



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:58 PM GMT

PDB ID : 6GSY
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 : 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
<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) : **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

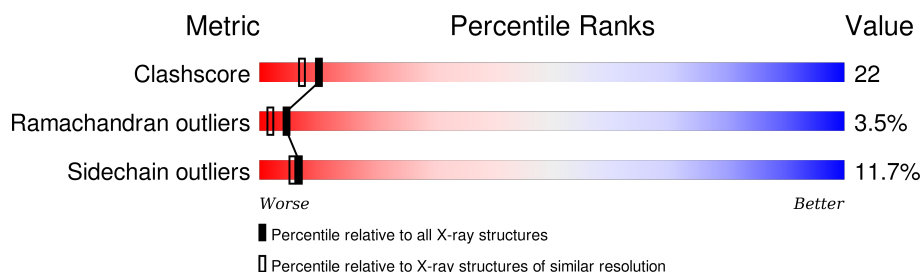
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.


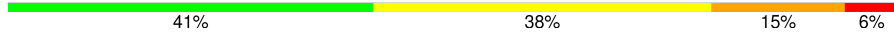
Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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 3 unique types of molecules in this entry. The entry contains 4074 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.

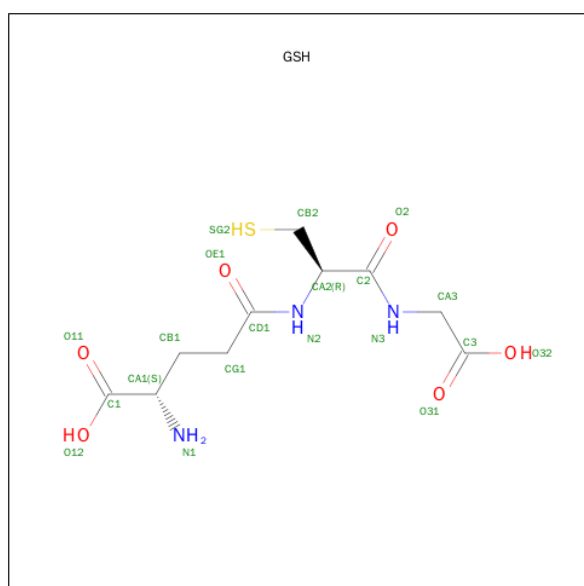
- Molecule 1 is a protein called MU CLASS GLUTATHIONE S-TRANSFERASE OF ISOENZYME 3-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1816	1177	303	325	11			
1	B	217	Total	C	N	O	S	0	0	0
			1816	1177	303	325	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	6	PHE	TYR	ENGINEERED	UNP P04905
B	6	PHE	TYR	ENGINEERED	UNP P04905

- Molecule 2 is GLUTATHIONE (three-letter code: GSH) (formula: $C_{10}H_{17}N_3O_6S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

- Molecule 3 is water.

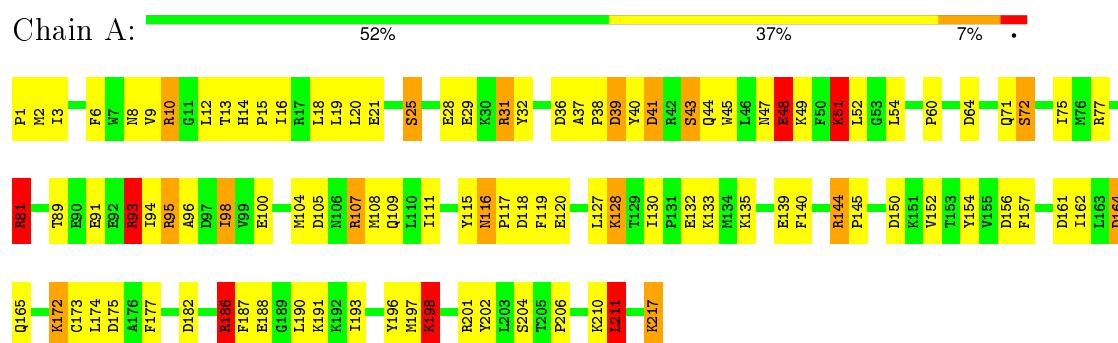
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	209	Total	O	0	0
			209	209		
3	B	193	Total	O	0	0
			193	193		

