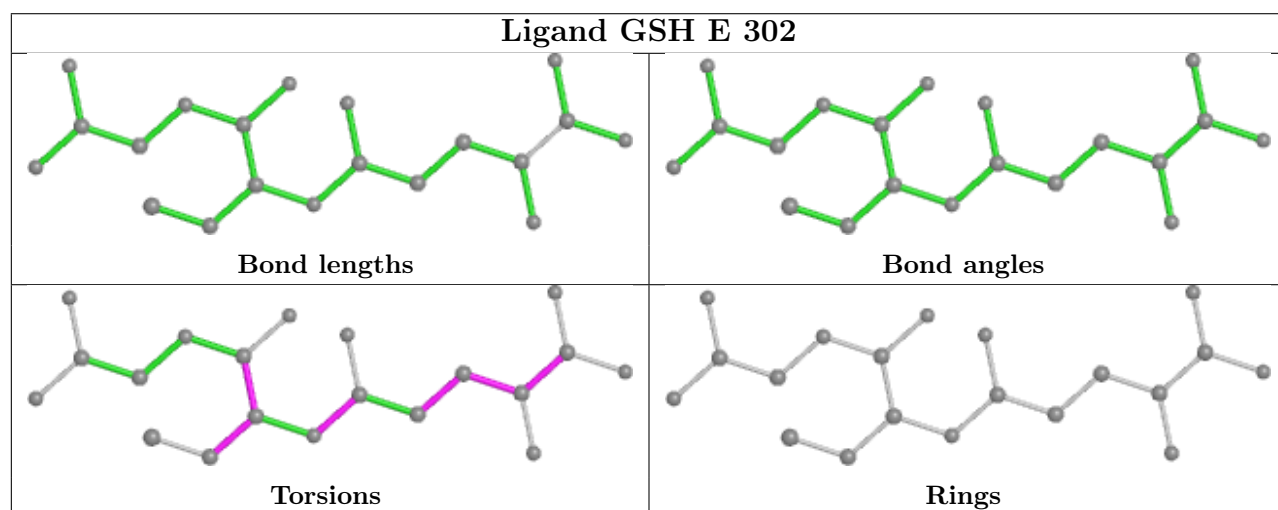
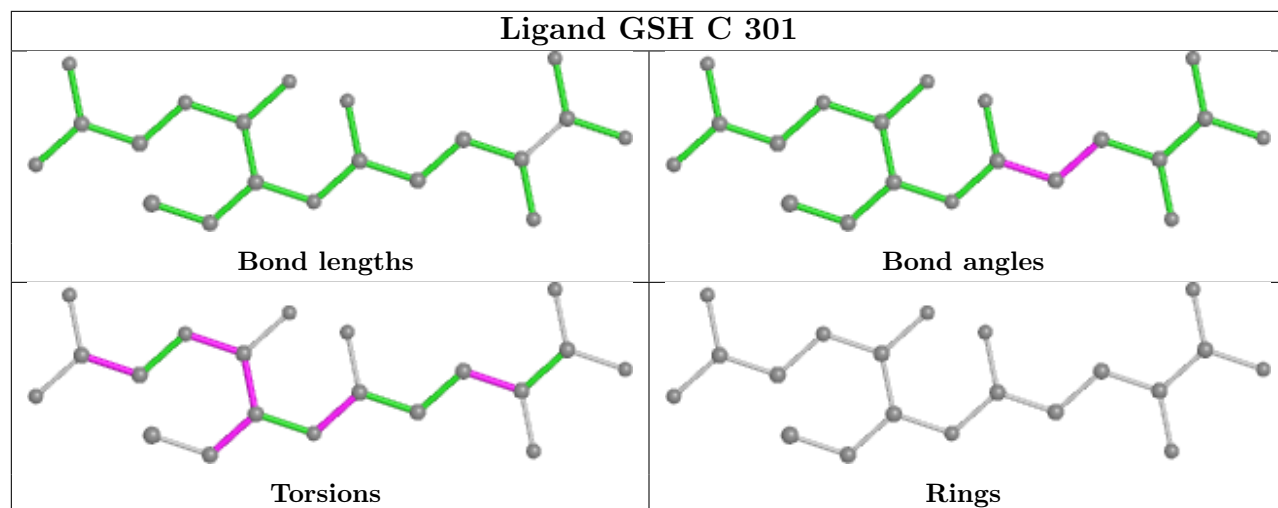
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

