

Full wwPDB X-ray Structure Validation Report (i)

Oct 23, 2021 – 11:47 AM EDT

PDB ID	:	1F75
Title	:	CRYSTAL STRUCTURE OF UNDECAPRENYL DIPHOSPHATE SYN-
		THASE FROM MICROCOCCUS LUTEUS B-P 26
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Deposited on		
Resolution	:	2.20 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 (1)) were used in the production of this report:

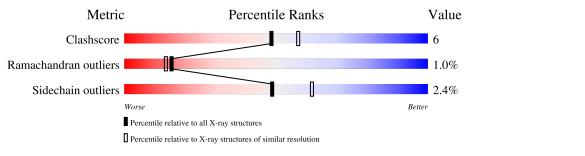
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.20 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	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)

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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	А	249	71%	16%	• 13%			
1	В	249	73%	12%	15%			



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3552 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.

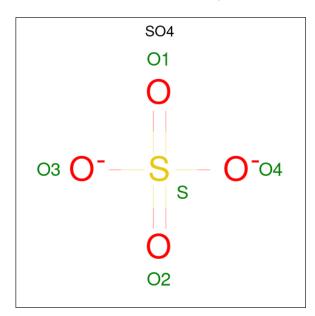
• Molecule 1 is a protein called UNDECAPRENYL PYROPHOSPHATE SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A 91'	217	Total	С	Ν	0	S	0	0	0
	211	1769	1140	294	328	$\overline{7}$	0	0	0	
1	В	212	Total	С	Ν	0	S	0	0	0
	D	212	1735	1120	287	321	7	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	54	ARG	LYS	engineered mutation	UNP 082827
В	54	ARG	LYS	engineered mutation	UNP 082827

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



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	А	1	Total 5	0 4	S 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	В	1	Total 5	0 4	S 1	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	19	Total O 19 19	0	0
3	В	19	Total O 19 19	0	0



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. 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: UNDECAPRENYL PYROPHOSPHATE SYNTHETASE



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	127.20Å 60.16 Å 75.70 Å	Depositor
a, b, c, α , β , γ	90.00° 105.69° 90.00°	Depositor
Resolution (Å)	50.00 - 2.20	Depositor
% Data completeness	(Not available) (50.00-2.20)	Depositor
(in resolution range)	(1100 available) (50.00 2.20)	Depositor
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.190 , 0.246	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3552	wwPDB-VP
Average B, all atoms $(Å^2)$	47.0	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: 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		lengths	Bond angles		
		RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.33	0/1808	0.54	0/2440	
1	В	0.32	0/1774	0.54	0/2393	
All	All	0.32	0/3582	0.54	0/4833	

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	А	1769	0	1761	30	0
1	В	1735	0	1728	18	0
2	А	5	0	0	0	0
2	В	5	0	0	0	0
3	А	19	0	0	0	0
3	В	19	0	0	1	0
All	All	3552	0	3489	42	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 6.

All (42) close contacts within the same asymmetric unit are listed below, sorted by their clash



magnitude.

