W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

Oct 23, 2021 – 03:30 PM EDT

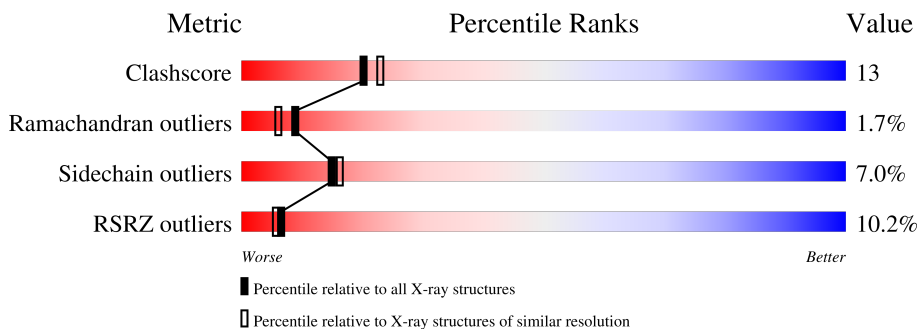|		
PDB ID	:	1EV4
Title	:	RAT GLUTATHIONE S-TRANSFERASE A1-1: MUTANT W21F/F220Y WITH GSO3 BOUND
Authors	:	Adman, E.T.; Le Trong, I.; Stenkamp, R.E.; Nieslanik, B.S.; Dietze, E.C.; Tai, G.; Ibarra, C.; Atkins, W.M.
Deposited on	:	2000-04-19
Resolution	:	2.20 Å(reported)

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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	Percentile Ranks	Value
Clashscore		13
Ramachandran outliers		1.7%
Sidechain outliers		7.0%
RSRZ outliers		10.2%

Worse   *Better*

▌ Percentile relative to all X-ray structures

◻ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
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 >=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 <=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	221	10% 68% 27% · ·
1	C	221	13% 57% 32% · 6%
1	D	221	7% 75% 22% · ·

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 5599 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 A1-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	221	Total	C	N	O	S	0	0	0
			1786	1152	299	324	11			
1	C	207	Total	C	N	O	S	0	0	0
			1664	1073	278	302	11			
1	D	221	Total	C	N	O	S	0	0	0
			1786	1152	299	324	11			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	21	PHE	TRP	engineered mutation	UNP P00502
A	96	SER	THR	conflict	UNP P00502
A	220	TYR	PHE	engineered mutation	UNP P00502
C	21	PHE	TRP	engineered mutation	UNP P00502
C	96	SER	THR	conflict	UNP P00502
C	220	TYR	PHE	engineered mutation	UNP P00502
D	21	PHE	TRP	engineered mutation	UNP P00502
D	96	SER	THR	conflict	UNP P00502
D	220	TYR	PHE	engineered mutation	UNP P00502

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

SO4

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0
2	D	1	Total 5	O 4	S 1	0	0

- Molecule 3 is GLUTATHIONE SULFONIC ACID (three-letter code: GTS) (formula: $C_{10}H_{17}N_3O_9S$).

GTS
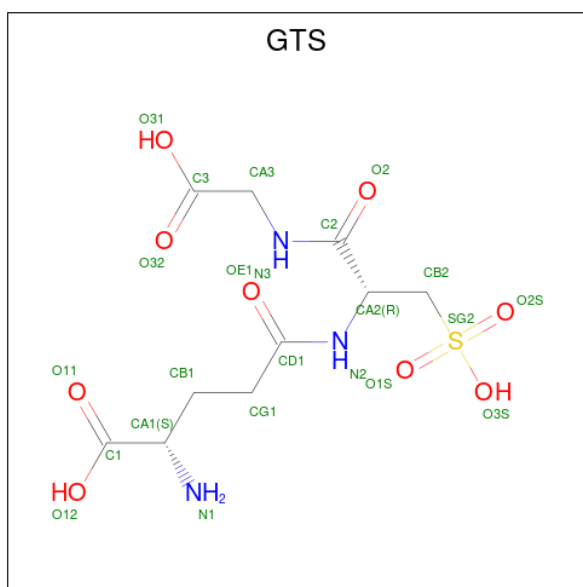
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 23	C 10	N 3	O 9	S 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	S	0	0
			23	10	3	9	1		
3	D	1	Total	C	N	O	S	0	0
			23	10	3	9	1		