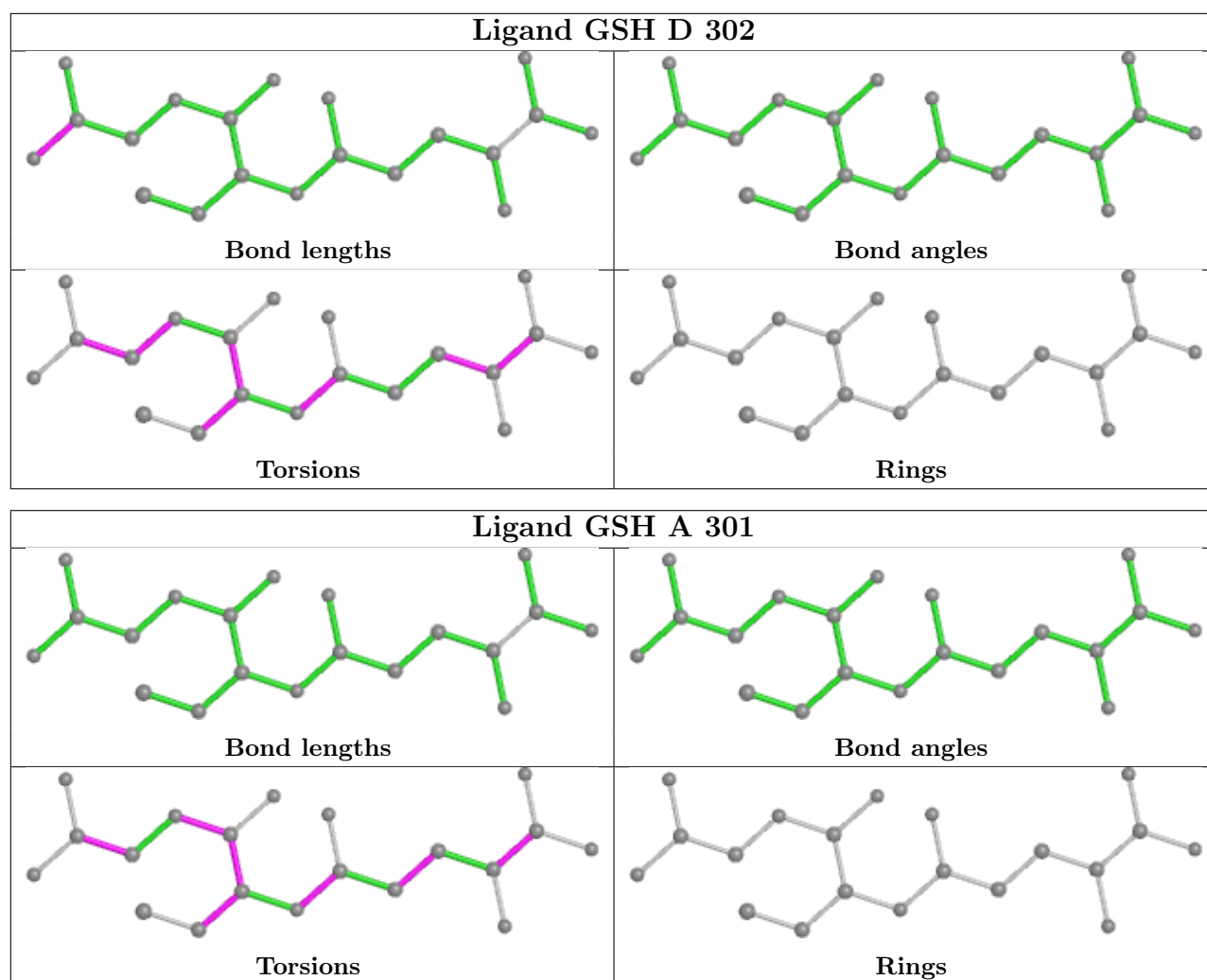
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 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	219/219 (100%)	-0.56	1 (0%) 91 90	5, 14, 37, 60	0
1	B	219/219 (100%)	-0.58	0 100 100	6, 17, 32, 60	0
1	C	219/219 (100%)	-0.69	1 (0%) 91 90	4, 13, 27, 59	0
1	D	219/219 (100%)	-0.73	0 100 100	5, 13, 26, 51	0
1	E	219/219 (100%)	-0.69	0 100 100	5, 14, 29, 36	0
1	F	219/219 (100%)	-0.70	0 100 100	5, 13, 29, 46	0
1	G	219/219 (100%)	-0.66	0 100 100	7, 13, 29, 46	0
1	H	219/219 (100%)	-0.67	0 100 100	5, 14, 30, 56	0
All	All	1752/1752 (100%)	-0.66	2 (0%) 95 95	4, 14, 30, 60	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	67	ASP	4.0
1	C	1	MET	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

