



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 07:09 AM GMT

PDB ID : 2E90
Title : S. cerevisiae geranylgeranyl pyrophosphate synthase in complex with magnesium, pyrophosphate and FPP
Authors : Guo, R.T.; Ko, T.P.; Chen, C.K.-M.; Jeng, W.Y.; Chang, T.H.; Liang, P.H.; Oldfield, E.; Wang, A.H.-J.
Deposited on : 2007-01-24
Resolution : 2.55 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

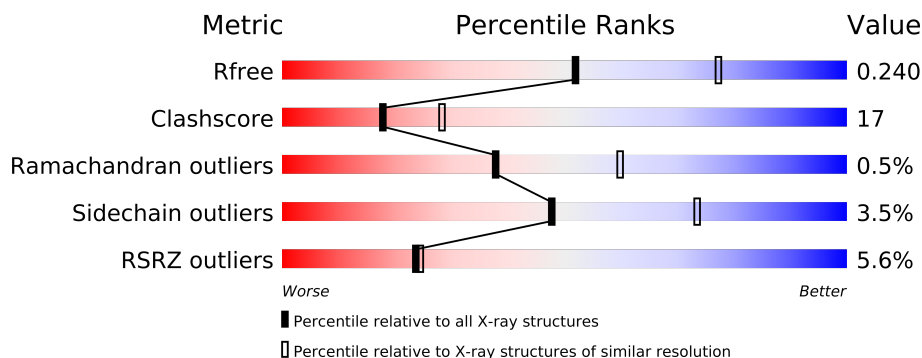
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.55 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}	66092	3413 (2.58-2.50)
Clashscore	79885	4284 (2.58-2.50)
Ramachandran outliers	78287	4193 (2.58-2.50)
Sidechain outliers	78261	4195 (2.58-2.50)
RSRZ outliers	66119	3414 (2.58-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	340	
1	B	340	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MG	A	1201	-	X
2	MG	A	1202	-	X
2	MG	B	1203	-	X
2	MG	B	1204	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5313 atoms, of which 0 are hydrogen and 0 are deuterium.

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 synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	0	0
			2440	1562	413	455	10			
1	B	304	Total	C	N	O	S	0	0	0
			2485	1595	423	457	10			

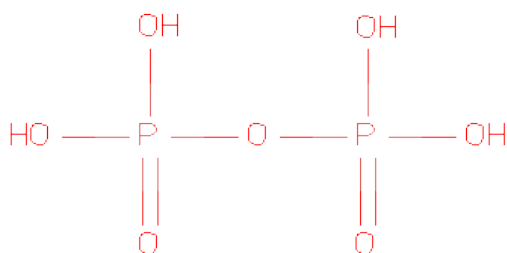
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	CLONING ARTIFACT	UNP Q12051
A	2	THR	-	CLONING ARTIFACT	UNP Q12051
A	3	LYS	-	CLONING ARTIFACT	UNP Q12051
A	4	ASN	-	CLONING ARTIFACT	UNP Q12051
A	5	LYS	-	CLONING ARTIFACT	UNP Q12051
B	1	MET	-	CLONING ARTIFACT	UNP Q12051
B	2	THR	-	CLONING ARTIFACT	UNP Q12051
B	3	LYS	-	CLONING ARTIFACT	UNP Q12051
B	4	ASN	-	CLONING ARTIFACT	UNP Q12051
B	5	LYS	-	CLONING ARTIFACT	UNP Q12051

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

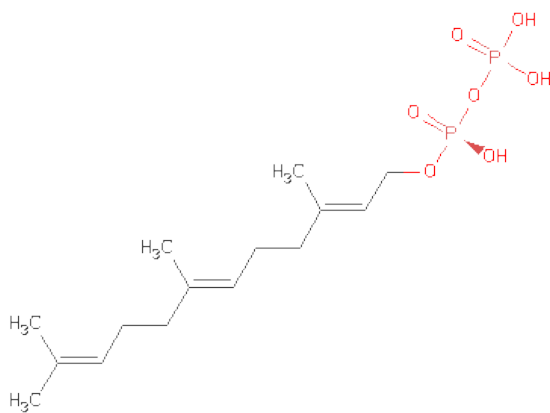
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		

