



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 10, 2020 – 02:32 AM BST

PDB ID : 6SXN
Title : Crystal structure of P212121 apo form of CrtE
Authors : Feng, Y.; Morgan, R.M.L.; Nixon, P.J.
Deposited on : 2019-09-26
Resolution : 2.66 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtrriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

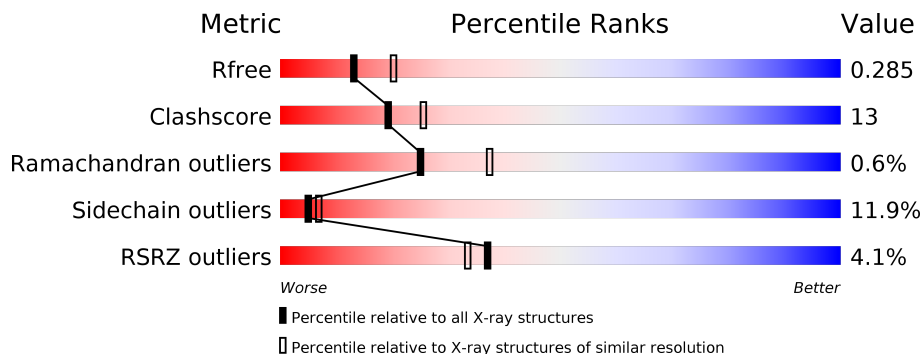
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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}	130704	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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 $\leq 5\%$. 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	267	
2	B	259	
3	C	258	
4	D	267	
5	E	259	
6	F	253	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11367 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 Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	267	1931	1221	328	371	11	0	0	0

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	THR	deletion	UNP B1XJV9
A	?	-	ALA	deletion	UNP B1XJV9
A	?	-	THR	deletion	UNP B1XJV9
A	?	-	ALA	deletion	UNP B1XJV9
A	?	-	GLU	deletion	UNP B1XJV9
A	?	-	GLU	deletion	UNP B1XJV9
A	?	-	LEU	deletion	UNP B1XJV9
A	?	-	GLY	deletion	UNP B1XJV9
A	?	-	LYS	deletion	UNP B1XJV9
A	?	-	THR	deletion	UNP B1XJV9
A	?	-	ALA	deletion	UNP B1XJV9
A	?	-	GLY	deletion	UNP B1XJV9
A	?	-	LYS	deletion	UNP B1XJV9
A	?	-	ASP	deletion	UNP B1XJV9
A	?	-	LEU	deletion	UNP B1XJV9
A	?	-	GLU	deletion	UNP B1XJV9
A	?	-	ALA	deletion	UNP B1XJV9
A	?	-	GLN	deletion	UNP B1XJV9
A	?	-	LYS	deletion	UNP B1XJV9
A	?	-	SER	deletion	UNP B1XJV9
A	?	-	LEU	deletion	UNP B1XJV9
A	?	-	TRP	deletion	UNP B1XJV9
A	?	-	GLY	deletion	UNP B1XJV9
A	?	-	ILE	deletion	UNP B1XJV9
A	?	-	GLU	deletion	UNP B1XJV9
A	300	ALA	-	expression tag	UNP B1XJV9