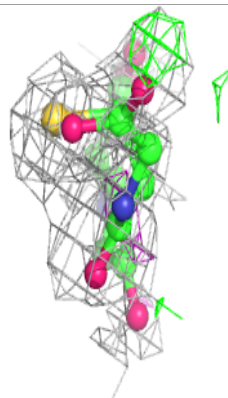
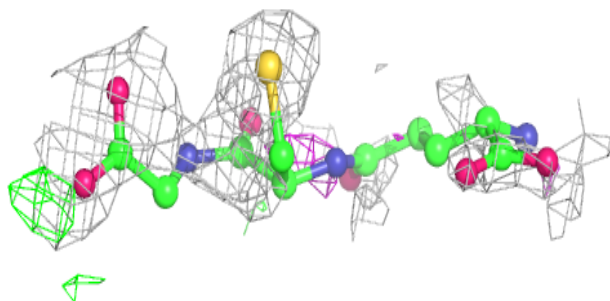
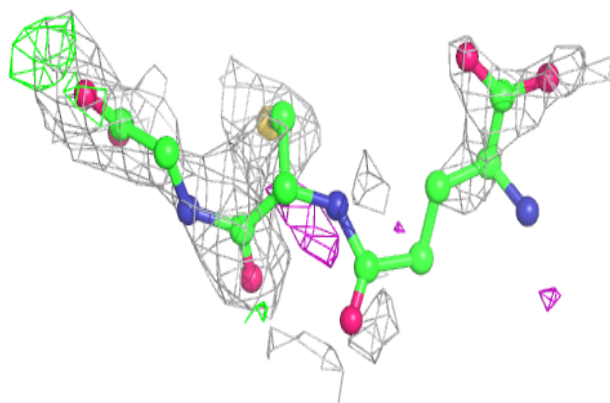
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	GSH	B	302	20/20	0.64	0.38	35,62,80,87	0
2	GSH	F	301	20/20	0.75	0.29	18,40,65,75	0
2	GSH	G	302	20/20	0.76	0.32	19,49,74,86	0
2	GSH	A	301	20/20	0.77	0.25	34,51,62,67	0
2	GSH	D	302	20/20	0.77	0.30	15,47,80,84	0
4	AG	E	301	1/1	0.77	0.11	134,134,134,134	0
2	GSH	H	301	20/20	0.81	0.23	26,40,53,55	0
4	AG	D	301	1/1	0.82	0.20	132,132,132,132	0
2	GSH	C	301	20/20	0.83	0.17	17,38,55,71	0
2	GSH	E	302	20/20	0.86	0.20	16,27,39,53	0
3	1PE	C	302	16/16	0.89	0.13	9,19,32,32	0
2	GSH	F	302	20/20	0.92	0.17	12,28,37,39	0
4	AG	G	301	1/1	0.93	0.07	88,88,88,88	0
2	GSH	B	301	20/20	0.95	0.13	10,22,38,44	0

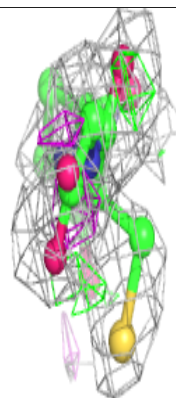
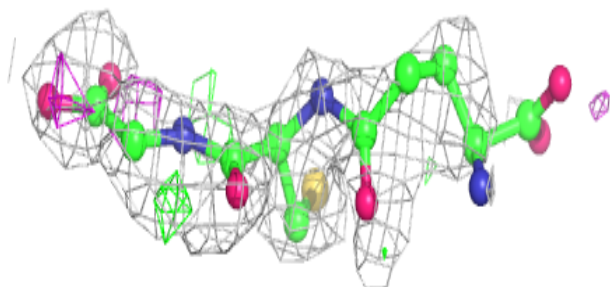
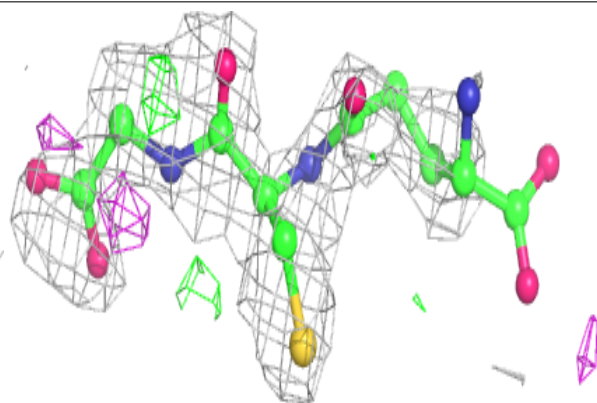
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around GSH B 302:**