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, α , β , γ	87.84Å 69.70Å 81.57Å 90.00° 105.38° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20	Depositor
% Data completeness (in resolution range)	88.3 (6.00-2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	GPRLSA	Depositor
R, R_{free}	0.168 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4074	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: GSH

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	1.08	0/1865	2.06	57/2513 (2.3%)
1	B	1.06	1/1865 (0.1%)	2.28	70/2513 (2.8%)
All	All	1.07	1/3730 (0.0%)	2.18	127/5026 (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
1	B	0	5
All	All	0	10

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	216	ASN	C-O	5.98	1.34	1.23

All (127) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	186	ARG	NE-CZ-NH1	28.20	134.40	120.30
1	B	107	ARG	NE-CZ-NH2	-23.90	108.35	120.30
1	B	81	ARG	NE-CZ-NH1	19.02	129.81	120.30
1	A	186	ARG	NE-CZ-NH1	17.81	129.20	120.30
1	B	95	ARG	CD-NE-CZ	16.98	147.38	123.60
1	B	186	ARG	CD-NE-CZ	16.27	146.38	123.60
1	A	201	ARG	CD-NE-CZ	16.07	146.10	123.60
1	B	186	ARG	NE-CZ-NH2	-15.44	112.58	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	186	ARG	CD-NE-CZ	15.34	145.08	123.60
1	A	77	ARG	NE-CZ-NH1	15.00	127.80	120.30
1	A	31	ARG	NE-CZ-NH1	14.52	127.56	120.30
1	B	77	ARG	NE-CZ-NH2	-14.10	113.25	120.30
1	A	107	ARG	NE-CZ-NH1	13.63	127.12	120.30
1	A	95	ARG	NE-CZ-NH1	12.87	126.74	120.30
1	B	67	ARG	NE-CZ-NH1	12.15	126.37	120.30
1	A	182	ASP	CB-CG-OD1	11.88	128.99	118.30
1	B	67	ARG	CD-NE-CZ	11.21	139.29	123.60
1	B	188	GLU	CA-CB-CG	11.08	137.78	113.40
1	B	191	LYS	CA-CB-CG	10.92	137.42	113.40
1	B	116	ASN	CA-CB-CG	10.41	136.31	113.40
1	B	107	ARG	NH1-CZ-NH2	10.24	130.67	119.40
1	A	188	GLU	OE1-CD-OE2	-9.64	111.73	123.30
1	B	61	TYR	CB-CG-CD1	9.62	126.77	121.00
1	B	40	TYR	CB-CG-CD1	9.58	126.75	121.00
1	B	144	ARG	NE-CZ-NH2	-9.34	115.63	120.30
1	A	77	ARG	NE-CZ-NH2	-9.26	115.67	120.30
1	B	186	ARG	NH1-CZ-NH2	-9.26	109.22	119.40
1	A	93	ARG	NE-CZ-NH2	-9.06	115.77	120.30
1	A	186	ARG	NH1-CZ-NH2	-9.05	109.44	119.40