- Molecule 2 is a protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	259	1884	1192	319	362	11	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	GLY	deletion	UNP B1XJV9
B	?	-	LYS	deletion	UNP B1XJV9
B	258	SER	ASP	conflict	UNP B1XJV9
B	?	-	ILE	deletion	UNP B1XJV9
B	?	-	THR	deletion	UNP B1XJV9
B	?	-	ALA	deletion	UNP B1XJV9
B	?	-	THR	deletion	UNP B1XJV9
B	?	-	ALA	deletion	UNP B1XJV9
B	?	-	GLU	deletion	UNP B1XJV9
B	?	-	GLU	deletion	UNP B1XJV9
B	?	-	LEU	deletion	UNP B1XJV9
B	?	-	GLY	deletion	UNP B1XJV9
B	?	-	LYS	deletion	UNP B1XJV9
B	?	-	THR	deletion	UNP B1XJV9
B	?	-	ALA	deletion	UNP B1XJV9
B	?	-	GLY	deletion	UNP B1XJV9
B	?	-	LYS	deletion	UNP B1XJV9
B	?	-	ASP	deletion	UNP B1XJV9
B	?	-	LEU	deletion	UNP B1XJV9
B	?	-	GLU	deletion	UNP B1XJV9
B	?	-	ALA	deletion	UNP B1XJV9
B	?	-	GLN	deletion	UNP B1XJV9
B	?	-	LYS	deletion	UNP B1XJV9
B	?	-	VAL	deletion	UNP B1XJV9
B	?	-	THR	deletion	UNP B1XJV9
B	?	-	TYR	deletion	UNP B1XJV9
B	?	-	PRO	deletion	UNP B1XJV9
B	?	-	SER	deletion	UNP B1XJV9
B	?	-	LEU	deletion	UNP B1XJV9
B	?	-	TRP	deletion	UNP B1XJV9
B	?	-	GLY	deletion	UNP B1XJV9
B	259	LEU	ILE	conflict	UNP B1XJV9

- Molecule 3 is a protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	258	1877	1187	318	361	11	0	0	0

There are 31 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	LYS	deletion	UNP B1XJV9
C	?	-	THR	deletion	UNP B1XJV9
C	?	-	ASP	deletion	UNP B1XJV9
C	?	-	VAL	deletion	UNP B1XJV9
C	?	-	ALA	deletion	UNP B1XJV9
C	?	-	ASP	deletion	UNP B1XJV9
C	?	-	ILE	deletion	UNP B1XJV9
C	?	-	THR	deletion	UNP B1XJV9
C	?	-	ALA	deletion	UNP B1XJV9
C	?	-	THR	deletion	UNP B1XJV9
C	?	-	ALA	deletion	UNP B1XJV9
C	?	-	GLU	deletion	UNP B1XJV9
C	?	-	GLU	deletion	UNP B1XJV9
C	?	-	LEU	deletion	UNP B1XJV9
C	?	-	GLY	deletion	UNP B1XJV9
C	?	-	LYS	deletion	UNP B1XJV9
C	?	-	THR	deletion	UNP B1XJV9
C	?	-	ALA	deletion	UNP B1XJV9
C	?	-	GLY	deletion	UNP B1XJV9
C	?	-	LYS	deletion	UNP B1XJV9
C	?	-	ASP	deletion	UNP B1XJV9
C	?	-	LEU	deletion	UNP B1XJV9
C	?	-	GLU	deletion	UNP B1XJV9
C	?	-	ALA	deletion	UNP B1XJV9
C	?	-	GLN	deletion	UNP B1XJV9
C	?	-	LYS	deletion	UNP B1XJV9
C	?	-	VAL	deletion	UNP B1XJV9
C	?	-	THR	deletion	UNP B1XJV9
C	?	-	TYR	deletion	UNP B1XJV9
C	?	-	PRO	deletion	UNP B1XJV9
C	?	-	SER	deletion	UNP B1XJV9

