



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 07:09 AM GMT

PDB ID : 2E92
Title : S. cerevisiae geranylgeranyl pyrophosphate synthase in complex with magnesium and BPH-261
Authors : Guo, R.T.; Ko, T.P.; Cao, R.; Jeng, W.Y.; Chen, C.K.-M.; Chang, T.H.; Liang, P.H.; Oldfield, E.; Wang, A.H.-J.
Deposited on : 2007-01-24
Resolution : 2.31 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>

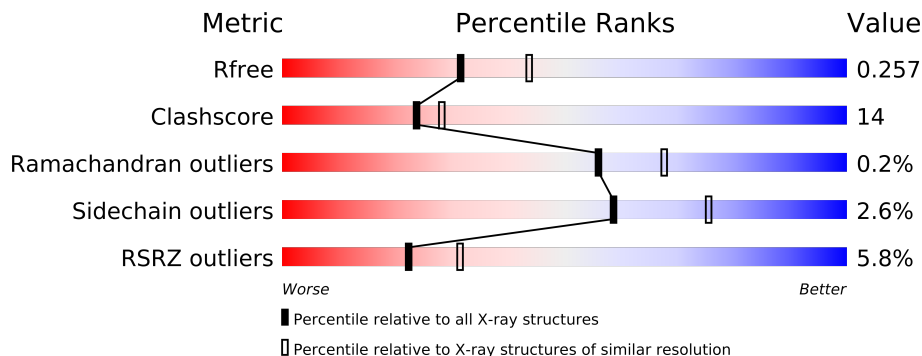
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.31 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	3293 (2.34-2.30)
Clashscore	79885	4097 (2.34-2.30)
Ramachandran outliers	78287	4055 (2.34-2.30)
Sidechain outliers	78261	4054 (2.34-2.30)
RSRZ outliers	66119	3294 (2.34-2.30)

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	1302	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5111 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	291	Total	C	N	O	S	0	0	0
			2373	1524	400	439	10			
1	B	290	Total	C	N	O	S	0	0	0
			2368	1523	400	435	10			

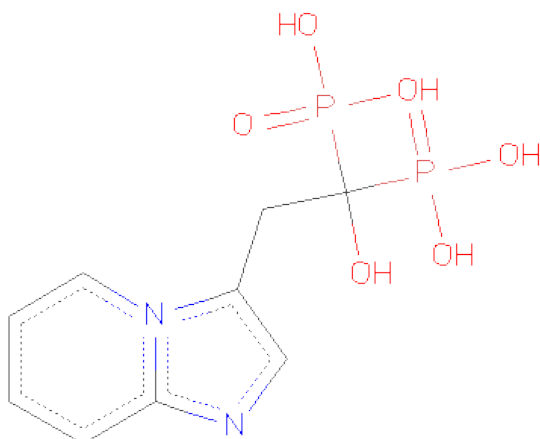
There are 10 discrepancies between the modelled and reference sequences:

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 (1-HYDROXY-2-IMIDAZO[1,2-A]PYRIDIN-3-YLETHANE-1,1-DIYL)BIS(P HOSPHONICACID) (three-letter code: M0N) (formula: C₉H₁₂N₂O₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			20	9	2	7	2		
3	B	1	Total	C	N	O	P	0	0
			20	9	2	7	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	150	Total	O	0	0
			150	150		
4	B	176	Total	O	0	0
			176	176		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.87Å 116.05Å 127.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.31 29.05 – 2.31	Depositor EDS
% Data completeness (in resolution range)	91.4 (30.00-2.31) 91.3 (29.05-2.31)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.31Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.199 , 0.257 0.199 , 0.257	Depositor DCC
R_{free} test set	1397 reflections (4.87%)	DCC
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.732	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 28714 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5111	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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: M0N, MG

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	0.69	0/2413	0.77	2/3265 (0.1%)
1	B	0.67	0/2408	0.74	0/3257
All	All	0.68	0/4821	0.76	2/6522 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	ARG	NE-CZ-NH1	6.21	123.41	120.30
1	A	269	ARG	NE-CZ-NH2	-6.08	117.26	120.30

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	2373	0	2421	74	0
1	B	2368	0	2425	66	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	20	0	8	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	20	0	8	5	0
4	A	150	0	0	3	0
4	B	176	0	0	3	0
All	All	5111	0	4862	138	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 14.