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

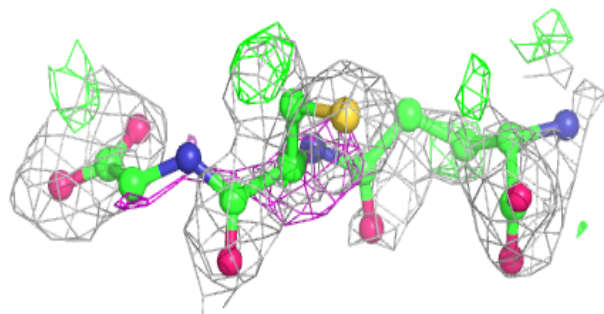
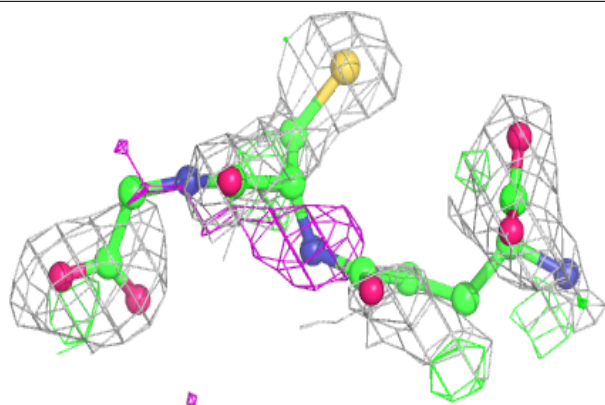
**Electron density around GSH F 301:**

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



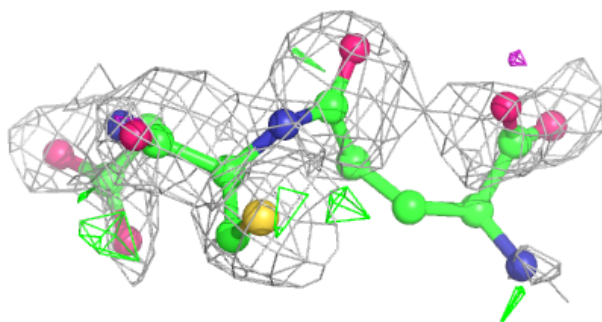
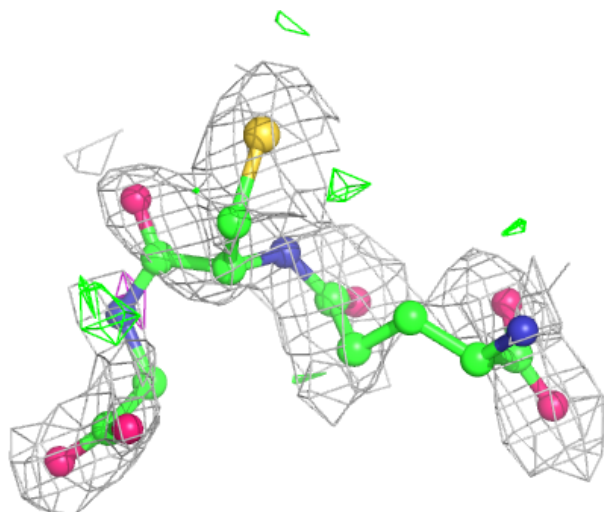
**Electron density around GSH G 302:**

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



**Electron density around GSH A 301:**

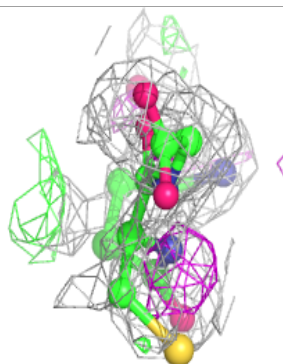
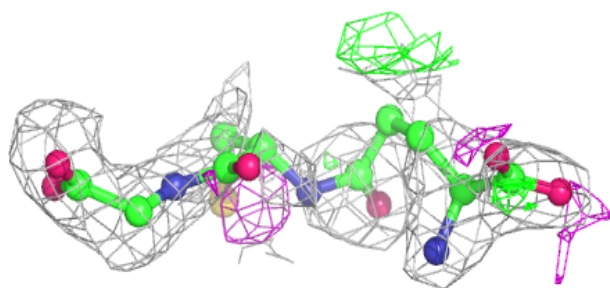
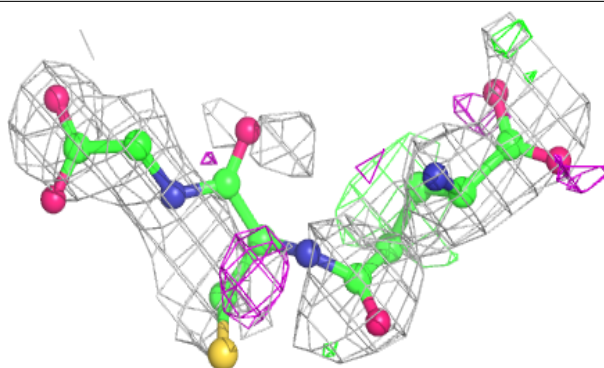
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



