Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:22 AM GMT

PDB ID	:	2GST
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
Authors	:	Ji, X.; Armstrong, R.N.; Gilliland, G.L.
Deposited on	:	1993-06-07
Resolution	:	1.80 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 (i) symbol.

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

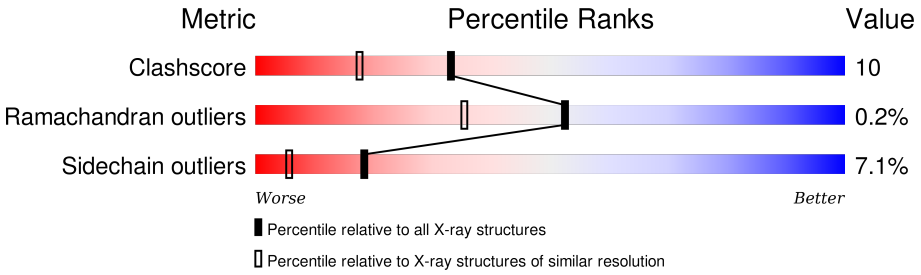
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

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 <=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	64% 28% 6%
1	B	217	68% 25% 6%

2 Entry composition

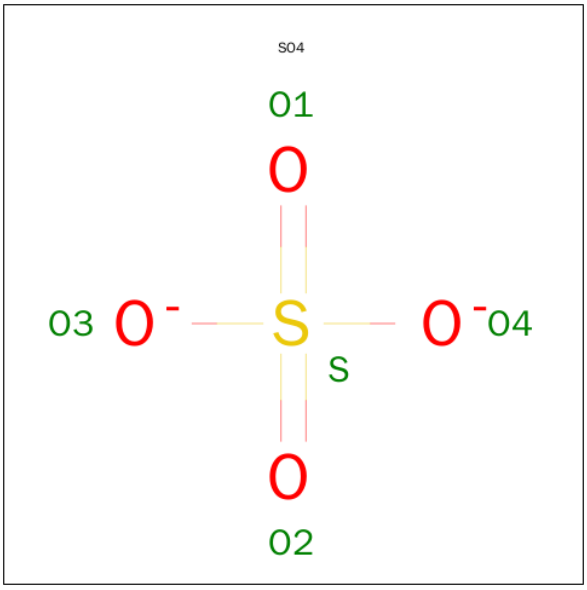
There are 4 unique types of molecules in this entry. The entry contains 4186 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

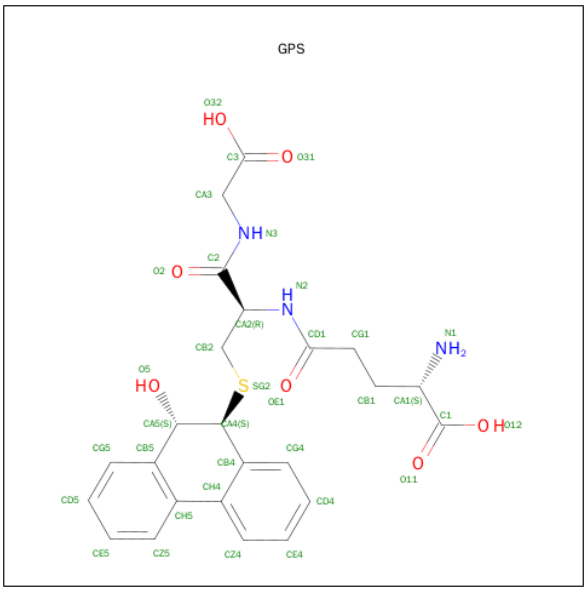
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	0	0
			1818	1177	303	327	11			
1	B	217	Total	C	N	O	S	0	0	0
			1818	1177	303	327	11			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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: $C_{24}H_{27}N_3O_7S$).