1	B	17	ARG	NE-CZ-NH1	9.05	124.83	120.30
1	A	31	ARG	NE-CZ-NH2	-8.89	115.85	120.30
1	B	201	ARG	CD-NE-CZ	8.88	136.04	123.60
1	B	88	GLU	CA-CB-CG	8.75	132.65	113.40
1	A	96	ALA	CB-CA-C	8.61	123.01	110.10
1	A	10	ARG	NE-CZ-NH1	8.45	124.53	120.30
1	B	116	ASN	CB-CA-C	8.43	127.25	110.40
1	B	77	ARG	NE-CZ-NH1	8.37	124.48	120.30
1	A	93	ARG	NE-CZ-NH1	8.15	124.37	120.30
1	A	39	ASP	CB-CG-OD1	-8.13	110.99	118.30
1	B	10	ARG	NE-CZ-NH2	7.98	124.29	120.30
1	B	81	ARG	NH1-CZ-NH2	-7.93	110.68	119.40
1	A	48	GLU	CA-CB-CG	7.89	130.76	113.40
1	B	61	TYR	CB-CG-CD2	-7.87	116.28	121.00
1	B	114	CYS	CB-CA-C	7.86	126.11	110.40
1	B	42	ARG	NE-CZ-NH1	7.81	124.20	120.30
1	B	202	TYR	CB-CG-CD2	-7.67	116.40	121.00
1	A	175	ASP	CB-CG-OD1	7.50	125.05	118.30
1	B	144	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	A	91	GLU	CG-CD-OE2	-7.32	103.66	118.30
1	B	115	TYR	CA-CB-CG	7.31	127.29	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	ARG	NE-CZ-NH2	-7.24	116.68	120.30
1	A	186	ARG	CA-CB-CG	7.08	128.99	113.40
1	B	170	GLU	OE1-CD-OE2	7.08	131.79	123.30
1	A	95	ARG	CD-NE-CZ	7.03	133.44	123.60
1	B	40	TYR	CB-CG-CD2	-7.01	116.79	121.00
1	A	81	ARG	NE-CZ-NH1	-6.97	116.81	120.30
1	A	64	ASP	CB-CG-OD1	6.92	124.53	118.30
1	B	24	ASP	CB-CG-OD1	6.91	124.52	118.30
1	A	156	ASP	CB-CG-OD1	-6.85	112.14	118.30
1	B	101	ASN	O-C-N	-6.77	111.87	122.70
1	B	42	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	A	77	ARG	CD-NE-CZ	6.64	132.89	123.60
1	A	64	ASP	CB-CG-OD2	-6.62	112.34	118.30
1	B	121	LYS	N-CA-CB	6.59	122.46	110.60
1	A	51	LYS	CA-CB-CG	6.51	127.73	113.40
1	B	77	ARG	CD-NE-CZ	6.43	132.60	123.60
1	B	70	THR	N-CA-CB	6.33	122.32	110.30
1	B	105	ASP	CB-CG-OD2	6.21	123.89	118.30
1	A	202	TYR	CB-CG-CD2	-6.18	117.29	121.00
1	B	93	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	B	157	PHE	CB-CG-CD2	-6.13	116.51	120.80
1	A	100	GLU	CG-CD-OE2	-6.13	106.04	118.30
1	B	106	ASN	N-CA-CB	6.13	121.63	110.60
1	B	91	GLU	OE1-CD-OE2	6.11	130.63	123.30
1	A	164	ASP	CB-CG-OD1	6.04	123.74	118.30
1	B	40	TYR	CA-CB-CG	6.03	124.85	113.40
1	B	202	TYR	CB-CG-CD1	6.01	124.60	121.00