- Molecule 3 is PYROPHOSPHATE (three-letter code: PPV) (formula: H₄O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			9	7	2		
3	B	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is FARNESYL DIPHOSPHATE (three-letter code: FPP) (formula: $C_{15}H_{28}O_7P_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	P	
			24	15	7	2	0
4	B	1	Total	C	O	P	
			24	15	7	2	0

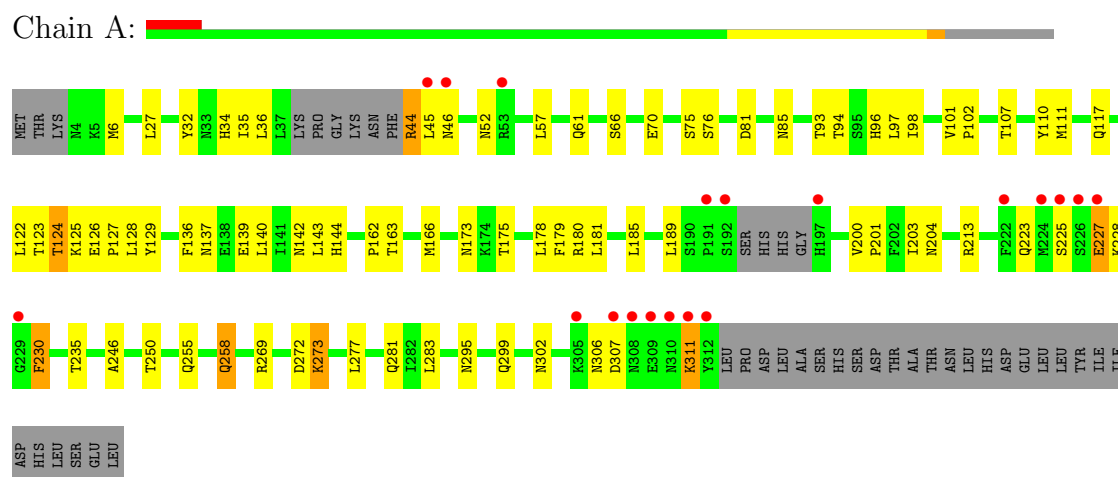
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	168	Total 168	O 168	0	0
5	B	150	Total 150	O 150	0	0

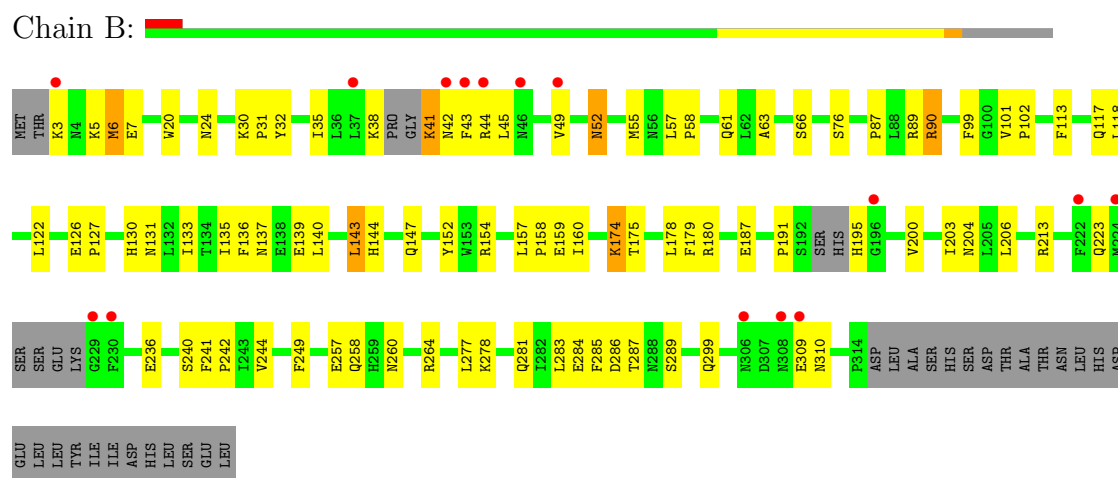
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 synthetase



