



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:07 AM UTC

PDB ID : 2GSQ / pdb_00002gsq
Title : GLUTATHIONE S-TRANSFERASE FROM SQUID DIGESTIVE GLAND
COMPLEXED WITH S-(3-IODOBENZYL)GLUTATHIONE
Authors : Ji, X.; Armstrong, R.N.; Gilliland, G.L.
Deposited on : 1995-04-14
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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

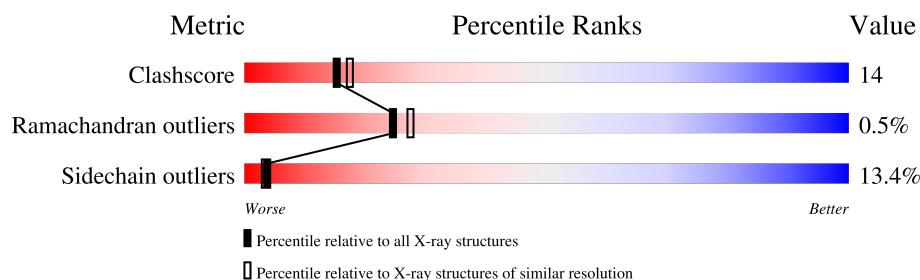
1 Overall quality at a glance

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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	202	 53% 36% 8%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1870 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.

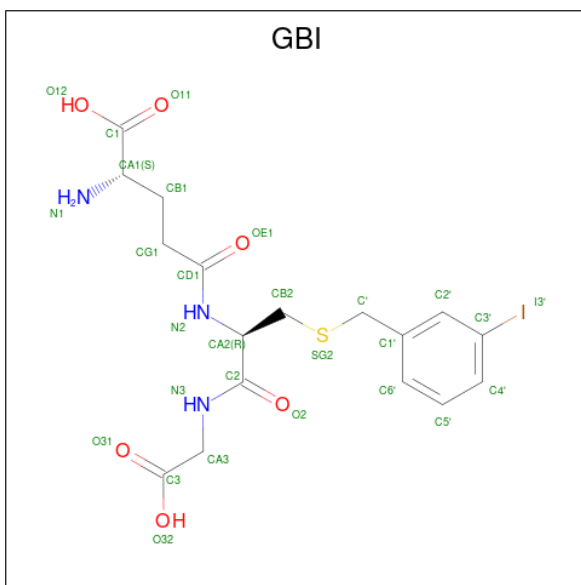
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	202	1598	1022	273	290	13	0	0	0

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is S-(3-IODOBENZYL)GLUTATHIONE (CCD ID: GBI) (formula: C₁₇H₂₂I₁N₃O₆S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 28	C 17	I 1	N 3	O 6	S 1	0	0
3	A	1	Total 28	C 17	I 1	N 3	O 6	S 1	0	0

- Molecule 4 is water.

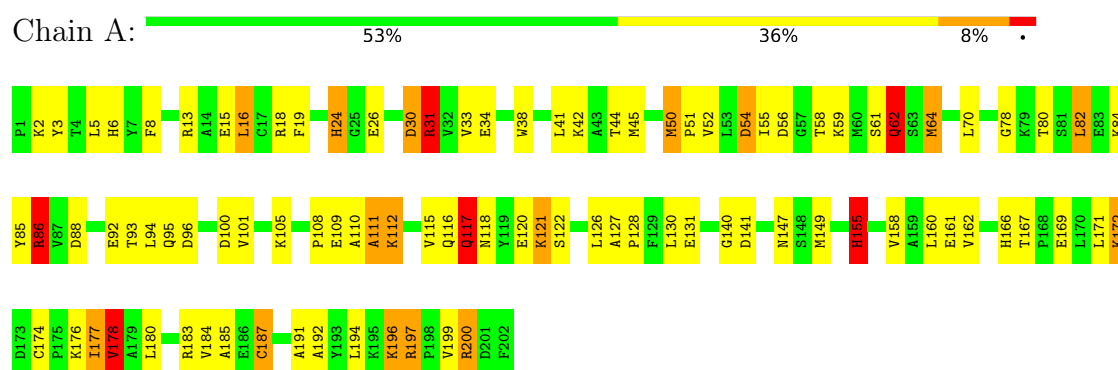
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	211	Total O 211 211	0	0

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. 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. 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.