1	A	202	TYR	CB-CG-CD1	6.00	124.60	121.00
1	A	107	ARG	NH1-CZ-NH2	-6.00	112.80	119.40
1	A	91	GLU	CG-CD-OE1	5.99	130.28	118.30
1	B	190	LEU	CA-CB-CG	5.97	129.03	115.30
1	A	202	TYR	CB-CA-C	5.90	122.19	110.40
1	B	36	ASP	CA-CB-CG	5.87	126.32	113.40
1	B	105	ASP	CA-CB-CG	5.87	126.31	113.40
1	A	47	ASN	N-CA-CB	5.86	121.15	110.60
1	B	67	ARG	NH1-CZ-NH2	-5.83	112.99	119.40
1	B	39	ASP	CB-CG-OD2	-5.79	113.08	118.30
1	A	154	TYR	CB-CG-CD2	-5.78	117.53	121.00
1	A	18	LEU	CA-CB-CG	5.73	128.48	115.30
1	B	164	ASP	CB-CG-OD1	5.72	123.45	118.30
1	B	89	THR	CA-CB-OG1	-5.70	97.04	109.00
1	A	182	ASP	CA-CB-CG	5.67	125.88	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	91	GLU	CG-CD-OE2	-5.64	107.02	118.30
1	A	211	LEU	N-CA-CB	-5.63	99.13	110.40
1	B	105	ASP	CB-CG-OD1	5.63	123.37	118.30
1	B	150	ASP	CB-CG-OD1	5.63	123.37	118.30
1	B	81	ARG	CD-NE-CZ	5.60	131.44	123.60
1	A	47	ASN	CA-C-O	-5.60	108.34	120.10
1	A	13	THR	CA-CB-OG1	-5.60	97.25	109.00
1	B	105	ASP	OD1-CG-OD2	-5.56	112.74	123.30
1	A	43	SER	CB-CA-C	-5.54	99.58	110.10
1	B	29	GLU	CG-CD-OE1	5.49	129.27	118.30
1	A	81	ARG	NE-CZ-NH2	-5.39	117.61	120.30
1	A	81	ARG	NH1-CZ-NH2	5.36	125.30	119.40
1	A	217	LYS	CA-CB-CG	5.36	125.19	113.40
1	B	29	GLU	CB-CA-C	-5.33	99.74	110.40
1	B	100	GLU	CG-CD-OE2	-5.25	107.81	118.30
1	B	120	GLU	CA-CB-CG	5.24	124.94	113.40
1	A	89	THR	CA-CB-OG1	-5.20	98.09	109.00
1	A	28	GLU	CG-CD-OE1	-5.19	107.92	118.30
1	B	173	CYS	CA-CB-SG	5.19	123.35	114.00
1	B	31	ARG	NE-CZ-NH1	-5.17	117.72	120.30
1	B	39	ASP	CB-CG-OD1	5.16	122.94	118.30
1	B	72	SER	N-CA-CB	5.15	118.22	110.50
1	A	182	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	B	97	ASP	CB-CG-OD1	5.13	122.92	118.30
1	A	177	PHE	CB-CG-CD1	5.12	124.38	120.80
1	A	198	LYS	N-CA-CB	5.12	119.81	110.60
1	A	157	PHE	CB-CG-CD1	5.11	124.38	120.80
1	A	152	VAL	CG1-CB-CG2	-5.09	102.75	110.90
1	A	72	SER	CA-CB-OG	5.09	124.93	111.20
1	B	201	ARG	NE-CZ-NH2	-5.09	117.76	120.30
1	A	39	ASP	OD1-CG-OD2	5.05	132.90	123.30
1	B	61	TYR	CA-CB-CG	5.03	122.95	113.40
1	B	69	ILE	CA-CB-CG1	5.02	120.53	111.00
1	B	202	TYR	CB-CA-C	5.01	120.41	110.40
1	A	40	TYR	N-CA-CB	-5.00	101.59	110.60