- Molecule 4 is a protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	267	1918	1212	327	368	11	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	ASP	deletion	UNP B1XJV9
D	?	-	ILE	deletion	UNP B1XJV9
D	?	-	THR	deletion	UNP B1XJV9
D	?	-	ALA	deletion	UNP B1XJV9
D	?	-	THR	deletion	UNP B1XJV9
D	?	-	ALA	deletion	UNP B1XJV9
D	?	-	GLU	deletion	UNP B1XJV9
D	?	-	GLU	deletion	UNP B1XJV9
D	?	-	LEU	deletion	UNP B1XJV9
D	?	-	GLY	deletion	UNP B1XJV9
D	?	-	LYS	deletion	UNP B1XJV9
D	?	-	THR	deletion	UNP B1XJV9
D	?	-	ALA	deletion	UNP B1XJV9
D	?	-	GLY	deletion	UNP B1XJV9
D	?	-	LYS	deletion	UNP B1XJV9
D	?	-	ASP	deletion	UNP B1XJV9
D	?	-	LEU	deletion	UNP B1XJV9
D	?	-	GLU	deletion	UNP B1XJV9
D	?	-	ALA	deletion	UNP B1XJV9
D	?	-	GLN	deletion	UNP B1XJV9
D	?	-	LYS	deletion	UNP B1XJV9
D	?	-	VAL	deletion	UNP B1XJV9

- Molecule 5 is a protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	E	259	1859	1175	320	353	11	0	0	0

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	GLY	deletion	UNP B1XJV9
E	?	-	LYS	deletion	UNP B1XJV9
E	?	-	THR	deletion	UNP B1XJV9
E	?	-	ASP	deletion	UNP B1XJV9
E	?	-	ASP	deletion	UNP B1XJV9
E	?	-	ASP	deletion	UNP B1XJV9
E	?	-	ILE	deletion	UNP B1XJV9
E	?	-	LEU	deletion	UNP B1XJV9
E	?	-	ASP	deletion	UNP B1XJV9

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Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	ILE	deletion	UNP B1XJV9
E	?	-	THR	deletion	UNP B1XJV9
E	?	-	ALA	deletion	UNP B1XJV9
E	?	-	THR	deletion	UNP B1XJV9
E	?	-	ALA	deletion	UNP B1XJV9
E	?	-	GLU	deletion	UNP B1XJV9
E	?	-	GLU	deletion	UNP B1XJV9
E	?	-	LEU	deletion	UNP B1XJV9
E	?	-	GLY	deletion	UNP B1XJV9
E	?	-	LYS	deletion	UNP B1XJV9
E	?	-	THR	deletion	UNP B1XJV9
E	?	-	ALA	deletion	UNP B1XJV9
E	?	-	GLY	deletion	UNP B1XJV9
E	?	-	LYS	deletion	UNP B1XJV9
E	?	-	ASP	deletion	UNP B1XJV9
E	?	-	LEU	deletion	UNP B1XJV9
E	?	-	GLU	deletion	UNP B1XJV9
E	?	-	ALA	deletion	UNP B1XJV9
E	?	-	GLN	deletion	UNP B1XJV9
E	?	-	LYS	deletion	UNP B1XJV9
E	?	-	VAL	deletion	UNP B1XJV9
E	?	-	THR	deletion	UNP B1XJV9
E	?	-	TYR	deletion	UNP B1XJV9
E	?	-	PRO	deletion	UNP B1XJV9
E	?	-	SER	deletion	UNP B1XJV9
E	?	-	LEU	deletion	UNP B1XJV9
E	?	-	TRP	deletion	UNP B1XJV9
E	?	-	GLY	deletion	UNP B1XJV9
E	?	-	ILE	deletion	UNP B1XJV9
E	?	-	GLU	deletion	UNP B1XJV9

- Molecule 6 is a protein called Geranylgeranyl pyrophosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	F	253	1852	1172	314	355	11	0	0	0

