



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 12:05 AM UTC

PDB ID : 3B61 / pdb_00003b61
Title : EmrE multidrug transporter, apo crystal form
Authors : Chang, G.; Chen, Y.J.
Deposited on : 2007-10-26
Resolution : 4.50 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
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

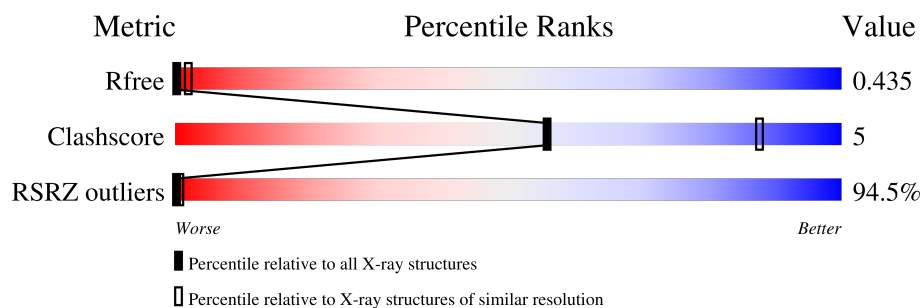
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.50 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	1127 (5.10-3.90)
Clashscore	190562	1002 (5.06-3.94)
RSRZ outliers	180081	1122 (5.10-3.90)

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	110	90% 97% .
1	B	110	92% 95% ..
1	C	110	95% 97% .
1	D	110	93% 95% ..
1	E	110	94% 97% .
1	F	110	87% 95% ..
1	G	110	95% 97% .

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Mol	Chain	Length	Quality of chain
1	H	110	91% 95%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 856 atoms, of which 0 are hydrogens and 0 are deuteriums.

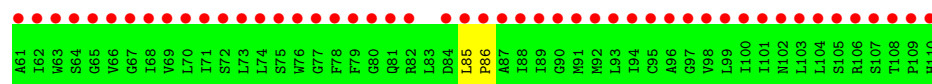
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Multidrug transporter emrE.

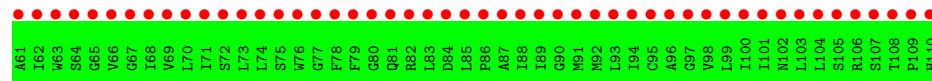
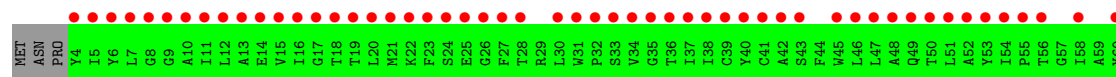
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
1	A	107	Total C 107 107	0	0	107
1	B	107	Total C 107 107	0	0	107
1	C	107	Total C 107 107	0	0	107
1	D	107	Total C 107 107	0	0	107
1	E	107	Total C 107 107	0	0	107
1	F	107	Total C 107 107	0	0	107
1	G	107	Total C 107 107	0	0	107
1	H	107	Total C 107 107	0	0	107

- Molecule 1: Multidrug transporter emrE

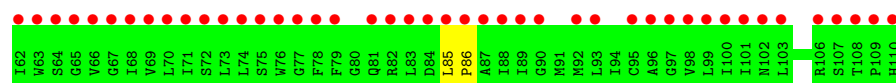
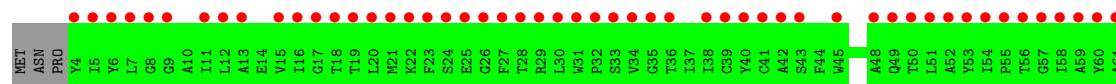
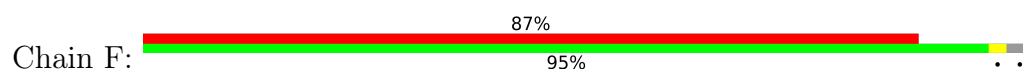




