wwPDB X-ray Structure Validation Summary Report (i)

Sep 17, 2023 – 05:28 PM EDT

PDB ID	:	4Z6M
Title	:	Structure of H200Q variant of Homoprotocatechuate 2,3-Dioxygenase from B.fuscum at 1.35 Ang resolution
Authors	:	Kovaleva, E.G.; Lipscomb, J.D.
Deposited on	:	2015-04-06
Resolution	:	1.35 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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:

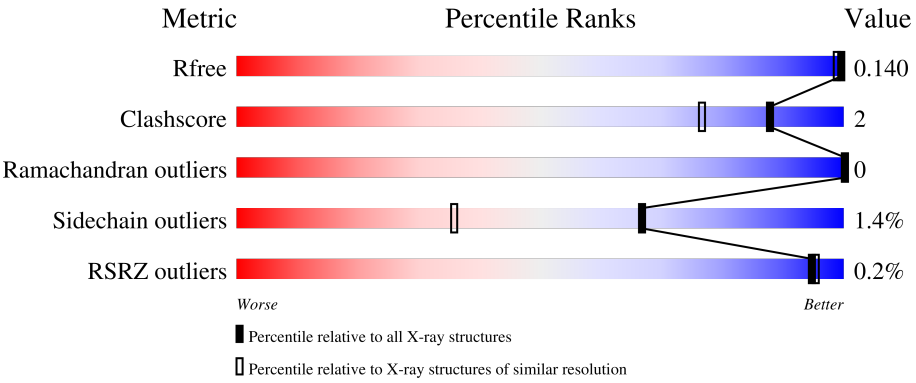
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 1.35 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	1509 (1.38-1.34)
Clashscore	141614	1551 (1.38-1.34)
Ramachandran outliers	138981	1530 (1.38-1.34)
Sidechain outliers	138945	1530 (1.38-1.34)
RSRZ outliers	127900	1487 (1.38-1.34)

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% 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	365	92% 6% ·
1	B	365	92% 7% ·
1	C	365	92% 7% ·
1	D	365	% 92% 6% ·

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 14005 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 Homoprotocatechuate 2,3-dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	S	0	12	0
			2993	1880	531	575	7			
1	B	359	Total	C	N	O	S	0	16	0
			3028	1900	539	581	8			
1	C	359	Total	C	N	O	S	0	12	0
			2991	1881	529	574	7			
1	D	359	Total	C	N	O	S	0	15	0
			3016	1897	530	581	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	200	GLN	HIS	engineered mutation	UNP Q45135
B	200	GLN	HIS	engineered mutation	UNP Q45135
C	200	GLN	HIS	engineered mutation	UNP Q45135
D	200	GLN	HIS	engineered mutation	UNP Q45135

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Fe	0	0
			1	1		
2	B	1	Total	Fe	0	0
			1	1		
2	C	1	Total	Fe	0	0
			1	1		
2	D	1	Total	Fe	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	B	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ca 1 1	0	0

- Molecule 5 is water.

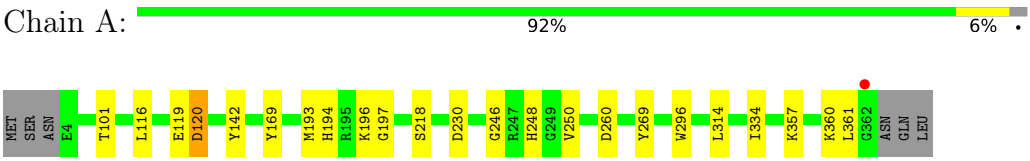
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	501	Total O 504 504	0	7
5	B	530	Total O 534 534	0	7
5	C	462	Total O 464 464	0	5
5	D	464	Total O 466 466	0	6

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 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.

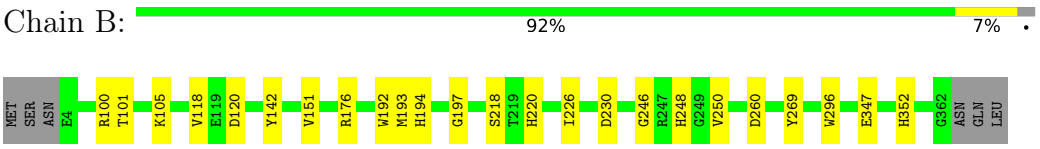
• Molecule 1: Homoprotocatechuate 2,3-dioxygenase

Chain A:



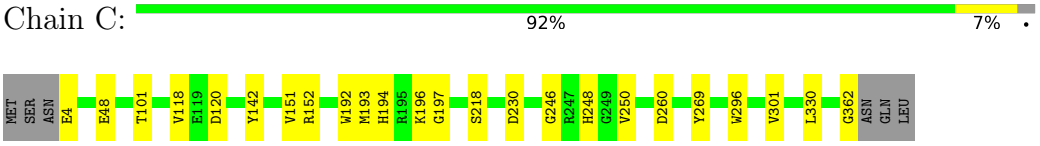
• Molecule 1: Homoprotocatechuate 2,3-dioxygenase

Chain B:



