



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 20, 2017 – 11:31 PM EST

PDB ID : 3GST  
Title : STRUCTURE OF THE XENOBIOTIC SUBSTRATE BINDING SITE OF A GLUTATHIONE S-TRANSFERASE AS REVEALED BY X-RAY CRYSTALLOGRAPHIC ANALYSIS OF PRODUCT COMPLEXES WITH THE DIASTEREOMERS OF 9-(S-GLUTATHIONYL)-10-HYDROXY-9, 10-DIHYDROPHENANTHRENE  
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Deposited on : unknown  
Resolution : 1.90 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20030345



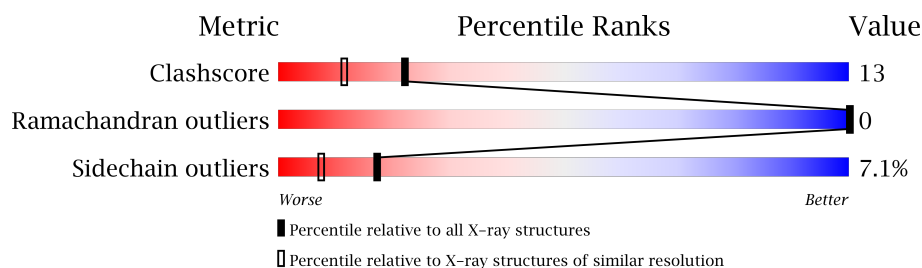
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	217	
1	B	217	



## 2 Entry composition [i](#)

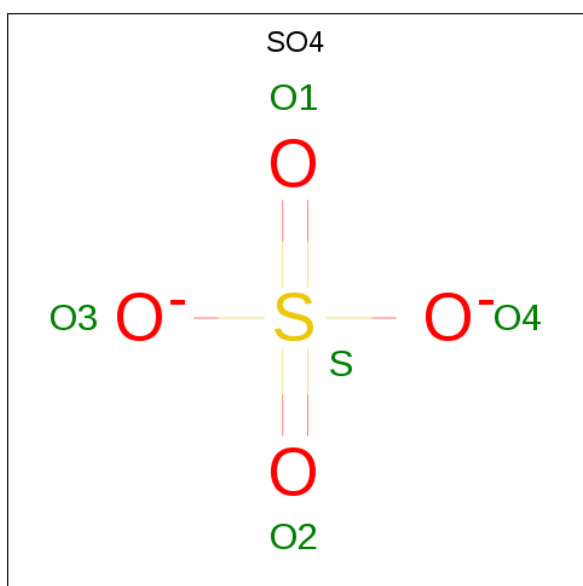
There are 4 unique types of molecules in this entry. The entry contains 4178 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
1	A	217	Total	C	N	O	S	0	0	0
			1818	1177	303	327	11			
1	B	217	Total	C	N	O	S	0	0	0
			1818	1177	303	327	11			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

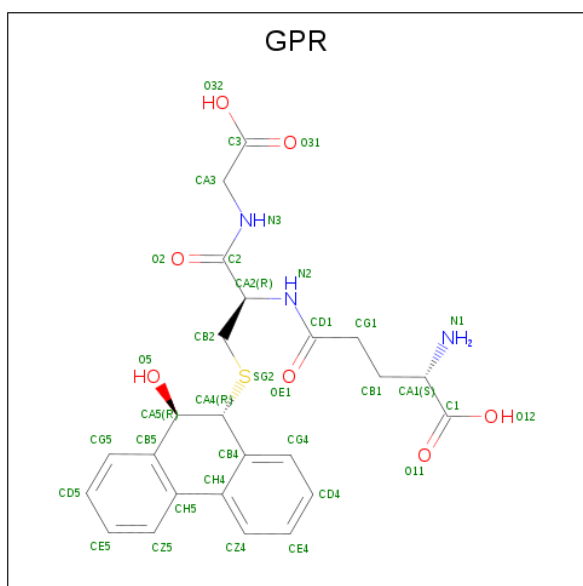
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is (9R,10R)-9-(S-GLUTATHIONYL)-10-HYDROXY-9,10-DIHYDROPHENANTHRENE (three-letter code: GPR) (formula:  $C_{24}H_{27}N_3O_7S$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0
			35	24	3	7	1	
3	B	1	Total	C	N	O	S	0
			35	24	3	7	1	