There are no chirality outliers.

All (10) planarity outliers are listed below:

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

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Mol	Chain	Res	Type	Group
1	A	81	ARG	Sidechain
1	A	93	ARG	Sidechain
1	A	98	ILE	Mainchain
1	B	107	ARG	Sidechain
1	B	144	ARG	Sidechain
1	B	154	TYR	Mainchain
1	B	186	ARG	Sidechain
1	B	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	1816	0	1805	53	3
1	B	1816	0	1805	108	4
2	A	20	0	15	1	0
2	B	20	0	14	2	0
3	A	209	0	0	4	0
3	B	193	0	0	11	0
All	All	4074	0	3639	159	4

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 22.

All (159) 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:10:ARG:HD3	1:B:207:ILE:HD13	1.29	1.09
1:B:170:GLU:HB3	1:B:173:CYS:HB3	1.31	1.06
1:A:48:GLU:HG3	1:A:52:LEU:HD21	1.45	0.96
1:A:135:LYS:NZ	1:A:139:GLU:OE2	2.03	0.92
1:B:95:ARG:NH2	1:B:144:ARG:HE	1.70	0.88
1:A:140:PHE:O	1:A:144:ARG:NH1	2.10	0.83
1:B:120:GLU:HA	1:B:123:LYS:HG3	1.61	0.80
1:B:144:ARG:HD3	1:B:149:GLY:HA2	1.65	0.77
1:B:110:LEU:HD13	1:B:166:TYR:HE1	1.49	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ARG:CD	1:B:149:GLY:HA2	2.18	0.73
1:B:135:LYS:HD2	1:B:139:GLU:OE2	1.90	0.72
1:B:40:TYR:HE2	1:B:211:LEU:H	1.38	0.72
1:A:105:ASP:O	1:A:109:GLN:HG3	1.90	0.71
1:B:124:PRO:O	1:B:127:LEU:HB2	1.93	0.69
1:B:40:TYR:CE2	1:B:211:LEU:HB2	2.27	0.69
1:B:10:ARG:HB3	1:B:207:ILE:HA	1.75	0.69
1:B:123:LYS:HB3	1:B:124:PRO:CD	2.24	0.67
1:B:123:LYS:HB3	1:B:124:PRO:HD2	1.77	0.67
1:A:16:ILE:HD13	1:A:75:ILE:HG21	1.78	0.66
1:B:113:LEU:HD22	1:B:126:PHE:CD2	2.31	0.65
1:B:191:LYS:HE2	1:B:191:LYS:O	1.96	0.65
1:A:95:ARG:NH2	1:A:144:ARG:NH1	2.45	0.65
1:A:41:ASP:OD1	1:A:43:SER:OG	2.11	0.65
1:A:193:ILE:O	1:A:197:MET:HG3	1.97	0.65
1:A:95:ARG:NH2	1:A:144:ARG:HH11	1.96	0.64
1:B:123:LYS:O	1:B:127:LEU:HD23	1.97	0.63
1:A:95:ARG:HH22	1:A:144:ARG:HH11	1.47	0.62
1:A:32:TYR:HD2	1:A:44:GLN:HG2	1.64	0.62
1:B:86:CYS:HB2	3:B:300:HOH:O	1.98	0.62
1:B:12:LEU:H	1:B:12:LEU:HD12	1.65	0.62
1:B:105:ASP:O	1:B:109:GLN:HG3	1.99	0.62
1:B:95:ARG:HH21	1:B:144:ARG:HE	1.46	0.61
1:B:124:PRO:O	1:B:127:LEU:N	2.27	0.61
1:B:172:LYS:O	1:B:174:LEU:N	2.33	0.61
1:B:40:TYR:CE2	1:B:210:LYS:HB2	2.36	0.61
1:A:95:ARG:HH22	1:A:144:ARG:NH1	1.99	0.61
1:B:110:LEU:HD13	1:B:166:TYR:CE1	2.33	0.60
1:A:21:GLU:HG3	1:A:196:TYR:CG	2.36	0.60
1:B:113:LEU:HD22	1:B:126:PHE:HD2	1.65	0.59
1:B:197:MET:HG2	1:B:202:TYR:CE1	2.37	0.59
1:B:40:TYR:HE2	1:B:210:LYS:HB2	1.67	0.59
1:B:2:MET:HE1	3:B:370:HOH:O	2.02	0.58
1:B:170:GLU:CB	1:B:173:CYS:HB3	2.22	0.58
