



Full wwPDB X-ray Structure Validation Report

Mar 1, 2014 – 02:11 AM GMT

PDB ID : 3LXZ
Title : Structure of probable Glutathione S-transferase(PP0183)from Pseudomonas putida
Authors : Ramagopal, U.A.; Toro, R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2010-02-25
Resolution : 1.76 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

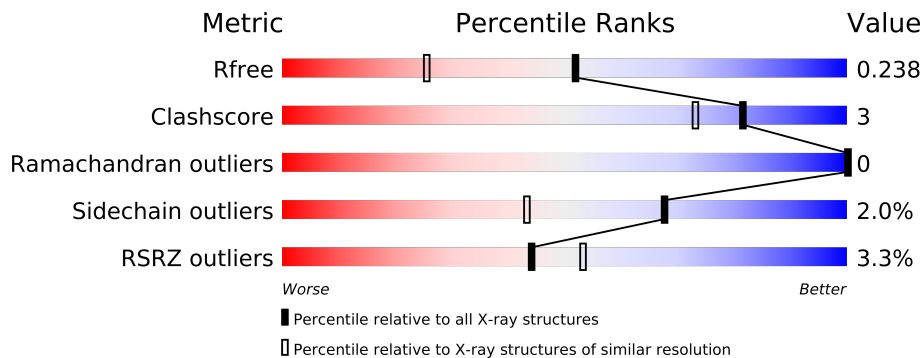
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.76 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))
R_{free}	66092	1134 (1.76-1.76)
Clashscore	79885	1304 (1.76-1.76)
Ramachandran outliers	78287	1288 (1.76-1.76)
Sidechain outliers	78261	1288 (1.76-1.76)
RSRZ outliers	66119	1135 (1.76-1.76)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	229	
1	B	229	
1	C	229	
1	D	229	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7538 atoms, of which 0 are hydrogen and 0 are deuterium.

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 family protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	218	Total	C	N	O	S	Se	0	1	0
			1699	1101	272	316	2	8			
1	B	217	Total	C	N	O	S	Se	0	5	0
			1729	1120	279	320	2	8			
1	C	218	Total	C	N	O	S	Se	0	6	0
			1747	1127	286	324	2	8			
1	D	217	Total	C	N	O	S	Se	0	3	0
			1710	1105	279	316	2	8			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	EXPRESSION TAG	UNP Q88RE7
A	1	SER	-	EXPRESSION TAG	UNP Q88RE7
A	221	GLU	-	EXPRESSION TAG	UNP Q88RE7
A	222	GLY	-	EXPRESSION TAG	UNP Q88RE7
A	223	HIS	-	EXPRESSION TAG	UNP Q88RE7
A	224	HIS	-	EXPRESSION TAG	UNP Q88RE7
A	225	HIS	-	EXPRESSION TAG	UNP Q88RE7
A	226	HIS	-	EXPRESSION TAG	UNP Q88RE7
A	227	HIS	-	EXPRESSION TAG	UNP Q88RE7
A	228	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	0	MSE	-	EXPRESSION TAG	UNP Q88RE7
B	1	SER	-	EXPRESSION TAG	UNP Q88RE7
B	221	GLU	-	EXPRESSION TAG	UNP Q88RE7
B	222	GLY	-	EXPRESSION TAG	UNP Q88RE7
B	223	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	224	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	225	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	226	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	227	HIS	-	EXPRESSION TAG	UNP Q88RE7
B	228	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	0	MSE	-	EXPRESSION TAG	UNP Q88RE7

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	SER	-	EXPRESSION TAG	UNP Q88RE7
C	221	GLU	-	EXPRESSION TAG	UNP Q88RE7
C	222	GLY	-	EXPRESSION TAG	UNP Q88RE7
C	223	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	224	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	225	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	226	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	227	HIS	-	EXPRESSION TAG	UNP Q88RE7
C	228	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	0	MSE	-	EXPRESSION TAG	UNP Q88RE7
D	1	SER	-	EXPRESSION TAG	UNP Q88RE7
D	221	GLU	-	EXPRESSION TAG	UNP Q88RE7
D	222	GLY	-	EXPRESSION TAG	UNP Q88RE7
D	223	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	224	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	225	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	226	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	227	HIS	-	EXPRESSION TAG	UNP Q88RE7
D	228	HIS	-	EXPRESSION TAG	UNP Q88RE7

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	148	Total O 148 148	0	0
2	B	181	Total O 181 181	0	0
2	C	170	Total O 170 170	0	0
2	D	154	Total O 154 154	0	0

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.