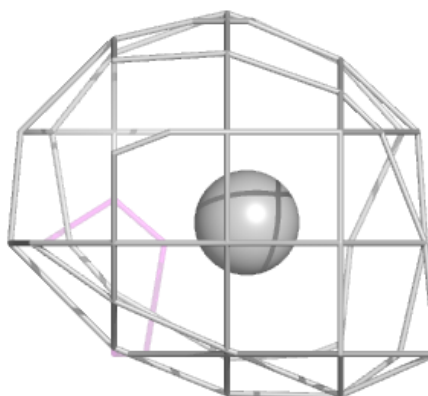
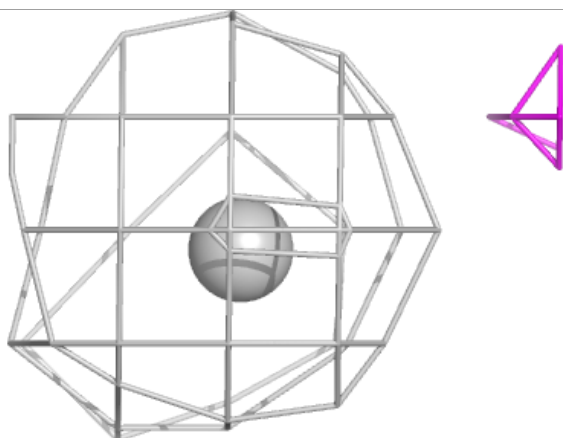
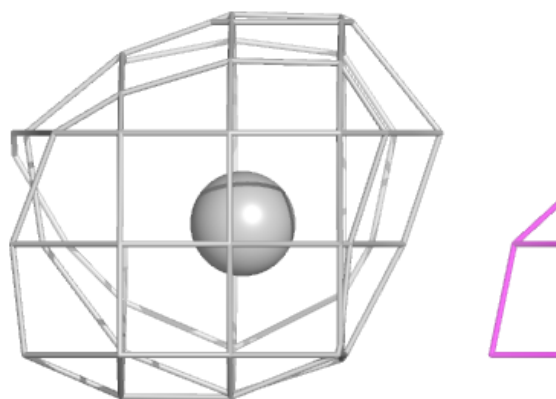


**Electron density around GSH D 302:**

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

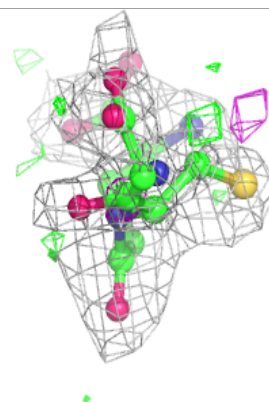
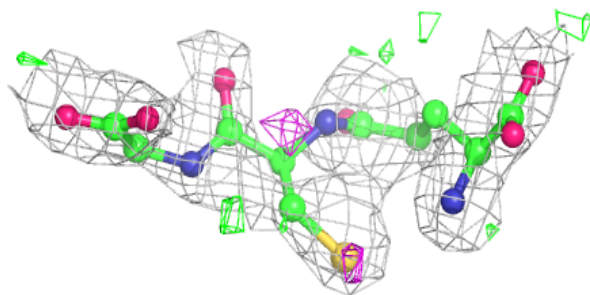
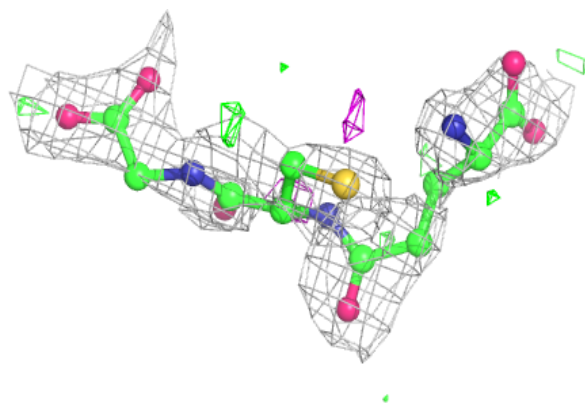
**Electron density around AG E 301:**