• Molecule 4 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total	O	0	0
			126	126		
4	C	56	Total	O	0	0
			56	56		
4	D	97	Total	O	0	0
			97	97		

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.
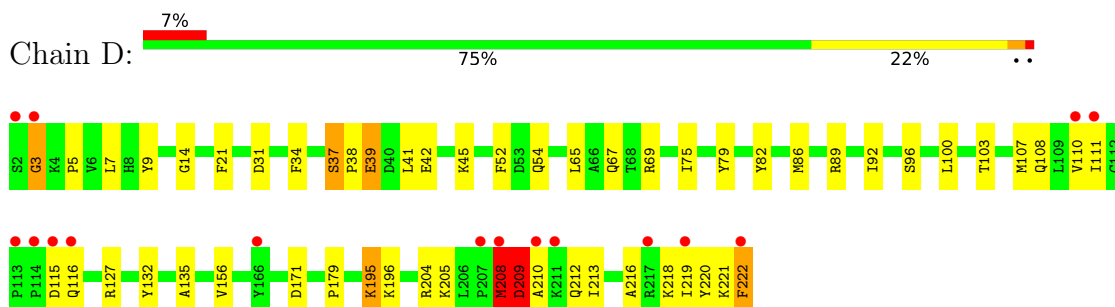
- Molecule 1: GLUTATHIONE S-TRANSFERASE A1-1
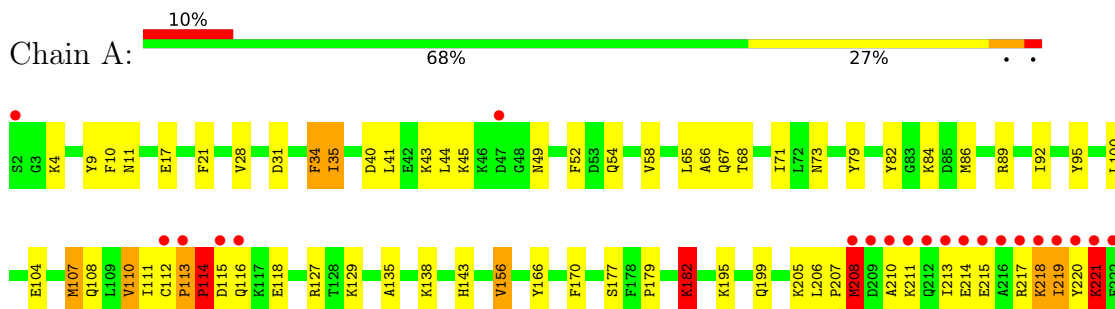
- Molecule 1: GLUTATHIONE S-TRANSFERASE A1-1
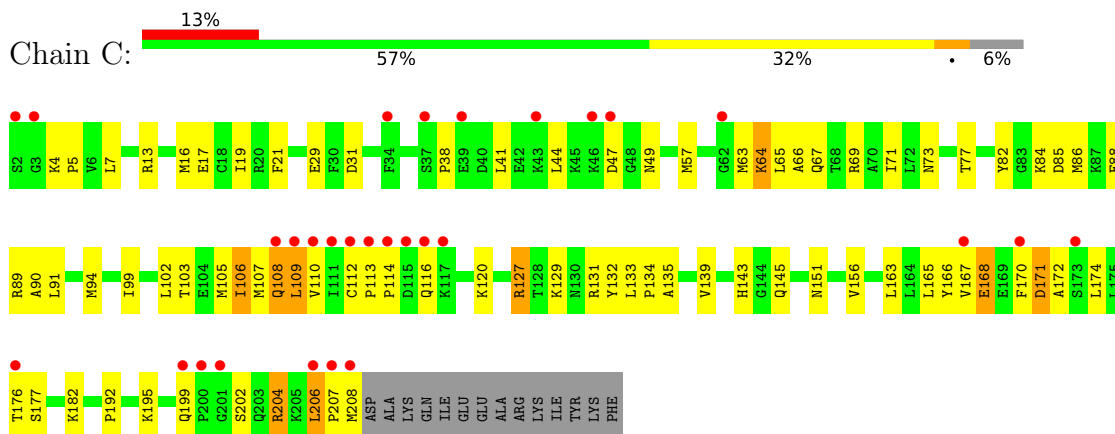
- Molecule 1: GLUTATHIONE S-TRANSFERASE A1-1