1:B:42:ARG:HD3	3:B:338:HOH:O	2.02	0.58
1:B:120:GLU:O	1:B:123:LYS:HB2	2.03	0.58
1:A:115:TYR:HE1	1:A:211:LEU:HB3	1.67	0.57
1:B:130:ILE:HD12	1:B:170:GLU:HG3	1.86	0.57
1:B:161:ASP:O	1:B:165:GLN:HG3	2.05	0.57
1:B:111:ILE:HG12	1:B:208:PHE:CZ	2.40	0.56
1:B:123:LYS:CB	1:B:124:PRO:HD2	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:VAL:HG12	1:A:206:PRO:HG2	1.87	0.56
1:B:151:LYS:HG2	1:B:152:VAL:O	2.05	0.56
1:B:8:ASN:H	1:B:8:ASN:HD22	1.53	0.55
1:A:172:LYS:HD3	1:A:172:LYS:N	2.21	0.55
1:A:145:PRO:HA	1:A:186:ARG:NH1	2.21	0.55
1:B:125:GLU:HA	1:B:128:LYS:HE2	1.88	0.55
1:B:142:GLY:O	1:B:179:ASN:ND2	2.40	0.55
1:A:21:GLU:HG3	1:A:196:TYR:CD1	2.42	0.55
1:B:8:ASN:N	1:B:8:ASN:HD22	2.05	0.54
1:B:100:GLU:OE1	1:B:154:TYR:OH	2.20	0.54
1:A:36:ASP:OD1	1:A:210:LYS:HD3	2.07	0.54
1:A:10:ARG:NH2	1:A:164:ASP:OD2	2.39	0.53
1:A:173:CYS:SG	1:A:174:LEU:CD1	2.97	0.53
1:B:8:ASN:ND2	3:B:305:HOH:O	2.42	0.53
1:B:214:TRP:O	1:B:216:ASN:N	2.42	0.53
1:B:107:ARG:O	1:B:111:ILE:HG13	2.09	0.53
1:B:130:ILE:O	1:B:133:LYS:N	2.41	0.52
1:A:132:GLU:HB2	3:A:365:HOH:O	2.09	0.52
1:B:92:GLU:HG2	1:B:148:ALA:O	2.09	0.52
1:B:40:TYR:HE2	1:B:211:LEU:N	2.07	0.52
1:B:39:ASP:O	1:B:40:TYR:C	2.48	0.52
1:A:10:ARG:HG3	1:A:14:HIS:HB2	1.91	0.52
1:B:12:LEU:HD21	2:B:218:GSH:HB12	1.91	0.52
1:A:6:PHE:O	1:A:31:ARG:HA	2.10	0.52
1:A:71:GLN:OE1	2:A:218:GSH:N1	2.43	0.51
1:B:191:LYS:C	1:B:191:LYS:HE2	2.31	0.51
1:B:111:ILE:HG12	1:B:208:PHE:CE2	2.46	0.51
1:A:12:LEU:HD13	1:A:72:SER:OG	2.12	0.50
1:A:20:LEU:O	1:A:25:SER:HB2	2.11	0.50
1:B:127:LEU:C	1:B:129:THR:H	2.13	0.50
1:B:120:GLU:HA	1:B:123:LYS:CG	2.37	0.50
1:B:106:ASN:ND2	1:B:137:TYR:OH	2.45	0.50
1:B:47:ASN:ND2	3:B:270:HOH:O	2.44	0.50
1:A:1:PRO:O	1:A:3:ILE:HD12	2.12	0.50
1:A:52:LEU:HB2	1:A:54:LEU:HG	1.93	0.50
1:A:37:ALA:HB1	1:A:38:PRO:HA	1.95	0.49
1:A:81:ARG:HE	1:B:90:GLU:CD	2.15	0.49
1:B:210:LYS:O	1:B:216:ASN:HB2	2.13	0.49
1:B:34:MET:HE2	1:B:40:TYR:HD2	1.77	0.49
1:B:125:GLU:N	1:B:128:LYS:HE2	2.27	0.48
1:B:59:LEU:H	2:B:218:GSH:HA31	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:LYS:HD3	3:A:229:HOH:O	2.12	0.48
1:B:207:ILE:HB	1:B:215:SER:HB3	1.96	0.48
1:A:104:MET:O	1:A:108:MET:HG2	2.14	0.48
1:A:107:ARG:O	1:A:111:ILE:HG13	2.14	0.48
1:B:94:ILE:O	1:B:98:ILE:HG13	2.14	0.47
1:A:94:ILE:O	1:A:98:ILE:HG13	2.15	0.47
1:A:45:TRP:CZ2	1:A:49:LYS:HG3	2.49	0.47
1:B:91:GLU:O	1:B:95:ARG:HG3	2.14	0.47
1:B:40:TYR:CE2	1:B:211:LEU:N	2.82	0.47
1:A:29:GLU:CD	1:A:31:ARG:HH21	2.18	0.47
1:B:116:ASN:OD1	1:B:117:PRO:HD2	2.16	0.46
1:B:86:CYS:HB2	3:B:322:HOH:O	2.16	0.46