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	272	Total	O	0	0
			272	272		
4	B	198	Total	O	0	0
			198	198		

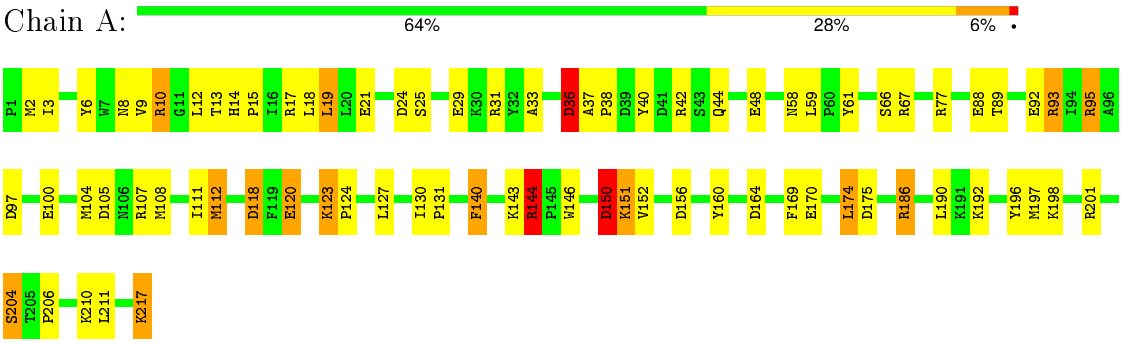
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

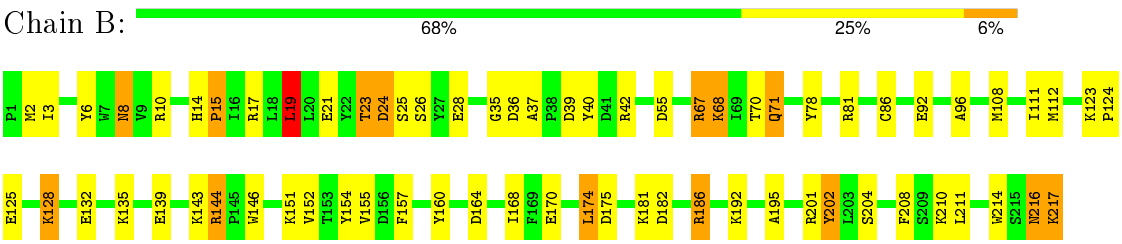
- Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain A:



- Molecule 1: GLUTATHIONE S-TRANSFERASE

Chain B:



4 Data and refinement statistics (i)

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	88.24Å 69.44Å 81.28Å 90.00° 106.01° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	GPRLSA	Depositor
R, R_{free}	0.160 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4186	wwPDB-VP
Average B, all atoms (Å^2)	17.0	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	1.17	3/1867 (0.2%)	1.88	51/2515 (2.0%)
1	B	1.13	4/1867 (0.2%)	1.89	41/2515 (1.6%)
All	All	1.15	7/3734 (0.2%)	1.89	92/5030 (1.8%)

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	1
1	B	0	2
All	All	0	3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	100	GLU	CD-OE1	-7.81	1.17	1.25
1	B	21	GLU	CD-OE1	-5.87	1.19	1.25
1	A	92	GLU	CD-OE1	-5.79	1.19	1.25
1	B	78	TYR	CB-CG	5.45	1.59	1.51
1	A	29	GLU	CD-OE1	-5.29	1.19	1.25
1	B	92	GLU	CD-OE1	-5.25	1.19	1.25
1	B	81	ARG	CZ-NH1	5.14	1.39	1.33