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	84.60Å 275.20Å 70.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.20 19.82 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.20) 77.2 (19.82-1.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.91 (at 1.80Å)	Xtriage
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.218 , 0.285 0.234 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	20.8	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 76.9	EDS
L-test for twinning[2]	$<|L|> = 0.49, <L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5599	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $<|L|>$, $<L^2>$ 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: SO4, GTS

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	0.66	0/1818	1.09	17/2440 (0.7%)
1	C	0.62	0/1694	1.01	5/2276 (0.2%)
1	D	0.61	0/1818	1.00	8/2440 (0.3%)
All	All	0.63	0/5330	1.03	30/7156 (0.4%)

There are no bond length outliers.

The worst 5 of 30 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	222	PHE	N-CA-CB	-7.39	97.29	110.60
1	A	182	LYS	CB-CG-CD	7.20	130.32	111.60
1	D	31	ASP	CA-CB-CG	7.00	128.80	113.40
1	A	218	LYS	CB-CG-CD	6.81	129.30	111.60
1	A	170	PHE	N-CA-CB	-6.71	98.52	110.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1786	0	1852	52	0
1	C	1664	0	1726	55	0
1	D	1786	0	1852	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	23	0	14	1	0
3	C	23	0	14	0	0
3	D	23	0	14	1	0
4	A	126	0	0	16	0
4	C	56	0	0	10	0
4	D	97	0	0	6	0
All	All	5599	0	5472	135	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 13.

The worst 5 of 135 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:PHE:HA	4:A:790:HOH:O	1.63	0.97
1:A:86:MET:HA	4:A:757:HOH:O	1.73	0.87
1:C:206:LEU:HB3	1:C:207:PRO:HD2	1.66	0.76
1:C:91:LEU:HA	4:C:735:HOH:O	1.85	0.75
1:A:73:ASN:HB3	4:A:455:HOH:O	1.87	0.75

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	219/221 (99%)	199 (91%)	16 (7%)	4 (2%)	8	5
1	C	205/221 (93%)	186 (91%)	17 (8%)	2 (1%)	15	14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	219/221 (99%)	203 (93%)	11 (5%)	5 (2%)	6	3
All	All	643/663 (97%)	588 (91%)	44 (7%)	11 (2%)	9	6

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	PRO
1	A	221	LYS
1	D	208	MET
1	C	171	ASP
1	D	210	ALA

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	195/195 (100%)	181 (93%)	14 (7%)	14	15
1	C	183/195 (94%)	168 (92%)	15 (8%)	11	11
1	D	195/195 (100%)	184 (94%)	11 (6%)	21	25
All	All	573/585 (98%)	533 (93%)	40 (7%)	15	16

5 of 40 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	204	ARG
1	D	116	GLN
1	D	21	PHE
1	D	82	TYR
1	D	204	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	49	ASN
1	D	108	GLN
1	D	199	GLN
1	D	193	ASN
1	C	193	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 ⓘ

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	GTS	A	230	-	15,22,22	0.75	0	19,30,30	1.51	3 (15%)
2	SO4	A	260	-	4,4,4	0.54	0	6,6,6	0.37	0
3	GTS	D	250	-	15,22,22	1.05	1 (6%)	19,30,30	1.58	3 (15%)
2	SO4	C	261	-	4,4,4	0.43	0	6,6,6	0.24	0
3	GTS	C	240	-	15,22,22	0.69	0	19,30,30	1.64	3 (15%)
2	SO4	D	262	-	4,4,4	0.42	0	6,6,6	0.25	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GTS	C	240	-	-	1/21/27/27	-
3	GTS	A	230	-	-	1/21/27/27	-
3	GTS	D	250	-	-	1/21/27/27	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	250	GTS	CB2-CA2	-2.49	1.51	1.53

