



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:30 PM GMT

PDB ID : 1B4P  
Title : CRYSTAL STRUCTURES OF CLASS MU CHIMERIC GST ISOENZYMES  
M1-2 AND M2-1  
Authors : Xiao, G.; Chen, J.; Armstrong, R.N.; Gilliland, G.L.  
Deposited on : 1998-12-26  
Resolution : 1.70 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.

---

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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

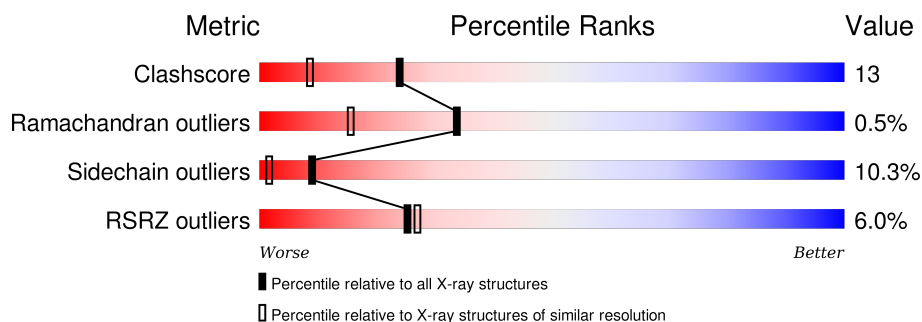
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)
RSRZ outliers	91569	3200 (1.70-1.70)

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.

Mol	Chain	Length	Quality of chain
1	A	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
2	SO4	A	223	-	-	-	X
3	GPS	A	219	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2148 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.

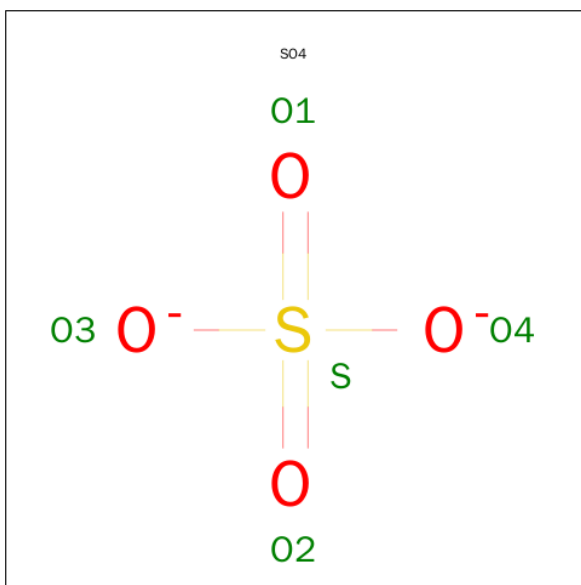
- Molecule 1 is a protein called PROTEIN (GLUTATHIONE S-TRANSFERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1813	1175	304	323	11			

There are 15 discrepancies between the modelled and reference sequences:

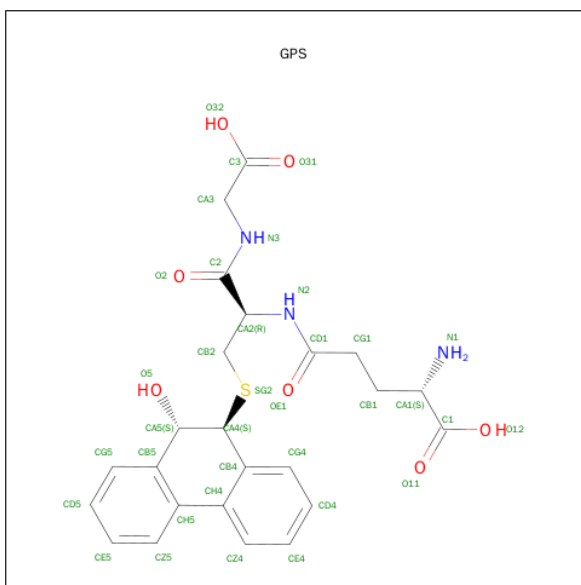
Chain	Residue	Modelled	Actual	Comment	Reference
A	3	ILE	THR	CONFLICT	UNP P08010
A	8	ASN	ASP	CONFLICT	UNP P08010
A	9	VAL	ILE	CONFLICT	UNP P08010
A	13	THR	ALA	CONFLICT	UNP P08010
A	15	PRO	ALA	CONFLICT	UNP P08010
A	19	LEU	PHE	CONFLICT	UNP P08010
A	25	SER	THR	CONFLICT	UNP P08010
A	29	GLU	ASP	CONFLICT	UNP P08010
A	31	ARG	LYS	CONFLICT	UNP P08010
A	33	ALA	SER	CONFLICT	UNP P08010
A	47	ASN	SER	CONFLICT	UNP P08010
A	67	ARG	HIS	CONFLICT	UNP P08010
A	76	MET	LEU	CONFLICT	UNP P08010
A	80	ALA	GLY	CONFLICT	UNP P08010
A	84	HIS	ASN	CONFLICT	UNP P08010