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



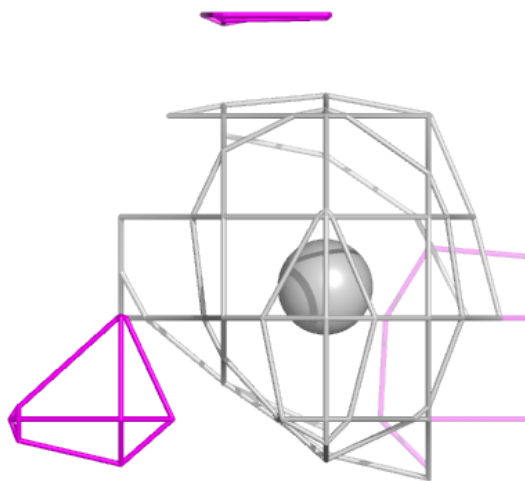
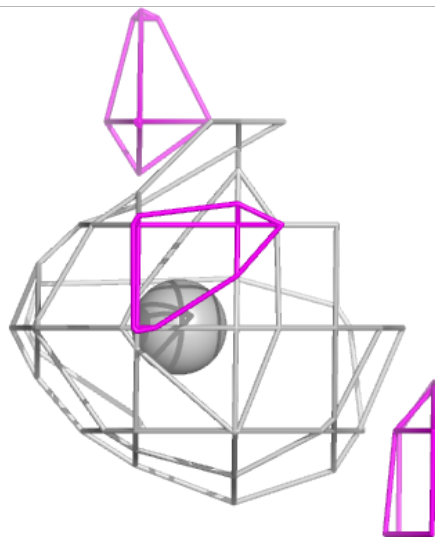
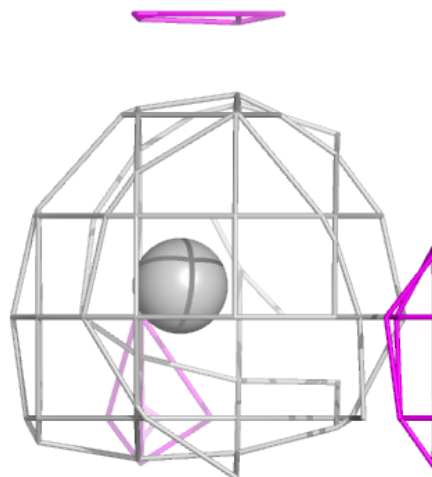
**Electron density around GSH H 301:**

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



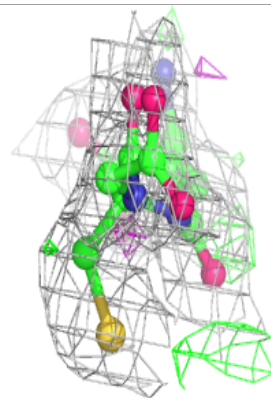
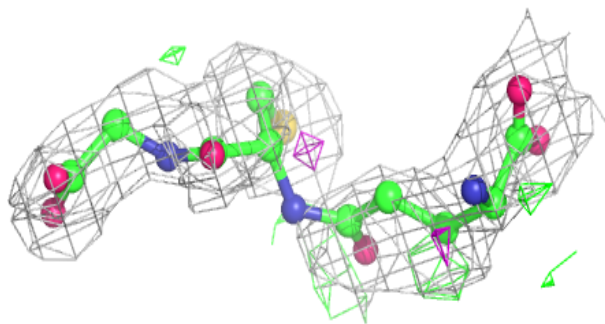
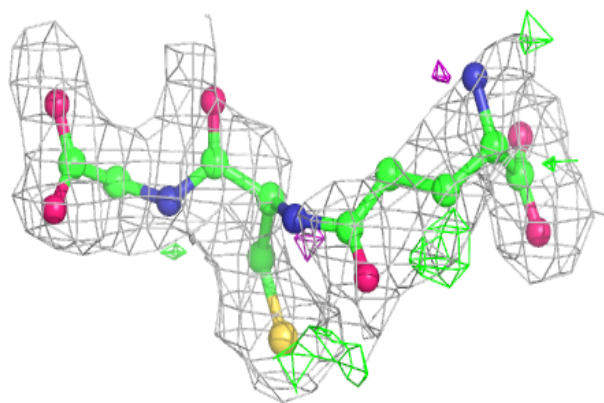
**Electron density around AG D 301:**

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



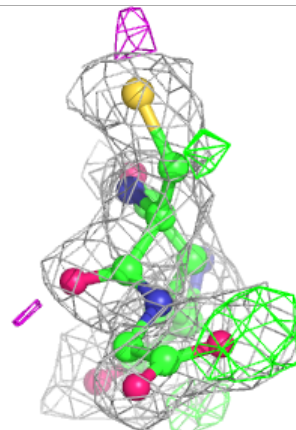
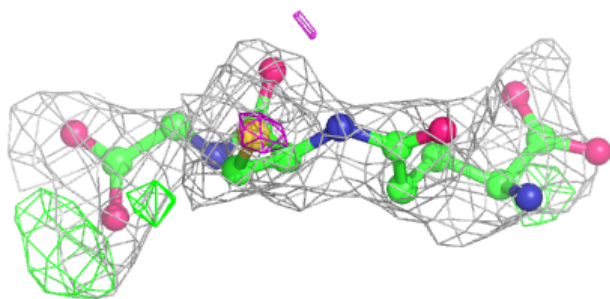
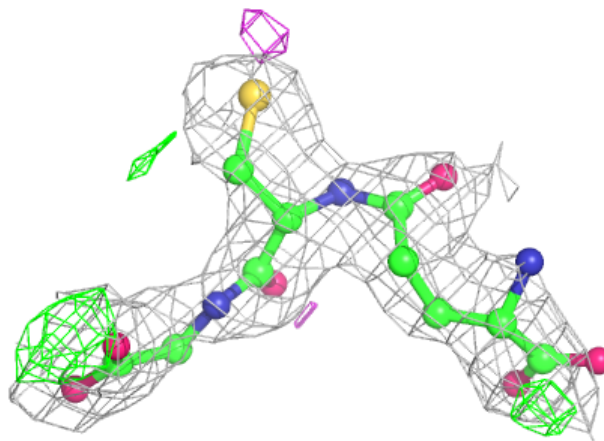
**Electron density around GSH C 301:**

