



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 7B53 / pdb_00007b53
Title : Crystal structure of MurE from E.coli
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Deposited on : 2020-12-03
Resolution : 1.75 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

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

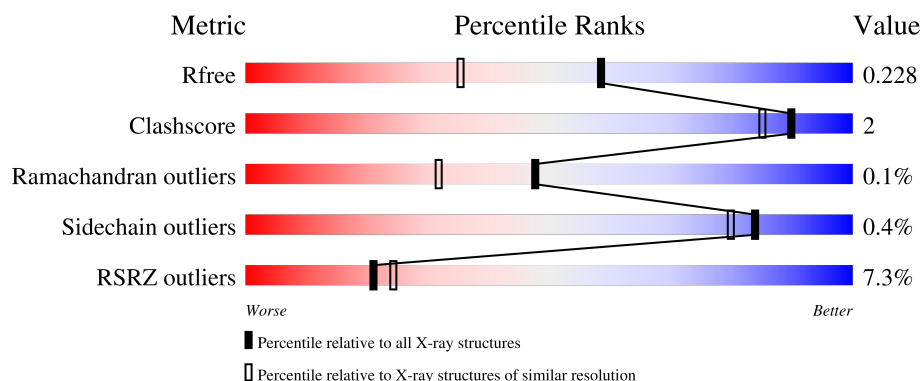
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	3183 (1.76-1.76)
Clashscore	190562	3299 (1.76-1.76)
Ramachandran outliers	187476	3274 (1.76-1.76)
Sidechain outliers	187428	3274 (1.76-1.76)
RSRZ outliers	180081	3183 (1.76-1.76)

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	496	5% 94% • •
1	B	496	9% 91% 5% •

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7842 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 UDP-N-acetylmuramoyl-L-alanyl-D-glutamate--2,6-diaminopimelate ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	487	Total	C	N	O	S	0	9	0
			3761	2349	681	710	21			
1	B	477	Total	C	N	O	S	0	7	0
			3642	2277	651	692	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P22188
B	0	SER	-	expression tag	UNP P22188

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

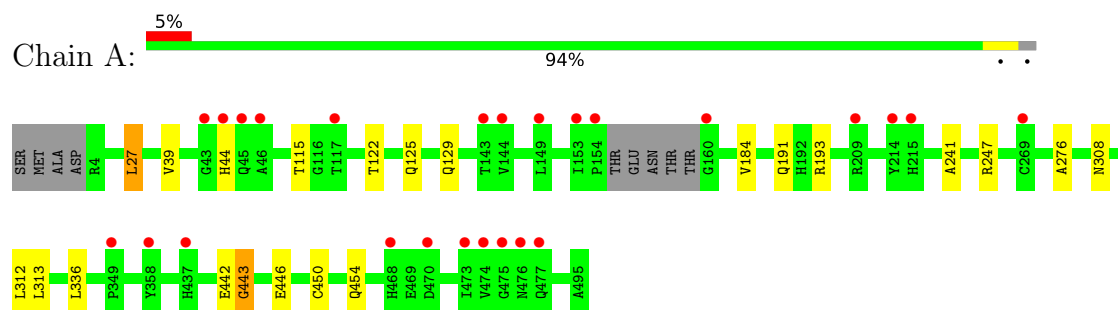
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	201	Total	O	0	0
			201	201		
3	B	222	Total	O	0	0
			222	222		

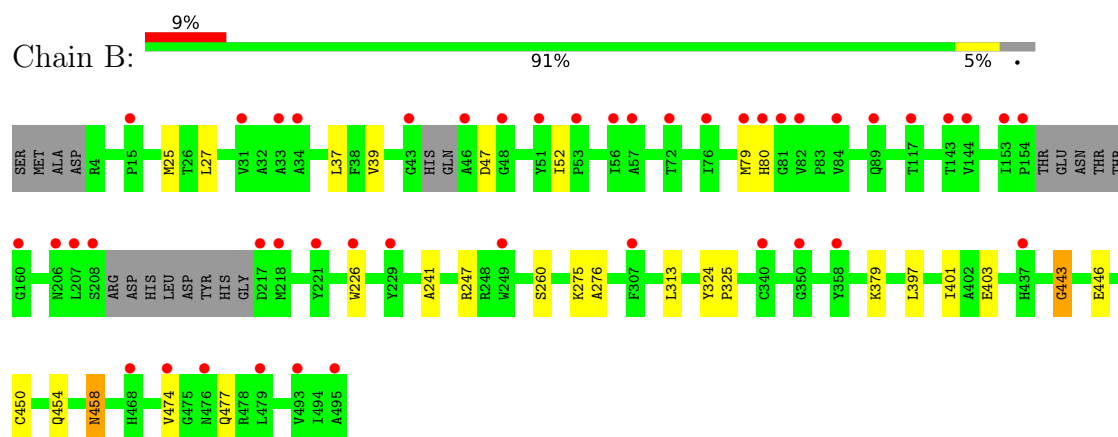
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: UDP-N-acetylmuramoyl-L-alanyl-D-glutamate--2,6-diaminopimelate ligase



