



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:58 PM GMT

PDB ID : 6GSW  
Title : FIRST-SPHERE AND SECOND-SPHERE ELECTROSTATIC EFFECTS IN  
THE ACTIVE SITE OF A CLASS MU GLUTATHIONE TRANSFERASE  
Authors : Xiao, G.; Ji, X.; Armstrong, R.N.; Gilliland, G.L.  
Deposited on : 1996-01-26  
Resolution : 1.85 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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

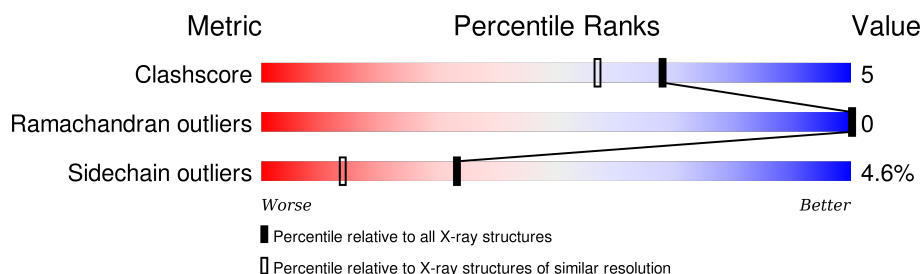
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)

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	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GPS	A	221	X	-	-	-
3	GPS	B	218	X	-	-	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4045 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.

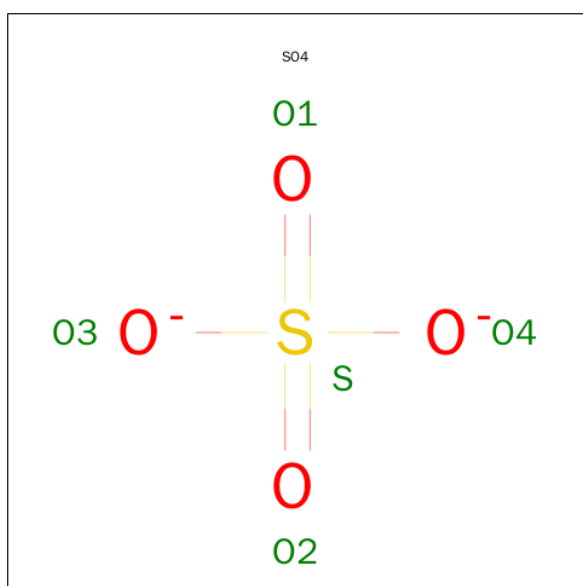
- Molecule 1 is a protein called MU CLASS GLUTATHIONE S-TRANSFERASE OF ISOENZYME 3-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1818	1178	303	326	11			
1	B	217	Total	C	N	O	S	0	0	0
			1818	1178	303	326	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	VAL	THR	ENGINEERED	UNP P04905
B	13	VAL	THR	ENGINEERED	UNP P04905

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



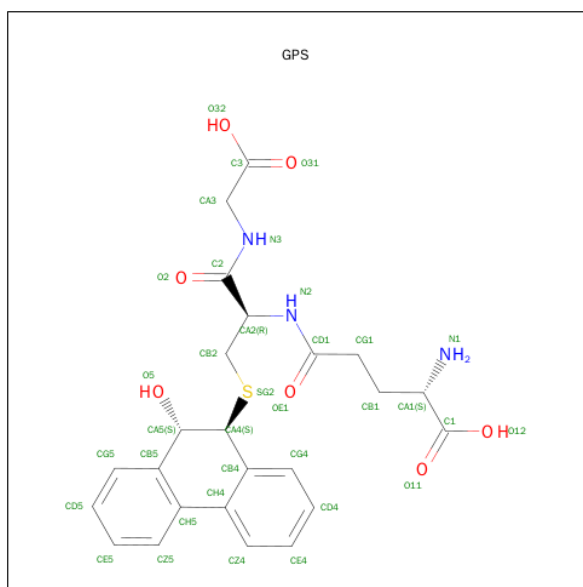
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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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		

- Molecule 3 is L-GAMMA-GLUTAMYL-S-[(9S,10S)-10-HYDROXY-9,10-DIHYDRO PHENANTHREN-9-YL]-L-CYSTEINYLGLYCINE (three-letter code: GPS) (formula: C<sub>24</sub>H<sub>27</sub>N<sub>3</sub>O<sub>7</sub>S).