Note EDS was not executed.

• Molecule 1: GLUTATHIONE S-TRANSFERASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	73.15Å 73.15Å 94.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.173 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1870	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

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: GBI, SO4

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	1.30	3/1632 (0.2%)	2.19	68/2200 (3.1%)

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	A	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	167	THR	CA-CB	5.94	1.61	1.53
1	A	78	GLY	N-CA	5.67	1.50	1.45
1	A	161	GLU	N-CA	5.13	1.52	1.46

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	ARG	CD-NE-CZ	20.31	152.83	124.40
1	A	200	ARG	CD-NE-CZ	12.71	142.20	124.40
1	A	30	ASP	CA-C-O	-9.93	109.38	120.70
1	A	96	ASP	CA-CB-CG	9.79	122.39	112.60
1	A	158	VAL	N-CA-C	9.57	119.61	110.42
1	A	197	ARG	NE-CZ-NH2	9.46	127.72	119.20
1	A	178	VAL	N-CA-CB	9.28	121.41	110.55
1	A	197	ARG	NE-CZ-NH1	-8.47	113.03	121.50
1	A	111	ALA	N-CA-C	-7.75	102.69	114.16
1	A	18	ARG	CD-NE-CZ	7.70	135.18	124.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	183	ARG	CD-NE-CZ	7.22	134.51	124.40
1	A	94	LEU	CA-C-N	6.79	130.06	120.28
1	A	94	LEU	C-N-CA	6.79	130.06	120.28
1	A	111	ALA	CA-C-O	6.76	125.95	118.52
1	A	172	LYS	CA-CB-CG	6.47	127.04	114.10
1	A	177	ILE	N-CA-C	-6.43	104.48	110.53
1	A	183	ARG	CA-CB-CG	6.29	126.69	114.10
1	A	160	LEU	N-CA-C	6.29	120.78	113.18
1	A	62	GLN	OE1-CD-NE2	-6.18	116.42	122.60
1	A	192	ALA	CA-C-N	6.17	128.86	120.54
1	A	192	ALA	C-N-CA	6.17	128.86	120.54
1	A	85	TYR	N-CA-C	-6.12	104.53	111.14
1	A	58	THR	CA-CB-OG1	-6.03	100.56	109.60
1	A	118	ASN	CA-C-N	6.00	128.24	120.44
1	A	118	ASN	C-N-CA	6.00	128.24	120.44
1	A	180	LEU	CA-C-O	-5.95	113.63	120.24
1	A	162	VAL	CA-CB-CG1	5.90	120.43	110.40
1	A	88	ASP	CA-CB-CG	5.87	118.47	112.60
1	A	183	ARG	NE-CZ-NH1	5.87	127.37	121.50
1	A	54	ASP	CA-C-O	-5.81	113.92	120.32
1	A	110	ALA	CA-C-N	5.76	130.09	120.70
1	A	110	ALA	C-N-CA	5.76	130.09	120.70
1	A	44	THR	O-C-N	5.75	128.94	122.27
1	A	13	ARG	NE-CZ-NH2	-5.72	114.05	119.20
1	A	100	ASP	N-CA-CB	5.66	119.01	110.30
1	A	117	GLN	CB-CA-C	5.65	121.91	110.38
1	A	178	VAL	CA-C-O	-5.65	115.18	121.17
1	A	185	ALA	CA-C-N	5.58	131.91	122.26
1	A	185	ALA	C-N-CA	5.58	131.91	122.26
1	A	15	GLU	N-CA-C	5.55	117.33	111.28
1	A	92	GLU	CG-CD-OE2	-5.46	105.85	118.40
1	A	50	MET	CG-SD-CE	-5.44	88.93	100.90
1	A	177	ILE	O-C-N	5.39	127.50	121.90
1	A	178	VAL	O-C-N	5.39	127.19	121.91
1	A	86	ARG	NE-CZ-NH1	5.38	126.88	121.50
1	A	122	SER	CA-C-O	-5.38	115.17	120.82
1	A	197	ARG	CB-CA-C	5.38	117.93	109.27
1	A	55	ILE	CA-C-O	-5.36	114.58	120.48
1	A	56	ASP	CA-C-N	-5.33	117.09	123.44
1	A	56	ASP	C-N-CA	-5.33	117.09	123.44
1	A	24	HIS	N-CA-C	-5.31	106.39	112.92
1	A	110	ALA	CA-C-O	5.30	128.09	120.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	GLY	N-CA-C	-5.29	106.36	112.29
1	A	18	ARG	NE-CZ-NH1	5.23	126.73	121.50
1	A	155	HIS	CA-CB-CG	-5.21	108.58	113.80
1	A	16	LEU	CA-C-O	-5.12	115.12	120.55
1	A	30	ASP	O-C-N	5.11	129.66	123.01
1	A	19	PHE	N-CA-C	5.11	116.93	111.36
1	A	199	VAL	CB-CA-C	5.11	118.46	111.31
1	A	19	PHE	CA-CB-CG	5.08	118.88	113.80
1	A	31	ARG	N-CA-CB	5.06	117.89	110.35
1	A	3	TYR	CA-C-O	-5.05	114.76	120.32
1	A	178	VAL	CA-CB-CG2	5.05	118.98	110.40
1	A	166	HIS	CA-C-N	5.04	128.09	122.74
1	A	166	HIS	C-N-CA	5.04	128.09	122.74
1	A	191	ALA	N-CA-C	-5.03	105.79	111.28
1	A	64	MET	CA-C-O	-5.01	115.11	120.42
1	A	187	CYS	N-CA-C	-5.01	100.81	109.58