A + a 1	A +	Interatomic	Clash	
Atom-1	Atom-2	distance (\AA)	overlap (Å)	
1:A:107:ASN:HD21	1:A:138:THR:H	1.31	0.78	
1:A:204:LEU:HB2	1:B:217:PHE:CE1	2.28	0.69	
1:A:107:ASN:ND2	1:A:138:THR:H	1.94	0.66	
1:A:154:ILE:O	1:A:158:VAL:HG23	2.00	0.61	
1:A:15:ILE:HD13	1:A:19:GLN:HG3	1.83	0.59	
1:A:219:PHE:HE1	1:B:219:PHE:HE1	1.51	0.57	
1:B:181:ASN:HB3	3:B:627:HOH:O	2.06	0.56	
1:A:23:HIS:HD2	1:A:190:PRO:O	1.90	0.54	
1:B:216:GLU:HG3	1:B:242:ARG:HH21	1.72	0.54	
1:A:15:ILE:HD11	1:A:17:ALA:HB3	1.90	0.54	
1:A:17:ALA:C	1:A:19:GLN:H	2.10	0.54	
1:B:157:ALA:O	1:B:161:ILE:HG13	2.08	0.53	
1:A:36:LYS:HE3	1:A:37:GLN:HE21	1.73	0.53	
1:B:197:ARG:HB3	1:B:219:PHE:HD2	1.76	0.50	
1:A:239:TYR:C	1:A:241:ASN:H	2.15	0.49	
1:A:102:GLU:O	1:A:106:LYS:HB2	2.13	0.49	
1:B:91:LEU:HB3	1:B:92:PRO:HD3	1.94	0.49	
1:B:39:LYS:HB2	1:B:39:LYS:NZ	2.27	0.49	
1:A:241:ASN:O	1:A:242:ARG:HB2	2.14	0.48	
1:A:38:LYS:HB3	1:A:40:MET:HE2	1.96	0.47	
1:B:86:ASN:O	1:B:90:LYS:HE3	2.14	0.47	
1:B:224:TRP:HB3	1:B:225:PRO:HD3	1.96	0.46	
1:A:204:LEU:HB2	1:B:217:PHE:CZ	2.51	0.45	
1:A:49:GLY:O	1:A:53:VAL:HG23	2.16	0.45	
1:B:234:GLN:O	1:B:238:ILE:HG13	2.16	0.45	
1:A:160:LEU:O	1:A:164:ARG:HG3	2.17	0.44	
1:B:216:GLU:CG	1:B:242:ARG:HH21	2.30	0.44	
1:A:38:LYS:HB3	1:A:40:MET:CE	2.47	0.44	
1:A:166:LYS:HD2	1:B:165:TYR:CE2	2.53	0.44	
1:A:34:TRP:CH2	1:A:48:GLU:HG3	2.53	0.44	
1:A:52:THR:HA	1:A:55:LYS:HE2	2.00	0.44	
1:A:17:ALA:O	1:A:19:GLN:N	2.49	0.43	
1:A:204:LEU:HB2	1:B:217:PHE:HE1	1.77	0.43	
1:A:15:ILE:HG21	1:A:19:GLN:HG3	1.99	0.43	
1:A:204:LEU:HD23	1:A:205:SER:N	2.34	0.42	
1:A:189:MET:HA	1:A:190:PRO:HD3	1.84	0.42	
1:A:165:TYR:CE2	1:B:166:LYS:HD2	2.56	0.41	
1:B:122:HIS:CD2	1:B:122:HIS:H	2.37	0.41	
1:A:197:ARG:HD3	1:A:201:GLU:O	2.20	0.41	
1:B:23:HIS:CE1	1:B:192:PRO:HG3	2.55	0.41	
1:A:201:GLU:HG2	1:A:203:ARG:HH21	1.86	0.41	

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:TRP:HB3	1:A:225:PRO:HD3	2.03	0.41

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	Analysed Favoured Allowed		Outliers	Percentiles
1	А	213/249~(86%)	207~(97%)	3~(1%)	3~(1%)	11 8
1	В	208/249~(84%)	200 (96%)	7 (3%)	1 (0%)	29 31
All	All	421/498 (84%)	407 (97%)	10 (2%)	4 (1%)	15 14

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	17	ALA
1	В	201	GLU
1	А	18	ALA
1	А	240	GLN

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	Analysed Rotameric		Percentiles	
1	А	192/221~(87%)	189~(98%)	3~(2%)	62 76	

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	Chain	Analysed	Rotameric	Outliers	Percentiles
1	В	189/221~(86%)	183~(97%)	6 (3%)	39 50
All	All	381/442 (86%)	372 (98%)	9(2%)	49 62

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All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	23	HIS
1	А	42	ARG
1	А	221	ASP
1	В	54	ARG
1	В	87	TYR
1	В	106	LYS
1	В	117	ASP
1	В	140	LEU
1	В	208	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	16	ASN
1	А	23	HIS
1	А	37	GLN
1	А	107	ASN
1	А	147	ASN
1	В	122	HIS
1	В	147	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 monosaccharides in this entry.



5.6 Ligand geometry (i)

2 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	Tuno	Chain	Dec	Tinle	В	ond leng	gths	В	ond ang	gles
IVIOI	Type	Chain	nes	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
2	SO4	В	902	-	4,4,4	1.93	1 (25%)	$6,\!6,\!6$	0.88	0
2	SO4	А	901	-	4,4,4	1.77	1 (25%)	$6,\!6,\!6$	0.92	0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
2	В	902	SO4	O1-S	3.43	1.64	1.46
2	А	901	SO4	O1-S	3.06	1.62	1.46

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.