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	252	Total	O	0	0
			252	252		
4	B	185	Total	O	0	0
			185	185		

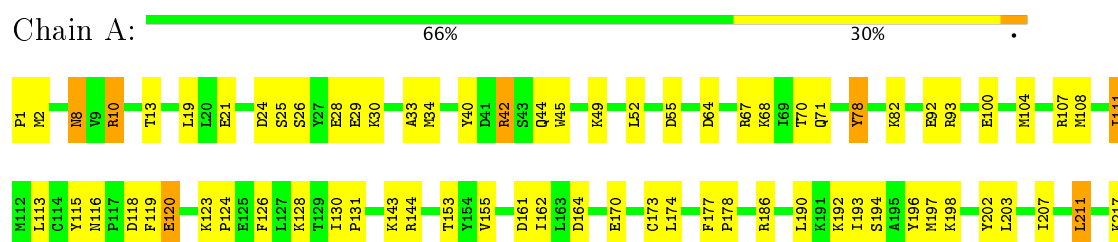


### 3 Residue-property plots [i](#)

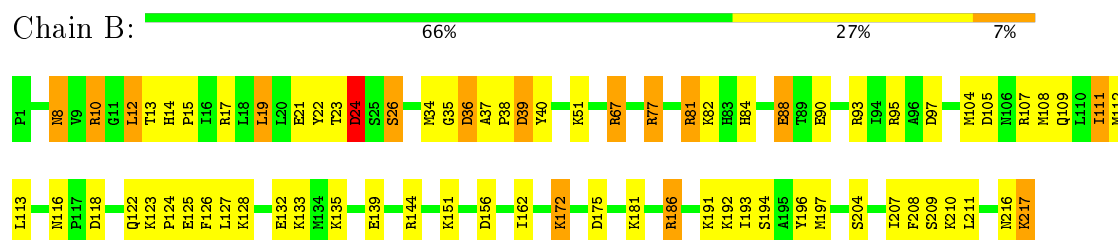
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

#### • Molecule 1: GLUTATHIONE S-TRANSFERASE



#### • 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	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.24Å 69.44Å 81.28Å 90.00° 106.01° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	GPRLSA	Depositor
R, $R_{free}$	0.159 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4178	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, GPR

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	1.06	0/1867	1.64	18/2515 (0.7%)
1	B	1.13	1/1867 (0.1%)	1.63	20/2515 (0.8%)
All	All	1.10	1/3734 (0.0%)	1.64	38/5030 (0.8%)

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	3
1	B	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	204	SER	CB-OG	6.08	1.50	1.42