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	31	ARG	Sidechain
1	A	86	ARG	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	1598	0	1620	44	0
2	A	5	0	0	1	0
3	A	56	0	40	1	0
4	A	211	0	0	4	0
All	All	1870	0	1660	45	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 14.

All (45) 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:109:GLU:HA	1:A:112:LYS:HE3	1.57	0.86
1:A:95:GLN:HE21	1:A:155:HIS:HE1	1.35	0.72
1:A:6:HIS:HB3	1:A:50:MET:CE	2.20	0.71
1:A:45:MET:HE2	1:A:52:VAL:HG21	1.74	0.68
1:A:117:GLN:HG3	1:A:121:LYS:HE2	1.78	0.65
1:A:116:GLN:O	1:A:120:GLU:HG3	1.98	0.64
1:A:131:GLU:OE1	1:A:177:ILE:HG12	1.98	0.63
1:A:131:GLU:OE1	1:A:174:CYS:HB3	2.00	0.61
1:A:80:THR:O	1:A:84:LYS:HG3	2.01	0.61
1:A:140:GLY:O	1:A:176:LYS:HG3	2.02	0.60
1:A:117:GLN:HG3	1:A:121:LYS:CE	2.33	0.58
1:A:95:GLN:NE2	1:A:155:HIS:HE1	2.01	0.56
1:A:93:THR:CG2	1:A:126:LEU:HD22	2.35	0.56
1:A:6:HIS:HB3	1:A:50:MET:HE1	1.89	0.55
1:A:16:LEU:HD13	1:A:155:HIS:HD2	1.73	0.54
1:A:16:LEU:HD13	1:A:155:HIS:CD2	2.44	0.52
1:A:196:LYS:HE2	4:A:463:HOH:O	2.09	0.52
1:A:38:TRP:CE2	1:A:42:LYS:HG3	2.44	0.51
1:A:116:GLN:NE2	1:A:120:GLU:OE2	2.40	0.51
1:A:8:PHE:HA	1:A:33:VAL:O	2.11	0.51
1:A:6:HIS:HB3	1:A:50:MET:HE2	1.93	0.50
1:A:31:ARG:NH2	4:A:496:HOH:O	2.44	0.50
1:A:45:MET:HE3	1:A:45:MET:HA	1.92	0.50
1:A:149:MET:HE1	1:A:187:CYS:SG	2.52	0.50
1:A:95:GLN:HE21	1:A:155:HIS:CE1	2.24	0.49
1:A:16:LEU:HD22	1:A:155:HIS:CD2	2.48	0.49
1:A:82:LEU:O	1:A:86:ARG:HG2	2.13	0.49
1:A:50:MET:HB3	1:A:51:PRO:HA	1.95	0.48
1:A:93:THR:HG23	1:A:126:LEU:HD22	1.95	0.48
1:A:38:TRP:CZ2	1:A:42:LYS:HG3	2.50	0.47
1:A:54:ASP:OD2	1:A:59:LYS:NZ	2.39	0.47
1:A:95:GLN:NE2	1:A:155:HIS:CE1	2.83	0.46
1:A:24:HIS:ND1	1:A:26:GLU:OE2	2.48	0.46
1:A:178:VAL:HG22	4:A:393:HOH:O	2.16	0.45
1:A:108:PRO:O	1:A:112:LYS:HE3	2.18	0.44
1:A:111:ALA:O	1:A:115:VAL:HG23	2.18	0.44
2:A:300:SO4:O4	3:A:203:GBI:I3'	3.05	0.44
1:A:64:MET:HE3	1:A:95:GLN:OE1	2.18	0.43
1:A:149:MET:HE2	1:A:184:VAL:HG22	2.01	0.43
1:A:174:CYS:O	1:A:178:VAL:HG13	2.18	0.43
1:A:54:ASP:OD1	1:A:54:ASP:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:34:GLU:HB3	4:A:443:HOH:O	2.21	0.41
1:A:127:ALA:N	1:A:128:PRO:CD	2.84	0.41
1:A:61:SER:O	1:A:62:GLN:HB2	2.20	0.40
1:A:169:GLU:CD	1:A:169:GLU:H	2.29	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	200/202 (99%)	190 (95%)	9 (4%)	1 (0%)	24 27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	GLN