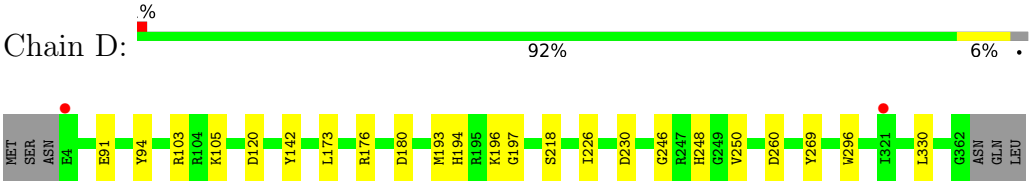
• Molecule 1: Homoprotocatechuate 2,3-dioxygenase

Chain C:



• Molecule 1: Homoprotocatechuate 2,3-dioxygenase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	110.48Å 152.09Å 96.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.09 – 1.35 48.09 – 1.35	Depositor EDS
% Data completeness (in resolution range)	95.9 (48.09-1.35) 95.9 (48.09-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.07 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.8.0069	Depositor
R, R_{free}	0.112 , 0.140 0.113 , 0.140	Depositor DCC
R_{free} test set	16910 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	9.7	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 44.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	14005	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FE2, CL, CA

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	0.47	0/3068	0.72	1/4162 (0.0%)
1	B	0.47	0/3103	0.75	1/4207 (0.0%)
1	C	0.47	0/3066	0.72	0/4161
1	D	0.43	0/3091	0.71	1/4193 (0.0%)
All	All	0.46	0/12328	0.73	3/16723 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	180	ASP	CB-CG-OD1	6.01	123.71	118.30
1	A	120	ASP	CB-CG-OD1	5.55	123.29	118.30
1	B	100	ARG	NE-CZ-NH1	5.00	122.80	120.30

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	2993	0	2833	13	0
1	B	3028	0	2868	13	0
1	C	2991	0	2838	15	0
1	D	3016	0	2854	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	1	0	0	0	0
5	A	504	0	0	0	0
5	B	534	0	0	1	0
5	C	464	0	0	0	0
5	D	466	0	0	0	0
All	All	14005	0	11393	46	0

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 2.

The worst 5 of 46 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:A:218[B]:SER:HA	1:A:269:TYR:O	2.08	0.54
1:C:48[A]:GLU:H	1:C:48[A]:GLU:CD	2.10	0.52
1:B:176:ARG:HD3	5:B:722:HOH:O	2.09	0.52
1:A:119:GLU:HG3	1:A:314:LEU:HG	1.91	0.52
1:B:218[A]:SER:HA	1:B:269:TYR:O	2.09	0.52

There are no symmetry-related clashes.

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	369/365 (101%)	359 (97%)	10 (3%)	0	100	100
1	B	373/365 (102%)	364 (98%)	9 (2%)	0	100	100
1	C	369/365 (101%)	361 (98%)	8 (2%)	0	100	100
1	D	372/365 (102%)	364 (98%)	8 (2%)	0	100	100
All	All	1483/1460 (102%)	1448 (98%)	35 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	318/313 (102%)	314 (99%)	4 (1%)	69	37
1	B	322/313 (103%)	318 (99%)	4 (1%)	71	42
1	C	318/313 (102%)	314 (99%)	4 (1%)	69	37
1	D	321/313 (103%)	316 (98%)	5 (2%)	62	30
All	All	1279/1252 (102%)	1262 (99%)	17 (1%)	67	37

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	176	ARG
1	D	260	ASP
1	B	260	ASP
1	C	4	GLU
1	C	120	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 9 ligands modelled in this entry, 9 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

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, 95^{th} 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	359/365 (98%)	-0.46	1 (0%) 94 95	6, 10, 20, 36	0
1	B	359/365 (98%)	-0.57	0 100 100	5, 9, 18, 31	0
1	C	359/365 (98%)	-0.35	0 100 100	6, 12, 22, 30	0
1	D	359/365 (98%)	-0.33	2 (0%) 89 91	6, 12, 25, 45	0
All	All	1436/1460 (98%)	-0.43	3 (0%) 95 95	5, 11, 22, 45	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	362	GLY	3.0
1	D	4	GLU	2.6
1	D	321	ILE	2.3

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 monosaccharides in this entry.

6.4 Ligands

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, 95^{th} 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	FE2	A	401	1/1	1.00	0.04	7,7,7,7	0
2	FE2	B	401	1/1	1.00	0.04	6,6,6,6	0
2	FE2	C	401	1/1	1.00	0.04	8,8,8,8	0
2	FE2	D	401	1/1	1.00	0.04	8,8,8,8	0
3	CL	A	402	1/1	1.00	0.04	10,10,10,10	0
3	CL	B	403	1/1	1.00	0.03	10,10,10,10	0
3	CL	C	402	1/1	1.00	0.03	13,13,13,13	0
3	CL	D	402	1/1	1.00	0.03	11,11,11,11	0
4	CA	B	402	1/1	1.00	0.03	14,14,14,14	1

6.5 Other polymers (i)

There are no such residues in this entry.