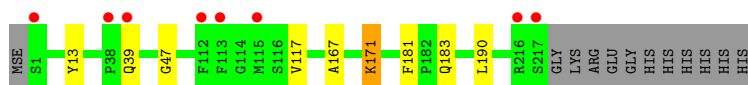
- Molecule 1: Glutathione S-transferase family protein

Chain A: 



- Molecule 1: Glutathione S-transferase family protein

Chain B: 



- Molecule 1: Glutathione S-transferase family protein

Chain C: 



- Molecule 1: Glutathione S-transferase family protein

Chain D: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.05Å 66.36Å 88.66Å 89.96° 85.27° 87.04°	Depositor
Resolution (Å)	50.00 – 1.76 30.58 – 1.76	Depositor EDS
% Data completeness (in resolution range)	90.6 (50.00-1.76) 90.6 (30.58-1.76)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 1.76Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.196 , 0.229 0.206 , 0.238	Depositor DCC
R_{free} test set	4901 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.6	EDS
Estimated twinning fraction	0.064 for -h,k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 97746 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7538	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.73	1/1726 (0.1%)	0.73	2/2318 (0.1%)
1	B	0.74	1/1755 (0.1%)	0.72	0/2355
1	C	0.75	0/1773	0.75	0/2377
1	D	0.75	1/1735 (0.1%)	0.72	0/2328
All	All	0.74	3/6989 (0.0%)	0.73	2/9378 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	205	GLU	CG-CD	5.22	1.59	1.51
1	A	159	CYS	CB-SG	-5.11	1.73	1.81
1	B	13	TYR	CD1-CE1	5.08	1.47	1.39

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	A	138	ARG	NE-CZ-NH1	5.26	122.93	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1699	0	1717	6	0
1	B	1729	0	1755	6	0
1	C	1747	0	1767	22	0
1	D	1710	0	1737	12	0
2	A	148	0	0	1	0
2	B	181	0	0	3	0
2	C	170	0	0	2	0
2	D	154	0	0	6	0
All	All	7538	0	6976	45	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (45) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:208:MSE:SE	1:C:212:MSE:HE3	2.21	0.90
1:C:211:PHE:CD2	1:C:212:MSE:HE2	2.08	0.87
1:C:211:PHE:CD2	1:C:212:MSE:CE	2.61	0.83
1:C:211:PHE:HD2	1:C:212:MSE:CE	2.02	0.71
1:C:47:GLY:HA2	2:C:245:HOH:O	1.93	0.68
1:D:126:ARG:NH1	1:D:175:ILE:HD11	2.08	0.68
1:C:211:PHE:HD2	1:C:212:MSE:HE2	1.57	0.65
1:A:126:ARG:HH11	1:A:175:ILE:HD11	1.61	0.64
1:D:126:ARG:HH12	1:D:175:ILE:HD11	1.64	0.62
1:C:211:PHE:CE2	1:C:212:MSE:HE1	2.36	0.61
1:C:214:MSE:HE3	1:C:215:ILE:CD1	2.31	0.61
1:A:185:LYS:HE2	2:A:443:HOH:O	2.01	0.61
1:B:47:GLY:HA2	2:B:281:HOH:O	2.02	0.58
1:D:118:GLU:HG3	2:D:274:HOH:O	2.05	0.56
1:C:214:MSE:HE3	1:C:215:ILE:HD13	1.88	0.56
1:C:211:PHE:CD2	1:C:212:MSE:HE1	2.38	0.55
1:D:47:GLY:HA2	2:D:261:HOH:O	2.06	0.55
1:A:126:ARG:NH1	1:A:175:ILE:HD11	2.23	0.54
1:B:167:ALA:O	1:B:171:LYS:HD3	2.08	0.53
1:B:171:LYS:N	1:B:171:LYS:HD3	2.23	0.52
1:C:166:ASN:HD21	1:C:177:PHE:N	2.08	0.52
1:A:185:LYS:HD3	2:D:248:HOH:O	2.10	0.52
1:C:46:ARG:HG3	1:C:46:ARG:O	2.11	0.50
1:D:183:GLN:CG	2:D:420:HOH:O	2.59	0.50
1:D:11:ASN:O	1:D:15:MSE:HG3	2.14	0.47
1:D:113:PHE:HB3	1:D:115:MSE:HE3	1.97	0.47
1:C:39[B]:GLN:H	1:C:39[B]:GLN:CD	2.19	0.46
1:B:190:LEU:C	1:B:190:LEU:HD23	2.37	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:6:GLY:HA3	1:A:13:TYR:CD1	2.53	0.44
1:C:141:ARG:NH1	2:C:360:HOH:O	2.51	0.44
1:C:190:LEU:HD23	1:C:190:LEU:C	2.39	0.43
1:C:44:SER:HB2	1:C:51:VAL:HG21	1.99	0.43
1:C:166:ASN:HD21	1:C:177:PHE:H	1.66	0.43
1:C:166:ASN:ND2	1:C:177:PHE:H	2.16	0.43
1:D:6:GLY:HA3	1:D:13:TYR:CD1	2.52	0.43
1:D:183:GLN:HG2	2:D:420:HOH:O	2.17	0.43
1:A:92:LEU:HD13	1:A:146:VAL:HG12	2.01	0.43
1:D:137:LYS:HE3	2:D:651:HOH:O	2.18	0.42
1:D:44:SER:HB2	1:D:51:VAL:HG21	2.02	0.42
1:B:183:GLN:HG2	2:B:319:HOH:O	2.19	0.41
1:B:171:LYS:H	1:B:171:LYS:HD3	1.86	0.41
1:C:18:LEU:HD23	1:C:156:LEU:HD22	2.02	0.41
1:C:45:PRO:HB2	1:D:135:THR:HG23	2.03	0.41
1:C:112:PHE:N	1:C:112:PHE:CD1	2.89	0.41
2:B:240:HOH:O	1:C:185:LYS:HD3	2.20	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	217/229 (95%)	213 (98%)	4 (2%)	0	100	100
1	B	220/229 (96%)	215 (98%)	5 (2%)	0	100	100
1	C	222/229 (97%)	219 (99%)	3 (1%)	0	100	100
1	D	218/229 (95%)	214 (98%)	4 (2%)	0	100	100
All	All	877/916 (96%)	861 (98%)	16 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	179/180 (99%)	175 (98%)	4 (2%)	64	38
1	B	184/180 (102%)	179 (97%)	5 (3%)	57	30
1	C	185/180 (103%)	182 (98%)	3 (2%)	75	55
1	D	181/180 (101%)	177 (98%)	4 (2%)	64	38
All	All	729/720 (101%)	713 (98%)	16 (2%)	68	38

