



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 18, 2018 – 04:08 AM EST

PDB ID : 5O7O
Title : The crystal structure of DfoC, the desferrioxamine biosynthetic pathway acetyltransferase/Non-Ribosomal Peptide Synthetase (NRPS)-Independent Siderophore (NIS) from the fire blight disease pathogen *Erwinia amylovora*
Authors : Salomone-Stagni, M.; Bartho, J.D.; Polsinelli, I.; Bellini, D.; Walsh, M.A.; Demitri, N.; Benini, S.
Deposited on : 2017-06-09
Resolution : 2.11 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

<http://wwpdb.org/validation/2016/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
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030736

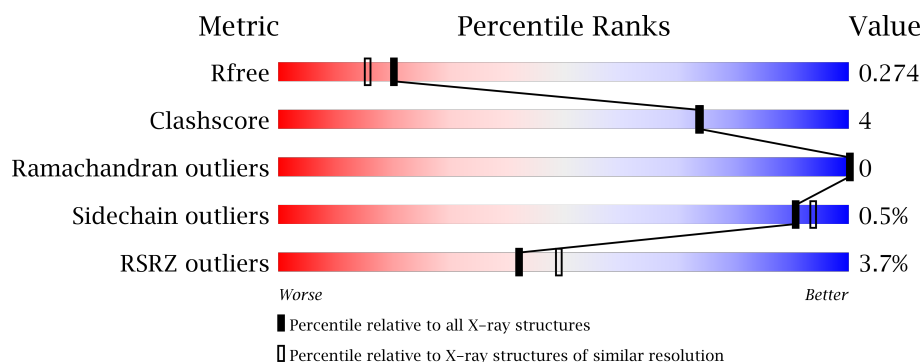
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.11 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}	100719	4988 (2.14-2.10)
Clashscore	112137	5557 (2.14-2.10)
Ramachandran outliers	110173	5504 (2.14-2.10)
Sidechain outliers	110143	5505 (2.14-2.10)
RSRZ outliers	101464	5021 (2.14-2.10)

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 $\leq 5\%$. 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	787	<div> <div style="width: 16%; background-color: red;"></div> <div style="width: 16%; background-color: orange;"></div> <div style="width: 16%; background-color: yellow;"></div> <div style="width: 16%; background-color: green;"></div> <div style="width: 81%; background-color: grey;"></div> </div>
1	B	787	<div> <div style="width: 16%; background-color: red;"></div> <div style="width: 16%; background-color: orange;"></div> <div style="width: 16%; background-color: yellow;"></div> <div style="width: 16%; background-color: green;"></div> <div style="width: 82%; background-color: grey;"></div> </div>
1	C	787	<div> <div style="width: 3%; background-color: red;"></div> <div style="width: 65%; background-color: green;"></div> <div style="width: 7%; background-color: yellow;"></div> <div style="width: 27%; background-color: grey;"></div> </div>
1	D	787	<div> <div style="width: 3%; background-color: red;"></div> <div style="width: 66%; background-color: green;"></div> <div style="width: 6%; background-color: yellow;"></div> <div style="width: 27%; background-color: grey;"></div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12009 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 Desferrioxamine siderophore biosynthesis protein dfoC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	0	0
			1179	754	206	210	9			
1	B	145	Total	C	N	O	S	0	0	0
			1173	752	204	207	10			
1	C	575	Total	C	N	O	S	0	0	0
			4616	2950	803	840	23			
1	D	575	Total	C	N	O	S	0	0	0
			4616	2950	802	841	23			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP D4I247
A	-1	ALA	-	expression tag	UNP D4I247
A	0	MET	-	expression tag	UNP D4I247
A	1	SER	-	expression tag	UNP D4I247
B	-2	GLY	-	expression tag	UNP D4I247
B	-1	ALA	-	expression tag	UNP D4I247
B	0	MET	-	expression tag	UNP D4I247
B	1	SER	-	expression tag	UNP D4I247
C	-2	GLY	-	expression tag	UNP D4I247
C	-1	ALA	-	expression tag	UNP D4I247
C	0	MET	-	expression tag	UNP D4I247
C	1	SER	-	expression tag	UNP D4I247
D	-2	GLY	-	expression tag	UNP D4I247
D	-1	ALA	-	expression tag	UNP D4I247
D	0	MET	-	expression tag	UNP D4I247
D	1	SER	-	expression tag	UNP D4I247