There are 39 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	?	-	THR	deletion	UNP B1XJV9
F	?	-	ASP	deletion	UNP B1XJV9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	?	-	VAL	deletion	UNP B1XJV9
F	?	-	ALA	deletion	UNP B1XJV9
F	?	-	VAL	deletion	UNP B1XJV9
F	?	-	ILE	deletion	UNP B1XJV9
F	?	-	LEU	deletion	UNP B1XJV9
F	?	-	ASP	deletion	UNP B1XJV9
F	?	-	ILE	deletion	UNP B1XJV9
F	?	-	THR	deletion	UNP B1XJV9
F	?	-	ALA	deletion	UNP B1XJV9
F	?	-	THR	deletion	UNP B1XJV9
F	?	-	ALA	deletion	UNP B1XJV9
F	?	-	GLU	deletion	UNP B1XJV9
F	?	-	GLU	deletion	UNP B1XJV9
F	?	-	LEU	deletion	UNP B1XJV9
F	?	-	GLY	deletion	UNP B1XJV9
F	?	-	LYS	deletion	UNP B1XJV9
F	?	-	THR	deletion	UNP B1XJV9
F	?	-	ALA	deletion	UNP B1XJV9
F	?	-	GLY	deletion	UNP B1XJV9
F	?	-	LYS	deletion	UNP B1XJV9
F	?	-	ASP	deletion	UNP B1XJV9
F	?	-	LEU	deletion	UNP B1XJV9
F	?	-	GLU	deletion	UNP B1XJV9
F	?	-	ALA	deletion	UNP B1XJV9
F	?	-	GLN	deletion	UNP B1XJV9
F	?	-	LYS	deletion	UNP B1XJV9
F	?	-	VAL	deletion	UNP B1XJV9
F	?	-	THR	deletion	UNP B1XJV9
F	?	-	TYR	deletion	UNP B1XJV9
F	?	-	PRO	deletion	UNP B1XJV9
F	?	-	SER	deletion	UNP B1XJV9
F	?	-	LEU	deletion	UNP B1XJV9
F	?	-	TRP	deletion	UNP B1XJV9
F	?	-	GLY	deletion	UNP B1XJV9
F	?	-	ILE	deletion	UNP B1XJV9
F	?	-	GLU	deletion	UNP B1XJV9
F	?	-	LYS	deletion	UNP B1XJV9

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	11	Total O 11 11	0	0

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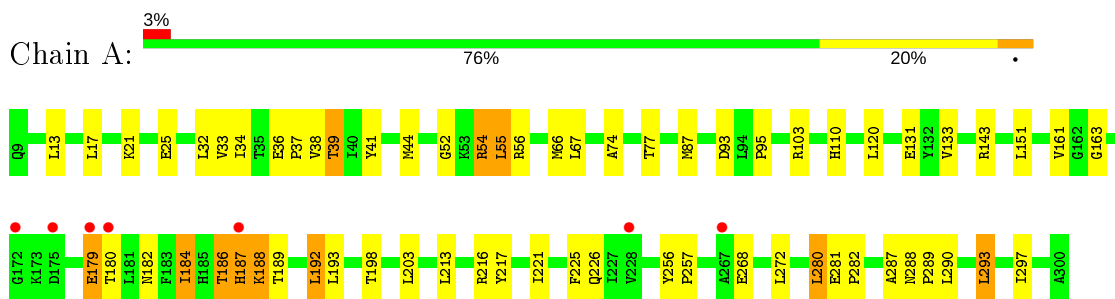
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	12	Total 12	O 12	0	0
7	C	5	Total 5	O 5	0	0
7	D	2	Total 2	O 2	0	0
7	E	9	Total 9	O 9	0	0
7	F	7	Total 7	O 7	0	0

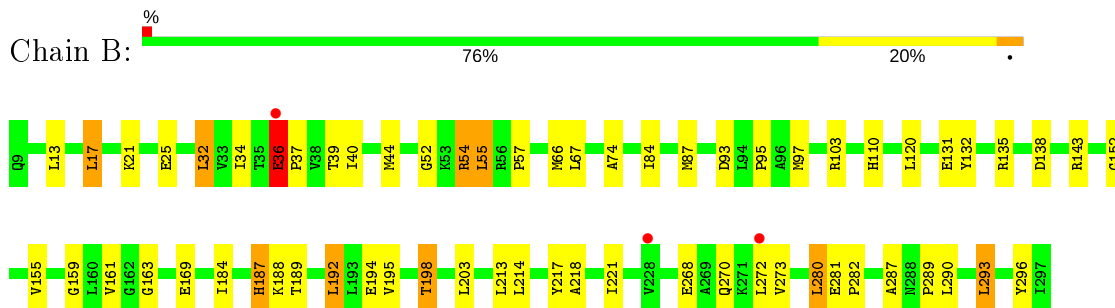
3 Residue-property plots [i](#)

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 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.