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	240	GTS	O3S-SG2-CB2	3.65	111.55	105.74
3	D	250	GTS	O1S-SG2-CB2	3.55	111.16	106.94
3	A	230	GTS	O3S-SG2-CB2	3.51	111.33	105.74
3	C	240	GTS	O1S-SG2-CB2	3.49	111.08	106.94
3	D	250	GTS	O3S-SG2-CB2	3.45	111.24	105.74

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	250	GTS	C3-CA3-N3-C2
3	C	240	GTS	C3-CA3-N3-C2
3	A	230	GTS	C3-CA3-N3-C2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	230	GTS	1	0
3	D	250	GTS	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 then 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 A 230

Bond lengths

Bond angles

Torsions

Rings

Ligand GTS D 250

Bond lengths

Bond angles

Torsions

Rings

Ligand GTS C 240

Bond lengths

Bond angles

Torsions

Rings
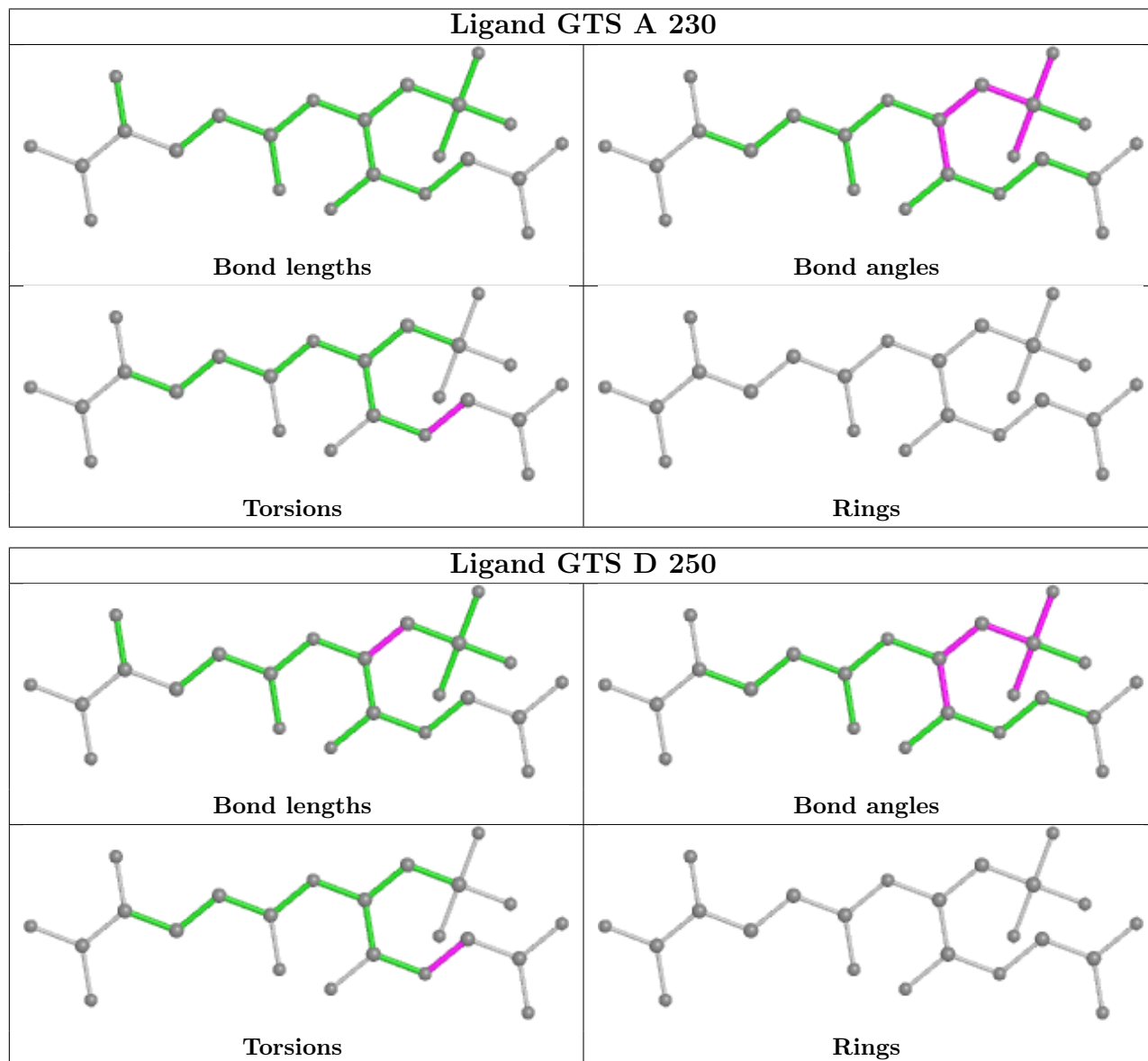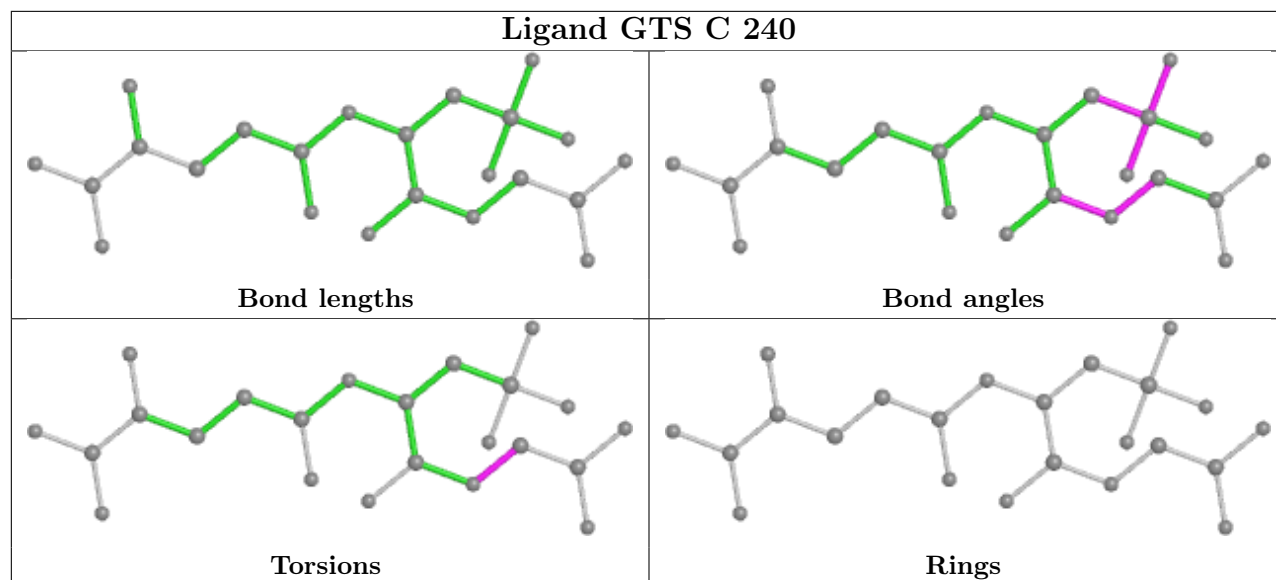
5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	221/221 (100%)	0.29	21 (9%)	8	7	16, 27, 95, 100	0
1	C	207/221 (93%)	0.70	29 (14%)	2	2	26, 47, 88, 100	0
1	D	221/221 (100%)	0.18	16 (7%)	15	14	23, 35, 81, 100	0
All	All	649/663 (97%)	0.38	66 (10%)	6	6	16, 38, 86, 100	0

The worst 5 of 66 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	ALA	25.1
1	A	219	ILE	10.4
1	A	222	PHE	9.7
1	C	2	SER	9.7
1	C	112	CYS	8.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

