



wwPDB X-ray Structure Validation Summary Report ⓘ

May 25, 2020 – 06:18 am BST

PDB ID : 4XGV
Title : Crystal structure of Escherichia coli Flavin trafficking protein, an FMN transferase
Authors : Tomchick, D.R.; Brautigam, C.A.; Deka, R.K.; Norgard, M.V.
Deposited on : 2015-01-02
Resolution : 1.88 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 ⓘ 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

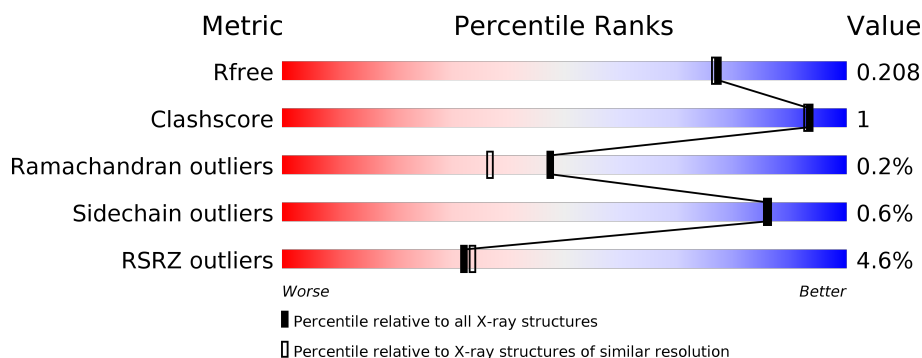
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	9470 (1.90-1.86)
Clashscore	141614	10282 (1.90-1.86)
Ramachandran outliers	138981	10152 (1.90-1.86)
Sidechain outliers	138945	10152 (1.90-1.86)
RSRZ outliers	127900	9303 (1.90-1.86)

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	340	<div> <div>4%</div> <div> <div></div> <div>86%</div> <div>9%</div> </div> </div>
1	B	340	<div> <div>5%</div> <div> <div></div> <div>87%</div> <div>9%</div> </div> </div>
1	C	340	<div> <div>4%</div> <div> <div></div> <div>88%</div> <div>9%</div> </div> </div>
1	D	340	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>10%</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 19989 atoms, of which 9603 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 FAD:protein FMN transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	309	Total	C	H	N	O	S	0	3	0
			4793	1511	2393	414	465	10			
1	B	309	Total	C	H	N	O	S	0	4	0
			4787	1510	2390	409	468	10			
1	C	309	Total	C	H	N	O	S	0	1	0
			4776	1506	2385	410	465	10			
1	D	305	Total	C	H	N	O	S	0	2	0
			4714	1485	2354	406	459	10			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP P0AB85
A	333	LEU	-	expression tag	UNP P0AB85
A	334	GLU	-	expression tag	UNP P0AB85
A	335	HIS	-	expression tag	UNP P0AB85
A	336	HIS	-	expression tag	UNP P0AB85
A	337	HIS	-	expression tag	UNP P0AB85
A	338	HIS	-	expression tag	UNP P0AB85
A	339	HIS	-	expression tag	UNP P0AB85
A	340	HIS	-	expression tag	UNP P0AB85
B	1	MET	-	initiating methionine	UNP P0AB85
B	333	LEU	-	expression tag	UNP P0AB85
B	334	GLU	-	expression tag	UNP P0AB85
B	335	HIS	-	expression tag	UNP P0AB85
B	336	HIS	-	expression tag	UNP P0AB85
B	337	HIS	-	expression tag	UNP P0AB85
B	338	HIS	-	expression tag	UNP P0AB85
B	339	HIS	-	expression tag	UNP P0AB85
B	340	HIS	-	expression tag	UNP P0AB85
C	1	MET	-	initiating methionine	UNP P0AB85
C	333	LEU	-	expression tag	UNP P0AB85
C	334	GLU	-	expression tag	UNP P0AB85

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Chain	Residue	Modelled	Actual	Comment	Reference
C	335	HIS	-	expression tag	UNP P0AB85
C	336	HIS	-	expression tag	UNP P0AB85
C	337	HIS	-	expression tag	UNP P0AB85
C	338	HIS	-	expression tag	UNP P0AB85
C	339	HIS	-	expression tag	UNP P0AB85
C	340	HIS	-	expression tag	UNP P0AB85
D	1	MET	-	initiating methionine	UNP P0AB85
D	333	LEU	-	expression tag	UNP P0AB85
D	334	GLU	-	expression tag	UNP P0AB85
D	335	HIS	-	expression tag	UNP P0AB85
D	336	HIS	-	expression tag	UNP P0AB85
D	337	HIS	-	expression tag	UNP P0AB85
D	338	HIS	-	expression tag	UNP P0AB85
D	339	HIS	-	expression tag	UNP P0AB85
D	340	HIS	-	expression tag	UNP P0AB85