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

- Molecule 4 is water.

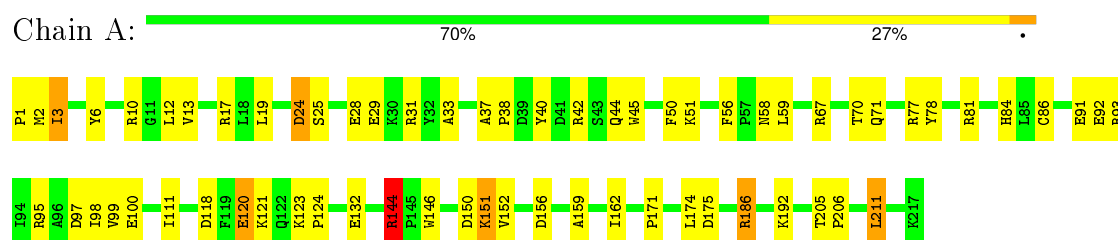
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	196	Total	O	0	0
			196	196		
4	B	128	Total	O	0	0
			128	128		

### 3 Residue-property plots

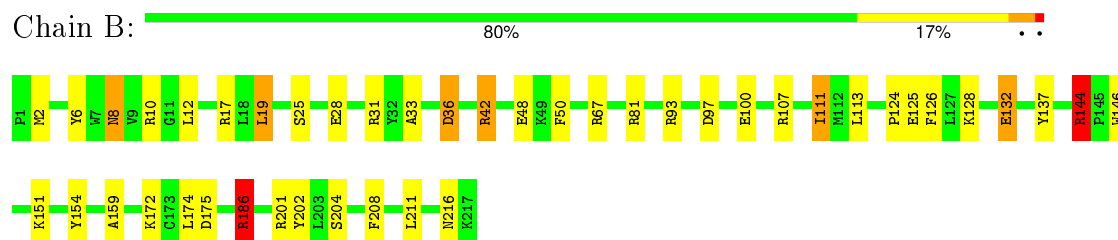
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

#### • Molecule 1: MU CLASS GLUTATHIONE S-TRANSFERASE OF ISOENZYME 3-3



#### • Molecule 1: MU CLASS GLUTATHIONE S-TRANSFERASE OF ISOENZYME 3-3



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.53 Å   69.28 Å   81.41 Å 90.00°   105.74°   90.00°	Depositor
Resolution (Å)	6.00 – 1.85	Depositor
% Data completeness (in resolution range)	84.2 (6.00-1.85)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	GPRLSA	Depositor
R, $R_{free}$	0.163 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4045	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	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, GPS

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	1.03	2/1867 (0.1%)	1.71	42/2515 (1.7%)
1	B	1.03	1/1867 (0.1%)	1.90	35/2515 (1.4%)
All	All	1.03	3/3734 (0.1%)	1.81	77/5030 (1.5%)

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
1	B	0	3
All	All	0	5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	100	GLU	CD-OE1	-6.82	1.18	1.25
1	A	100	GLU	CD-OE1	-5.87	1.19	1.25
1	A	29	GLU	CD-OE1	-5.42	1.19	1.25