- Molecule 1: Geranylgeranyl pyrophosphate synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.77Å 115.94Å 126.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.14 – 2.55 34.14 – 2.56	Depositor EDS
% Data completeness (in resolution range)	93.5 (34.14-2.55) 94.1 (34.14-2.56)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.54Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.185 , 0.239 0.185 , 0.240	Depositor DCC
R_{free} test set	1040 reflections (4.89%)	DCC
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.424	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 41.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 21259 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5313	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 5.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, FPP, PPV

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	0.53	0/2482	0.67	0/3358
1	B	0.52	0/2529	0.67	0/3419
All	All	0.52	0/5011	0.67	0/6777

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2440	0	2483	94	0
1	B	2485	0	2535	87	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	9	0	0	0	0
3	B	9	0	0	1	0
4	A	24	0	25	4	0
4	B	24	0	25	14	0
5	A	168	0	0	5	0
5	B	150	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5313	0	5068	173	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 17.

All (173) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:299:GLN:HE21	1:B:3:LYS:HD3	1.17	1.09
1:A:163:THR:H	1:A:166:MET:HE3	1.19	1.08
1:B:143:LEU:HD21	4:B:1102:FPP:H92	1.47	0.96
1:B:143:LEU:HD11	4:B:1102:FPP:H7	1.47	0.96
1:A:52:ASN:HD21	1:A:57:LEU:H	1.09	0.93
1:A:123:THR:HG22	1:A:125:LYS:H	1.31	0.93
1:B:180:ARG:HH11	1:B:204:ASN:HD21	1.17	0.92
1:B:52:ASN:HD21	1:B:57:LEU:H	1.14	0.90
1:A:258:GLN:HE21	1:A:258:GLN:N	1.69	0.90
1:B:213:ARG:HD3	5:B:559:HOH:O	1.75	0.87
1:A:123:THR:HG21	1:A:128:LEU:HD23	1.54	0.87
1:A:258:GLN:NE2	1:A:258:GLN:H	1.73	0.86
1:A:258:GLN:HE21	1:A:258:GLN:H	0.88	0.83
1:B:55:MET:HE1	1:B:187:GLU:HG2	1.61	0.80
1:A:273:LYS:H	1:A:273:LYS:CD	1.92	0.80
1:A:272:ASP:HA	1:A:273:LYS:HE3	1.62	0.80
1:B:38:LYS:HE2	3:B:1002:PPV:O12	1.82	0.79
1:A:52:ASN:ND2	1:A:57:LEU:H	1.81	0.78
1:B:180:ARG:HH11	1:B:204:ASN:ND2	1.82	0.76
1:A:163:THR:N	1:A:166:MET:HE3	1.97	0.76