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	186	ARG	NE-CZ-NH1	23.33	131.97	120.30
1	B	186	ARG	NE-CZ-NH2	-20.81	109.89	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	186	ARG	NE-CZ-NH1	14.71	127.66	120.30
1	B	144	ARG	NE-CZ-NH1	14.24	127.42	120.30
1	B	144	ARG	NE-CZ-NH2	-13.48	113.56	120.30
1	A	17	ARG	NE-CZ-NH1	13.24	126.92	120.30
1	B	42	ARG	NE-CZ-NH1	13.07	126.83	120.30
1	B	10	ARG	NE-CZ-NH2	-12.91	113.84	120.30
1	A	10	ARG	NE-CZ-NH2	-12.20	114.20	120.30
1	A	144	ARG	NE-CZ-NH1	11.87	126.23	120.30
1	B	186	ARG	CD-NE-CZ	11.22	139.31	123.60
1	B	201	ARG	NE-CZ-NH1	10.87	125.73	120.30
1	B	17	ARG	NE-CZ-NH2	-9.99	115.30	120.30
1	B	144	ARG	CD-NE-CZ	9.51	136.92	123.60
1	A	144	ARG	CD-NE-CZ	9.30	136.62	123.60
1	A	105	ASP	CB-CG-OD1	9.14	126.53	118.30
1	B	42	ARG	NE-CZ-NH2	-8.94	115.83	120.30
1	A	95	ARG	CD-NE-CZ	8.94	136.11	123.60
1	B	78	TYR	CB-CG-CD2	-8.86	115.69	121.00
1	A	175	ASP	CB-CG-OD2	-8.70	110.47	118.30
1	A	42	ARG	NE-CZ-NH1	8.66	124.63	120.30
1	A	59	LEU	CB-CG-CD1	-8.56	96.44	111.00
1	A	17	ARG	NE-CZ-NH2	-8.44	116.08	120.30
1	A	67	ARG	NE-CZ-NH2	-8.39	116.10	120.30
1	A	156	ASP	CB-CG-OD2	-8.35	110.78	118.30
1	B	6	TYR	CB-CG-CD1	-8.31	116.02	121.00
1	A	31	ARG	NE-CZ-NH1	7.94	124.27	120.30
1	A	186	ARG	CD-NE-CZ	7.90	134.66	123.60
1	B	10	ARG	NE-CZ-NH1	7.88	124.24	120.30
1	B	17	ARG	NH1-CZ-NH2	7.77	127.95	119.40
1	A	42	ARG	NE-CZ-NH2	-7.69	116.46	120.30
1	A	118	ASP	CB-CG-OD2	-7.65	111.41	118.30
1	A	186	ARG	NE-CZ-NH2	-7.62	116.49	120.30
1	B	157	PHE	CB-CG-CD2	-7.60	115.48	120.80
1	B	17	ARG	NE-CZ-NH1	-7.33	116.64	120.30
1	A	42	ARG	CD-NE-CZ	7.28	133.79	123.60
1	A	105	ASP	CB-CG-OD2	-7.21	111.81	118.30
1	B	19	LEU	CB-CG-CD2	7.14	123.14	111.00
1	B	201	ARG	NH1-CZ-NH2	-6.84	111.88	119.40
1	B	201	ARG	CD-NE-CZ	6.83	133.16	123.60
1	B	202	TYR	CB-CG-CD2	-6.63	117.02	121.00
1	B	182	ASP	CB-CG-OD2	-6.59	112.37	118.30
1	B	24	ASP	CB-CG-OD1	6.41	124.07	118.30
1	A	150	ASP	CA-CB-CG	-6.38	99.36	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	78	TYR	CG-CD2-CE2	-6.35	116.22	121.30
1	B	154	TYR	CB-CG-CD2	-6.34	117.20	121.00