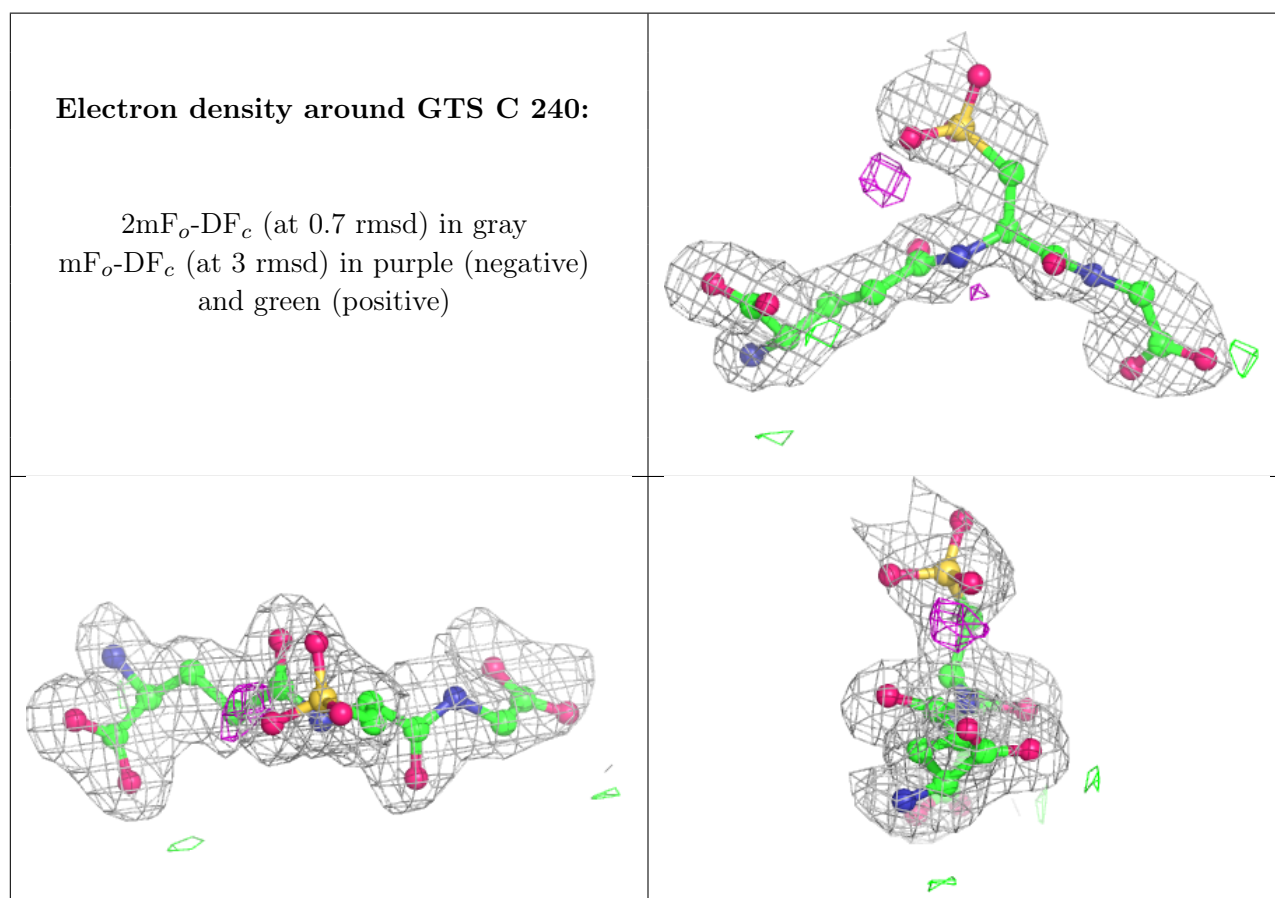There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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, 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($\mathring{A}^2$)	Q<0.9
3	GTS	C	240	23/23	0.84	0.15	29,60,80,83	0
3	GTS	A	230	23/23	0.90	0.13	16,33,51,59	0
3	GTS	D	250	23/23	0.90	0.12	26,39,56,58	0
2	SO4	D	262	5/5	0.93	0.12	81,81,83,85	0
2	SO4	C	261	5/5	0.94	0.14	74,74,75,77	0
2	SO4	A	260	5/5	0.95	0.11	40,42,47,49	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.
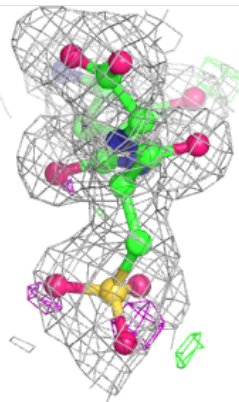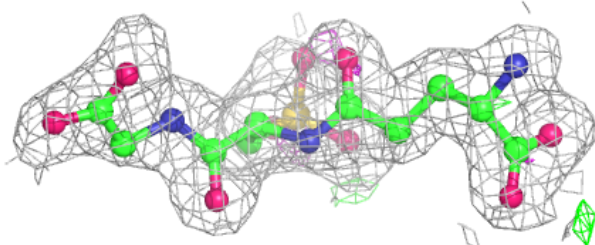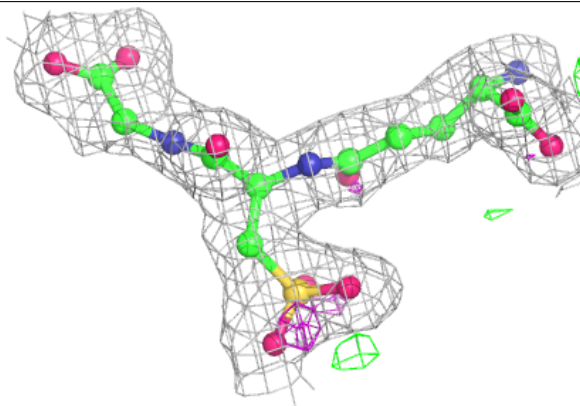