- 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 5	O 4	S 1	0	0
2	A	1	Total 5	O 4	S 1	0	0
2	A	1	Total 5	O 4	S 1	0	0

- 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	A	1	Total	C	N	O	S	0	0
			35	24	3	7	1		

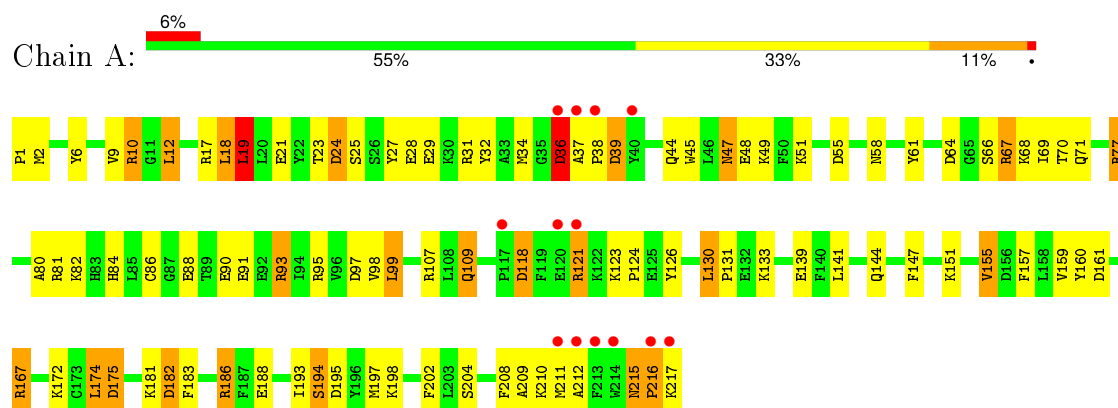
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	250	Total	O	0	0
			250	250		

### 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.

- Molecule 1: PROTEIN (GLUTATHIONE S-TRANSFERASE)



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.59Å 82.55Å 79.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 1.70 45.21 – 1.66	Depositor EDS
% Data completeness (in resolution range)	88.6 (6.00-1.70) 88.7 (45.21-1.66)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.14 (at 1.66Å)	Xtriage
Refinement program	GPRLSA	Depositor
R, $R_{free}$	0.179 , (Not available) 0.173 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	14.8	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 80.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 25872 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2148	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.40% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.375 respectively for untwinned datasets, and 0.333, 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, 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	1.17	0/1863	2.05	63/2509 (2.5%)

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	5

There are no bond length outliers.