All (138) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:163:THR:H	1:A:166:MET:HE3	0.98	1.09
1:B:143:LEU:HD11	3:B:902:M0N:H5	1.24	1.08
1:A:180:ARG:HH11	1:A:204:ASN:HD21	1.01	0.97
1:A:163:THR:H	1:A:166:MET:CE	1.82	0.93
1:A:139:GLU:HB2	1:A:181:LEU:HD12	1.50	0.92
1:A:258:GLN:HE21	1:A:258:GLN:H	0.92	0.92
1:A:163:THR:N	1:A:166:MET:HE3	1.84	0.91
1:A:123:THR:HG21	1:A:128:LEU:HD23	1.51	0.90
1:A:180:ARG:HH11	1:A:204:ASN:ND2	1.69	0.88
1:A:52:ASN:HD21	1:A:57:LEU:H	1.21	0.86
1:A:258:GLN:HE21	1:A:258:GLN:N	1.73	0.86
1:B:52:ASN:HD21	1:B:57:LEU:H	1.23	0.86
1:A:123:THR:HG22	1:A:125:LYS:H	1.40	0.84
1:A:143:LEU:HD11	3:A:901:M0N:H5	1.59	0.83
1:A:76:SER:HB3	3:A:901:M0N:H4	1.66	0.78
1:B:157:LEU:HD12	1:B:158:PRO:HA	1.64	0.78
1:B:180:ARG:HH11	1:B:204:ASN:HD21	1.31	0.77
1:A:258:GLN:H	1:A:258:GLN:NE2	1.78	0.73
1:A:143:LEU:HD11	3:A:901:M0N:C5	2.20	0.71
1:B:143:LEU:HD11	3:B:902:M0N:C5	2.13	0.71
1:A:94:THR:HG21	1:A:96:HIS:CE1	2.27	0.70
1:B:55:MET:CE	1:B:187:GLU:HA	2.23	0.68
3:A:901:M0N:O2	4:A:723:HOH:O	2.12	0.68
1:B:258:GLN:HG3	1:B:286:ASP:OD2	1.93	0.68
1:A:45:LEU:O	1:A:49:VAL:HG23	1.93	0.68
1:B:143:LEU:CD1	3:B:902:M0N:H5	2.15	0.68
1:B:180:ARG:HH11	1:B:204:ASN:ND2	1.93	0.67
1:B:258:GLN:OE1	1:B:282:ILE:HD13	1.94	0.66
1:A:200:VAL:HB	1:A:201:PRO:HD3	1.78	0.66
3:A:901:M0N:O6	4:A:723:HOH:O	2.14	0.66
1:A:52:ASN:ND2	1:A:57:LEU:H	1.94	0.65
1:A:148:GLY:C	1:B:28:ILE:HD11	2.17	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:52:ASN:ND2	1:B:57:LEU:H	1.96	0.63
1:A:76:SER:HB3	3:A:901:M0N:C4	2.29	0.62
1:B:126:GLU:HA	1:B:126:GLU:OE2	1.99	0.62
1:A:123:THR:HG22	1:A:124:THR:H	1.65	0.62
1:B:55:MET:O	1:B:191:PRO:HD2	1.99	0.61
1:B:52:ASN:HD22	1:B:55:MET:HB2	1.65	0.61
1:A:105:ILE:HD11	1:B:105:ILE:HD11	1.82	0.60
1:B:55:MET:HE2	1:B:187:GLU:HA	1.84	0.59
1:A:123:THR:CG2	1:A:128:LEU:HD23	2.27	0.58
1:A:44:ARG:HG3	1:A:44:ARG:HH11	1.69	0.58
1:A:94:THR:CG2	1:A:96:HIS:CE1	2.87	0.58
1:B:120:SER:HB3	1:B:129:TYR:CE1	2.38	0.57
1:B:42:ASN:N	4:B:691:HOH:O	2.38	0.57
1:A:283:LEU:HB3	1:A:289:SER:HB2	1.86	0.57
1:A:163:THR:HG23	1:A:166:MET:CE	2.37	0.55