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ARG	NE-CZ-NH1	24.44	132.52	120.30
1	B	81	ARG	NE-CZ-NH1	14.06	127.33	120.30
1	A	93	ARG	CD-NE-CZ	13.93	143.10	123.60
1	B	95	ARG	NE-CZ-NH1	12.69	126.64	120.30
1	B	17	ARG	NE-CZ-NH1	12.52	126.56	120.30
1	A	93	ARG	NE-CZ-NH2	-12.26	114.17	120.30
1	B	144	ARG	CD-NE-CZ	12.14	140.60	123.60
1	A	107	ARG	NE-CZ-NH2	11.86	126.23	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	ARG	NE-CZ-NH2	-11.79	114.40	120.30
1	B	144	ARG	NE-CZ-NH1	11.13	125.87	120.30
1	B	81	ARG	NE-CZ-NH2	-10.90	114.85	120.30
1	A	10	ARG	NE-CZ-NH2	-10.12	115.24	120.30
1	B	156	ASP	CB-CG-OD2	-10.03	109.28	118.30
1	B	24	ASP	CB-CG-OD2	-9.30	109.93	118.30
1	A	161	ASP	CB-CG-OD1	8.23	125.70	118.30
1	B	105	ASP	CB-CG-OD1	7.78	125.30	118.30
1	A	202	TYR	CB-CG-CD2	-7.37	116.58	121.00
1	A	120	GLU	OE1-CD-OE2	-7.17	114.69	123.30
1	B	144	ARG	NE-CZ-NH2	-7.17	116.72	120.30
1	A	120	GLU	CG-CD-OE1	7.13	132.56	118.30
1	B	10	ARG	NE-CZ-NH2	-6.71	116.94	120.30
1	B	26	SER	CB-CA-C	6.47	122.40	110.10
1	A	29	GLU	OE1-CD-OE2	6.29	130.85	123.30
1	A	107	ARG	NE-CZ-NH1	-6.01	117.30	120.30
1	B	84	HIS	CA-CB-CG	-5.86	103.64	113.60
1	B	39	ASP	CB-CG-OD2	5.76	123.48	118.30
1	B	93	ARG	NE-CZ-NH2	-5.57	117.52	120.30
1	A	93	ARG	NH1-CZ-NH2	-5.54	113.31	119.40
1	B	90	GLU	CG-CD-OE1	5.52	129.35	118.30
1	B	77	ARG	CD-NE-CZ	-5.51	115.88	123.60
1	A	164	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	B	97	ASP	CB-CG-OD1	5.23	123.00	118.30
1	B	77	ARG	NE-CZ-NH2	-5.19	117.70	120.30
1	A	92	GLU	CG-CD-OE2	-5.17	107.97	118.30
1	A	78	TYR	CB-CG-CD1	-5.08	117.95	121.00
1	A	55	ASP	CB-CG-OD2	-5.08	113.73	118.30
1	A	42	ARG	CD-NE-CZ	-5.05	116.53	123.60
1	A	100	GLU	OE1-CD-OE2	-5.04	117.25	123.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	144	ARG	Sidechain
1	A	186	ARG	Sidechain
1	A	42	ARG	Sidechain
1	B	186	ARG	Sidechain



## 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	1818	0	1805	36	0
1	B	1818	0	1805	62	0
2	A	25	0	0	0	0
2	B	10	0	0	2	0
3	A	35	0	25	0	0
3	B	35	0	25	2	0
4	A	252	0	0	2	0
4	B	185	0	0	4	0
All	All	4178	0	3660	98	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.