Electron density around GTS C 240:

$2mF_o$-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 A 230:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
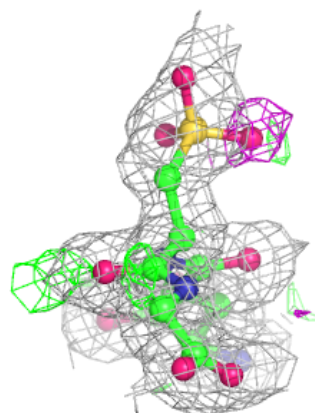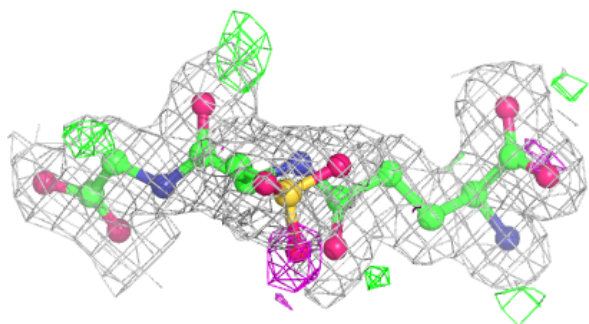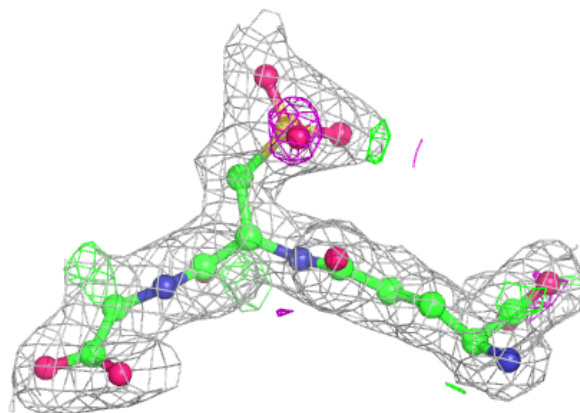
Electron density around GTS D 250:

$2mF_o$-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.