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



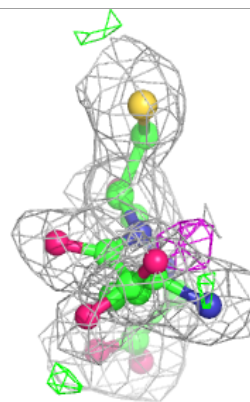
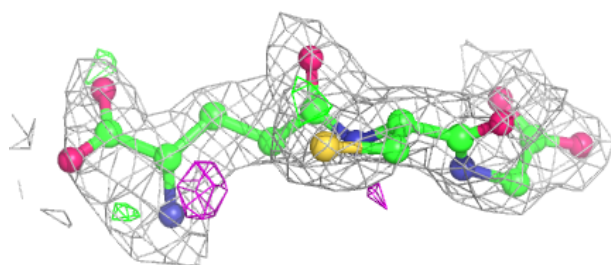
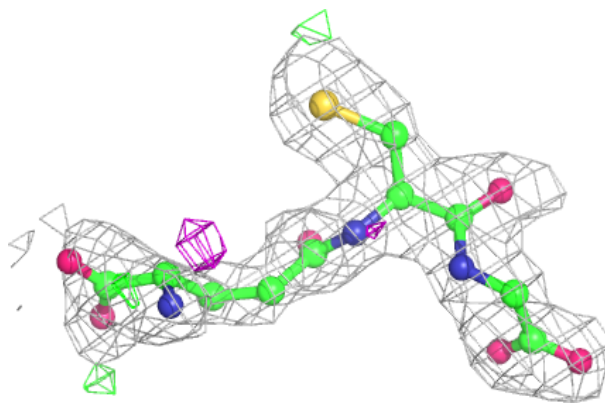
**Electron density around GSH E 302:**

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



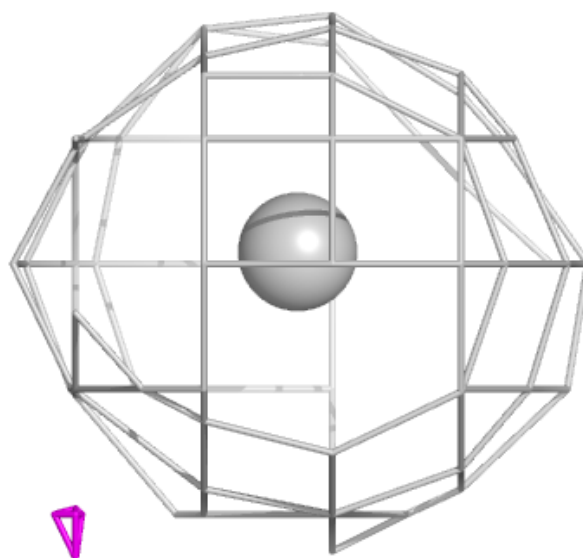
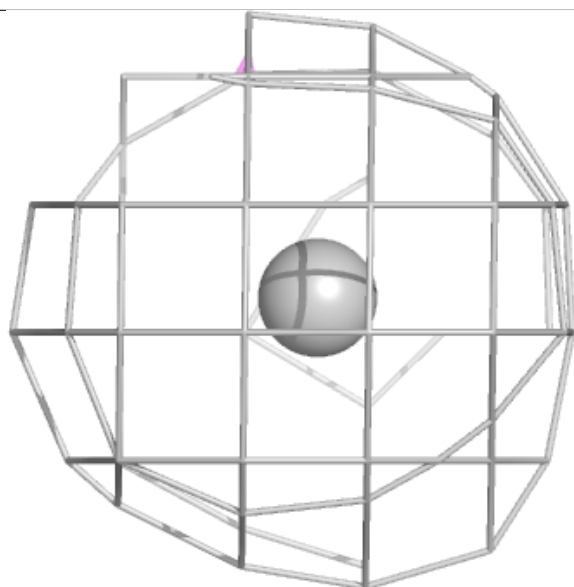
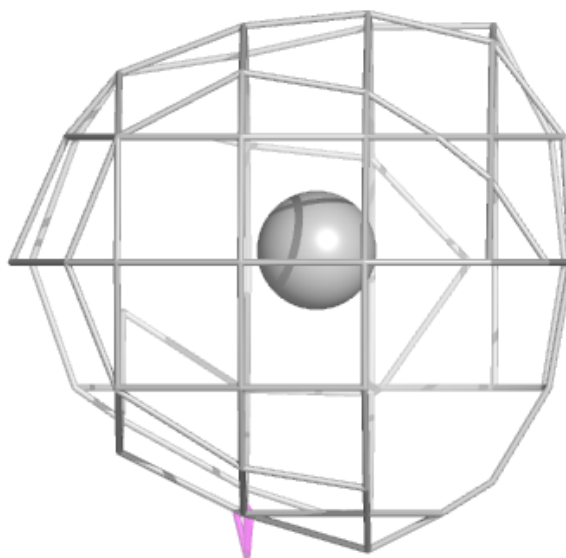
**Electron density around GSH F 302:**