All (98) 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:108:MET:O	1:B:112:MET:HG3	1.69	0.92
1:A:1:PRO:HG2	1:A:28:GLU:OE1	1.75	0.87
1:A:34:MET:CE	1:A:40:TYR:HB3	2.05	0.87
1:B:123:LYS:HB3	1:B:124:PRO:HD3	1.64	0.79
1:A:34:MET:HE3	1:A:40:TYR:HB3	1.64	0.78
1:A:52:LEU:O	1:A:68:LYS:NZ	2.16	0.77
1:B:88:GLU:HG2	1:B:151:LYS:NZ	2.00	0.77
1:B:36:ASP:N	1:B:36:ASP:OD1	2.18	0.76
1:A:124:PRO:O	1:A:128:LYS:HE3	1.90	0.71
1:B:35:GLY:O	1:B:40:TYR:HD1	1.74	0.69
1:B:40:TYR:CE2	1:B:211:LEU:HD13	2.28	0.69
1:B:172:LYS:HD2	1:B:175:ASP:OD2	1.93	0.67
1:A:34:MET:HE2	1:A:40:TYR:HB3	1.75	0.67
1:B:34:MET:HE3	1:B:209:SER:HB2	1.77	0.66
1:B:88:GLU:HG2	1:B:151:LYS:HZ1	1.59	0.66
1:B:109:GLN:OE1	1:B:133:LYS:HE2	1.96	0.65
1:B:40:TYR:CD2	1:B:211:LEU:HD13	2.34	0.63
1:B:34:MET:CE	1:B:209:SER:HB2	2.31	0.61
1:A:104:MET:O	1:A:108:MET:HG2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ASP:OD2	1:B:192:LYS:HE2	2.03	0.59
1:B:24:ASP:OD2	1:B:192:LYS:CE	2.52	0.57
1:A:64:ASP:OD2	1:A:82:LYS:NZ	2.37	0.57
1:B:38:PRO:HG2	1:B:39:ASP:N	2.20	0.57
1:B:125:GLU:OE2	1:B:128:LYS:NZ	2.36	0.57
1:B:67:ARG:NH2	4:B:287:HOH:O	2.37	0.56
1:B:175:ASP:OD1	1:B:181:LYS:NZ	2.30	0.56
1:A:116:ASN:OD1	1:A:118:ASP:HB2	2.06	0.55
1:B:22:TYR:O	1:B:192:LYS:HD2	2.07	0.54
1:B:116:ASN:OD1	1:B:118:ASP:HB2	2.08	0.53
1:A:113:LEU:HD22	1:A:126:PHE:CG	2.44	0.53
1:B:40:TYR:CE2	1:B:211:LEU:CD1	2.91	0.53
1:B:35:GLY:O	1:B:40:TYR:HA	2.09	0.53
1:B:19:LEU:HD22	1:B:23:THR:HG23	1.90	0.53
1:B:111:ILE:HG12	1:B:208:PHE:CE1	2.43	0.52
1:B:186:ARG:HD3	2:B:218:SO4:O1	2.09	0.52
1:B:135:LYS:NZ	4:B:289:HOH:O	2.42	0.52
1:A:21:GLU:HG3	1:A:196:TYR:CD1	2.44	0.52
1:B:14:HIS:N	1:B:15:PRO:CD	2.73	0.51
1:A:45:TRP:CZ2	1:A:49:LYS:HG3	2.46	0.51
1:B:8:ASN:H	1:B:8:ASN:HD22	1.57	0.51
1:A:10:ARG:HB3	1:A:207:ILE:HA	1.93	0.51
1:B:111:ILE:HG22	1:B:112:MET:N	2.26	0.51
1:A:190:LEU:HB2	1:A:193:ILE:HG12	1.94	0.50
1:A:111:ILE:HD13	1:A:115:TYR:CE1	2.48	0.49
1:B:193:ILE:O	1:B:197:MET:HG3	2.13	0.49
1:A:153:THR:OG1	1:A:155:VAL:HG22	2.13	0.48
1:B:217:LYS:NZ	4:B:352:HOH:O	2.43	0.48
1:B:135:LYS:NZ	1:B:139:GLU:OE2	2.41	0.48
1:B:108:MET:O	1:B:112:MET:CG	2.52	0.48
1:A:115:TYR:CE2	1:A:211:LEU:HD23	2.49	0.48
1:B:8:ASN:H	1:B:8:ASN:ND2	2.12	0.47
1:B:123:LYS:HB3	1:B:124:PRO:CD	2.39	0.47