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	ARG	CD-NE-CZ	16.63	146.89	123.60
1	A	10	ARG	NE-CZ-NH2	-15.27	112.66	120.30
1	A	77	ARG	NE-CZ-NH1	12.00	126.30	120.30
1	A	77	ARG	NE-CZ-NH2	-11.52	114.54	120.30
1	A	67	ARG	NE-CZ-NH2	-11.00	114.80	120.30
1	A	186	ARG	NE-CZ-NH2	-10.84	114.88	120.30
1	A	186	ARG	NE-CZ-NH1	10.80	125.70	120.30
1	A	29	GLU	CG-CD-OE2	-9.62	99.05	118.30
1	A	139	GLU	OE1-CD-OE2	9.30	134.46	123.30
1	A	6	TYR	CB-CG-CD2	9.04	126.42	121.00
1	A	81	ARG	CD-NE-CZ	9.00	136.20	123.60
1	A	29	GLU	CG-CD-OE1	8.98	136.26	118.30
1	A	107	ARG	NE-CZ-NH2	-8.77	115.92	120.30
1	A	93	ARG	CD-NE-CZ	8.46	135.44	123.60
1	A	61	TYR	CB-CG-CD2	-8.44	115.94	121.00
1	A	95	ARG	NE-CZ-NH1	8.13	124.37	120.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	81	ARG	NE-CZ-NH1	-8.06	116.27	120.30
1	A	121	ARG	NE-CZ-NH2	7.93	124.27	120.30
1	A	18	LEU	CB-CG-CD2	7.53	123.80	111.00
1	A	17	ARG	NE-CZ-NH1	7.14	123.87	120.30
1	A	39	ASP	CB-CG-OD2	-6.97	112.02	118.30
1	A	93	ARG	NE-CZ-NH1	6.88	123.74	120.30
1	A	77	ARG	CD-NE-CZ	6.72	133.01	123.60
1	A	18	LEU	CA-CB-CG	6.70	130.71	115.30
1	A	34	MET	CA-CB-CG	-6.37	102.46	113.30
1	A	19	LEU	CB-CG-CD2	6.33	121.76	111.00
1	A	194	SER	N-CA-CB	6.31	119.97	110.50
1	A	107	ARG	NH1-CZ-NH2	6.27	126.29	119.40
1	A	36	ASP	CB-CG-OD1	6.21	123.89	118.30
1	A	161	ASP	CB-CG-OD1	6.12	123.81	118.30
1	A	130	LEU	CB-CA-C	6.07	121.74	110.20
1	A	98	VAL	O-C-N	6.06	132.39	122.70
1	A	55	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	24	ASP	CB-CG-OD2	-6.00	112.90	118.30
1	A	90	GLU	CG-CD-OE2	-5.91	106.47	118.30
1	A	36	ASP	CB-CA-C	5.90	122.19	110.40
1	A	95	ARG	NE-CZ-NH2	-5.85	117.38	120.30
1	A	155	VAL	O-C-N	5.85	132.05	122.70
1	A	28	GLU	N-CA-CB	5.84	121.12	110.60
1	A	31	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	182	ASP	CB-CG-OD1	5.78	123.50	118.30
1	A	97	ASP	O-C-N	-5.67	113.64	122.70
1	A	31	ARG	NE-CZ-NH1	-5.60	117.50	120.30
1	A	141	LEU	CB-CG-CD2	-5.58	101.51	111.00
1	A	186	ARG	N-CA-CB	5.52	120.54	110.60
1	A	107	ARG	NE-CZ-NH1	-5.52	117.54	120.30
1	A	91	GLU	CG-CD-OE1	5.44	129.18	118.30
1	A	32	TYR	CB-CG-CD1	5.42	124.25	121.00
1	A	32	TYR	CB-CG-CD2	-5.41	117.75	121.00
1	A	175	ASP	CB-CG-OD2	-5.39	113.44	118.30
1	A	39	ASP	CB-CA-C	5.27	120.94	110.40
1	A	195	ASP	CB-CG-OD2	-5.26	113.56	118.30
1	A	19	LEU	CA-C-O	-5.23	109.12	120.10
1	A	157	PHE	CB-CG-CD2	-5.22	117.15	120.80
1	A	58	ASN	O-C-N	5.21	131.04	122.70
1	A	6	TYR	CG-CD2-CE2	5.20	125.46	121.30
1	A	160	TYR	CB-CG-CD1	-5.17	117.90	121.00
1	A	80	ALA	CB-CA-C	5.16	117.83	110.10

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	ARG	NH1-CZ-NH2	5.13	125.05	119.40
1	A	61	TYR	CZ-CE2-CD2	-5.12	115.19	119.80
1	A	183	PHE	CB-CG-CD2	-5.05	117.26	120.80
1	A	1	PRO	CA-N-CD	-5.05	104.43	111.50
1	A	69	ILE	CB-CG1-CD1	5.04	128.01	113.90

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	182	ASP	Mainchain
1	A	186	ARG	Sidechain
1	A	188	GLU	Mainchain
1	A	67	ARG	Sidechain
1	A	77	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	1813	0	1800	49	0
2	A	15	0	0	1	0
3	A	70	0	52	7	0
4	A	250	0	0	4	0
All	All	2148	0	1852	49	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 (49) 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:212:ALA:HB3	1:A:216:PRO:HB3	1.51	0.91
1:A:25:SER:O	3:A:219:GPS:N1	2.07	0.86
1:A:68:LYS:NZ	2:A:222:SO4:O2	2.20	0.74