1	A	160	TYR	CB-CG-CD2	-6.33	117.20	121.00
1	A	61	TYR	CD1-CE1-CZ	-6.25	114.18	119.80
1	A	29	GLU	OE1-CD-OE2	6.21	130.76	123.30
1	B	78	TYR	CD1-CG-CD2	6.02	124.52	117.90
1	A	107	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	B	3	ILE	O-C-N	5.98	132.26	122.70
1	A	127	LEU	O-C-N	5.93	132.18	122.70
1	A	77	ARG	NE-CZ-NH2	-5.92	117.34	120.30
1	A	120	GLU	CA-CB-CG	5.91	126.39	113.40
1	A	169	PHE	CB-CG-CD1	-5.77	116.76	120.80
1	A	48	GLU	O-C-N	5.75	131.90	122.70
1	B	28	GLU	CA-CB-CG	5.73	126.01	113.40
1	A	61	TYR	CB-CG-CD1	-5.71	117.57	121.00
1	A	164	ASP	CB-CG-OD2	-5.71	113.16	118.30
1	B	15	PRO	O-C-N	5.69	131.81	122.70
1	B	17	ARG	O-C-N	-5.64	113.67	122.70
1	B	68	LYS	O-C-N	5.62	131.68	122.70
1	A	93	ARG	CD-NE-CZ	5.57	131.40	123.60
1	A	156	ASP	O-C-N	5.57	131.61	122.70
1	B	164	ASP	CB-CG-OD2	-5.56	113.29	118.30
1	A	19	LEU	CA-CB-CG	5.50	127.94	115.30
1	A	36	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	A	97	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	A	10	ARG	CD-NE-CZ	-5.44	115.99	123.60
1	A	58	ASN	N-CA-CB	-5.40	100.88	110.60
1	B	195	ALA	N-CA-CB	-5.38	102.57	110.10
1	A	10	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	A	156	ASP	OD1-CG-OD2	5.27	133.32	123.30
1	A	140	PHE	CB-CG-CD1	-5.25	117.13	120.80
1	B	23	THR	O-C-N	5.23	131.07	122.70
1	B	160	TYR	CG-CD2-CE2	5.22	125.47	121.30
1	A	204	SER	CA-CB-OG	-5.20	97.15	111.20
1	B	23	THR	CA-C-O	-5.17	109.25	120.10
1	A	174	LEU	CB-CG-CD1	5.16	119.77	111.00
1	A	197	MET	CA-CB-CG	-5.16	104.53	113.30
1	B	81	ARG	O-C-N	5.12	130.89	122.70
1	A	170	GLU	CG-CD-OE2	5.10	128.49	118.30
1	A	6	TYR	CB-CG-CD1	-5.09	117.95	121.00
1	B	55	ASP	CB-CG-OD2	-5.09	113.72	118.30
1	A	18	LEU	CB-CA-C	5.05	119.79	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	144	ARG	NH1-CZ-NH2	-5.05	113.85	119.40
1	B	164	ASP	CB-CG-OD1	5.05	122.84	118.30
1	B	78	TYR	CA-CB-CG	-5.03	103.84	113.40
1	B	81	ARG	NE-CZ-NH1	-5.01	117.79	120.30
1	A	190	LEU	O-C-N	5.00	130.71	122.70
1	A	61	TYR	CG-CD2-CE2	-5.00	117.30	121.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