1:B:143:LEU:HD21	4:B:1102:FPP:C9	2.16	0.75
1:A:44:ARG:O	1:A:46:ASN:N	2.20	0.75
1:A:299:GLN:NE2	1:B:3:LYS:HD3	1.98	0.74
1:B:52:ASN:ND2	1:B:57:LEU:H	1.86	0.72
1:B:147:GLN:HA	1:B:147:GLN:HE21	1.55	0.71
1:A:52:ASN:HD21	1:A:57:LEU:N	1.88	0.71
1:B:89:ARG:HH11	1:B:90:ARG:HH21	1.38	0.71
1:B:159:GLU:HG2	1:B:160:ILE:HG13	1.73	0.70
1:B:157:LEU:HD12	1:B:158:PRO:HA	1.73	0.70
1:A:123:THR:CG2	1:A:128:LEU:HD23	2.22	0.69
1:A:175:THR:HG22	1:A:179:PHE:HE2	1.57	0.69
1:A:225:SER:HB3	1:A:228:LYS:HD3	1.75	0.69
1:B:55:MET:HE2	1:B:187:GLU:HA	1.77	0.67
1:A:6:MET:HE1	1:B:299:GLN:CB	2.25	0.66
1:B:258:GLN:HG2	1:B:286:ASP:OD2	1.95	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:126:GLU:HA	1:B:126:GLU:OE2	1.95	0.66
1:A:6:MET:HE1	1:B:299:GLN:HB2	1.78	0.66
1:A:273:LYS:H	1:A:273:LYS:CE	2.09	0.65
1:B:41:LYS:HE2	5:B:688:HOH:O	1.96	0.65
1:B:143:LEU:HD11	4:B:1102:FPP:H12	1.79	0.65
1:A:273:LYS:H	1:A:273:LYS:HE3	1.61	0.65
1:B:45:LEU:HD11	5:B:462:HOH:O	1.98	0.64
1:A:299:GLN:CB	1:B:6:MET:HE3	2.28	0.64
4:B:1102:FPP:H7	4:B:1102:FPP:H12	1.80	0.64
1:A:180:ARG:HH11	1:A:204:ASN:HD21	1.43	0.64
1:B:143:LEU:HD13	4:B:1102:FPP:H142	1.79	0.64
1:B:174:LYS:HE3	4:B:1102:FPP:H42	1.80	0.64
1:A:44:ARG:NH1	5:A:424:HOH:O	2.31	0.63
1:B:41:LYS:HD3	1:B:43:PHE:HB2	1.79	0.63
1:A:307:ASP:OD1	1:A:311:LYS:HG2	1.98	0.63
1:B:140:LEU:O	1:B:144:HIS:HD2	1.82	0.62
1:A:225:SER:C	1:A:227:GLU:H	2.02	0.62
1:A:66:SER:O	1:A:70:GLU:HG2	2.01	0.61
1:B:147:GLN:NE2	1:B:147:GLN:HA	2.14	0.61
1:A:123:THR:HG22	1:A:124:THR:H	1.65	0.61
1:A:94:THR:HG23	5:A:515:HOH:O	2.01	0.61
1:B:41:LYS:HG2	1:B:43:PHE:H	1.66	0.60
1:B:52:ASN:HD21	1:B:57:LEU:N	1.93	0.60
1:B:89:ARG:HH11	1:B:90:ARG:NH2	1.97	0.60
1:A:225:SER:HB3	1:A:228:LYS:CG	2.33	0.59
1:A:122:LEU:HD13	1:A:189:LEU:HD11	1.82	0.59
1:A:123:THR:HG22	1:A:124:THR:N	2.18	0.59
4:B:1102:FPP:C7	4:B:1102:FPP:H12	2.32	0.58
1:A:81:ASP:HB3	1:A:94:THR:HG21	1.86	0.57
1:B:143:LEU:HD21	4:B:1102:FPP:C11	2.35	0.57
1:A:162:PRO:HA	1:A:166:MET:HE1	1.86	0.57
1:B:55:MET:CE	1:B:187:GLU:HA	2.34	0.56
1:A:27:LEU:HD22	1:B:152:TYR:CZ	2.41	0.56
1:A:139:GLU:HA	1:A:139:GLU:OE2	2.05	0.56
1:B:45:LEU:O	1:B:49:VAL:HG23	2.06	0.56
1:A:175:THR:O	1:A:179:PHE:HD2	1.89	0.56
1:B:38:LYS:NZ	1:B:38:LYS:HB3	2.20	0.55
1:A:180:ARG:HD3	1:A:204:ASN:HD21	1.72	0.55
1:B:240:SER:O	1:B:244:VAL:HG23	2.07	0.55