1:A:161:ASP:O	1:A:165:GLN:HG3	2.16	0.46
1:B:12:LEU:HB3	1:B:107:ARG:HD2	1.98	0.45
1:A:10:ARG:CG	1:A:14:HIS:HB2	2.46	0.45
1:B:208:PHE:O	1:B:209:SER:O	2.34	0.45
1:B:40:TYR:OH	1:B:211:LEU:HA	2.17	0.45
1:B:209:SER:O	1:B:216:ASN:HB2	2.16	0.45
1:B:127:LEU:C	1:B:129:THR:N	2.69	0.45
1:A:111:ILE:HG23	1:A:115:TYR:CD2	2.52	0.45
1:B:194:SER:HA	3:B:401:HOH:O	2.17	0.45
1:B:130:ILE:O	1:B:131:PRO:C	2.55	0.45
1:B:123:LYS:O	1:B:124:PRO:C	2.54	0.45
1:A:2:MET:C	1:A:3:ILE:HD12	2.36	0.45
1:B:172:LYS:O	1:B:175:ASP:N	2.27	0.45
1:B:7:TRP:HB3	3:B:305:HOH:O	2.18	0.44
1:A:109:GLN:OE1	1:A:133:LYS:HE2	2.17	0.44
1:B:134:MET:HE1	1:B:166:TYR:HD2	1.83	0.44
1:A:145:PRO:O	1:A:186:ARG:NH1	2.39	0.44
1:A:128:LYS:NZ	3:A:419:HOH:O	2.41	0.44
1:B:181:LYS:HB3	3:B:393:HOH:O	2.17	0.44
1:B:130:ILE:HG13	1:B:130:ILE:H	1.61	0.44
1:B:176:ALA:O	1:B:178:PRO:HD3	2.17	0.44
1:A:173:CYS:SG	1:A:174:LEU:HD13	2.58	0.44
1:B:123:LYS:NZ	1:B:169:PHE:HZ	2.16	0.43
1:B:8:ASN:H	1:B:8:ASN:ND2	2.15	0.43
1:A:14:HIS:N	1:A:15:PRO:CD	2.81	0.43
1:B:116:ASN:CG	1:B:117:PRO:HD3	2.39	0.43
1:A:116:ASN:HA	1:A:117:PRO:HD2	1.76	0.43
1:B:172:LYS:C	1:B:174:LEU:N	2.70	0.43
1:B:178:PRO:HD2	3:B:266:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:GLU:HA	1:B:123:LYS:HB2	2.00	0.43
1:A:162:ILE:HD12	1:A:162:ILE:HA	1.76	0.43
1:B:204:SER:OG	1:B:205:THR:HG23	2.19	0.43
1:B:120:GLU:C	1:B:122:GLN:H	2.22	0.43
1:B:134:MET:HE3	1:B:166:TYR:HB3	2.01	0.42
1:B:104:MET:O	1:B:108:MET:HG2	2.19	0.42
1:B:212:ALA:HB3	1:B:216:ASN:HB3	2.02	0.42
1:A:187:PHE:O	1:A:190:LEU:HD12	2.20	0.42
1:B:162:ILE:HD12	1:B:162:ILE:HA	1.88	0.42
1:B:172:LYS:C	1:B:174:LEU:H	2.23	0.42
1:A:12:LEU:HD11	1:A:60:PRO:CD	2.49	0.42
1:B:123:LYS:NZ	1:B:169:PHE:CZ	2.86	0.42
1:A:71:GLN:NE2	1:B:105:ASP:OD2	2.34	0.42
1:B:184:LEU:O	1:B:188:GLU:HG2	2.19	0.42
1:B:208:PHE:O	1:B:216:ASN:HA	2.19	0.41
1:B:1:PRO:HG3	3:B:350:HOH:O	2.21	0.41
1:B:4:LEU:HB3	1:B:29:GLU:HG2	2.02	0.41
1:A:127:LEU:HA	1:A:130:ILE:HD12	2.03	0.41
1:B:111:ILE:O	1:B:112:MET:C	2.59	0.41
1:B:125:GLU:HA	1:B:128:LYS:HG3	2.02	0.41
1:B:48:GLU:O	1:B:52:LEU:HG	2.20	0.41
1:B:195:ALA:HA	1:B:198:LYS:HE3	2.03	0.41
1:B:119:PHE:CG	1:B:213:GLN:HG3	2.56	0.41
1:B:115:TYR:O	1:B:116:ASN:HB2	2.21	0.41
1:B:214:TRP:O	1:B:215:SER:C	2.59	0.40
1:B:210:LYS:HA	1:B:216:ASN:O	2.21	0.40
1:A:198:LYS:HB2	3:A:395:HOH:O	2.22	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:LYS:CD	1:B:118:ASP:OD1[3_455]	1.30	0.90
1:B:67:ARG:NH2	1:B:67:ARG:NH2[2_555]	1.83	0.37
1:A:191:LYS:CE	1:B:118:ASP:OD1[3_455]	2.01	0.19
1:A:191:LYS:NZ	1:B:118:ASP:CG[3_455]	2.18	0.02