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

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1818	0	1805	37	1
1	B	1818	0	1805	39	1
2	A	10	0	0	0	0
3	A	35	0	25	2	0
3	B	35	0	25	2	0
4	A	272	0	0	4	0
4	B	198	0	0	4	0
All	All	4186	0	3660	76	1

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 10.

All (76) 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:2:MET:CE	1:B:25:SER:HB3	2.04	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASP:OD2	1:A:192:LYS:HE3	1.76	0.85
1:B:2:MET:HE1	1:B:25:SER:HB3	1.59	0.83
1:B:111:ILE:HD11	3:B:218:GPS:HD4	1.69	0.73
1:B:86:CYS:SG	4:B:335:HOH:O	2.46	0.72
1:B:37:ALA:HB2	1:B:40:TYR:CZ	2.26	0.70
1:A:144:ARG:HH11	1:A:144:ARG:HG3	1.57	0.70
1:B:125:GLU:OE1	1:B:128:LYS:NZ	2.26	0.69
1:B:2:MET:HE2	1:B:26:SER:O	1.92	0.69
1:A:111:ILE:HD11	3:A:220:GPS:HD4	1.77	0.66
1:B:125:GLU:HA	1:B:128:LYS:HE3	1.76	0.65
1:B:2:MET:HE3	1:B:25:SER:HB3	1.78	0.65
1:A:150:ASP:HB3	1:A:151:LYS:HE3	1.80	0.63
1:B:125:GLU:O	1:B:128:LYS:HG3	1.99	0.62
1:A:3:ILE:HD12	1:A:3:ILE:N	2.15	0.61
1:B:123:LYS:HB3	1:B:124:PRO:HD3	1.84	0.59
1:B:67:ARG:NH2	4:B:396:HOH:O	2.36	0.59
1:B:68:LYS:NZ	4:B:364:HOH:O	2.36	0.59
1:A:2:MET:C	1:A:3:ILE:HD12	2.24	0.58
1:A:33:ALA:H	1:A:44:GLN:NE2	2.03	0.57
1:B:111:ILE:HD11	3:B:218:GPS:CD4	2.36	0.56
1:A:44:GLN:NE2	4:A:356:HOH:O	2.38	0.56
1:B:96:ALA:HA	1:B:155:VAL:HG11	1.87	0.56
1:B:8:ASN:H	1:B:8:ASN:HD22	1.54	0.55
1:B:143:LYS:NZ	4:B:411:HOH:O	2.38	0.55
1:A:111:ILE:HD11	3:A:220:GPS:CD4	2.39	0.53
1:B:135:LYS:HE3	1:B:139:GLU:OE2	2.09	0.53
1:B:35:GLY:O	1:B:40:TYR:HA	2.10	0.52
1:A:95:ARG:NH2	1:A:144:ARG:HE	2.07	0.52
1:A:144:ARG:NH1	1:A:144:ARG:HG3	2.19	0.52
1:A:140:PHE:O	1:A:144:ARG:NH2	2.43	0.52
1:A:40:TYR:CZ	1:A:211:LEU:HD13	2.45	0.52
1:A:21:GLU:HG3	1:A:196:TYR:CD1	2.45	0.51
1:A:40:TYR:CE2	1:A:211:LEU:HD13	2.46	0.51
1:B:37:ALA:HB2	1:B:40:TYR:OH	2.10	0.51
1:A:198:LYS:NZ	4:A:410:HOH:O	2.43	0.51
1:A:36:ASP:OD1	1:A:210:LYS:NZ	2.33	0.50
1:B:175:ASP:OD1	1:B:181:LYS:HE3	2.12	0.50
1:A:143:LYS:HA	1:A:143:LYS:HE2	1.94	0.49
1:B:175:ASP:OD1	1:B:181:LYS:CE	2.60	0.49
1:A:2:MET:CE	1:A:25:SER:HB3	2.43	0.48
1:B:8:ASN:N	1:B:8:ASN:HD22	2.10	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:VAL:HG12	1:A:206:PRO:HG2	1.96	0.47
1:A:24:ASP:OD2	1:A:192:LYS:CE	2.56	0.47
1:B:146:TRP:CE2	1:B:152:VAL:HG22	2.49	0.47
1:A:21:GLU:HG3	1:A:196:TYR:CG	2.51	0.46
1:A:108:MET:O	1:A:112:MET:HG2	2.16	0.46
1:A:201:ARG:NH2	4:A:415:HOH:O	2.48	0.46
1:A:95:ARG:CZ	1:A:144:ARG:HE	2.29	0.46
1:B:174:LEU:HB3	1:B:181:LYS:HG3	1.98	0.46
1:B:108:MET:O	1:B:112:MET:HG2	2.15	0.46
1:B:125:GLU:OE1	1:B:128:LYS:CE	2.65	0.45
1:A:89:THR:O	1:A:93:ARG:HG3	2.16	0.45
1:A:151:LYS:NZ	4:A:379:HOH:O	2.50	0.45
1:A:146:TRP:CE2	1:A:152:VAL:HG22	2.52	0.45
1:B:216:ASN:C	1:B:216:ASN:HD22	2.20	0.44
1:B:202:TYR:CZ	1:B:204:SER:HB3	2.53	0.44
1:B:168:ILE:HG22	1:B:214:TRP:HZ2	1.82	0.43
1:A:10:ARG:HD2	1:A:204:SER:O	2.19	0.43
1:A:14:HIS:N	1:A:15:PRO:CD	2.82	0.43
1:B:123:LYS:N	1:B:124:PRO:CD	2.81	0.43
1:A:217:LYS:HB2	1:A:217:LYS:HE2	1.46	0.43
1:A:13:THR:O	1:A:13:THR:HG22	2.18	0.43
1:B:2:MET:HE1	1:B:25:SER:CB	2.40	0.42
1:A:37:ALA:HB1	1:A:38:PRO:HA	2.01	0.42
1:B:8:ASN:H	1:B:8:ASN:ND2	2.17	0.42
1:B:19:LEU:HD22	1:B:23:THR:HG23	2.01	0.42
1:B:111:ILE:HG12	1:B:208:PHE:HE1	1.85	0.41
1:B:216:ASN:ND2	1:B:217:LYS:HG2	2.35	0.41
1:B:24:ASP:CG	1:B:192:LYS:NZ	2.73	0.41
1:A:36:ASP:HA	1:A:210:LYS:HD3	2.03	0.41
1:B:70:THR:O	1:B:71:GLN:HB2	2.20	0.41
1:B:14:HIS:N	1:B:15:PRO:CD	2.83	0.41
1:A:130:ILE:N	1:A:131:PRO:CD	2.83	0.41
1:A:123:LYS:HB3	1:A:124:PRO:HD3	2.03	0.41
1:A:104:MET:O	1:A:108:MET:HG2	2.21	0.40