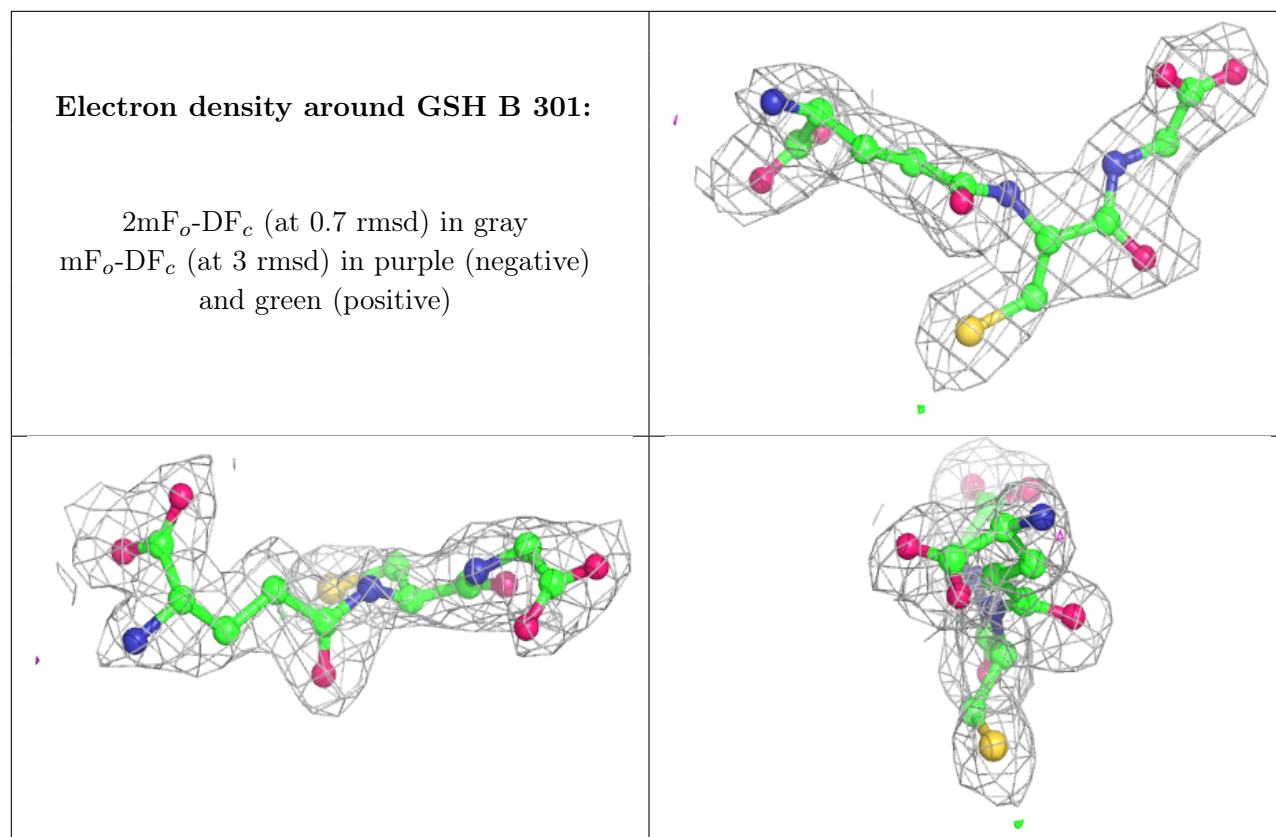
$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around AG G 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.