

Full wwPDB X-ray Structure Validation Report (i)

Jan 27, 2024 – 02:30 PM EST

PDB ID : 1AW9

Title : STRUCTURE OF GLUTATHIONE S-TRANSFERASE III IN APO FORM

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Deposited on : 1997-10-13

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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity: 4.02b-467

Xtriage (Phenix) : NOT EXECUTED EDS : NOT EXECUTED

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

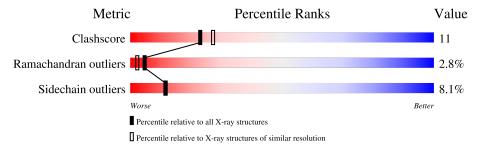
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY\ DIFFRACTION$

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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain		
1	A	216	75%	21%	•



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2485 atoms, of which 657 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 III.

Mol	Chain	Residues		Atoms			ZeroOcc	AltConf	Trace		
1	A	216	Total 1997	C 1083	H 335	N 278	O 300	S 1	5	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	48	ALA	THR	conflict	UNP Q9ZP62

• Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total Cd 5 5	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	A	161	Total 483	H 322	O 161	0	0

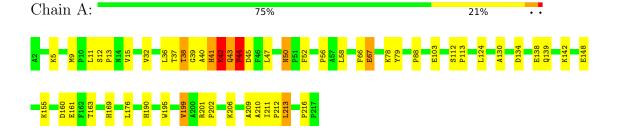


3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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 III





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants	97.77Å 97.77Å 116.45Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 - 2.20	Depositor
% Data completeness	(Not available) (8.00-2.20)	Depositor
(in resolution range)	(1100 available) (0.00 2.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.197 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2485	wwPDB-VP
Average B, all atoms (Å ²)	21.0	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: CD

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).

Mal	Chain	Bond	lengths	Bo	nd angles
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.67	0/1710	0.80	2/2343 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}(^{o})$
1	A	134	ASP	CB-CG-OD1	-5.95	112.94	118.30
1	A	134	ASP	CB-CG-OD2	5.09	122.88	118.30

There are no chirality outliers.

There are no planarity outliers.

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	1662	335	1677	38	1
2	A	5	0	0	0	0
3	A	161	322	0	6	1
All	All	1828	657	1677	38	2

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 11.

All (38) close contacts within the same asymmetric unit are listed below, sorted by their clash



magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	${\rm distance}(\mathring{\rm A})$	overlap (Å)
1:A:44:PRO:HB3	1:A:47:LEU:HD12	1.59	0.83
1:A:5:LYS:HE2	1:A:32:VAL:HG21	1.61	0.81
1:A:44:PRO:HA	3:A:375:HOH:O	1.89	0.70
1:A:42:LYS:HD3	1:A:42:LYS:H	1.59	0.67
1:A:37:THR:HB	3:A:311:HOH:O	1.97	0.65
1:A:148:GLU:OE2	1:A:190:HIS:HD2	1.78	0.65
1:A:9:MET:HE2	3:A:287:HOH:O	2.04	0.56
1:A:155:LYS:HZ3	1:A:160:ASP:HA	1.71	0.55
1:A:112:SER:OG	1:A:113:PRO:HD3	2.09	0.53
1:A:155:LYS:NZ	1:A:160:ASP:HA	2.24	0.52
1:A:13:PRO:HD2	3:A:246:HOH:O	2.10	0.52
1:A:41:HIS:O	1:A:43:GLN:HG3	2.10	0.52
1:A:195:TRP:O	1:A:199:VAL:HG23	2.10	0.51
1:A:78:LYS:HD3	1:A:79:TYR:CE2	2.45	0.51
1:A:50:ASN:ND2	1:A:52:PHE:H	2.09	0.51
1:A:202:PRO:HD2	3:A:298:HOH:O	2.12	0.50
1:A:42:LYS:HD3	1:A:42:LYS:N	2.28	0.48
1:A:88:PRO:HD2	1:A:163:THR:HG22	1.95	0.47
1:A:66:PHE:O	1:A:67:GLU:HB2	2.14	0.47
1:A:37:THR:O	1:A:39:GLY:N	2.45	0.47
1:A:210:ALA:HB1	3:A:273:HOH:O	2.16	0.46
1:A:155:LYS:NZ	1:A:155:LYS:HB2	2.31	0.46
1:A:11:LEU:HG	1:A:212:PRO:HG2	1.97	0.46
1:A:138:GLU:O	1:A:142:LYS:HG2	2.16	0.45
1:A:211:ILE:CG2	1:A:213:LEU:HD22	2.47	0.45
1:A:40:ALA:C	1:A:42:LYS:N	2.71	0.44
1:A:42:LYS:HE2	1:A:42:LYS:HB2	1.81	0.43
1:A:11:LEU:HD23	1:A:211:ILE:CG2	2.48	0.43
1:A:40:ALA:O	1:A:42:LYS:N	2.52	0.43
1:A:103:GLU:HB2	1:A:169:HIS:CE1	2.53	0.42
1:A:5:LYS:CE	1:A:32:VAL:HG21	2.39	0.42
1:A:206:LYS:O	1:A:209:ALA:HB3	2.20	0.42
1:A:37:THR:C	1:A:39:GLY:H	2.22	0.41
1:A:47:LEU:HD23	1:A:47:LEU:HA	1.82	0.41
1:A:201:ARG:HA	1:A:202:PRO:HD3	1.90	0.41
1:A:15:VAL:HG22	1:A:56:PRO:HB3	2.03	0.40
1:A:40:ALA:C	1:A:42:LYS:H	2.23	0.40
1:A:43:GLN:HA	1:A:44:PRO:HD2	1.63	0.40

All (2) 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	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	Clash overlap (Å)
1:A:130:ALA:H	1:A:139:GLN:HE22[2_655]	1.35	0.25
3:A:301:HOH:H1	3:A:318:HOH:O[2_655]	1.50	0.10

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	214/216 (99%)	200 (94%)	8 (4%)	6 (3%)	5 2

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	THR
1	A	44	PRO
1	A	41	HIS
1	A	42	LYS
1	A	67	GLU
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	173/173 (100%)	159 (92%)	14 (8%)	11 12

All (14) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Tuno
MIOI	Chain	nes	Type
1	A	12	SER
1	A	36	LEU
1	A	38	THR
1	A	42	LYS
1	A	43	GLN
1	A	44	PRO
1	A	45	ASP
1	A	50	ASN
1	A	58	LEU
1	A	124	LEU
1	A	161	GLU
1	A	176	LEU
1	A	199	VAL
1	A	213	LEU

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

Mol	Chain	Res	Type
1	A	50	ASN
1	A	139	GLN
1	A	190	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.



There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

