



Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:14 pm BST

PDB ID : 3WBX
Title : Crystal structure of Gox0644 at apoform
Authors : Yuan, Y.A.; Wang, C.
Deposited on : 2013-05-22
Resolution : 2.40 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

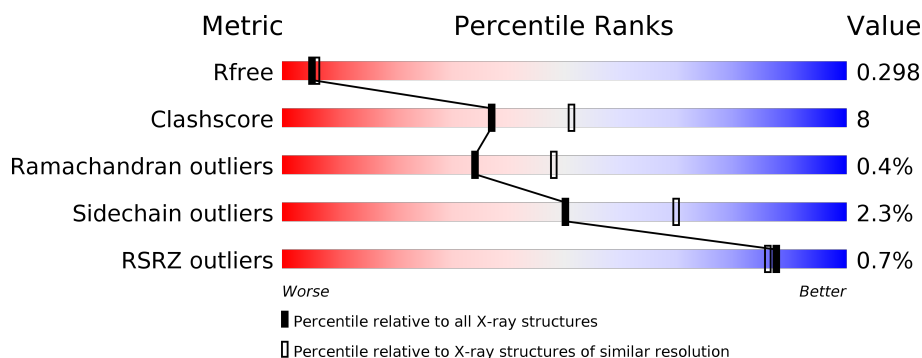
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 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 $\leq 5\%$. 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	282	
1	B	282	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4600 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 Putative 2,5-diketo-D-gluconic acid reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2183	1377	392	408	6			
1	B	271	Total	C	N	O	S	0	0	0
			2183	1377	392	408	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q5FT75
A	-1	SER	-	EXPRESSION TAG	UNP Q5FT75
A	0	HIS	-	EXPRESSION TAG	UNP Q5FT75
B	-2	GLY	-	EXPRESSION TAG	UNP Q5FT75
B	-1	SER	-	EXPRESSION TAG	UNP Q5FT75
B	0	HIS	-	EXPRESSION TAG	UNP Q5FT75

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	113	Total O 113 113	0	0
3	B	101	Total O 101 101	0	0

- Molecule 1: Putative 2,5-diketo-D-gluconic acid reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	55.88Å 75.78Å 125.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.40 31.36 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.9 (50.00-2.40) 98.0 (31.36-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.15 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.224 , 0.300 0.224 , 0.298	Depositor DCC
R_{free} test set	1084 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4600	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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

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	0.55	0/2232	0.65	0/3020
1	B	0.54	0/2232	0.62	0/3020
All	All	0.54	0/4464	0.64	0/6040

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2183	0	2161	31	0
1	B	2183	0	2161	37	0
2	A	5	0	0	0	0
2	B	15	0	0	0	0
3	A	113	0	0	1	1
3	B	101	0	0	2	0
All	All	4600	0	4322	68	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 8.