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	172/172 (100%)	149 (87%)	23 (13%)	4 3

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	5	LEU
1	A	30	ASP
1	A	41	LEU
1	A	70	LEU
1	A	82	LEU
1	A	86	ARG
1	A	101	VAL
1	A	105	LYS
1	A	112	LYS
1	A	117	GLN
1	A	121	LYS
1	A	130	LEU
1	A	141	ASP
1	A	147	ASN
1	A	155	HIS
1	A	171	LEU
1	A	172	LYS
1	A	178	VAL
1	A	194	LEU
1	A	196	LYS
1	A	197	ARG
1	A	200	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	95	GLN
1	A	117	GLN
1	A	147	ASN
1	A	155	HIS

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	GBI	A	203	-	27,28,28	1.17	3 (11%)	33,36,36	1.28	4 (12%)
2	SO4	A	300	-	4,4,4	0.78	0	6,6,6	1.10	1 (16%)
3	GBI	A	204	-	27,28,28	0.92	0	33,36,36	1.27	4 (12%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GBI	A	203	-	-	2/28/28/28	0/1/1/1
3	GBI	A	204	-	-	14/28/28/28	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	203	GBI	O12-C1	-2.74	1.21	1.30
3	A	203	GBI	O32-C3	-2.71	1.21	1.30
3	A	203	GBI	C'-SG2	2.10	1.85	1.82

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	204	GBI	CB2-SG2-C'	-3.01	94.95	101.24
3	A	203	GBI	C2-CA2-N2	-2.93	103.17	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	203	GBI	CB2-SG2-C'	-2.86	95.27	101.24
3	A	204	GBI	C4'-C3'-I3'	-2.71	115.64	119.69
3	A	204	GBI	C2'-C3'-I3'	2.59	123.04	119.38
3	A	203	GBI	C1'-C'-SG2	-2.44	108.24	113.97
2	A	300	SO4	O4-S-O1	-2.27	97.70	109.56
3	A	204	GBI	CB2-CA2-C2	2.11	114.22	109.46
3	A	203	GBI	C4'-C3'-I3'	-2.02	116.68	119.69