1:B:55:MET:HE1	1:B:187:GLU:HA	1.87	0.55
1:A:180:ARG:NH1	1:A:204:ASN:ND2	2.49	0.55
1:A:123:THR:HG22	1:A:124:THR:N	2.22	0.55
1:B:215:ASP:HB3	1:B:240:SER:HB2	1.88	0.54
1:B:119:VAL:HG11	1:B:133:ILE:HD13	1.90	0.53
1:A:246:ALA:HB2	1:A:283:LEU:HD22	1.89	0.53
1:B:180:ARG:HD3	1:B:204:ASN:HD21	1.74	0.53
1:A:145:ARG:HB3	1:B:20:TRP:CZ3	2.44	0.53
1:A:140:LEU:O	1:A:144:HIS:HD2	1.91	0.53
1:A:52:ASN:C	1:A:54:VAL:H	2.12	0.52
1:B:75:SER:OG	1:B:111:MET:HB2	2.09	0.52
1:B:183:LEU:HD11	1:B:199:LEU:HB2	1.90	0.52
1:A:47:LEU:O	1:A:51:ILE:HG13	2.10	0.52
1:A:223:GLN:HE21	1:A:281:GLN:HE22	1.58	0.51
1:B:5:LYS:O	1:B:9:LYS:HG3	2.10	0.51
1:B:230:PHE:O	1:B:276:LYS:NZ	2.42	0.51
1:A:299:GLN:HG2	4:A:569:HOH:O	2.11	0.51
1:B:46:ASN:O	1:B:50:GLN:HG2	2.11	0.51
1:A:180:ARG:O	1:A:184:ARG:HG3	2.10	0.50
1:B:260:ASN:O	1:B:264:ARG:HG3	2.11	0.50
1:A:145:ARG:HD3	1:B:20:TRP:CD2	2.47	0.50
1:A:5:LYS:O	1:A:9:LYS:HG3	2.12	0.50
1:B:47:LEU:O	1:B:51:ILE:HG13	2.12	0.49
1:A:132:LEU:HD11	1:A:189:LEU:HD21	1.93	0.49
1:A:281:GLN:HA	1:A:281:GLN:NE2	2.27	0.49
1:A:126:GLU:N	1:A:127:PRO:CD	2.75	0.49
1:B:128:LEU:O	1:B:132:LEU:HG	2.11	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:32:TYR:CD1	1:B:111:MET:HG3	2.48	0.48
1:A:234:ILE:HG12	1:A:279:LEU:HD23	1.94	0.48
1:A:255:GLN:HB3	1:A:258:GLN:NE2	2.28	0.48
3:B:902:M0N:H1	3:B:902:M0N:O	2.14	0.48
1:A:143:LEU:HD11	3:A:901:M0N:C6	2.44	0.47
1:B:147:GLN:HE21	1:B:174:LYS:HG3	1.78	0.47
1:B:34:HIS:HE1	1:B:93:THR:O	1.97	0.47
1:B:87:PRO:HA	1:B:97:LEU:HD11	1.96	0.47
1:B:116:MET:O	1:B:119:VAL:HB	2.14	0.47
1:B:132:LEU:HD11	1:B:189:LEU:HD21	1.96	0.47
1:A:139:GLU:OE2	1:A:139:GLU:HA	2.15	0.47
1:A:284:GLU:HB2	1:A:290:LEU:HD12	1.97	0.46
1:B:167:TYR:C	1:B:167:TYR:CD2	2.89	0.46
1:A:44:ARG:HG3	1:A:44:ARG:NH1	2.30	0.46
1:B:52:ASN:C	1:B:54:VAL:H	2.18	0.46
1:B:239:LEU:HD22	1:B:239:LEU:N	2.31	0.45
1:B:101:VAL:HB	1:B:102:PRO:HD3	1.97	0.45
1:B:52:ASN:HD21	1:B:57:LEU:N	2.02	0.45
1:A:302:ASN:HB3	1:A:306:ASN:HD22	1.80	0.45
1:B:249:PHE:CD2	1:B:287:THR:HG22	2.52	0.45
1:A:86:ALA:O	1:A:94:THR:HG21	2.17	0.45