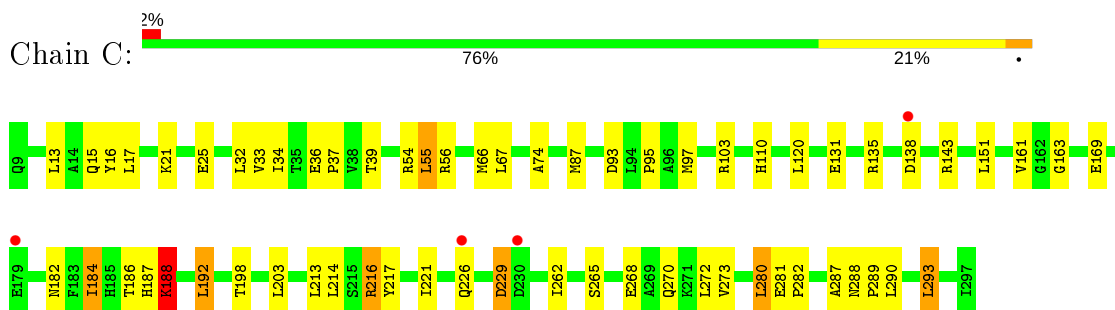
- Molecule 1: Geranylgeranyl pyrophosphate synthase



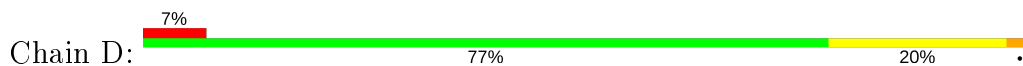
- Molecule 2: Geranylgeranyl pyrophosphate synthase

