



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 05:12 pm BST

PDB ID : 3QFI  
Title : X-ray crystal structure of transcriptional regulator (EF0465) from *Enterococcus faecalis*, Northeast Structural Genomics Consortium Target EfR190  
Authors : Forouhar, F.; Neely, H.; Seetharaman, J.; Wang, H.; Ciccocanti, C.; Mao, L.; Acton, T.B.; Xiao, R.; Everett, J.K.; Montelione, G.T.; Tong, L.; Hun, J.F.; Northeast Structural Genomics Consortium (NESG)  
Deposited on : 2011-01-21  
Resolution : 2.71 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](mailto: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.

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

MolProbity	:	4.02b-467
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

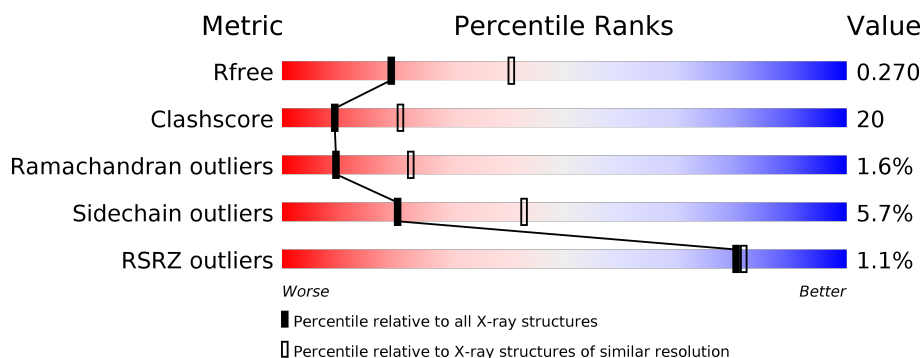
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.71 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	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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  $\geq 3$ , 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	301	<div> <div></div> <div> <div></div> <div>54%</div> <div>30%</div> <div>•</div> <div>13%</div> </div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2095 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 Transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	261	Total	C	N	O	S	0	0	0
			2057	1288	345	415	9			

There are 9 discrepancies between the modelled and reference sequences:

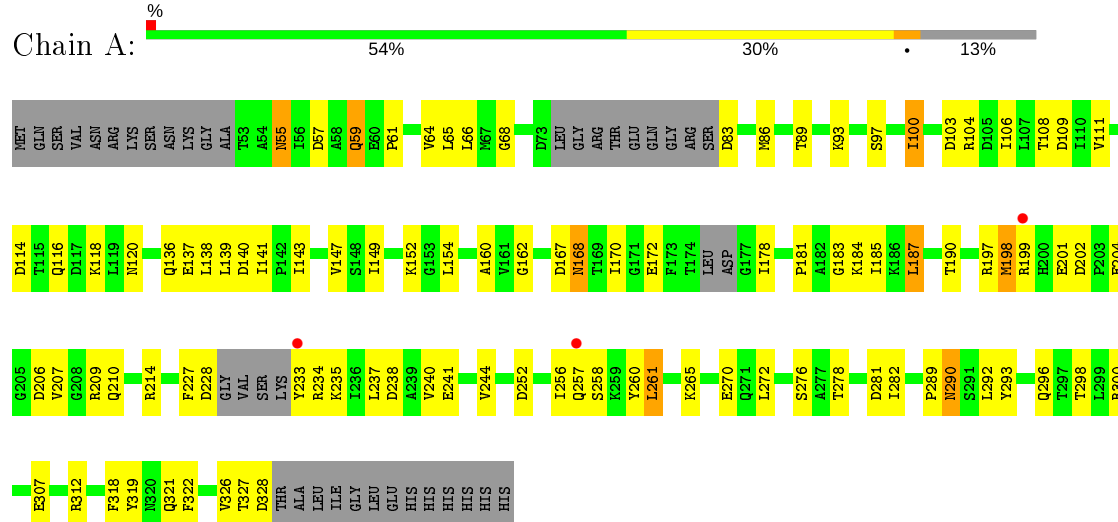
Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	-	INITIATING METHIONINE	UNP Q838I2
A	334	LEU	-	EXPRESSION TAG	UNP Q838I2
A	335	GLU	-	EXPRESSION TAG	UNP Q838I2
A	336	HIS	-	EXPRESSION TAG	UNP Q838I2
A	337	HIS	-	EXPRESSION TAG	UNP Q838I2
A	338	HIS	-	EXPRESSION TAG	UNP Q838I2
A	339	HIS	-	EXPRESSION TAG	UNP Q838I2
A	340	HIS	-	EXPRESSION TAG	UNP Q838I2
A	341	HIS	-	EXPRESSION TAG	UNP Q838I2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	38	Total	O	0	0
			38	38		