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	186	ARG	NE-CZ-NH1	27.59	134.10	120.30
1	B	186	ARG	CD-NE-CZ	25.79	159.70	123.60
1	B	186	ARG	NE-CZ-NH2	-19.07	110.76	120.30
1	A	77	ARG	NE-CZ-NH1	13.85	127.22	120.30
1	B	144	ARG	NE-CZ-NH2	-12.78	113.91	120.30
1	B	42	ARG	NE-CZ-NH1	12.61	126.60	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	77	ARG	NE-CZ-NH2	-12.12	114.24	120.30
1	B	42	ARG	NE-CZ-NH2	-12.01	114.30	120.30
1	B	10	ARG	NE-CZ-NH2	-11.44	114.58	120.30
1	A	186	ARG	NE-CZ-NH1	10.85	125.72	120.30
1	A	42	ARG	NE-CZ-NH1	10.70	125.65	120.30
1	B	154	TYR	CB-CG-CD2	-9.75	115.15	121.00
1	B	17	ARG	NE-CZ-NH2	-9.59	115.51	120.30
1	A	144	ARG	NE-CZ-NH2	9.01	124.81	120.30
1	A	81	ARG	NE-CZ-NH2	-8.73	115.94	120.30
1	B	17	ARG	NE-CZ-NH1	8.67	124.63	120.30
1	B	93	ARG	NE-CZ-NH2	-8.35	116.12	120.30
1	A	156	ASP	CB-CG-OD2	-8.26	110.87	118.30
1	A	31	ARG	NE-CZ-NH2	-8.12	116.24	120.30
1	A	118	ASP	CB-CG-OD2	-8.09	111.02	118.30
1	A	67	ARG	NE-CZ-NH1	7.96	124.28	120.30
1	A	120	GLU	CB-CG-CD	7.92	135.59	114.20
1	A	67	ARG	NE-CZ-NH2	-7.73	116.44	120.30
1	B	107	ARG	NE-CZ-NH1	7.65	124.13	120.30
1	B	93	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	B	97	ASP	CB-CG-OD1	7.23	124.81	118.30
1	A	24	ASP	CB-CG-OD1	7.20	124.78	118.30
1	B	28	GLU	N-CA-CB	6.88	122.98	110.60
1	A	42	ARG	CD-NE-CZ	6.51	132.72	123.60
1	A	144	ARG	CD-NE-CZ	6.48	132.67	123.60
1	B	137	TYR	CB-CG-CD1	-6.38	117.17	121.00
1	A	175	ASP	CB-CG-OD2	-6.38	112.56	118.30
1	B	81	ARG	NE-CZ-NH2	-6.35	117.13	120.30
1	B	31	ARG	NE-CZ-NH2	6.33	123.46	120.30
1	A	56	PHE	CB-CG-CD2	-6.30	116.39	120.80
1	B	28	GLU	CA-CB-CG	6.23	127.10	113.40
1	A	97	ASP	CB-CG-OD2	-6.19	112.73	118.30
1	A	17	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	92	GLU	OE1-CD-OE2	6.00	130.50	123.30
1	A	99	VAL	CA-CB-CG2	-6.00	101.90	110.90
1	B	50	PHE	CB-CG-CD1	-5.94	116.64	120.80
1	B	6	TYR	CB-CG-CD2	5.83	124.50	121.00
1	A	58	ASN	N-CA-CB	-5.79	100.18	110.60
1	B	33	ALA	N-CA-CB	-5.77	102.03	110.10
1	A	42	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	B	186	ARG	NH1-CZ-NH2	-5.70	113.13	119.40
1	A	132	GLU	CG-CD-OE2	-5.69	106.91	118.30
1	A	81	ARG	NE-CZ-NH1	5.69	123.14	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	121	LYS	CA-CB-CG	-5.64	100.99	113.40
1	A	93	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	B	211	LEU	CA-CB-CG	5.53	128.02	115.30
1	A	120	GLU	CG-CD-OE1	5.51	129.33	118.30
1	A	6	TYR	CB-CG-CD2	5.51	124.31	121.00
1	A	3	ILE	O-C-N	5.48	131.46	122.70
1	B	137	TYR	CB-CG-CD2	5.47	124.28	121.00
1	A	211	LEU	CA-CB-CG	5.43	127.79	115.30
1	B	97	ASP	CB-CG-OD2	-5.43	113.42	118.30
1	B	201	ARG	CD-NE-CZ	5.42	131.18	123.60
1	B	36	ASP	CB-CG-OD2	-5.41	113.43	118.30
1	B	48	GLU	CA-CB-CG	5.39	125.26	113.40
1	A	159	ALA	O-C-N	5.38	131.31	122.70
1	A	24	ASP	CB-CG-OD2	-5.34	113.50	118.30
1	A	45	TRP	O-C-N	5.31	131.19	122.70
1	B	19	LEU	CB-CG-CD2	5.30	120.00	111.00
1	A	98	ILE	O-C-N	5.29	131.17	122.70
1	B	132	GLU	CB-CG-CD	5.29	128.48	114.20
1	B	28	GLU	O-C-N	5.21	131.04	122.70
1	A	84	HIS	CA-CB-CG	-5.16	104.83	113.60
1	A	144	ARG	NH1-CZ-NH2	-5.14	113.74	119.40
1	A	59	LEU	CB-CG-CD1	-5.12	102.29	111.00
1	A	10	ARG	NE-CZ-NH2	-5.12	117.74	120.30
1	A	50	PHE	CB-CG-CD1	-5.11	117.22	120.80
1	B	154	TYR	CB-CG-CD1	5.07	124.04	121.00
1	B	159	ALA	O-C-N	5.06	130.80	122.70
1	B	201	ARG	NE-CZ-NH1	5.04	122.82	120.30
1	A	78	TYR	CB-CG-CD1	-5.03	117.98	121.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	144	ARG	Sidechain
1	A	186	ARG	Sidechain
1	B	144	ARG	Sidechain
1	B	42	ARG	Sidechain
1	B	67	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	1807	20	0
1	B	1818	0	1807	13	0
2	A	15	0	0	1	0
3	A	35	0	25	1	0
3	B	35	0	25	1	0
4	A	196	0	0	3	0
4	B	128	0	0	1	0
All	All	4045	0	3664	35	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 5.