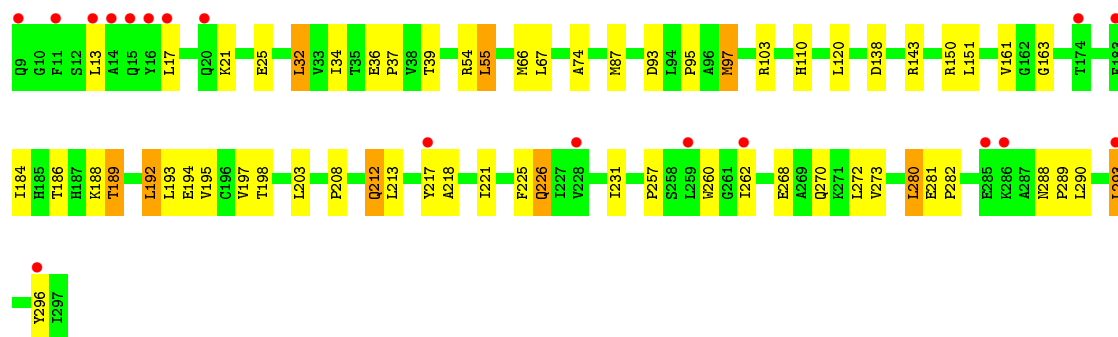


- Molecule 3: Geranylgeranyl pyrophosphate synthase

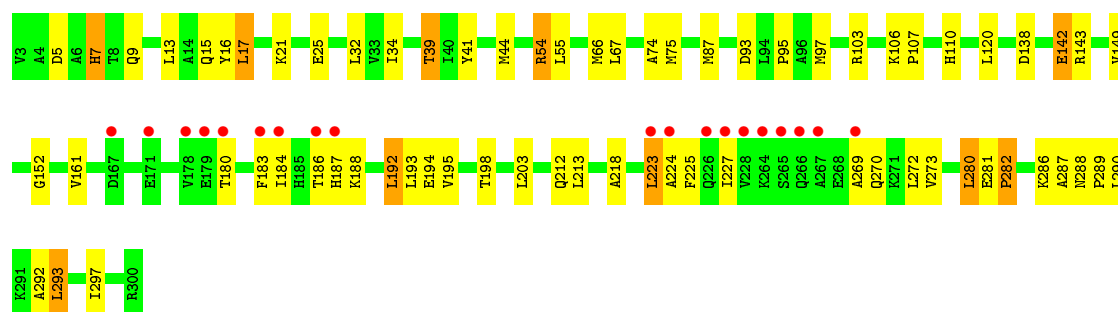
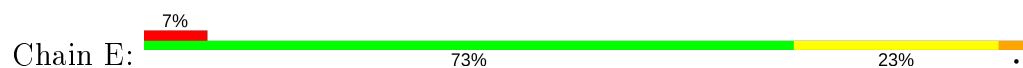


- Molecule 4: Geranylgeranyl pyrophosphate synthase

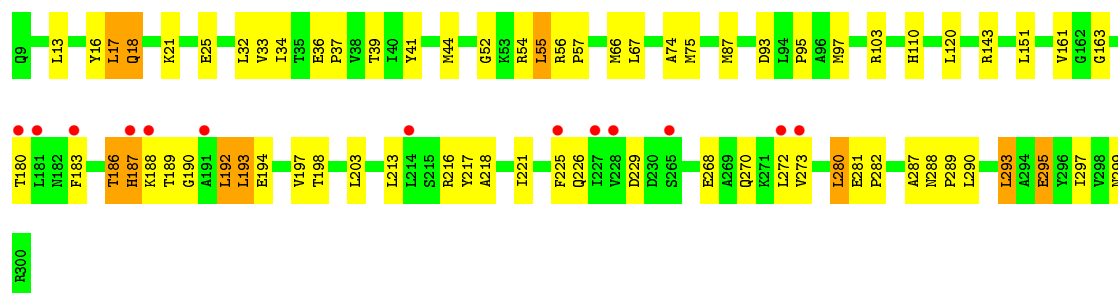




- Molecule 5: Geranylgeranyl pyrophosphate synthase



- Molecule 6: Geranylgeranyl pyrophosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.56Å 122.97Å 134.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.02 – 2.66 67.92 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.2 (68.02-2.66) 99.3 (67.92-2.66)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 2.65Å)	Xtrriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.228 , 0.285 0.230 , 0.285	Depositor DCC
R_{free} test set	2407 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	57.0	Xtrriage
Anisotropy	0.178	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11367	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.56% 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	0.49	0/1954	0.78	0/2656
2	B	0.48	0/1906	0.78	0/2589
3	C	0.52	1/1900 (0.1%)	0.83	2/2581 (0.1%)
4	D	0.45	0/1943	0.79	1/2644 (0.0%)
5	E	0.50	0/1882	0.83	2/2559 (0.1%)
6	F	0.48	0/1875	0.79	0/2547
All	All	0.49	1/11460 (0.0%)	0.80	5/15576 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	188	LYS	C-O	5.75	1.34	1.23

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	216	ARG	CG-CD-NE	10.49	133.82	111.80
5	E	186	THR	CB-CA-C	7.30	131.31	111.60
4	D	212	GLN	CB-CA-C	5.62	121.65	110.40
3	C	229	ASP	CB-CA-C	5.32	121.04	110.40
5	E	212	GLN	CB-CA-C	5.16	120.73	110.40

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	1931	0	1891	53	0
2	B	1884	0	1853	49	0
3	C	1877	0	1849	47	0
4	D	1918	0	1867	53	0
5	E	1859	0	1816	52	3
6	F	1852	0	1839	56	3
7	A	11	0	0	4	0
7	B	12	0	0	1	0
7	C	5	0	0	1	0
7	D	2	0	0	1	0
7	E	9	0	0	2	0
7	F	7	0	0	1	0
All	All	11367	0	11115	289	3

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 13.