1:A:223:GLN:NE2	1:A:281:GLN:HE22	2.14	0.45
1:B:119:VAL:HG11	1:B:133:ILE:CD1	2.47	0.44
1:A:280:ILE:HG22	1:A:281:GLN:HE21	1.81	0.44
1:B:238:LYS:NZ	3:B:902:M0N:O6	2.51	0.44
1:A:52:ASN:HD21	1:A:57:LEU:N	2.01	0.44
1:A:293:THR:O	1:A:297:ILE:HG13	2.17	0.44
1:A:203:ILE:HD12	1:A:203:ILE:HA	1.78	0.44
1:B:233:ASP:CG	1:B:238:LYS:HD3	2.39	0.44
1:A:52:ASN:C	1:A:54:VAL:N	2.70	0.44
1:A:180:ARG:HD3	1:A:204:ASN:HD21	1.83	0.43
1:B:55:MET:CE	1:B:190:SER:HB2	2.48	0.43
3:A:901:M0N:O	3:A:901:M0N:H1	2.17	0.43
1:B:34:HIS:CE1	1:B:93:THR:O	2.71	0.43
1:B:122:LEU:HD13	1:B:189:LEU:HD11	2.01	0.43
1:A:238:LYS:HB2	1:A:238:LYS:HE3	1.81	0.43
1:A:244:VAL:O	1:A:248:ASN:ND2	2.50	0.43
1:B:60:ASP:HB2	4:B:549:HOH:O	2.18	0.43
1:A:245:HIS:HD2	1:A:289:SER:OG	2.02	0.43
1:B:5:LYS:O	1:B:5:LYS:HG2	2.19	0.43
1:A:149:LEU:HD23	1:A:149:LEU:HA	1.90	0.42
1:B:32:TYR:O	1:B:35:ILE:HG22	2.19	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:32:TYR:CD1	1:A:111:MET:HG3	2.55	0.42
1:A:101:VAL:N	1:A:102:PRO:CD	2.83	0.42
1:B:126:GLU:CA	1:B:126:GLU:OE2	2.67	0.42
1:B:249:PHE:HD2	1:B:287:THR:HG22	1.85	0.41
1:B:55:MET:HE2	1:B:190:SER:HB2	2.01	0.41
1:A:92:GLN:HG3	1:A:93:THR:N	2.36	0.41
1:B:282:ILE:O	1:B:286:ASP:HB3	2.20	0.41
1:B:203:ILE:HD12	1:B:203:ILE:HA	1.85	0.41
1:A:148:GLY:CA	1:B:28:ILE:HD11	2.51	0.41
1:A:116:MET:CE	1:B:116:MET:SD	3.08	0.41
1:B:157:LEU:CD1	1:B:158:PRO:HA	2.43	0.41
1:A:302:ASN:HB3	1:A:306:ASN:ND2	2.36	0.41
1:A:89:ARG:O	1:A:90:ARG:HB2	2.20	0.41
1:A:241:PHE:HB3	1:A:242:PRO:HD3	2.01	0.41
1:B:133:ILE:HA	1:B:133:ILE:HD13	1.91	0.40
1:B:200:VAL:O	1:B:203:ILE:HG22	2.21	0.40
1:A:90:ARG:HD3	1:A:90:ARG:HA	1.89	0.40
1:B:42:ASN:ND2	4:B:574:HOH:O	2.55	0.40
1:A:42:ASN:O	1:A:46:ASN:N	2.54	0.40
1:A:60:ASP:O	1:A:63:ALA:HB3	2.21	0.40
1:A:235:THR:O	1:A:269:ARG:HA	2.21	0.40
1:B:154:ARG:HD2	1:B:236:GLU:O	2.21	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	283/340 (83%)	274 (97%)	9 (3%)	0	100	100
1	B	282/340 (83%)	274 (97%)	7 (2%)	1 (0%)	43	52
All	All	565/680 (83%)	548 (97%)	16 (3%)	1 (0%)	56	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	59	LYS