All (35) 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:125:GLU:OE1	1:B:128:LYS:NZ	2.14	0.76
1:A:24:ASP:OD2	1:A:192:LYS:HE2	1.98	0.63
1:A:95:ARG:NH2	1:A:144:ARG:HE	1.98	0.61
1:B:125:GLU:HA	1:B:128:LYS:HE3	1.83	0.59
1:A:33:ALA:H	1:A:44:GLN:NE2	2.00	0.58
1:A:40:TYR:CE2	1:A:211:LEU:HD13	2.40	0.55
1:A:2:MET:CE	1:A:25:SER:HB3	2.40	0.52
3:A:221:GPS:HB2	3:A:221:GPS:HG4	1.93	0.49
1:B:8:ASN:HD22	1:B:8:ASN:N	2.09	0.49
1:B:2:MET:HE1	1:B:25:SER:HB3	1.95	0.49
1:B:202:TYR:CE2	1:B:204:SER:HB3	2.48	0.48
1:A:146:TRP:CE2	1:A:152:VAL:HG22	2.48	0.48
1:B:144:ARG:NH2	4:B:320:HOH:O	2.47	0.48
1:A:51:LYS:CE	2:A:220:SO4:O4	2.61	0.48
1:B:146:TRP:CH2	1:B:186:ARG:HG2	2.48	0.48
1:A:86:CYS:HB2	4:A:324:HOH:O	2.14	0.47
1:A:205:THR:HB	1:A:206:PRO:HA	1.96	0.47
1:B:2:MET:CE	1:B:25:SER:HB3	2.45	0.47
1:A:123:LYS:HB3	1:A:124:PRO:HD3	1.96	0.46
1:B:111:ILE:HG12	1:B:208:PHE:CE1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:THR:O	1:A:71:GLN:HB2	2.17	0.45
1:B:113:LEU:HD22	1:B:126:PHE:CG	2.52	0.44
1:B:8:ASN:H	1:B:8:ASN:HD22	1.67	0.42
1:A:171:PRO:HB3	4:A:277:HOH:O	2.18	0.42
1:B:202:TYR:CZ	1:B:204:SER:HB3	2.54	0.41
1:A:91:GLU:HG3	4:A:326:HOH:O	2.20	0.41
1:A:1:PRO:HG2	1:A:28:GLU:OE2	2.21	0.41
1:A:37:ALA:HB1	1:A:38:PRO:HA	2.03	0.41
1:A:37:ALA:HB2	1:A:40:TYR:CZ	2.56	0.41
3:B:218:GPS:CG4	3:B:218:GPS:HB2	2.51	0.41
1:A:162:ILE:HD12	1:A:162:ILE:HA	1.92	0.40
1:A:150:ASP:CB	1:A:151:LYS:HE3	2.52	0.40
1:A:2:MET:HE3	1:A:25:SER:HB3	2.03	0.40
1:B:172:LYS:HD2	1:B:175:ASP:OD2	2.21	0.40
1:A:3:ILE:HG12	1:A:28:GLU:HG3	2.04	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	215/217 (99%)	212 (99%)	3 (1%)	0	100	100
1	B	215/217 (99%)	212 (99%)	3 (1%)	0	100	100
All	All	430/434 (99%)	424 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	197/197 (100%)	190 (96%)	7 (4%)	42	21
1	B	197/197 (100%)	186 (94%)	11 (6%)	26	9
All	All	394/394 (100%)	376 (95%)	18 (5%)	33	14