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	204	GBI	CG1-CD1-N2-CA2
3	A	204	GBI	N2-CA2-CB2-SG2
3	A	204	GBI	OE1-CD1-N2-CA2
3	A	204	GBI	O31-C3-CA3-N3
3	A	204	GBI	O2-C2-CA2-CB2
3	A	204	GBI	O32-C3-CA3-N3
3	A	204	GBI	N3-C2-CA2-CB2
3	A	203	GBI	C1-CA1-CB1-CG1
3	A	204	GBI	SG2-C'-C1'-C2'
3	A	204	GBI	SG2-C'-C1'-C6'
3	A	204	GBI	O12-C1-CA1-N1
3	A	203	GBI	O12-C1-CA1-N1
3	A	204	GBI	C3-CA3-N3-C2
3	A	204	GBI	C2-CA2-CB2-SG2
3	A	204	GBI	N1-CA1-CB1-CG1
3	A	204	GBI	O11-C1-CA1-N1

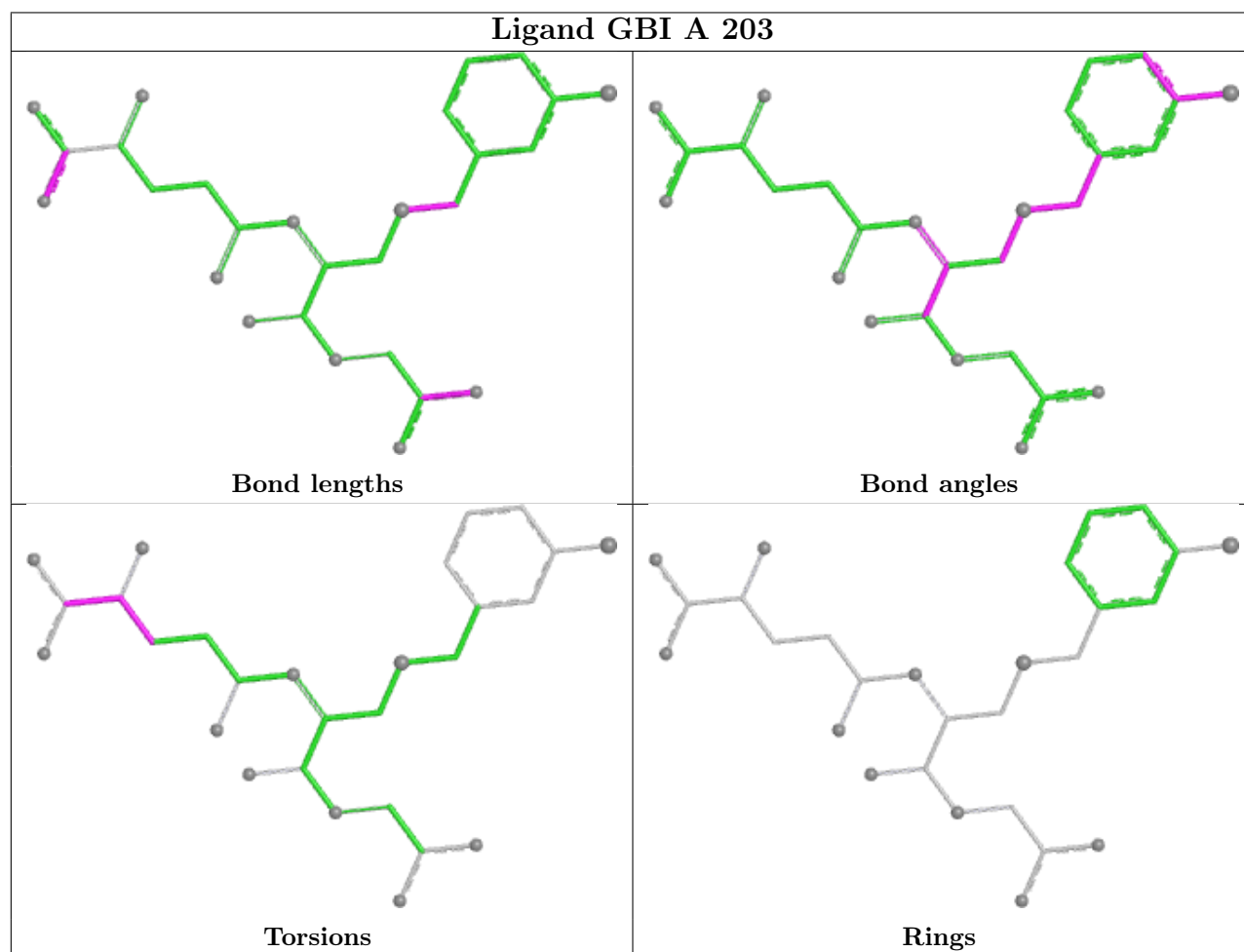
There are no ring outliers.