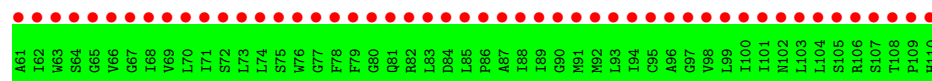
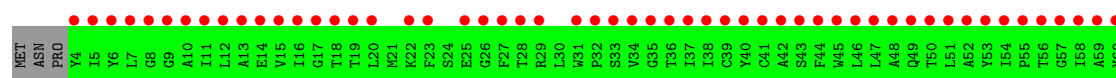
• Molecule 1: Multidrug transporter emrE



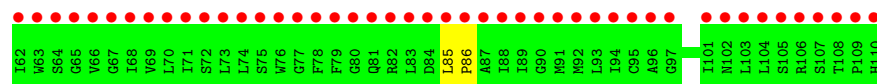
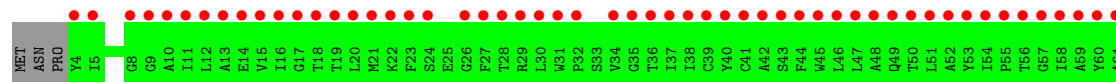
• Molecule 1: Multidrug transporter emrE



• Molecule 1: Multidrug transporter emrE



• Molecule 1: Multidrug transporter emrE



4 Data and refinement statistics

Property	Value	Source
Space group	F 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	181.00Å 239.20Å 284.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 4.50 19.99 – 4.50	Depositor EDS
% Data completeness (in resolution range)	75.8 (19.99-4.50) 74.8 (19.99-4.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 4.54Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.318 , 0.362 0.410 , 0.435	Depositor DCC
R_{free} test set	1369 reflections (9.89%)	wwPDB-VP
Wilson B-factor (Å ²)	187.7	Xtriage
Anisotropy	0.196	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	85.45 , 999.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.69	EDS
Total number of atoms	856	wwPDB-VP
Average B, all atoms (Å ²)	229.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.05% 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).

There are no protein, RNA or DNA chains available to summarize Z scores of covalent bonds and angles.

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	107	0	0	0	0
1	B	107	0	0	1	0
1	C	107	0	0	0	0
1	D	107	0	0	1	0
1	E	107	0	0	0	0
1	F	107	0	0	1	0
1	G	107	0	0	0	0
1	H	107	0	0	1	0
All	All	856	0	0	4	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 5.

All (4) 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:H:85:LEU:CA	1:H:86:PRO:CA	2.64	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:LEU:CA	1:D:86:PRO:CA	2.64	0.75
1:F:85:LEU:CA	1:F:86:PRO:CA	2.65	0.75
1:B:85:LEU:CA	1:B:86:PRO:CA	2.66	0.73

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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 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 ⓘ

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	107/110 (97%)	9.23	99 (92%) 0 1	82, 228, 336, 336	0
1	B	107/110 (97%)	7.52	101 (94%) 0 1	62, 240, 336, 336	0
1	C	107/110 (97%)	8.39	104 (97%) 0 0	76, 222, 336, 336	0
1	D	107/110 (97%)	8.58	102 (95%) 0 0	61, 209, 336, 336	0
1	E	107/110 (97%)	8.62	103 (96%) 0 0	56, 221, 336, 336	0
1	F	107/110 (97%)	7.30	96 (89%) 0 1	74, 209, 336, 336	0
1	G	107/110 (97%)	7.90	104 (97%) 0 0	48, 229, 336, 336	0
1	H	107/110 (97%)	7.80	100 (93%) 0 1	37, 241, 336, 336	0
All	All	856/880 (97%)	8.17	809 (94%) 0 1	37, 227, 336, 336	0

The worst 5 of 809 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	VAL	46.6
1	A	28	THR	39.4
1	E	84	ASP	35.3
1	A	102	ASN	34.5
1	D	14	GLU	34.3

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.