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	30	Total 30	O 30	0	0
2	B	32	Total 32	O 32	0	0
2	C	195	Total 195	O 195	0	0
2	D	168	Total 168	O 168	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	75.17Å 156.35Å 93.81Å 90.00° 95.57° 90.00°	Depositor
Resolution (Å)	78.18 – 2.11 78.17 – 2.11	Depositor EDS
% Data completeness (in resolution range)	99.8 (78.18-2.11) 99.8 (78.17-2.11)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.221 , 0.264 0.235 , 0.274	Depositor DCC
R_{free} test set	6119 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	44.9	Xtriage
Anisotropy	0.106	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12009	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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

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	0.94	1/1211 (0.1%)	0.93	4/1638 (0.2%)
1	B	0.90	1/1205 (0.1%)	1.02	8/1629 (0.5%)
1	C	0.87	4/4736 (0.1%)	0.93	19/6440 (0.3%)
1	D	0.83	2/4736 (0.0%)	0.90	14/6440 (0.2%)
All	All	0.86	8/11888 (0.1%)	0.92	45/16147 (0.3%)

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	B	0	1
1	C	0	1
1	D	0	1
All	All	0	3

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	649	MET	C-N	11.38	1.60	1.34
1	C	268	SER	C-N	-11.35	1.07	1.34
1	B	163	ILE	C-N	-11.08	1.08	1.34
1	C	245	GLY	C-N	10.91	1.59	1.34
1	C	246	SER	C-N	-9.44	1.12	1.34

The worst 5 of 45 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	490	MET	CB-CA-C	-15.29	79.83	110.40
1	B	163	ILE	O-C-N	-13.47	101.15	122.70
1	B	10	GLY	C-N-CA	10.81	148.73	121.70
1	B	10	GLY	O-C-N	-9.72	107.15	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	GLY	C-N-CA	9.54	145.55	121.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	163	ILE	Mainchain
1	C	200	GLY	Peptide
1	D	650	LYS	Mainchain

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	1179	0	1142	18	0
1	B	1173	0	1139	14	0
1	C	4616	0	4499	36	0
1	D	4616	0	4499	30	0
2	A	30	0	0	2	0
2	B	32	0	0	0	0
2	C	195	0	0	4	0
2	D	168	0	0	2	0
All	All	12009	0	11279	87	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 4.

The worst 5 of 87 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:139:VAL:HG12	1:A:141:PRO:HD3	1.46	0.93
1:D:361:TYR:OH	1:D:367:GLU:OE2	1.91	0.88
1:A:94:ALA:O	1:A:96:GLN:NE2	2.17	0.77
1:A:140:GLU:OE1	1:A:165:MET:CE	2.35	0.75
1:A:89:ILE:HD12	1:A:140:GLU:HB3	1.73	0.69

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	142/787 (18%)	136 (96%)	6 (4%)	0	100	100
1	B	141/787 (18%)	136 (96%)	5 (4%)	0	100	100
1	C	571/787 (73%)	551 (96%)	20 (4%)	0	100	100
1	D	571/787 (73%)	552 (97%)	19 (3%)	0	100	100
All	All	1425/3148 (45%)	1375 (96%)	50 (4%)	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	126/670 (19%)	126 (100%)	0	100	100
1	B	125/670 (19%)	123 (98%)	2 (2%)	68	73
1	C	488/670 (73%)	487 (100%)	1 (0%)	94	96
1	D	488/670 (73%)	485 (99%)	3 (1%)	89	92
All	All	1227/2680 (46%)	1221 (100%)	6 (0%)	91	94

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	490	MET
1	D	351	ASN
1	D	204	THR
1	B	110	THR
1	D	341	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	522	GLN
1	C	624	HIS
1	D	631	HIS
1	C	541	HIS
1	C	582	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	3
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	269:LEU	C	270:THR	N	1.18
1	C	246:SER	C	247:GLU	N	1.12
1	B	163:ILE	C	164:ASP	N	1.08
1	C	268:SER	C	269:LEU	N	1.07

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	146/787 (18%)	0.30	7 (4%) 31 37	36, 51, 70, 95	0
1	B	145/787 (18%)	0.40	11 (7%) 15 19	37, 55, 74, 84	0
1	C	575/787 (73%)	0.20	9 (1%) 72 76	34, 49, 73, 104	0
1	D	575/787 (73%)	0.25	27 (4%) 32 38	33, 50, 75, 115	0
All	All	1441/3148 (45%)	0.25	54 (3%) 42 49	33, 50, 74, 115	0

The worst 5 of 54 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	499	TYR	7.8
1	C	494	ARG	5.5
1	D	236	HIS	5.3
1	D	495	GLY	5.1
1	A	110	THR	5.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.