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%)	203 (94%)	11 (5%)	1 (0%)	34	35
1	B	215/217 (99%)	183 (85%)	18 (8%)	14 (6%)	1	0
All	All	430/434 (99%)	386 (90%)	29 (7%)	15 (4%)	4	2

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	39	ASP
1	B	122	GLN
1	B	123	LYS
1	B	173	CYS
1	B	209	SER
1	B	212	ALA
1	B	215	SER
1	B	38	PRO
1	B	40	TYR
1	B	117	PRO
1	B	124	PRO
1	B	125	GLU
1	A	119	PHE
1	B	71	GLN
1	B	116	ASN

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%)	179 (91%)	18 (9%)	12	11
1	B	197/197 (100%)	169 (86%)	28 (14%)	4	3
All	All	394/394 (100%)	348 (88%)	46 (12%)	7	5

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	19	LEU
1	A	25	SER
1	A	39	ASP
1	A	41	ASP
1	A	48	GLU
1	A	51	LYS
1	A	93	ARG
1	A	116	ASN
1	A	118	ASP
1	A	120	GLU
1	A	128	LYS
1	A	150	ASP
1	A	172	LYS
1	A	198	LYS
1	A	204	SER
1	A	211	LEU
1	A	217	LYS
1	B	8	ASN
1	B	12	LEU
1	B	19	LEU
1	B	26	SER
1	B	36	ASP
1	B	40	TYR
1	B	51	LYS
1	B	69	ILE
1	B	108	MET
1	B	115	TYR
1	B	119	PHE
1	B	122	GLN
1	B	123	LYS
1	B	128	LYS
1	B	129	THR
1	B	132	GLU
1	B	151	LYS

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Mol	Chain	Res	Type
1	B	169	PHE
1	B	173	CYS
1	B	175	ASP
1	B	178	PRO
1	B	186	ARG
1	B	191	LYS
1	B	192	LYS
1	B	203	LEU
1	B	209	SER
1	B	211	LEU
1	B	215	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	ASN
1	B	8	ASN
1	B	47	ASN
1	B	106	ASN
1	B	109	GLN
1	B	116	ASN
1	B	122	GLN
1	B	167	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

2 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	GSH	A	218	-	13,19,19	1.41	3 (23%)	15,24,24	2.74	6 (40%)
2	GSH	B	218	-	13,19,19	1.77	3 (23%)	15,24,24	3.33	8 (53%)

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	GSH	A	218	-	-	0/18/24/24	0/0/0/0
2	GSH	B	218	-	-	0/18/24/24	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	218	GSH	C2-N3	-3.76	1.25	1.33
2	B	218	GSH	CG1-CD1	-2.44	1.46	1.51
2	A	218	GSH	C2-N3	-2.44	1.28	1.33
2	A	218	GSH	CG1-CD1	-2.34	1.46	1.51
2	B	218	GSH	CA2-C2	2.25	1.59	1.52
2	A	218	GSH	CA3-N3	2.26	1.50	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	218	GSH	CA2-CB2-SG2	-6.38	106.32	114.16
2	B	218	GSH	O2-C2-CA2	-6.06	106.96	120.36
2	A	218	GSH	OE1-CD1-CG1	-4.24	114.66	121.98
2	B	218	GSH	C2-CA2-N2	-2.55	104.07	111.26
2	B	218	GSH	CA2-CB2-SG2	-2.53	111.05	114.16
2	B	218	GSH	OE1-CD1-CG1	-2.34	117.95	121.98
2	A	218	GSH	CB1-CA1-N1	2.32	117.10	110.52
2	B	218	GSH	CB1-CG1-CD1	2.38	118.92	113.27
2	A	218	GSH	CG1-CD1-N2	3.14	120.95	115.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	218	GSH	CB1-CG1-CD1	3.56	121.73	113.27
2	B	218	GSH	CB2-CA2-N2	3.59	116.45	111.40
2	A	218	GSH	CB2-CA2-N2	3.71	116.60	111.40
2	B	218	GSH	CA3-N3-C2	6.05	130.65	122.34
2	B	218	GSH	O2-C2-N3	7.05	137.22	123.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	218	GSH	1	0
2	B	218	GSH	2	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.