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	13	VAL
1	A	19	LEU
1	A	111	ILE
1	A	120	GLU
1	A	151	LYS
1	A	174	LEU
1	B	8	ASN
1	B	12	LEU
1	B	19	LEU
1	B	36	ASP
1	B	111	ILE
1	B	124	PRO
1	B	132	GLU
1	B	151	LYS
1	B	174	LEU
1	B	186	ARG
1	B	216	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
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 [i](#)

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

5 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	1.09	0	6,6,6	0.30	0
2	SO4	A	219	-	4,4,4	1.28	0	6,6,6	0.37	0
2	SO4	A	220	-	4,4,4	1.43	0	6,6,6	0.58	0
3	GPS	A	221	-	29,37,37	1.23	3 (10%)	35,51,51	1.85	13 (37%)
3	GPS	B	218	-	29,37,37	1.36	4 (13%)	35,51,51	1.62	7 (20%)

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
3	GPS	A	221	-	2/2/8/10	0/21/43/43	0/3/3/3
3	GPS	B	218	-	2/2/8/10	0/21/43/43	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	218	GPS	CB1-CG1	-2.20	1.45	1.52
3	B	218	GPS	C2-N3	-2.15	1.29	1.33
3	A	221	GPS	C2-N3	-2.03	1.29	1.33
3	B	218	GPS	CZ5-CH5	2.07	1.43	1.39
3	A	221	GPS	CZ5-CH5	2.18	1.43	1.39
3	B	218	GPS	CG4-CB4	2.90	1.43	1.39
3	A	221	GPS	CG4-CB4	2.91	1.43	1.39

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	221	GPS	CA3-N3-C2	-3.80	117.12	122.34
3	B	218	GPS	OE1-CD1-CG1	-3.27	116.34	121.98
3	A	221	GPS	O5-CA5-CB5	-3.02	104.13	110.42
3	B	218	GPS	O5-CA5-CB5	-2.96	104.27	110.42
3	B	218	GPS	CE4-CD4-CG4	-2.73	116.19	120.19
3	A	221	GPS	O2-C2-CA2	-2.69	114.41	120.36
3	A	221	GPS	CG5-CB5-CA5	-2.63	114.66	120.75
3	B	218	GPS	CG5-CB5-CA5	-2.59	114.75	120.75
3	A	221	GPS	CH5-CH4-CB4	-2.54	115.81	119.78
3	A	221	GPS	CE4-CD4-CG4	-2.53	116.48	120.19
3	A	221	GPS	CE5-CZ5-CH5	-2.27	116.38	120.33
3	A	221	GPS	CZ4-CH4-CB4	2.04	120.65	118.41
3	A	221	GPS	O2-C2-N3	2.32	127.74	123.08
3	A	221	GPS	CZ5-CH5-CB5	2.34	120.98	118.41
3	A	221	GPS	CB2-CA2-N2	2.39	116.70	110.83
3	B	218	GPS	CB1-CG1-CD1	2.44	119.08	113.27
3	B	218	GPS	CD4-CE4-CZ4	2.45	123.77	120.19
3	B	218	GPS	CG5-CB5-CH5	2.45	122.73	119.41
3	A	221	GPS	CB1-CG1-CD1	3.22	120.91	113.27
3	A	221	GPS	CD4-CE4-CZ4	3.52	125.33	120.19

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	221	GPS	CA5
3	A	221	GPS	CA4
3	B	218	GPS	CA5
3	B	218	GPS	CA4

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	220	SO4	1	0
3	A	221	GPS	1	0
3	B	218	GPS	1	0

## 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](#)

EDS was not executed - this section will therefore be empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.