1:B:8:ASN:N	1:B:8:ASN:HD22	2.12	0.47
1:B:81:ARG:HD3	2:B:219:SO4:O4	2.15	0.47
1:A:8:ASN:N	1:A:8:ASN:HD22	2.13	0.47
1:A:2:MET:CE	1:A:25:SER:HB3	2.45	0.46
1:B:123:LYS:HE2	1:B:127:LEU:HD11	1.97	0.46
1:A:33:ALA:H	1:A:44:GLN:NE2	2.14	0.46
1:B:113:LEU:HD22	1:B:126:PHE:CG	2.50	0.46
1:B:12:LEU:HB3	1:B:107:ARG:HD3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:21:GLU:HG3	1:B:196:TYR:CD1	2.51	0.46
1:B:38:PRO:CG	1:B:39:ASP:N	2.79	0.46
1:B:111:ILE:HG13	3:B:220:GPR:CE5	2.46	0.46
1:B:123:LYS:CB	1:B:124:PRO:HD3	2.42	0.46
1:B:194:SER:HA	1:B:197:MET:HE2	1.97	0.45
1:B:77:ARG:HH11	1:B:77:ARG:HD3	1.57	0.45
1:B:37:ALA:HB2	1:B:40:TYR:CZ	2.51	0.45
1:A:177:PHE:HA	1:A:178:PRO:HD2	1.83	0.45
1:B:88:GLU:CG	1:B:151:LYS:NZ	2.77	0.45
1:A:78:TYR:OH	1:A:82:LYS:NZ	2.36	0.45
1:B:24:ASP:OD2	1:B:192:LYS:NZ	2.49	0.45
1:B:88:GLU:CG	1:B:151:LYS:HZ1	2.26	0.44
1:A:13:THR:O	1:A:13:THR:HG22	2.17	0.44
1:B:10:ARG:HB3	1:B:207:ILE:HA	2.00	0.44
1:B:111:ILE:HG13	3:B:220:GPR:CD5	2.48	0.44
1:B:124:PRO:O	1:B:128:LYS:HG3	2.19	0.43
1:A:70:THR:O	1:A:71:GLN:HB2	2.19	0.43
1:B:38:PRO:HG2	1:B:39:ASP:OD2	2.19	0.43
1:A:8:ASN:H	1:A:8:ASN:HD22	1.65	0.42
1:B:104:MET:O	1:B:108:MET:HG2	2.19	0.42
1:B:111:ILE:CG2	1:B:112:MET:N	2.81	0.42
1:B:122:GLN:HB3	1:B:122:GLN:HE21	1.67	0.42
1:B:34:MET:HE2	1:B:34:MET:HB3	1.58	0.42
1:A:123:LYS:N	1:A:124:PRO:CD	2.81	0.42
1:A:193:ILE:HG22	1:A:197:MET:CE	2.50	0.42
1:A:24:ASP:OD2	1:A:192:LYS:HE3	2.20	0.42
1:B:82:LYS:HE3	4:B:327:HOH:O	2.20	0.42
1:A:194:SER:O	1:A:198:LYS:HD2	2.19	0.41
1:A:2:MET:HG2	1:A:64:ASP:HB2	2.03	0.41
1:A:119:PHE:CZ	1:A:123:LYS:HD3	2.56	0.41
1:A:162:ILE:HD12	1:A:162:ILE:HA	1.87	0.41
1:B:209:SER:O	1:B:216:ASN:HB2	2.21	0.41
1:B:40:TYR:CE1	1:B:210:LYS:HB3	2.55	0.41
1:A:130:ILE:N	1:A:131:PRO:CD	2.84	0.41
1:A:192:LYS:HD3	4:A:286:HOH:O	2.21	0.41
1:B:13:THR:HG22	1:B:13:THR:O	2.21	0.40
1:A:30:LYS:NZ	4:A:318:HOH:O	2.42	0.40
1:A:193:ILE:CG2	1:A:197:MET:CE	2.99	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	215/217 (99%)	212 (99%)	3 (1%)	0	100	100
1	B	215/217 (99%)	211 (98%)	4 (2%)	0	100	100
All	All	430/434 (99%)	423 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	197/197 (100%)	184 (93%)	13 (7%)	19	9
1	B	197/197 (100%)	182 (92%)	15 (8%)	15	6
All	All	394/394 (100%)	366 (93%)	28 (7%)	17	7