- Molecule 1: Transcriptional regulator



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.21Å 54.54Å 99.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.87 – 2.71 29.01 – 2.71	Depositor EDS
% Data completeness (in resolution range)	82.0 (19.87-2.71) 85.3 (29.01-2.71)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.95 (at 2.72Å)	Xtriage
Refinement program	CNS 1.2 & Xtalview	Depositor
R, $R_{free}$	0.201 , 0.264 0.207 , 0.270	Depositor DCC
$R_{free}$ test set	688 reflections (10.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.2	Xtriage
Anisotropy	0.808	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 51.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	2095	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.42	0/2084	0.59	1/2814 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	290	ASN	CB-CA-C	6.06	122.52	110.40

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	2057	0	2028	81	0
2	A	38	0	0	6	0
All	All	2095	0	2028	81	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 20.

All (81) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance ( $\text{\AA}$ )	Clash overlap ( $\text{\AA}$ )
1:A:136:GLN:HE21	1:A:143:ILE:H	1.17	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:THR:HG21	1:A:138:LEU:HD13	1.56	0.85
1:A:252:ASP:O	1:A:256:ILE:HG12	1.77	0.84
1:A:55:ASN:HD21	1:A:57:ASP:HB3	1.50	0.76
1:A:326:VAL:HG12	1:A:328:ASP:H	1.60	0.67
1:A:228:ASP:HB3	1:A:233:TYR:HB2	1.76	0.66
1:A:228:ASP:HB2	1:A:234:ARG:NE	2.10	0.65
1:A:59:GLN:HG3	1:A:59:GLN:O	1.95	0.64
1:A:100:ILE:HG22	1:A:100:ILE:O	1.97	0.64
1:A:59:GLN:HA	1:A:260:TYR:OH	1.98	0.62
1:A:204:GLU:HB3	1:A:207:VAL:HG23	1.81	0.62
1:A:108:THR:CG2	1:A:138:LEU:HD13	2.30	0.61
1:A:104:ARG:NH2	1:A:199:ARG:HH22	1.99	0.61
1:A:199:ARG:HH21	1:A:209:ARG:CZ	2.14	0.60
1:A:293:TYR:OH	1:A:307:GLU:HG2	2.03	0.59
1:A:55:ASN:C	1:A:55:ASN:HD22	2.06	0.59
1:A:136:GLN:NE2	1:A:143:ILE:H	1.95	0.58
1:A:136:GLN:HE21	1:A:143:ILE:N	1.96	0.57
1:A:104:ARG:HB2	1:A:120:ASN:HB2	1.85	0.57
1:A:65:LEU:HD12	1:A:86:MET:O	2.05	0.57
1:A:228:ASP:HB2	1:A:234:ARG:CD	2.35	0.56
1:A:59:GLN:HA	1:A:260:TYR:CZ	2.42	0.55
1:A:234:ARG:H	1:A:234:ARG:CD	2.19	0.55
1:A:100:ILE:CG2	1:A:100:ILE:O	2.54	0.55
1:A:61:PRO:HD3	1:A:93:LYS:HE3	1.88	0.55
1:A:149:ILE:HG23	1:A:244:VAL:HG22	1.91	0.53
1:A:272:LEU:N	2:A:512:HOH:O	2.35	0.53
1:A:327:THR:HG22	1:A:327:THR:O	2.09	0.53
1:A:233:TYR:O	1:A:237:LEU:HD23	2.09	0.52
1:A:178:ILE:HD13	1:A:190:THR:HG22	1.90	0.52
1:A:234:ARG:HB2	2:A:521:HOH:O	2.09	0.52
1:A:197:ARG:HH11	1:A:197:ARG:HG2	1.75	0.52
1:A:89:THR:O	1:A:97:SER:HA	2.09	0.52
1:A:270:GLU:OE1	1:A:298:THR:HG21	2.10	0.52
1:A:66:LEU:HD22	1:A:147:VAL:HG11	1.92	0.51
1:A:238:ASP:C	1:A:240:VAL:H	2.14	0.50
1:A:168:ASN:ND2	1:A:170:ILE:H	2.09	0.50
1:A:183:GLY:O	1:A:185:ILE:HG13	2.12	0.50
1:A:103:ASP:HB3	1:A:106:ILE:HG13	1.95	0.49
1:A:278:THR:HG23	1:A:282:ILE:O	2.13	0.49