1:B:147:GLN:CA	1:B:147:GLN:HE21	2.16	0.55
1:B:3:LYS:O	1:B:7:GLU:HG3	2.07	0.55
1:A:180:ARG:HH11	1:A:204:ASN:ND2	2.05	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:225:SER:HB3	1:A:228:LYS:CD	2.37	0.54
1:A:302:ASN:O	1:A:306:ASN:HB2	2.09	0.53
1:B:41:LYS:HG2	1:B:42:ASN:N	2.24	0.53
1:B:101:VAL:HB	1:B:102:PRO:HD3	1.91	0.53
1:A:250:THR:HG22	1:A:255:GLN:NE2	2.23	0.53
1:A:227:GLU:HG3	5:A:563:HOH:O	2.09	0.52
1:B:203:ILE:HD12	1:B:206:LEU:HD23	1.92	0.52
1:A:299:GLN:HB3	1:B:6:MET:HE3	1.91	0.52
1:B:249:PHE:CD2	1:B:287:THR:HG22	2.45	0.52
1:B:139:GLU:HA	1:B:139:GLU:OE2	2.10	0.51
1:A:94:THR:CG2	1:A:96:HIS:CD2	2.94	0.51
1:A:299:GLN:HB2	1:B:6:MET:HE3	1.93	0.51
1:A:140:LEU:HD13	1:B:113:PHE:CE2	2.46	0.51
1:B:87:PRO:HD2	5:B:710:HOH:O	2.11	0.51
1:A:162:PRO:HA	1:A:166:MET:CE	2.41	0.50
1:A:75:SER:OG	1:A:111:MET:HB2	2.11	0.50
1:B:20:TRP:CH2	1:B:24:ASN:HB3	2.47	0.50
1:A:140:LEU:HA	1:A:143:LEU:HD12	1.92	0.50
1:B:131:ASN:O	1:B:135:ILE:HG13	2.12	0.50
1:A:273:LYS:N	1:A:273:LYS:CD	2.70	0.50
1:A:273:LYS:H	1:A:273:LYS:HD3	1.73	0.49
1:B:200:VAL:O	1:B:203:ILE:HG22	2.12	0.49
1:A:101:VAL:N	1:A:102:PRO:CD	2.75	0.49
1:B:143:LEU:HD23	1:B:144:HIS:CD2	2.48	0.49
1:A:230:PHE:N	1:A:230:PHE:CD2	2.81	0.48
1:A:225:SER:CB	1:A:228:LYS:HD3	2.41	0.48
1:B:32:TYR:O	1:B:35:ILE:HG22	2.14	0.48
1:B:55:MET:O	1:B:191:PRO:HD2	2.13	0.48
1:B:143:LEU:HD11	4:B:1102:FPP:C12	2.43	0.48
1:B:61:GLN:NE2	1:B:122:LEU:O	2.46	0.47
1:A:295:ASN:HD21	1:B:3:LYS:NZ	2.12	0.47
1:A:93:THR:HB	1:A:98:ILE:HD11	1.95	0.47
1:A:213:ARG:NH1	5:A:698:HOH:O	2.47	0.47
1:B:30:LYS:HB3	1:B:99:PHE:CE2	2.48	0.47
1:A:94:THR:HG21	1:A:96:HIS:CD2	2.49	0.47
1:A:277:LEU:O	1:A:281:GLN:HG2	2.13	0.47
1:A:126:GLU:HB2	1:A:127:PRO:HD3	1.96	0.47
1:A:163:THR:HG23	1:A:166:MET:CE	2.45	0.47
1:A:175:THR:HG22	1:A:179:PHE:CE2	2.43	0.47
1:B:278:LYS:O	1:B:281:GLN:HG2	2.14	0.47
1:A:200:VAL:HB	1:A:201:PRO:HD3	1.96	0.47
1:B:63:ALA:O	1:B:66:SER:HB3	2.14	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:163:THR:H	1:A:166:MET:CE	2.07	0.46
1:A:136:PHE:CE2	1:A:140:LEU:HD12	2.50	0.46
1:B:58:PRO:HG2	1:B:61:GLN:HB2	1.97	0.46
4:B:1102:FPP:O2B	4:B:1102:FPP:C1	2.64	0.46
1:B:241:PHE:HB3	1:B:242:PRO:HD3	1.96	0.46
1:A:181:LEU:HD12	1:A:181:LEU:O	2.16	0.46
1:A:137:ASN:HD21	1:B:117:GLN:CA	2.29	0.46