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:ALA:HB2	3:A:218:GPS:CD5	2.18	0.73
1:A:21:GLU:HA	3:A:219:GPS:O12	1.91	0.70
1:A:208:PHE:O	1:A:216:PRO:HA	1.94	0.68
1:A:37:ALA:HB1	1:A:38:PRO:HA	1.74	0.67
1:A:86:CYS:O	1:A:151:LYS:HE2	1.96	0.66
1:A:21:GLU:OE1	3:A:219:GPS:O11	2.16	0.64
1:A:36:ASP:OD1	1:A:217:LYS:NZ	2.33	0.61
1:A:36:ASP:OD2	1:A:210:LYS:NZ	2.33	0.61
1:A:202:PHE:CE2	1:A:204:SER:HB3	2.36	0.61
1:A:210:LYS:HA	1:A:216:PRO:HB2	1.82	0.60
1:A:9:VAL:HG21	3:A:218:GPS:HE5	1.83	0.59
1:A:118:ASP:OD1	1:A:121:ARG:HD2	2.05	0.57
1:A:12:LEU:C	1:A:12:LEU:HD12	2.27	0.55
1:A:216:PRO:CD	1:A:217:LYS:H	2.27	0.48
1:A:2:MET:HE1	1:A:25:SER:HB3	1.95	0.48
1:A:99:LEU:HD13	1:A:159:VAL:HG21	1.95	0.47
1:A:70:THR:O	1:A:71:GLN:HB2	2.14	0.47
1:A:194:SER:HA	1:A:197:MET:HE2	1.97	0.46
1:A:45:TRP:CZ2	1:A:49:LYS:HG3	2.50	0.46
1:A:37:ALA:CB	1:A:38:PRO:HA	2.40	0.46
1:A:121:ARG:HG2	1:A:121:ARG:HH11	1.79	0.46
1:A:82:LYS:HE2	4:A:347:HOH:O	2.16	0.46
1:A:175:ASP:OD1	1:A:181:LYS:NZ	2.41	0.45
1:A:93:ARG:NH1	4:A:496:HOH:O	2.50	0.44
1:A:86:CYS:HB2	4:A:393:HOH:O	2.17	0.44
1:A:174:LEU:HB3	1:A:181:LYS:HG2	1.99	0.44
1:A:99:LEU:HD11	1:A:147:PHE:CD2	2.52	0.44
1:A:123:LYS:N	1:A:124:PRO:HD2	2.33	0.44
1:A:121:ARG:NH1	1:A:121:ARG:HG2	2.33	0.43
1:A:215:ASN:N	1:A:216:PRO:HD3	2.32	0.43
1:A:10:ARG:HD2	1:A:204:SER:O	2.18	0.43
1:A:198:LYS:HE2	1:A:198:LYS:HB2	1.81	0.43
1:A:88:GLU:OE1	1:A:151:LYS:NZ	2.52	0.42
1:A:2:MET:HG2	1:A:64:ASP:HB2	2.00	0.42
1:A:193:ILE:O	1:A:197:MET:HG3	2.19	0.42
1:A:2:MET:CE	1:A:25:SER:HB3	2.49	0.42
1:A:109:GLN:HE21	1:A:109:GLN:HB3	1.39	0.41
1:A:19:LEU:HD22	1:A:23:THR:HG23	2.01	0.41
1:A:209:ALA:HB2	3:A:218:GPS:CE5	2.50	0.41
1:A:27:TYR:HE1	3:A:219:GPS:HO12	1.69	0.41
1:A:144:GLN:HB2	1:A:144:GLN:HE21	1.50	0.40

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:LEU:N	1:A:131:PRO:CD	2.84	0.40
1:A:181:LYS:HE2	4:A:420:HOH:O	2.21	0.40
1:A:47:ASN:O	1:A:51:LYS:NZ	2.32	0.40
1:A:212:ALA:CB	1:A:216:PRO:HB3	2.36	0.40
1:A:126:TYR:CE1	1:A:133:LYS:NZ	2.89	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%)	205 (95%)	9 (4%)	1 (0%)	34 15

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	216	PRO

### 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	195/195 (100%)	175 (90%)	20 (10%)	9 2

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	18	LEU
1	A	19	LEU
1	A	24	ASP
1	A	36	ASP
1	A	39	ASP
1	A	44	GLN
1	A	47	ASN
1	A	48	GLU
1	A	66	SER
1	A	84	HIS
1	A	99	LEU
1	A	109	GLN
1	A	118	ASP
1	A	155	VAL
1	A	167	ARG
1	A	172	LYS
1	A	174	LEU
1	A	211	MET
1	A	215	ASN

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

Mol	Chain	Res	Type
1	A	44	GLN
1	A	47	ASN
1	A	109	GLN
1	A	144	GLN
1	A	215	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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$
3	GPS	A	218	-	29,37,37	1.64	9 (31%)	35,51,51	1.87	8 (22%)
3	GPS	A	219	-	29,37,37	1.87	8 (27%)	35,51,51	2.50	13 (37%)
2	SO4	A	221	-	4,4,4	1.02	0	6,6,6	0.23	0
2	SO4	A	222	-	4,4,4	1.23	0	6,6,6	0.33	0
2	SO4	A	223	-	4,4,4	1.08	0	6,6,6	0.17	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	GPS	A	218	-	-	0/21/43/43	0/3/3/3
3	GPS	A	219	-	-	0/21/43/43	0/3/3/3
2	SO4	A	221	-	-	0/0/0/0	0/0/0/0
2	SO4	A	222	-	-	0/0/0/0	0/0/0/0
2	SO4	A	223	-	-	0/0/0/0	0/0/0/0