- Molecule 1: UDP-N-acetylmuramoyl-L-alanyl-D-glutamate--2,6-diaminopimelate ligase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.63Å 58.94Å 74.23Å 97.32° 91.54° 105.17°	Depositor
Resolution (Å)	56.34 – 1.75 56.34 – 1.75	Depositor EDS
% Data completeness (in resolution range)	96.5 (56.34-1.75) 96.5 (56.34-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.195 , 0.220 0.203 , 0.228	Depositor DCC
R_{free} test set	4440 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	28.2	Xtriage
Anisotropy	0.223	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 32.6	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	0.96	EDS
Total number of atoms	7842	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.73% 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

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

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	0.54	0/3830	0.81	1/5197 (0.0%)
1	B	0.54	0/3708	0.80	1/5032 (0.0%)
All	All	0.54	0/7538	0.80	2/10229 (0.0%)

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	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	443	GLY	N-CA-C	-5.36	100.47	113.18
1	A	443	GLY	N-CA-C	-5.08	101.15	113.18

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	442	GLU	Peptide

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	3761	0	3711	12	0
1	B	3642	0	3602	15	0
2	A	8	0	12	0	0
2	B	8	0	12	1	0
3	A	201	0	0	0	0
3	B	222	0	0	2	0
All	All	7842	0	7337	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 2.

All (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:27:LEU:HD12	1:A:39:VAL:HB	1.84	0.59
1:B:27:LEU:HD23	1:B:39:VAL:HB	1.88	0.56
1:B:79:MET:HE3	1:B:80:HIS:CD2	2.42	0.54
1:A:443:GLY:HA3	1:A:446:GLU:OE1	2.08	0.54
1:A:450:CYS:SG	1:A:454:GLN:NE2	2.82	0.53
1:B:443:GLY:HA3	1:B:446:GLU:OE1	2.08	0.52
1:B:25:MET:HE3	1:B:37:LEU:HD23	1.92	0.50
1:B:450:CYS:SG	1:B:454:GLN:NE2	2.86	0.49
1:B:379:LYS:HG2	2:B:502:EDO:H12	1.95	0.48
1:B:403:GLU:O	3:B:601:HOH:O	2.21	0.46
1:B:474:VAL:O	1:B:477:GLN:HG2	2.14	0.46
1:A:276:ALA:HB2	1:A:313:LEU:HD11	1.98	0.46
1:A:115:THR:HG22	1:A:184:VAL:HB	1.99	0.44
1:A:241:ALA:O	1:A:247[A]:ARG:HG3	2.18	0.44
1:B:458:ASN:HD22	1:B:458:ASN:H	1.66	0.44
1:A:125[B]:GLN:CG	1:A:129:GLN:HE21	2.30	0.44
1:B:276:ALA:HB2	1:B:313:LEU:HD11	1.99	0.44
1:A:122:THR:HG21	1:A:308:ASN:ND2	2.33	0.43
1:A:191:GLN:OE1	1:A:193:ARG:NH2	2.48	0.43
1:B:226:TRP:NE1	3:B:608:HOH:O	2.52	0.43
1:B:241:ALA:O	1:B:247:ARG:HG3	2.19	0.43
1:A:241:ALA:O	1:A:247[B]:ARG:HG3	2.19	0.42
1:B:260:SER:O	1:B:275:LYS:HA	2.20	0.42
1:B:397:LEU:O	1:B:401:ILE:HG12	2.19	0.42
1:A:312[B]:LEU:HD11	1:A:336:LEU:CD1	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:HIS:O	1:A:44:HIS:CG	2.74	0.41
1:B:324:TYR:HA	1:B:325:PRO:HD3	1.97	0.40