All (1) 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:88:GLU:OE2	1:B:217:LYS:NZ[3_455]	2.05	0.15

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%)	3 (1%)	1 (0%)	34 17
All	All	430/434 (99%)	423 (98%)	6 (1%)	1 (0%)	52 35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	71	GLN

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%)	183 (93%)	14 (7%)	18 6
1	B	197/197 (100%)	183 (93%)	14 (7%)	18 6
All	All	394/394 (100%)	366 (93%)	28 (7%)	18 6

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

Mol	Chain	Res	Type
1	A	8	ASN
1	A	12	LEU
1	A	19	LEU
1	A	36	ASP
1	A	66	SER

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Mol	Chain	Res	Type
1	A	112	MET
1	A	118	ASP
1	A	120	GLU
1	A	123	LYS
1	A	144	ARG
1	A	150	ASP
1	A	151	LYS
1	A	174	LEU
1	A	217	LYS
1	B	8	ASN
1	B	19	LEU
1	B	36	ASP
1	B	39	ASP
1	B	67	ARG
1	B	128	LYS
1	B	132	GLU
1	B	151	LYS
1	B	170	GLU
1	B	174	LEU
1	B	210	LYS
1	B	211	LEU
1	B	216	ASN
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	44	GLN
1	B	8	ASN
1	B	122	GLN
1	B	216	ASN

5.3.3 RNA (i)

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)

4 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	1.43	0	6,6,6	0.42	0
2	SO4	A	219	-	4,4,4	1.12	0	6,6,6	1.01	1 (16%)
3	GPS	A	220	-	29,37,37	1.40	5 (17%)	35,51,51	1.74	9 (25%)
3	GPS	B	218	-	29,37,37	1.35	4 (13%)	35,51,51	1.52	8 (22%)

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
3	GPS	A	220	-	-	0/21/43/43	0/3/3/3
3	GPS	B	218	-	-	0/21/43/43	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	218	GPS	CH5-CH4	-4.03	1.39	1.47
3	A	220	GPS	CH5-CH4	-2.60	1.42	1.47
3	A	220	GPS	CH5-CB5	-2.31	1.37	1.40
3	B	218	GPS	OE1-CD1	2.23	1.27	1.23
3	A	220	GPS	O5-CA5	2.32	1.47	1.42
3	B	218	GPS	CB2-CA2	2.34	1.59	1.53
3	B	218	GPS	CG5-CB5	2.40	1.43	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	220	GPS	CG5-CB5	2.88	1.43	1.39
3	A	220	GPS	CG4-CB4	3.02	1.43	1.39

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	220	GPS	O2-C2-CA2	-5.06	109.18	120.36
3	A	220	GPS	CD4-CG4-CB4	-2.97	117.16	121.02
3	A	220	GPS	CB4-CA4-CA5	-2.69	103.75	109.89
3	B	218	GPS	O2-C2-CA2	-2.65	114.50	120.36
3	B	218	GPS	CD4-CG4-CB4	-2.58	117.66	121.02
3	A	220	GPS	CB2-CA2-N2	-2.53	104.59	110.83
3	A	220	GPS	C2-CA2-N2	-2.41	104.46	111.26
3	A	220	GPS	CG1-CD1-N2	-2.26	112.15	115.83
3	A	220	GPS	O5-CA5-CB5	-2.22	105.80	110.42
3	A	220	GPS	CB1-CA1-N1	-2.18	104.33	110.52
3	B	218	GPS	CB2-CA2-N2	-2.18	105.46	110.83
3	B	218	GPS	CB4-CA4-CA5	-2.05	105.21	109.89
3	B	218	GPS	OE1-CD1-N2	-2.04	119.54	123.01
2	A	219	SO4	O4-S-O3	2.02	117.19	108.98
3	B	218	GPS	O2-C2-N3	2.20	127.49	123.08
3	B	218	GPS	CA3-N3-C2	2.46	125.73	122.34
3	B	218	GPS	CZ4-CH4-CB4	2.60	121.26	118.41
3	A	220	GPS	O2-C2-N3	3.17	129.44	123.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

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

5.7 Other polymers (i)

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 will therefore be empty.

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

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

6.3 Carbohydrates

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

6.4 Ligands

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

6.5 Other polymers

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