1:A:136:GLN:HG2	1:A:143:ILE:HG12	1.96	0.48
1:A:206:ASP:HB3	1:A:210:GLN:HE21	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:ASP:H	1:A:234:ARG:NH1	2.11	0.48
1:A:139:LEU:HB2	1:A:141:ILE:HG12	1.96	0.48
1:A:68:GLY:O	1:A:83:ASP:HB3	2.14	0.47
1:A:106:ILE:O	1:A:118:LYS:HA	2.14	0.47
1:A:167:ASP:OD2	1:A:184:LYS:HE3	2.15	0.46
1:A:181:PRO:HG2	1:A:185:ILE:HG12	1.98	0.45
1:A:109:ASP:OD1	1:A:116:GLN:HG3	2.16	0.45
1:A:281:ASP:HB3	2:A:524:HOH:O	2.16	0.45
1:A:152:LYS:HD3	2:A:535:HOH:O	2.16	0.45
1:A:234:ARG:H	1:A:234:ARG:HD2	1.82	0.45
1:A:318:PHE:O	1:A:321:GLN:HG2	2.16	0.45
1:A:140:ASP:O	1:A:300:ARG:NH1	2.49	0.45
1:A:154:LEU:HD12	1:A:154:LEU:O	2.16	0.45
1:A:111:VAL:HG21	1:A:137:GLU:HG3	1.99	0.45
1:A:240:VAL:HG22	1:A:244:VAL:HG23	1.98	0.45
1:A:66:LEU:HA	1:A:147:VAL:HG13	1.99	0.44
1:A:162:GLY:N	2:A:518:HOH:O	2.50	0.44
1:A:238:ASP:HA	1:A:241:GLU:HG2	2.00	0.44
1:A:238:ASP:C	1:A:240:VAL:N	2.72	0.44
1:A:289:PRO:HD3	1:A:319:TYR:CD2	2.53	0.44
1:A:235:LYS:O	1:A:235:LYS:HD3	2.18	0.43
1:A:265:LYS:NZ	2:A:528:HOH:O	2.51	0.43
1:A:197:ARG:O	1:A:198:MET:HB2	2.18	0.43
1:A:296:GLN:O	1:A:300:ARG:HG3	2.18	0.43
1:A:234:ARG:HG2	1:A:257:GLN:OE1	2.19	0.42
1:A:66:LEU:HD22	1:A:147:VAL:CG1	2.49	0.42
1:A:202:ASP:C	1:A:204:GLU:H	2.23	0.42
1:A:321:GLN:HG3	1:A:322:PHE:CD2	2.54	0.42
1:A:234:ARG:HD2	1:A:234:ARG:N	2.34	0.41
1:A:237:LEU:O	1:A:240:VAL:HG12	2.20	0.41
1:A:199:ARG:HH21	1:A:209:ARG:NE	2.18	0.41
1:A:261:LEU:HD12	1:A:261:LEU:H	1.86	0.41
1:A:278:THR:HA	1:A:282:ILE:O	2.21	0.41
1:A:289:PRO:HD3	1:A:319:TYR:CG	2.56	0.41
1:A:187:LEU:HA	1:A:187:LEU:HD12	1.89	0.41
1:A:293:TYR:OH	1:A:307:GLU:CG	2.68	0.41
1:A:210:GLN:O	1:A:214:ARG:HG3	2.20	0.41
1:A:160:ALA:CB	1:A:237:LEU:HD13	2.50	0.41
1:A:204:GLU:O	1:A:207:VAL:HG22	2.21	0.40

There are no symmetry-related clashes.



## 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	253/301 (84%)	226 (89%)	23 (9%)	4 (2%)	9	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	MET
1	A	227	PHE
1	A	258	SER
1	A	172	GLU

### 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	227/260 (87%)	214 (94%)	13 (6%)	20	43

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	59	GLN
1	A	64	VAL
1	A	100	ILE
1	A	114	ASP
1	A	168	ASN
1	A	187	LEU

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Mol	Chain	Res	Type
1	A	201	GLU
1	A	261	LEU
1	A	276	SER
1	A	290	ASN
1	A	292	LEU
1	A	312	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	ASN
1	A	59	GLN
1	A	91	ASN
1	A	120	ASN
1	A	133	ASN
1	A	136	GLN
1	A	150	ASN
1	A	168	ASN
1	A	210	GLN
1	A	243	ASN
1	A	285	GLN
1	A	296	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](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	261/301 (86%)	-0.11	3 (1%)	80 82	2, 23, 49, 59	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	257	GLN	2.4
1	A	199	ARG	2.4
1	A	233	TYR	2.1

### 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.