Full wwPDB X-ray Structure Validation Report (i)

May 13, 2020 – 11:11 am BST

PDB ID	:	1FF3
Title	:	STRUCTURE OF THE PEPTIDE METHIONINE SULFOXIDE REDUCTASE FROM ESCHERICHIA COLI
Authors	:	Tete-Favier, F.; Cobessi, D.; Boschi-Muller, S.; Azza, S.; Branlant, G.; Aubry, A.
Deposited on	:	2000-07-25
Resolution	:	1.90 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 (i) symbol.

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

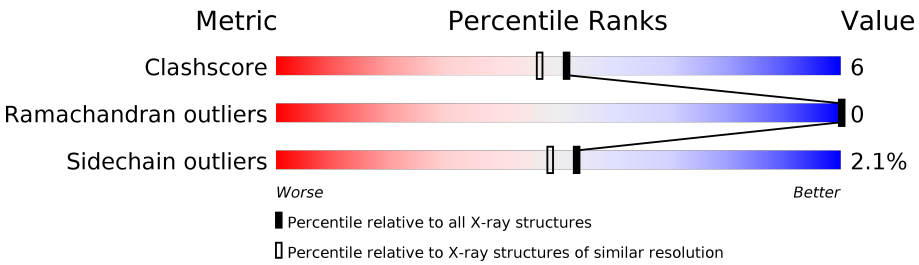
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

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 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	211	90% 9% •
1	B	211	86% 5% • 8%
1	C	211	73% 13% • 12%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5153 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.

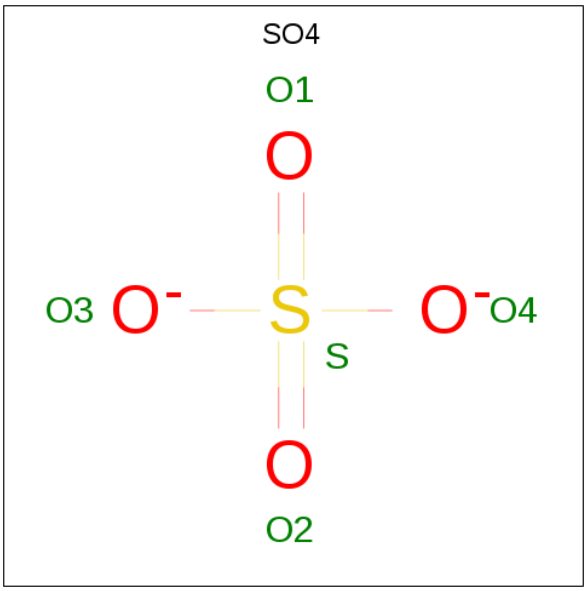
- Molecule 1 is a protein called PEPTIDE METHIONINE SULFOXIDE REDUCTASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	As	C	N	O	S			
1	A	211	1637	1	1033	280	313	10	0	0	0
			Total	As	C	N	O	S			
1	B	194	1521	1	958	263	291	8	0	0	0
			Total	As	C	N	O	S			
1	C	185	1445	1	909	249	278	8	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	CAS	CYS	MODIFIED RESIDUE	UNP P0A744
B	51	CAS	CYS	MODIFIED RESIDUE	UNP P0A744
C	51	CAS	CYS	MODIFIED RESIDUE	UNP P0A744

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	A	1	5	4	1	0	0
			Total	O	S		
2	B	1	5	4	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	229	229	229	0	0
			Total	O		
3	B	216	216	216	0	0
			Total	O		
3	C	85	85	85	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

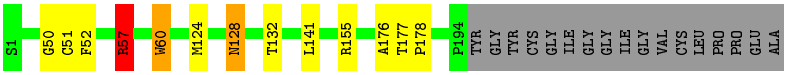
Note EDS was not executed.