The worst 5 of 289 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)
2:B:36:GLU:HB3	2:B:37:PRO:CD	1.52	1.30
2:B:36:GLU:HB3	2:B:37:PRO:HD2	1.28	1.14
4:D:66:MET:HE3	4:D:290:LEU:HD12	1.21	1.14
6:F:186:THR:CG2	6:F:226:GLN:NE2	2.11	1.13
6:F:186:THR:HG23	6:F:226:GLN:NE2	1.64	1.10

All (3) 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)
5:E:180:THR:CB	6:F:187:HIS:NE2[1_655]	1.70	0.50
5:E:180:THR:N	6:F:187:HIS:NE2[1_655]	2.12	0.08
5:E:180:THR:CA	6:F:187:HIS:NE2[1_655]	2.18	0.02

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	259/267 (97%)	249 (96%)	10 (4%)	0	100	100
2	B	251/259 (97%)	237 (94%)	12 (5%)	2 (1%)	19	29
3	C	252/258 (98%)	240 (95%)	11 (4%)	1 (0%)	34	48
4	D	263/267 (98%)	247 (94%)	14 (5%)	2 (1%)	19	29
5	E	253/259 (98%)	237 (94%)	12 (5%)	4 (2%)	9	14
6	F	247/253 (98%)	239 (97%)	8 (3%)	0	100	100
All	All	1525/1563 (98%)	1449 (95%)	67 (4%)	9 (1%)	25	37

5 of 9 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	36	GLU
3	C	262	ILE
4	D	262	ILE
2	B	296	TYR
4	D	231	ILE

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	187/211 (89%)	164 (88%)	23 (12%)	4	6
2	B	184/205 (90%)	164 (89%)	20 (11%)	6	9
3	C	184/203 (91%)	160 (87%)	24 (13%)	4	5
4	D	183/211 (87%)	164 (90%)	19 (10%)	7	10
5	E	176/203 (87%)	156 (89%)	20 (11%)	5	8
6	F	184/199 (92%)	159 (86%)	25 (14%)	3	5
All	All	1098/1232 (89%)	967 (88%)	131 (12%)	5	7

5 of 131 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	C	198	THR
4	D	87	MET
6	F	189	THR
3	C	216	ARG
3	C	293	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
4	D	153	GLN
5	E	7	HIS
6	F	226	GLN
4	D	164	GLN
4	D	288	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](#)

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	B	3
1	A	3
5	E	2
3	C	2
6	F	2
4	D	1

The worst 5 of 13 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	235:THR	C	254:VAL	N	15.68
1	D	232:LEU	C	255:THR	N	14.53
1	C	232:LEU	C	259:LEU	N	12.09
1	A	257:PRO	C	264:LYS	N	12.08
1	E	228:VAL	C	264:LYS	N	11.77

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	267/267 (100%)	0.09	7 (2%) 56 52	32, 61, 127, 191	0
2	B	259/259 (100%)	0.04	3 (1%) 79 77	29, 55, 109, 156	0
3	C	258/258 (100%)	0.04	4 (1%) 72 69	30, 52, 119, 165	0
4	D	267/267 (100%)	0.38	18 (6%) 17 14	41, 79, 137, 178	0
5	E	259/259 (100%)	0.24	19 (7%) 15 12	31, 59, 158, 239	0
6	F	253/253 (100%)	0.25	13 (5%) 28 25	34, 64, 144, 192	0
All	All	1563/1563 (100%)	0.17	64 (4%) 37 33	29, 63, 132, 239	0

The worst 5 of 64 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
5	E	183	PHE	6.4
4	D	9	GLN	6.1
6	F	228	VAL	5.6
5	E	184	ILE	5.2
5	E	226	GLN	5.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.