1:A:143:LEU:HD13	4:A:1101:FPP:H142	1.98	0.45
1:A:140:LEU:O	1:A:144:HIS:HD2	1.99	0.45
1:B:260:ASN:O	1:B:264:ARG:HG3	2.16	0.45
1:A:101:VAL:HB	1:A:102:PRO:HD3	1.99	0.45
1:A:129:TYR:CE2	1:B:130:HIS:CD2	3.05	0.45
1:A:142:ASN:O	1:A:173:ASN:HB3	2.17	0.45
1:A:225:SER:C	1:A:227:GLU:N	2.69	0.45
1:B:154:ARG:HD2	1:B:236:GLU:O	2.16	0.45
1:B:175:THR:O	1:B:179:PHE:HD2	2.00	0.45
1:B:76:SER:CB	4:B:1102:FPP:C12	2.95	0.45
1:A:94:THR:CG2	1:A:96:HIS:HD2	2.29	0.44
1:A:223:GLN:HE21	1:A:281:GLN:HE22	1.65	0.44
1:B:133:ILE:HD13	1:B:133:ILE:HA	1.85	0.44
1:B:5:LYS:HA	1:B:5:LYS:HD2	1.67	0.44
1:A:32:TYR:O	1:A:35:ILE:HG22	2.17	0.44
1:A:117:GLN:N	1:B:137:ASN:HD21	2.16	0.44
1:B:38:LYS:HZ2	1:B:38:LYS:HB3	1.81	0.44
1:A:225:SER:HB3	1:A:228:LYS:HG3	1.98	0.43
1:B:30:LYS:HB2	1:B:31:PRO:HD3	2.00	0.43
1:B:76:SER:HB3	4:B:1102:FPP:C12	2.48	0.43
1:B:118:LEU:HD23	1:B:118:LEU:HA	1.80	0.43
1:B:41:LYS:HD2	1:B:44:ARG:N	2.34	0.43
1:A:181:LEU:HD12	1:A:185:LEU:HG	2.00	0.43
1:A:235:THR:O	1:A:269:ARG:HA	2.19	0.43
1:B:136:PHE:CE2	1:B:140:LEU:HD12	2.53	0.43
1:B:283:LEU:HG	1:B:289:SER:HB2	2.01	0.43
1:A:140:LEU:HD21	4:A:1101:FPP:H153	2.01	0.42
4:A:1101:FPP:H12	4:A:1101:FPP:C7	2.50	0.42
1:B:309:GLU:O	1:B:310:ASN:HB2	2.19	0.42
1:A:246:ALA:HB2	1:A:283:LEU:HD22	2.02	0.42
1:A:94:THR:HG22	5:A:701:HOH:O	2.19	0.42
1:A:203:ILE:HD12	1:A:203:ILE:HA	1.88	0.42
1:B:140:LEU:O	1:B:144:HIS:CD2	2.67	0.42
1:B:204:ASN:HA	1:B:204:ASN:HD22	1.66	0.41
1:A:34:HIS:HE1	1:A:93:THR:O	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:B:1102:FPP:H11	4:B:1102:FPP:O2B	2.19	0.41
1:A:302:ASN:HB3	1:A:306:ASN:HD22	1.85	0.41
1:B:180:ARG:NH1	1:B:204:ASN:ND2	2.58	0.41
1:B:52:ASN:HA	1:B:52:ASN:HD22	1.63	0.41
1:B:223:GLN:HE22	1:B:284:GLU:CG	2.34	0.41
1:A:36:LEU:HD23	1:A:36:LEU:HA	1.87	0.41
1:A:97:LEU:HD23	1:A:97:LEU:HA	1.93	0.40
1:A:281:GLN:HA	1:A:281:GLN:NE2	2.37	0.40
1:A:107:THR:O	1:A:110:TYR:HB3	2.21	0.40
1:A:273:LYS:HD3	1:A:273:LYS:N	2.35	0.40
1:A:140:LEU:CD2	4:A:1101:FPP:H153	2.51	0.40
1:B:285:PHE:N	1:B:285:PHE:CD2	2.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	293/340 (86%)	279 (95%)	13 (4%)	1 (0%)	50	71
1	B	296/340 (87%)	286 (97%)	8 (3%)	2 (1%)	30	49
All	All	589/680 (87%)	565 (96%)	21 (4%)	3 (0%)	38	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	B	174	LYS
1	B	127	PRO