- Molecule 1: PEPTIDE METHIONINE SULFOXIDE REDUCTASE

Chain A:

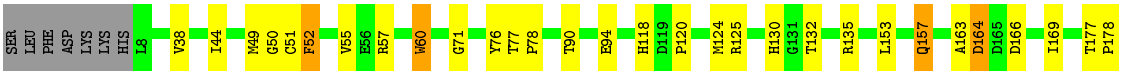
- Molecule 1: PEPTIDE METHIONINE SULFOXIDE REDUCTASE

Chain B:



- Molecule 1: PEPTIDE METHIONINE SULFOXIDE REDUCTASE

Chain C:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α, β, γ	102.50Å 102.50Å 292.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) (30.00-1.90)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.195 , 0.218	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5153	wwPDB-VP
Average B, all atoms ($Å^2$)	24.0	wwPDB-VP

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: CAS, SO4

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	0.31	0/1677	0.57	0/2289
1	B	0.32	0/1557	0.63	1/2126 (0.0%)
1	C	0.29	0/1478	0.54	0/2020
All	All	0.31	0/4712	0.58	1/6435 (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	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	57	ARG	NE-CZ-NH1	6.82	123.71	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	57	ARG	Sidechain

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	1637	0	1524	15	0
1	B	1521	0	1414	10	0
1	C	1445	0	1337	24	1
2	A	15	0	0	0	0
2	B	5	0	0	0	0
3	A	229	0	0	1	0
3	B	216	0	0	2	0
3	C	85	0	0	0	0
All	All	5153	0	4275	49	1

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 6.

All (49) 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:51:CAS:CE2	1:A:51:CAS:AS	2.23	1.47
1:B:51:CAS:CE2	1:B:51:CAS:AS	2.23	1.46
1:A:51:CAS:CE1	1:A:51:CAS:AS	2.24	1.45
1:C:51:CAS:AS	1:C:51:CAS:CE2	2.24	1.45
1:C:51:CAS:CE1	1:C:51:CAS:AS	2.25	1.45
1:B:51:CAS:CE1	1:B:51:CAS:AS	2.25	1.45
1:C:157:GLN:HG2	1:C:169:ILE:HD12	1.62	0.79
1:B:128:ASN:ND2	3:B:724:HOH:O	2.18	0.75
1:C:124:MET:SD	1:C:132:THR:HG22	2.34	0.68
1:B:57:ARG:HG3	1:B:60:TRP:CZ2	2.31	0.66
1:C:38:VAL:HG21	1:C:44:ILE:HD11	1.78	0.66
1:A:37:ASN:OD1	1:A:38:VAL:N	2.36	0.59
1:B:141:LEU:HD23	1:B:176:ALA:HB3	1.87	0.56
1:A:200:ILE:HG23	1:A:200:ILE:O	2.06	0.55
1:B:124:MET:SD	1:B:132:THR:HG22	2.48	0.54
1:C:124:MET:HE2	1:C:135:ARG:HB3	1.91	0.53
1:B:177:THR:HB	1:B:178:PRO:HD2	1.92	0.51
1:B:177:THR:HB	1:B:178:PRO:CD	2.41	0.51
1:A:71:GLY:HA3	1:A:179:PHE:CZ	2.46	0.51
1:C:118:HIS:O	1:C:120:PRO:HD3	2.11	0.50
1:C:163:ALA:O	1:C:164:ASP:HB2	2.11	0.50
1:C:157:GLN:HE21	1:C:157:GLN:CA	2.25	0.49
1:C:125:ARG:HH11	1:C:125:ARG:HG2	1.78	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:ARG:HD2	1:C:130:HIS:CD2	2.47	0.49
1:C:50:GLY:O	1:C:51:CAS:C	2.61	0.48
1:C:177:THR:HB	1:C:178:PRO:CD	2.44	0.48
1:A:177:THR:HB	1:A:178:PRO:CD	2.44	0.47
1:C:124:MET:HE2	1:C:135:ARG:CB	2.44	0.47
1:A:193:ASN:HB3	1:A:195:TYR:CE2	2.50	0.47
1:C:52:PHE:CD2	1:C:94:GLU:HB3	2.50	0.46
1:C:76:TYR:O	1:C:78:PRO:HD3	2.16	0.46
1:A:125:ARG:HB2	1:A:130:HIS:HA	1.97	0.45
1:B:50:GLY:O	1:B:51:CAS:C	2.64	0.44
1:C:71:GLY:HA3	1:C:179:PHE:CZ	2.52	0.44
1:A:177:THR:HB	1:A:178:PRO:HD2	2.00	0.44
1:C:177:THR:HB	1:C:178:PRO:HD2	2.01	0.43
1:A:50:GLY:O	1:A:51:CAS:C	2.66	0.43
1:A:6:LYS:HG2	3:A:699:HOH:O	2.19	0.43
1:C:153:LEU:C	1:C:153:LEU:HD23	2.39	0.43
1:C:57:ARG:HA	1:C:60:TRP:CD2	2.53	0.43
1:A:153:LEU:HD23	1:A:153:LEU:C	2.41	0.42
1:B:155:ARG:HD3	3:B:706:HOH:O	2.20	0.41
1:C:125:ARG:HB2	1:C:130:HIS:HA	2.02	0.41
1:C:49:MET:HG3	1:C:55:VAL:HG22	2.03	0.41
1:A:54:GLY:HA3	1:A:197:TYR:CE2	2.55	0.41
1:A:57:ARG:HA	1:A:60:TRP:CD2	2.55	0.41
1:A:33:HIS:HB3	1:A:66:TYR:CZ	2.56	0.40
1:C:124:MET:O	1:C:125:ARG:HB3	2.22	0.40
1:C:77:THR:OG1	1:C:90:THR:HA	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:166:ASP:OD1	1:C:166:ASP:OD1[8_555]	2.01	0.19