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	492/496 (99%)	485 (99%)	7 (1%)	0	100	100
1	B	476/496 (96%)	467 (98%)	8 (2%)	1 (0%)	43	27
All	All	968/992 (98%)	952 (98%)	15 (2%)	1 (0%)	48	32

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	ASP

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	388/387 (100%)	387 (100%)	1 (0%)	86	83
1	B	377/387 (97%)	375 (100%)	2 (0%)	81	75
All	All	765/774 (99%)	762 (100%)	3 (0%)	84	80

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LEU
1	B	52	ILE
1	B	458	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	268	ASN
1	A	302	HIS
1	A	308	ASN
1	A	454	GLN
1	A	468	HIS
1	B	118	ASN
1	B	280	ASN
1	B	454	GLN
1	B	458	ASN

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](#)

4 ligands are modelled in this entry.

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	$\# Z > 2$
2	EDO	A	501	-	3,3,3	0.08	0	2,2,2	0.14	0
2	EDO	B	502	-	3,3,3	0.05	0	2,2,2	0.17	0
2	EDO	B	501	-	3,3,3	0.05	0	2,2,2	0.08	0
2	EDO	A	502	-	3,3,3	0.12	0	2,2,2	0.06	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
2	EDO	A	501	-	-	0/1/1/1	-
2	EDO	B	502	-	-	0/1/1/1	-
2	EDO	B	501	-	-	0/1/1/1	-
2	EDO	A	502	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	502	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	502	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

6.1 Protein, DNA and RNA chains

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	487/496 (98%)	0.48	25 (5%)	33 39	11, 34, 58, 108	9 (1%)
1	B	477/496 (96%)	0.56	45 (9%)	14 17	13, 33, 71, 105	7 (1%)
All	All	964/992 (97%)	0.52	70 (7%)	21 24	11, 34, 67, 108	16 (1%)

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	144	VAL	6.3
1	B	46	ALA	5.2
1	B	144	VAL	4.5
1	B	160	GLY	4.1
1	A	44	HIS	3.8
1	A	470	ASP	3.7
1	A	358	TYR	3.7
1	A	215	HIS	3.5
1	B	474	VAL	3.4
1	B	154	PRO	3.4
1	B	217	ASP	3.4
1	A	474	VAL	3.3
1	A	476	ASN	3.3
1	A	477	GLN	3.3
1	B	358	TYR	3.2
1	B	79	MET	3.1
1	A	475	GLY	3.0
1	B	81	GLY	3.0
1	A	143	THR	3.0
1	B	495	ALA	3.0
1	A	468	HIS	3.0
1	B	153	ILE	3.0
1	B	207	LEU	3.0
1	A	154	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	226	TRP	2.9
1	B	43	GLY	2.9
1	B	56	ILE	2.8
1	B	143	THR	2.8
1	A	45	GLN	2.8
1	B	82	VAL	2.7
1	A	473	ILE	2.7
1	A	160	GLY	2.7
1	B	80	HIS	2.6
1	B	340	CYS	2.6
1	B	57	ALA	2.6
1	B	72	THR	2.5
1	B	307	PHE	2.5
1	A	269[A]	CYS	2.5
1	B	208	SER	2.5
1	A	349	PRO	2.5
1	B	437	HIS	2.5
1	B	206	ASN	2.5
1	A	209	ARG	2.4
1	A	153	ILE	2.4
1	B	229	TYR	2.4
1	B	476	ASN	2.4
1	B	493	VAL	2.3
1	B	48	GLY	2.3
1	B	117	THR	2.3
1	A	149	LEU	2.3
1	A	117	THR	2.3
1	B	218	MET	2.3
1	B	33	ALA	2.2
1	A	43	GLY	2.2
1	B	15	PRO	2.2
1	A	214	TYR	2.2
1	B	51	TYR	2.2
1	B	221	TYR	2.2
1	B	84	VAL	2.1
1	B	479	LEU	2.1
1	B	249	TRP	2.1
1	B	53	PRO	2.1
1	B	350	GLY	2.1
1	A	46	ALA	2.0
1	B	34	ALA	2.0
1	B	31	VAL	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	468	HIS	2.0
1	B	89	GLN	2.0
1	B	76	ILE	2.0
1	A	437	HIS	2.0

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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	EDO	B	502	4/4	0.71	0.22	49,58,65,67	0
2	EDO	A	502	4/4	0.79	0.24	38,38,42,51	0
2	EDO	A	501	4/4	0.80	0.16	41,42,53,56	0
2	EDO	B	501	4/4	0.89	0.20	45,46,49,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.