All (68) 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:168:LEU:HB3	1:B:190:SER:HB3	1.45	0.98
1:A:222:ARG:HE	1:A:226:GLN:HE21	1.08	0.96
1:A:226:GLN:HE22	1:A:252:LEU:H	1.13	0.96
1:B:226:GLN:HE22	1:B:252:LEU:H	1.06	0.93
1:B:222:ARG:HE	1:B:226:GLN:HE21	1.21	0.87
1:B:226:GLN:HE22	1:B:252:LEU:N	1.83	0.75
1:B:226:GLN:NE2	1:B:252:LEU:H	1.86	0.71
1:A:168:LEU:HB3	1:A:190:SER:HB3	1.72	0.70
1:A:222:ARG:HE	1:A:226:GLN:NE2	1.85	0.69
1:B:175:ARG:NH2	1:B:261:GLU:O	2.25	0.69
1:B:168:LEU:HB3	1:B:190:SER:CB	2.20	0.68
1:A:222:ARG:NE	1:A:226:GLN:HE21	1.87	0.67
1:B:66:LYS:O	1:B:69:GLU:HG2	1.96	0.66
1:B:146:GLU:H	1:B:149:HIS:HD2	1.46	0.63
1:B:62:GLU:HG3	1:B:100:LEU:HD22	1.80	0.63
1:A:25:ILE:HD12	1:A:245:LEU:HD12	1.83	0.61
1:B:215:THR:H	1:B:218:GLN:HE21	1.49	0.60
1:A:226:GLN:HE22	1:A:252:LEU:N	1.93	0.59
1:B:165:GLN:HA	1:B:189:GLU:O	2.02	0.59
1:B:215:THR:H	1:B:218:GLN:NE2	2.01	0.58
1:A:199:VAL:HG22	1:A:205:ILE:HG13	1.85	0.57
1:B:181:HIS:HD2	1:B:188:THR:OG1	1.88	0.57
1:B:224:HIS:HB3	1:B:231:VAL:HG22	1.88	0.55
1:B:222:ARG:NE	1:B:226:GLN:HE21	1.98	0.54
1:B:171:ASP:OD2	1:B:198:ARG:HD3	2.08	0.54
1:B:125:TRP:O	1:B:129:VAL:HG23	2.09	0.53
1:B:69:GLU:O	1:B:70:ASP:HB2	2.09	0.53
1:A:215:THR:H	1:A:218:GLN:HE21	1.57	0.53
1:A:99:ARG:NH1	3:A:509:HOH:O	2.41	0.53
1:A:25:ILE:HG23	1:A:225:LEU:HD21	1.90	0.52
1:B:168:LEU:CB	1:B:190:SER:HB3	2.29	0.52
1:A:164:ASN:HD22	1:A:181:HIS:CE1	2.27	0.52
1:A:164:ASN:HD22	1:A:181:HIS:HE1	1.57	0.52
1:A:238:PRO:HA	1:A:241:LEU:HB2	1.92	0.51
1:A:66:LYS:O	1:A:69:GLU:HB2	2.11	0.51
1:B:234:LYS:O	1:B:235:SER:HB2	2.11	0.50
1:A:231:VAL:HG22	1:A:233:PRO:HD3	1.94	0.50
1:A:25:ILE:HG22	1:A:231:VAL:HG13	1.95	0.49
1:A:88:ASP:O	1:A:92:ARG:HG3	2.12	0.49
1:A:28:GLY:O	1:A:234:LYS:HB3	2.14	0.48
1:A:74:ILE:O	1:A:103:ARG:NH2	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:TRP:CH2	1:A:191:TRP:HB2	2.49	0.47
1:B:188:THR:HB	1:B:229:LEU:HD22	1.97	0.47
1:A:80:LEU:HD13	1:A:90:THR:HG23	1.97	0.47
1:A:14:ILE:CG1	1:A:24:GLN:HG3	2.45	0.46
1:A:14:ILE:HG12	1:A:24:GLN:HG3	1.96	0.46
1:B:192:ARG:HD3	1:B:196:LYS:HG3	1.97	0.46
1:B:14:ILE:HG12	1:B:24:GLN:HG3	1.97	0.46
1:B:27:LEU:HD13	1:B:241:LEU:HD21	1.97	0.45
1:B:29:VAL:C	1:B:30:TRP:HD1	2.19	0.45
1:A:237:ASN:HD22	1:A:240:ARG:NE	2.14	0.45
1:B:146:GLU:H	1:B:149:HIS:CD2	2.31	0.44
1:A:25:ILE:CD1	1:A:245:LEU:HD12	2.47	0.44
1:A:27:LEU:HD13	1:A:241:LEU:HD21	1.99	0.44
1:A:75:PHE:HA	1:A:107:ASP:OD1	2.18	0.43
1:B:144:ASN:O	1:B:274:PRO:HB3	2.19	0.43
1:B:123:GLU:HB3	3:B:491:HOH:O	2.18	0.43
1:B:238:PRO:HA	1:B:241:LEU:HB2	2.01	0.43
1:B:55:ALA:HB3	1:B:58:TYR:CE1	2.54	0.41
1:B:66:LYS:HG2	1:B:102:ARG:NH2	2.35	0.41
1:A:33:PRO:HA	1:A:34:PRO:HD2	1.83	0.41
1:A:33:PRO:HG2	1:A:36:GLU:HG2	2.03	0.41
1:B:208:ILE:O	1:B:212:HIS:HD2	2.04	0.41
1:A:237:ASN:HA	1:A:238:PRO:HD2	1.92	0.41
1:B:50:ARG:HA	1:B:74:ILE:HD12	2.03	0.41
1:B:57:LEU:HD22	1:B:81:TRP:CZ3	2.56	0.41
1:B:237:ASN:HB2	3:B:406:HOH:O	2.21	0.40
1:B:258:GLN:O	1:B:262:GLN: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)
3:A:401:HOH:O	3:A:512:HOH:O[1_455]	2.19	0.01

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	269/282 (95%)	263 (98%)	5 (2%)	1 (0%)	34	48
1	B	269/282 (95%)	262 (97%)	6 (2%)	1 (0%)	34	48
All	All	538/564 (95%)	525 (98%)	11 (2%)	2 (0%)	34	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	235	SER
1	A	34	PRO

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	235/244 (96%)	230 (98%)	5 (2%)	53	72
1	B	235/244 (96%)	229 (97%)	6 (3%)	46	66
All	All	470/488 (96%)	459 (98%)	11 (2%)	50	70

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	LEU
1	A	190	SER
1	A	198	ARG
1	A	235	SER
1	A	240	ARG
1	B	60	ASN
1	B	76	LEU
1	B	152	ARG
1	B	190	SER
1	B	198	ARG

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Mol	Chain	Res	Type
1	B	239	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	20	HIS
1	A	118	GLN
1	A	181	HIS
1	A	218	GLN
1	A	226	GLN
1	A	262	GLN
1	B	60	ASN
1	B	85	GLN
1	B	118	GLN
1	B	120	GLN
1	B	149	HIS
1	B	181	HIS
1	B	218	GLN
1	B	226	GLN

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	B	303	-	4,4,4	0.25	0	6,6,6	0.18	0
2	SO4	B	301	-	4,4,4	0.18	0	6,6,6	0.31	0
2	SO4	B	302	-	4,4,4	0.14	0	6,6,6	0.26	0
2	SO4	A	301	-	4,4,4	0.23	0	6,6,6	0.23	0

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](#)

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	271/282 (96%)	-0.35	1 (0%) 92 91	11, 22, 38, 45	1 (0%)
1	B	271/282 (96%)	-0.10	3 (1%) 80 79	13, 26, 40, 46	1 (0%)
All	All	542/564 (96%)	-0.23	4 (0%) 87 86	11, 23, 38, 46	2 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	236	VAL	7.2
1	A	236	VAL	3.7
1	B	9	GLU	2.4
1	B	30	TRP	2.4

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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	SO4	B	303	5/5	0.71	0.29	53,53,54,54	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SO4	A	301	5/5	0.90	0.16	53,54,54,55	0
2	SO4	B	301	5/5	0.98	0.13	34,35,35,36	0
2	SO4	B	302	5/5	0.99	0.10	25,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.