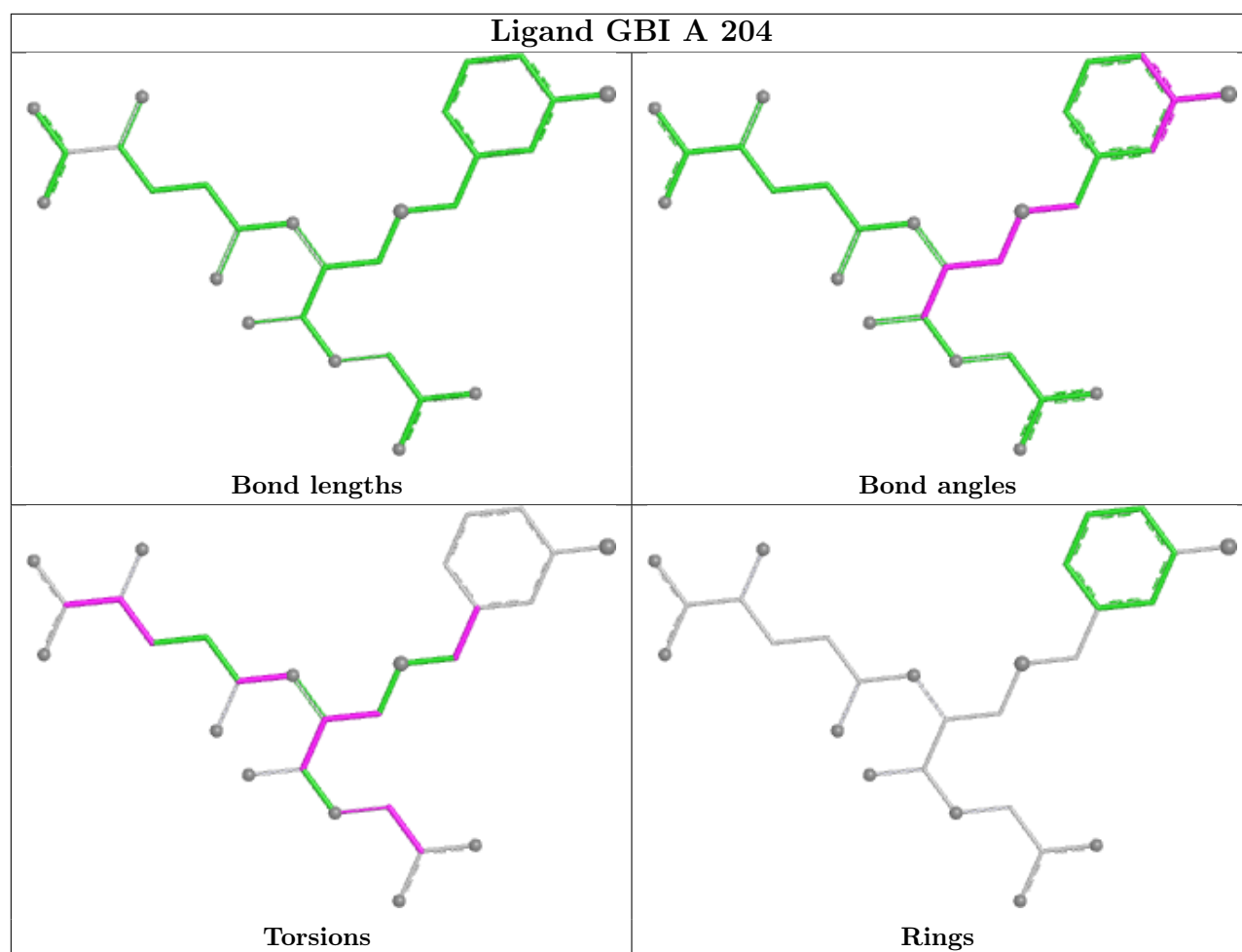
2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	203	GBI	1	0
2	A	300	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be

highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.