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0
3	C	1	Total Na 1 1	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	B	1	Total	C	H	O	0	0
			10	2	6	2		
4	B	1	Total	C	H	O	0	0
			10	2	6	2		
4	B	1	Total	C	H	O	0	0
			10	2	6	2		
4	B	1	Total	C	H	O	0	0
			10	2	6	2		
4	C	1	Total	C	H	O	0	0
			10	2	6	2		
4	D	1	Total	C	H	O	0	0
			10	2	6	2		
4	D	1	Total	C	H	O	0	0
			10	2	6	2		
4	D	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	179	Total	O	0	0
			179	179		
6	B	194	Total	O	0	0
			194	194		
6	C	192	Total	O	0	0
			192	192		
6	D	211	Total	O	0	0
			211	211		

His	His
S329	
GLU	
LYS	
ASN	
LEU	
GLU	
HIS	
HIS	
HIS	
HIS	
HIS	

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.20Å 70.64Å 85.96Å 75.70° 72.45° 69.28°	Depositor
Resolution (Å)	31.51 – 1.88 31.51 – 1.88	Depositor EDS
% Data completeness (in resolution range)	96.4 (31.51-1.88) 96.4 (31.51-1.88)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.88Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.163 , 0.201 0.176 , 0.208	Depositor DCC
R_{free} test set	4544 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 49.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for h,h-k,h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19989	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, CA, EDO, ACT

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	0.30	0/2457	0.50	0/3330
1	B	0.29	0/2459	0.49	0/3336
1	C	0.30	0/2439	0.49	0/3306
1	D	0.32	0/2412	0.51	0/3270
All	All	0.30	0/9767	0.50	0/13242

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 [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	2400	2393	2386	7	0
1	B	2397	2390	2381	8	0
1	C	2391	2385	2382	5	0
1	D	2360	2354	2349	7	0
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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	1	0	0	0	0
4	A	16	24	24	1	0
4	B	20	30	30	0	0
4	C	4	6	6	0	0
4	D	12	18	18	1	0
5	A	4	3	3	0	0
6	A	179	0	0	3	0
6	B	194	0	0	2	0
6	C	192	0	0	1	0
6	D	211	0	0	2	0
All	All	10386	9603	9579	27	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 1.

The worst 5 of 27 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:222:VAL:N	6:A:604:HOH:O	2.33	0.60
1:A:148:SER:O	6:A:601:HOH:O	2.16	0.59
1:A:247:GLY:N	6:A:606:HOH:O	2.37	0.56
1:C:28:ARG:NH1	6:C:601:HOH:O	2.39	0.56
1:B:134:LYS:NZ	6:B:607:HOH:O	2.41	0.54

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	306/340 (90%)	299 (98%)	5 (2%)	2 (1%)	22	11
1	B	309/340 (91%)	304 (98%)	5 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	304/340 (89%)	299 (98%)	5 (2%)	0	100	100
1	D	301/340 (88%)	296 (98%)	5 (2%)	0	100	100
All	All	1220/1360 (90%)	1198 (98%)	20 (2%)	2 (0%)	47	37

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	312	GLY

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	262/288 (91%)	258 (98%)	4 (2%)	65	59
1	B	264/288 (92%)	264 (100%)	0	100	100
1	C	262/288 (91%)	261 (100%)	1 (0%)	91	90
1	D	259/288 (90%)	258 (100%)	1 (0%)	91	90
All	All	1047/1152 (91%)	1041 (99%)	6 (1%)	86	86

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

Mol	Chain	Res	Type
1	A	149	HIS
1	D	242	TYR
1	A	241	ASN
1	A	72	SER
1	C	252	HIS

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

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

Of 20 ligands modelled in this entry, 6 are monoatomic - leaving 14 for Mogul analysis.