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	208/211 (99%)	203 (98%)	5 (2%)	0	100	100
1	B	191/211 (90%)	188 (98%)	3 (2%)	0	100	100
1	C	182/211 (86%)	178 (98%)	4 (2%)	0	100	100
All	All	581/633 (92%)	569 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	167/167 (100%)	164 (98%)	3 (2%)	59	55
1	B	156/167 (93%)	153 (98%)	3 (2%)	57	53
1	C	147/167 (88%)	143 (97%)	4 (3%)	44	38
All	All	470/501 (94%)	460 (98%)	10 (2%)	53	48

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	PHE
1	A	60	TRP
1	A	125	ARG
1	B	52	PHE
1	B	60	TRP
1	B	128	ASN
1	C	52	PHE
1	C	60	TRP
1	C	157	GLN
1	C	164	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	GLN
1	B	19	ASN
1	B	128	ASN
1	C	61	GLN
1	C	130	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

3 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	CAS	A	51	1	5,8,9	6.75	2 (40%)	1,9,11	0.34	0
1	CAS	B	51	1	5,8,9	6.86	2 (40%)	1,9,11	0.15	0
1	CAS	C	51	1,3	5,8,9	6.90	2 (40%)	1,9,11	0.16	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
1	CAS	A	51	1	-	0/0/7/9	-
1	CAS	B	51	1	-	0/0/7/9	-
1	CAS	C	51	1,3	-	0/0/7/9	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	51	CAS	AS-CE1	11.18	2.25	1.96

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	51	CAS	AS-CE1	11.15	2.25	1.96
1	A	51	CAS	AS-CE1	10.98	2.24	1.96
1	C	51	CAS	AS-CE2	10.61	2.24	1.96
1	B	51	CAS	AS-CE2	10.47	2.23	1.96
1	A	51	CAS	AS-CE2	10.30	2.23	1.96

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	51	CAS	3	0
1	B	51	CAS	3	0
1	C	51	CAS	3	0

5.5 Carbohydrates (i)

There are no carbohydrates 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	SO4	A	534	-	4,4,4	0.23	0	6,6,6	0.07	0
2	SO4	B	536	-	4,4,4	0.24	0	6,6,6	0.06	0
2	SO4	A	535	-	4,4,4	0.28	0	6,6,6	0.08	0
2	SO4	A	533	-	4,4,4	0.26	0	6,6,6	0.09	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.