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	180[A]	ASP
1	A	180[B]	ASP
1	A	181	PHE
1	B	39	GLN
1	B	117[A]	VAL
1	B	117[B]	VAL
1	B	171	LYS
1	B	181	PHE
1	C	181	PHE
1	C	189	GLN
1	C	214	MSE
1	D	120	LEU
1	D	171	LYS
1	D	181	PHE
1	D	189	GLN

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

Mol	Chain	Res	Type
1	B	39	GLN
1	C	166	ASN
1	C	189	GLN
1	D	189	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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, 95th 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 ²)	Q<0.9
1	A	218/229 (95%)	0.05	6 (2%)	50	60	14, 21, 34, 47	0
1	B	217/229 (94%)	-0.06	8 (3%)	39	48	13, 18, 33, 47	0
1	C	218/229 (95%)	-0.03	8 (3%)	39	48	11, 18, 32, 49	0
1	D	217/229 (94%)	0.04	7 (3%)	45	54	13, 19, 34, 55	0
All	All	870/916 (94%)	0.00	29 (3%)	44	53	11, 19, 33, 55	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	115	MSE	4.2
1	D	217	SER	4.0
1	C	218	GLY	3.8
1	B	1	SER	3.7
1	A	1	SER	3.7
1	C	112	PHE	3.4
1	D	114	GLY	3.3
1	B	217	SER	3.1
1	C	113	PHE	3.0
1	C	39[A]	GLN	3.0
1	C	217	SER	2.9
1	D	1	SER	2.8
1	B	112	PHE	2.7
1	C	216	ARG	2.7
1	A	215	ILE	2.6
1	C	1	SER	2.6
1	B	115	MSE	2.5
1	D	113	PHE	2.5
1	B	216	ARG	2.4
1	D	216	ARG	2.3
1	B	38	PRO	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	113	PHE	2.3
1	B	113	PHE	2.3
1	C	38	PRO	2.2
1	A	115	MSE	2.1
1	A	39	GLN	2.1
1	D	38	PRO	2.1
1	B	39	GLN	2.0
1	A	216	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.