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	$\# Z > 2$
4	EDO	B	505	-	3,3,3	0.43	0	2,2,2	0.32	0
4	EDO	B	504	-	3,3,3	0.49	0	2,2,2	0.21	0
4	EDO	C	503	-	3,3,3	0.46	0	2,2,2	0.34	0
4	EDO	A	504	-	3,3,3	0.48	0	2,2,2	0.36	0
4	EDO	D	504	-	3,3,3	0.40	0	2,2,2	0.25	0
4	EDO	D	503	-	3,3,3	0.40	0	2,2,2	0.39	0
4	EDO	A	505	-	3,3,3	0.52	0	2,2,2	0.23	0
4	EDO	B	503	-	3,3,3	0.45	0	2,2,2	0.41	0
4	EDO	D	502	-	3,3,3	0.47	0	2,2,2	0.34	0
4	EDO	A	506	-	3,3,3	0.44	0	2,2,2	0.35	0
4	EDO	B	506	-	3,3,3	0.41	0	2,2,2	0.44	0
5	ACT	A	507	-	1,3,3	1.42	0	0,3,3	0.00	-
4	EDO	B	502	-	3,3,3	0.44	0	2,2,2	0.29	0
4	EDO	A	503	-	3,3,3	0.53	0	2,2,2	0.11	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	B	505	-	-	1/1/1/1	-
4	EDO	B	504	-	-	0/1/1/1	-
4	EDO	C	503	-	-	0/1/1/1	-
4	EDO	A	504	-	-	0/1/1/1	-
4	EDO	D	504	-	-	0/1/1/1	-
4	EDO	D	503	-	-	0/1/1/1	-
4	EDO	A	505	-	-	0/1/1/1	-
4	EDO	B	503	-	-	0/1/1/1	-
4	EDO	D	502	-	-	0/1/1/1	-
4	EDO	A	506	-	-	1/1/1/1	-
4	EDO	B	506	-	-	0/1/1/1	-
4	EDO	B	502	-	-	1/1/1/1	-
4	EDO	A	503	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	502	EDO	O1-C1-C2-O2
4	B	505	EDO	O1-C1-C2-O2
4	A	503	EDO	O1-C1-C2-O2
4	A	506	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	504	EDO	1	0
4	A	503	EDO	1	0

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 [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	309/340 (90%)	-0.10	13 (4%) 36 37	14, 25, 63, 97	0
1	B	309/340 (90%)	-0.07	16 (5%) 27 28	15, 25, 63, 86	0
1	C	309/340 (90%)	-0.06	13 (4%) 36 37	15, 27, 59, 96	0
1	D	305/340 (89%)	-0.20	15 (4%) 29 31	13, 22, 63, 101	0
All	All	1232/1360 (90%)	-0.11	57 (4%) 32 34	13, 25, 63, 101	0

The worst 5 of 57 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	222	VAL	8.6
1	D	13	THR	6.0
1	C	60	TYR	5.4
1	A	120	GLN	5.0
1	A	215	PRO	4.8

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(\AA^2)	Q<0.9
4	EDO	D	504	4/4	0.76	0.22	49,68,76,82	0
4	EDO	B	504	4/4	0.81	0.18	35,42,50,50	0
4	EDO	A	506	4/4	0.83	0.20	42,50,56,58	0
4	EDO	A	503	4/4	0.84	0.15	33,44,54,65	0
5	ACT	A	507	4/4	0.86	0.19	39,46,47,50	0
4	EDO	B	505	4/4	0.88	0.12	45,54,58,61	0
4	EDO	B	506	4/4	0.89	0.13	28,38,45,47	0
4	EDO	D	503	4/4	0.92	0.18	33,39,47,48	0
4	EDO	A	504	4/4	0.92	0.13	32,38,40,49	0
3	NA	C	502	1/1	0.94	0.15	31,31,31,31	0
3	NA	A	502	1/1	0.94	0.21	36,36,36,36	0
4	EDO	C	503	4/4	0.94	0.10	19,30,39,46	0
4	EDO	A	505	4/4	0.94	0.09	32,39,40,40	0
4	EDO	D	502	4/4	0.94	0.10	24,29,30,30	0
4	EDO	B	502	4/4	0.96	0.09	21,30,46,55	0
4	EDO	B	503	4/4	0.96	0.12	20,32,44,53	0
2	CA	A	501	1/1	0.99	0.04	20,20,20,20	0
2	CA	C	501	1/1	0.99	0.03	26,26,26,26	0
2	CA	B	501	1/1	0.99	0.02	22,22,22,22	0
2	CA	D	501	1/1	1.00	0.05	20,20,20,20	0

6.5 Other polymers [i](#)

There are no such residues in this entry.