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	19	LEU
1	A	26	SER
1	A	67	ARG
1	A	111	ILE
1	A	120	GLU
1	A	143	LYS
1	A	170	GLU

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Mol	Chain	Res	Type
1	A	173	CYS
1	A	174	LEU
1	A	203	LEU
1	A	211	LEU
1	A	217	LYS
1	B	8	ASN
1	B	12	LEU
1	B	19	LEU
1	B	24	ASP
1	B	26	SER
1	B	36	ASP
1	B	51	LYS
1	B	67	ARG
1	B	88	GLU
1	B	111	ILE
1	B	132	GLU
1	B	162	ILE
1	B	172	LYS
1	B	191	LYS
1	B	217	LYS

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

Mol	Chain	Res	Type
1	A	8	ASN
1	A	44	GLN
1	B	8	ASN
1	B	122	GLN

### 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 carbohydrates in this entry.



## 5.6 Ligand geometry

9 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
2	SO4	A	218	-	4,4,4	0.90	0	6,6,6	0.33	0
2	SO4	A	219	-	4,4,4	0.86	0	6,6,6	0.23	0
2	SO4	A	220	-	4,4,4	0.91	0	6,6,6	0.33	0
2	SO4	A	221	-	4,4,4	0.95	0	6,6,6	0.45	0
2	SO4	A	222	-	4,4,4	0.89	0	6,6,6	0.33	0
3	GPR	A	223	-	27,37,37	1.13	1 (3%)	34,51,51	1.59	7 (20%)
2	SO4	B	218	-	4,4,4	0.76	0	6,6,6	0.38	0
2	SO4	B	219	-	4,4,4	1.10	0	6,6,6	0.38	0
3	GPR	B	220	-	27,37,37	1.63	7 (25%)	34,51,51	1.87	10 (29%)

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	SO4	A	218	-	-	0/0/0/0	0/0/0/0
2	SO4	A	219	-	-	0/0/0/0	0/0/0/0
2	SO4	A	220	-	-	0/0/0/0	0/0/0/0
2	SO4	A	221	-	-	0/0/0/0	0/0/0/0
2	SO4	A	222	-	-	0/0/0/0	0/0/0/0
3	GPR	A	223	-	-	0/21/43/43	0/3/3/3
2	SO4	B	218	-	-	0/0/0/0	0/0/0/0
2	SO4	B	219	-	-	0/0/0/0	0/0/0/0
3	GPR	B	220	-	-	0/21/43/43	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	220	GPR	C2-N3	-3.09	1.27	1.33
3	B	220	GPR	CH4-CB4	2.02	1.43	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	223	GPR	CG4-CB4	2.15	1.42	1.39
3	B	220	GPR	CZ5-CH5	2.31	1.43	1.39
3	B	220	GPR	CA2-N2	2.51	1.51	1.45
3	B	220	GPR	CA3-N3	2.66	1.51	1.46
3	B	220	GPR	CG4-CB4	3.15	1.43	1.39
3	B	220	GPR	O5-CA5	3.54	1.49	1.42

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	220	GPR	O5-CA5-CB5	-3.90	102.60	110.47
3	A	223	GPR	OE1-CD1-CG1	-3.54	115.35	122.01
3	B	220	GPR	CG1-CB1-CA1	-3.33	106.07	113.84
3	A	223	GPR	CA3-N3-C2	-3.04	118.28	122.39
3	B	220	GPR	O2-C2-CA2	-2.97	114.03	120.43
3	B	220	GPR	OE1-CD1-CG1	-2.95	116.47	122.01
3	A	223	GPR	CG1-CB1-CA1	-2.95	106.96	113.84
3	B	220	GPR	CA2-CB2-SG2	-2.89	105.87	113.12
3	A	223	GPR	O5-CA5-CB5	-2.71	105.01	110.47
3	A	223	GPR	CA5-CA4-SG2	-2.61	102.78	110.94
3	B	220	GPR	CA2-N2-CD1	-2.14	115.97	121.62
3	A	223	GPR	OE1-CD1-N2	2.02	126.46	122.97
3	B	220	GPR	CA3-N3-C2	2.07	125.19	122.39
3	B	220	GPR	CG5-CB5-CH5	2.14	122.23	119.42
3	B	220	GPR	O2-C2-N3	2.56	128.00	123.07
3	A	223	GPR	CB1-CG1-CD1	2.57	119.00	113.18
3	B	220	GPR	OE1-CD1-N2	3.80	129.53	122.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	218	SO4	1	0
2	B	219	SO4	1	0
3	B	220	GPR	2	0



## 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

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.