5.3.2 Protein sidechains ⓘ

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	280/317 (88%)	269 (96%)	11 (4%)	43	69
1	B	284/317 (90%)	275 (97%)	9 (3%)	51	77
All	All	564/634 (89%)	544 (96%)	20 (4%)	48	73

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ARG
1	A	61	GLN
1	A	76	SER
1	A	85	ASN
1	A	124	THR
1	A	178	LEU
1	A	227	GLU
1	A	230	PHE
1	A	258	GLN
1	A	273	LYS
1	A	311	LYS
1	B	6	MET
1	B	41	LYS
1	B	52	ASN
1	B	90	ARG
1	B	143	LEU
1	B	178	LEU
1	B	195	HIS
1	B	257	GLU
1	B	277	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	HIS
1	A	52	ASN
1	A	61	GLN
1	A	85	ASN
1	A	92	GLN
1	A	96	HIS
1	A	109	ASN

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Mol	Chain	Res	Type
1	A	137	ASN
1	A	144	HIS
1	A	204	ASN
1	A	211	GLN
1	A	223	GLN
1	A	255	GLN
1	A	258	GLN
1	A	260	ASN
1	A	281	GLN
1	A	295	ASN
1	A	298	ASN
1	A	299	GLN
1	A	306	ASN
1	B	33	ASN
1	B	52	ASN
1	B	67	GLN
1	B	92	GLN
1	B	96	HIS
1	B	106	ASN
1	B	109	ASN
1	B	121	GLN
1	B	130	HIS
1	B	137	ASN
1	B	144	HIS
1	B	147	GLN
1	B	195	HIS
1	B	204	ASN
1	B	211	GLN
1	B	223	GLN
1	B	255	GLN
1	B	281	GLN
1	B	288	ASN
1	B	298	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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
3	PPV	A	1001	-	8,8,8	1.80	1 (12%)	13,13,13	2.13	1 (7%)
4	FPP	A	1101	2	23,23,23	2.05	8 (34%)	31,31,31	1.36	5 (16%)
3	PPV	B	1002	-	8,8,8	1.99	3 (37%)	13,13,13	2.12	1 (7%)
4	FPP	B	1102	2	23,23,23	2.05	8 (34%)	31,31,31	1.30	4 (12%)

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
3	PPV	A	1001	-	-	0/6/6/6	0/0/0/0
4	FPP	A	1101	2	-	0/25/25/25	0/0/0/0
3	PPV	B	1002	-	-	0/6/6/6	0/0/0/0
4	FPP	B	1102	2	-	0/25/25/25	0/0/0/0