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	218	GPS	CH5-CH4	-2.80	1.42	1.47
3	A	219	GPS	CH5-CH4	-2.67	1.42	1.47
3	A	219	GPS	CH5-CB5	-2.55	1.36	1.40
3	A	218	GPS	CG1-CD1	-2.35	1.46	1.51
3	A	218	GPS	CH5-CB5	-2.11	1.37	1.40
3	A	218	GPS	CE4-CD4	2.10	1.43	1.38
3	A	218	GPS	CB2-CA2	2.11	1.59	1.53
3	A	218	GPS	O5-CA5	2.18	1.47	1.42
3	A	218	GPS	CA3-N3	2.26	1.50	1.46
3	A	219	GPS	CA1-N1	2.29	1.58	1.48

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	218	GPS	CD1-N2	2.31	1.38	1.34
3	A	219	GPS	CG5-CB5	2.75	1.43	1.39
3	A	219	GPS	OE1-CD1	2.79	1.29	1.23
3	A	218	GPS	CG5-CB5	2.88	1.43	1.39
3	A	219	GPS	CZ5-CH5	3.08	1.45	1.39
3	A	219	GPS	CG4-CB4	3.79	1.44	1.39
3	A	219	GPS	O5-CA5	4.76	1.52	1.42

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	219	GPS	O2-C2-CA2	-7.32	104.17	120.36
3	A	219	GPS	CA2-N2-CD1	-4.77	109.41	121.58
3	A	218	GPS	CA3-N3-C2	-4.36	116.34	122.34
3	A	218	GPS	CB4-CA4-CA5	-4.25	100.19	109.89
3	A	218	GPS	O2-C2-CA2	-3.88	111.78	120.36
3	A	219	GPS	CB2-CA2-N2	-3.55	102.09	110.83
3	A	219	GPS	CA2-CB2-SG2	-3.31	105.11	112.97
3	A	219	GPS	CB1-CG1-CD1	-3.18	105.70	113.27
3	A	218	GPS	CH5-CH4-CB4	-3.11	114.92	119.78
3	A	218	GPS	CD5-CG5-CB5	-3.04	117.06	121.02
3	A	219	GPS	OE1-CD1-N2	-3.03	117.87	123.01
3	A	219	GPS	O5-CA5-CB5	-2.80	104.59	110.42
3	A	219	GPS	CB1-CA1-N1	-2.47	103.50	110.52
3	A	218	GPS	C2-CA2-N2	-2.35	104.63	111.26
3	A	219	GPS	CD4-CG4-CB4	-2.32	118.00	121.02
3	A	219	GPS	CB4-CA4-CA5	-2.27	104.71	109.89
3	A	218	GPS	CG5-CB5-CA5	-2.08	115.94	120.75
3	A	219	GPS	CH5-CH4-CB4	-2.03	116.61	119.78
3	A	219	GPS	CG1-CD1-N2	2.86	120.50	115.83
3	A	218	GPS	O2-C2-N3	2.93	128.96	123.08
3	A	219	GPS	O2-C2-N3	6.61	136.35	123.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	218	GPS	3	0
3	A	219	GPS	4	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	222	SO4	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](#)

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	217/217 (100%)	0.01	13 (5%) 25 27	7, 18, 36, 41	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	216	PRO	9.3
1	A	217	LYS	5.2
1	A	212	ALA	4.6
1	A	37	ALA	4.3
1	A	214	TRP	4.0
1	A	213	PHE	3.9
1	A	36	ASP	3.4
1	A	117	PRO	3.2
1	A	211	MET	3.2
1	A	121	ARG	2.5
1	A	40	TYR	2.3
1	A	38	PRO	2.1
1	A	120	GLU	2.1

### 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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	A	223	5/5	0.80	0.37	7.71	42,43,43,43	5
3	GPS	A	219	35/35	0.74	0.20	3.09	29,35,42,43	0
3	GPS	A	218	35/35	0.96	0.09	0.30	9,18,28,29	0
2	SO4	A	222	5/5	0.68	0.34	-	38,38,39,39	4
2	SO4	A	221	5/5	0.87	0.46	-	41,41,41,41	5

## 6.5 Other polymers [i](#)

There are no such residues in this entry.