



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 10:28 AM UTC

PDB ID : 2QR4 / pdb_00002qr4
Title : Crystal structure of oligoendopeptidase-F from *Enterococcus faecium*
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Deposited on : 2007-07-27
Resolution : 2.50 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

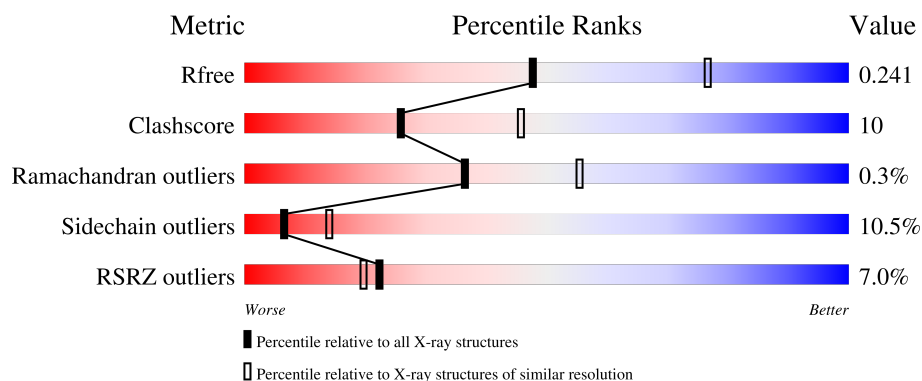
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	587	3% 68% 18% • 11%
1	B	587	9% 65% 17% • 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8549 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 Peptidase M3B, oligoendopeptidase F.

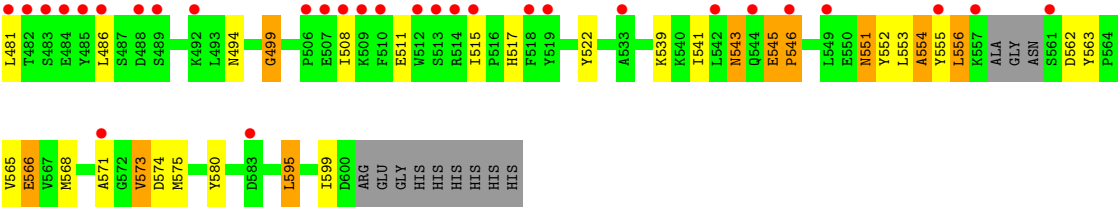
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	521	Total	C	N	O	Se	0	4	0
			4265	2723	700	832	10			
1	B	506	Total	C	N	O	Se	0	2	0
			4096	2612	675	800	9			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MSE	-	expression tag	UNP Q3XYC8
A	24	SER	-	expression tag	UNP Q3XYC8
A	25	LEU	-	expression tag	UNP Q3XYC8
A	602	GLU	-	expression tag	UNP Q3XYC8
A	603	GLY	-	expression tag	UNP Q3XYC8
A	604	HIS	-	expression tag	UNP Q3XYC8
A	605	HIS	-	expression tag	UNP Q3XYC8
A	606	HIS	-	expression tag	UNP Q3XYC8
A	607	HIS	-	expression tag	UNP Q3XYC8
A	608	HIS	-	expression tag	UNP Q3XYC8
A	609	HIS	-	expression tag	UNP Q3XYC8
B	23	MSE	-	expression tag	UNP Q3XYC8
B	24	SER	-	expression tag	UNP Q3XYC8
B	25	LEU	-	expression tag	UNP Q3XYC8
B	602	GLU	-	expression tag	UNP Q3XYC8
B	603	GLY	-	expression tag	UNP Q3XYC8
B	604	HIS	-	expression tag	UNP Q3XYC8
B	605	HIS	-	expression tag	UNP Q3XYC8
B	606	HIS	-	expression tag	UNP Q3XYC8
B	607	HIS	-	expression tag	UNP Q3XYC8
B	608	HIS	-	expression tag	UNP Q3XYC8
B	609	HIS	-	expression tag	UNP Q3XYC8

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	99	Total 99	O 99	0	0
2	B	89	Total 89	O 89	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	133.14Å 133.14Å 171.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.04 – 2.50 39.04 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.04-2.50) 99.9 (39.04-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.51Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.188 , 0.244 0.186 , 0.241	Depositor DCC
R_{free} test set	2726 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 50.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8549	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.86% 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](#)

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	0.53	0/4351	0.84	5/5872 (0.1%)
1	B	0.67	12/4176 (0.3%)	0.87	9/5632 (0.2%)
All	All	0.60	12/8527 (0.1%)	0.86	14/11504 (0.1%)

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	436	THR	CB-OG1	-9.60	1.28	1.43
1	B	438	LYS	C-O	9.29	1.35	1.24
1	B	437	GLU	CD-OE1	8.12	1.40	1.25
1	B	437	GLU	CD-OE2	7.69	1.40	1.25
1	B	438	LYS	CA-C	7.45	1.62	1.52

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	499	GLY	CA-C-N	9.60	130.26	119.32
1	B	499	GLY	C-N-CA	9.60	130.26	119.32
1	B	436	THR	CA-CB-OG1	7.47	120.80	109.60
1	A	478	GLY	CA-C-N	-7.19	117.41	122.59
1	A	478	GLY	C-N-CA	-7.19	117.41	122.59

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	4265	0	4074	72	0
1	B	4096	0	3863	87	0
2	A	99	0	0	5	0
2	B	89	0	0	3	0
All	All	8549	0	7937	155	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 10.

The worst 5 of 155 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:228:GLN:NE2	1:B:184:VAL:HG12	1.62	1.11
1:B:438:LYS:H	1:B:438:LYS:HD2	1.35	0.90
1:A:533:ALA:HB2	1:A:575:MSE:HE3	1.53	0.90
1:A:228:GLN:HE22	1:B:184:VAL:HG12	1.37	0.90
1:B:463:GLN:HE22	1:B:494:ASN:HD22	1.20	0.86

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	517/587 (88%)	505 (98%)	11 (2%)	1 (0%)	43	63
1	B	498/587 (85%)	476 (96%)	20 (4%)	2 (0%)	30	49
All	All	1015/1174 (86%)	981 (97%)	31 (3%)	3 (0%)	36	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	313	GLU

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Mol	Chain	Res	Type
1	B	546	PRO
1	B	339	MSE

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	454/504 (90%)	408 (90%)	46 (10%)	7	15
1	B	430/504 (85%)	383 (89%)	47 (11%)	6	13
All	All	884/1008 (88%)	791 (90%)	93 (10%)	6	14

5 of 93 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	110	VAL
1	B	311	LEU
1	B	146	VAL
1	B	187	THR
1	B	386	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	472	HIS
1	B	521	ASN
1	B	140	HIS
1	B	448	ASN
1	B	53	GLN

5.3.3 RNA ⓘ

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 oligosaccharides 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](#)

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	511/587 (87%)	0.29	15 (2%) 53 49	17, 45, 68, 123	4 (0%)
1	B	496/587 (84%)	0.68	55 (11%) 10 8	20, 45, 66, 91	2 (0%)
All	All	1007/1174 (85%)	0.48	70 (6%) 22 20	17, 45, 68, 123	6 (0%)

The worst 5 of 70 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	415	ILE	6.0
1	B	400	TYR	5.1
1	B	312	GLY	5.1
1	A	316	ILE	4.7
1	B	519	TYR	4.3

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 oligosaccharides 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.