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1101	FPP	C6-C7	-3.84	1.39	1.50
4	A	1101	FPP	C11-C12	-3.83	1.39	1.50
4	B	1102	FPP	C6-C7	-3.82	1.39	1.50
4	B	1102	FPP	C11-C12	-3.81	1.39	1.50
4	A	1101	FPP	C7-C8	3.28	1.39	1.32
4	B	1102	FPP	C7-C8	3.27	1.39	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1101	FPP	C2-C3	3.19	1.39	1.32
4	B	1102	FPP	C2-C3	3.19	1.39	1.32
4	A	1101	FPP	PB-O1B	3.09	1.61	1.51
4	B	1102	FPP	PB-O1B	3.05	1.61	1.51
3	A	1001	PPV	P2-O22	3.04	1.61	1.51
3	B	1002	PPV	P2-O22	2.96	1.61	1.51
4	A	1101	FPP	C1-C2	-2.85	1.39	1.49
4	B	1102	FPP	C1-C2	-2.83	1.39	1.49
4	B	1102	FPP	PA-O1A	2.66	1.61	1.51
4	A	1101	FPP	PA-O1A	2.65	1.61	1.51
4	B	1102	FPP	C12-C13	2.23	1.39	1.32
4	A	1101	FPP	C12-C13	2.23	1.39	1.32
3	B	1002	PPV	P1-OPP	2.19	1.64	1.60
3	B	1002	PPV	P1-O21	2.03	1.62	1.54

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1002	PPV	P2-OPP-P1	-7.38	110.04	131.68
3	A	1001	PPV	P2-OPP-P1	-7.34	110.15	131.68
4	B	1102	FPP	PA-O3A-PB	-3.24	122.18	131.68
4	A	1101	FPP	C10-C8-C9	3.22	120.28	115.39
4	A	1101	FPP	PA-O3A-PB	-3.06	122.70	131.68
4	B	1102	FPP	C4-C3-C5	3.04	120.01	115.39
4	A	1101	FPP	C4-C3-C5	3.02	119.98	115.39
4	B	1102	FPP	C10-C8-C9	2.76	119.58	115.39
4	B	1102	FPP	C15-C13-C14	2.15	120.11	114.62
4	A	1101	FPP	O2A-PA-O3A	2.14	115.28	105.14
4	A	1101	FPP	C15-C13-C14	2.12	120.03	114.62

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	299/340 (87%)	-0.05	19 (6%) 19 19	16, 30, 80, 105	0
1	B	304/340 (89%)	-0.04	15 (4%) 28 30	15, 34, 78, 106	0
All	All	603/680 (88%)	-0.05	34 (5%) 24 24	15, 32, 78, 106	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	308	ASN	6.6
1	A	226	SER	6.5
1	A	225	SER	6.2
1	A	227	GLU	5.5
1	A	224	MET	5.2
1	B	3	LYS	4.4
1	A	309	GLU	4.3
1	B	42	ASN	4.2
1	A	312	TYR	4.0
1	B	229	GLY	3.8
1	B	196	GLY	3.8
1	A	310	ASN	3.7
1	B	43	PHE	3.7
1	A	311	LYS	3.3
1	B	224	MET	3.1
1	A	229	GLY	3.0
1	B	222	PHE	3.0
1	A	305	LYS	2.9
1	B	230	PHE	2.8
1	A	307	ASP	2.6
1	A	222	PHE	2.6
1	B	49	VAL	2.5
1	B	309	GLU	2.5
1	A	53	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	191	PRO	2.4
1	A	46	ASN	2.3
1	A	192	SER	2.3
1	B	46	ASN	2.2
1	B	308	ASN	2.2
1	B	37	LEU	2.2
1	B	306	ASN	2.1
1	A	45	LEU	2.1
1	B	44	ARG	2.0
1	A	197	HIS	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MG	A	1201	1/1	0.23	5.00	7,7,7,7	0
2	MG	A	1202	1/1	0.17	4.58	22,22,22,22	0
2	MG	B	1203	1/1	0.24	4.46	19,19,19,19	0
2	MG	B	1204	1/1	0.30	2.50	9,9,9,9	0
4	FPP	B	1102	24/24	0.24	1.42	41,44,47,47	0
3	PPV	A	1001	9/9	0.18	0.91	86,87,93,93	0
4	FPP	A	1101	24/24	0.17	0.88	36,40,44,46	0
3	PPV	B	1002	9/9	0.19	-0.38	101,102,105,105	0

6.5 Other polymers ⓘ

There are no such residues in this entry.