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	272/317 (86%)	267 (98%)	5 (2%)	71	86
1	B	271/317 (86%)	262 (97%)	9 (3%)	50	66
All	All	543/634 (86%)	529 (97%)	14 (3%)	59	75

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ASP
1	A	79	ILE
1	A	94	THR
1	A	158	PRO
1	A	258	GLN
1	B	26	SER
1	B	28	ILE
1	B	52	ASN
1	B	90	ARG
1	B	92	GLN
1	B	147	GLN
1	B	271	SER
1	B	281	GLN
1	B	288	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	61	GLN
1	A	85	ASN
1	A	109	ASN
1	A	121	GLN
1	A	130	HIS

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Mol	Chain	Res	Type
1	A	137	ASN
1	A	144	HIS
1	A	164	GLN
1	A	204	ASN
1	A	211	GLN
1	A	245	HIS
1	A	255	GLN
1	A	258	GLN
1	A	281	GLN
1	A	298	ASN
1	A	306	ASN
1	B	52	ASN
1	B	109	ASN
1	B	121	GLN
1	B	137	ASN
1	B	147	GLN
1	B	204	ASN
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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	M0N	A	901	2	21,21,21	2.85	11 (52%)	34,34,34	1.79	10 (29%)
3	M0N	B	902	2	21,21,21	2.88	10 (47%)	34,34,34	1.70	10 (29%)

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	M0N	A	901	2	-	1/23/23/23	0/0/2/2
3	M0N	B	902	2	-	1/23/23/23	0/0/2/2

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	M0N	P1-O3	6.66	1.61	1.50
3	B	902	M0N	P1-O3	6.65	1.61	1.50
3	B	902	M0N	P2-O6	6.49	1.61	1.50
3	A	901	M0N	P2-O6	6.39	1.61	1.50
3	A	901	M0N	P2-O5	3.52	1.61	1.54
3	B	902	M0N	C8-C9	-3.51	1.53	1.55
3	B	902	M0N	P1-O1	3.47	1.61	1.54
3	B	902	M0N	P2-O5	3.47	1.61	1.54
3	A	901	M0N	C8-C9	-3.35	1.53	1.55
3	A	901	M0N	P1-O1	3.31	1.61	1.54
3	A	901	M0N	P1-O2	-3.29	1.48	1.54
3	A	901	M0N	P2-O4	-3.22	1.48	1.54
3	B	902	M0N	P1-O2	-3.18	1.48	1.54
3	B	902	M0N	P2-O4	-3.12	1.48	1.54
3	A	901	M0N	P2-C9	2.47	1.87	1.85
3	B	902	M0N	C3-C4	2.45	1.39	1.35
3	B	902	M0N	P1-C9	2.40	1.87	1.85
3	A	901	M0N	C3-C4	2.28	1.39	1.35
3	B	902	M0N	P2-C9	2.26	1.86	1.85
3	A	901	M0N	C8-C2	2.17	1.54	1.50
3	A	901	M0N	C1-C2	-2.14	1.32	1.37

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	M0N	N1-C7-N2	-4.22	107.40	113.07
3	B	902	M0N	N1-C7-N2	-3.84	107.91	113.07
3	A	901	M0N	O5-P2-O6	-3.38	105.58	113.14
3	A	901	M0N	O1-P1-O3	-3.32	105.72	113.14
3	B	902	M0N	C9-C8-C2	-3.30	111.30	115.79
3	B	902	M0N	O5-P2-O6	-3.23	105.92	113.14
3	B	902	M0N	O4-P2-C9	2.96	112.94	105.98
3	A	901	M0N	O4-P2-C9	2.95	112.92	105.98
3	A	901	M0N	P1-C9-P2	-2.67	108.06	112.78
3	A	901	M0N	C9-C8-C2	-2.64	112.21	115.79
3	B	902	M0N	C2-N2-C7	-2.57	104.77	108.11
3	A	901	M0N	O2-P1-O1	2.54	115.17	107.93
3	B	902	M0N	O2-P1-C9	2.51	111.88	105.98
3	A	901	M0N	O2-P1-C9	2.47	111.78	105.98
3	B	902	M0N	O6-P2-C9	-2.32	103.95	110.04
3	B	902	M0N	O3-P1-C9	-2.31	103.96	110.04
3	B	902	M0N	O1-P1-O3	-2.22	108.17	113.14
3	A	901	M0N	O4-P2-O5	2.16	114.08	107.93
3	B	902	M0N	O4-P2-O5	2.10	113.91	107.93
3	A	901	M0N	C2-N2-C7	-2.08	105.41	108.11

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	M0N	C9-C8-C2-N2
3	B	902	M0N	C9-C8-C2-N2

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	291/340 (85%)	0.13	18 (6%) 20 29	27, 42, 74, 99	0
1	B	290/340 (85%)	0.17	16 (5%) 24 33	30, 47, 77, 95	0
All	All	581/680 (85%)	0.15	34 (5%) 22 31	27, 45, 77, 99	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	42	ASN	5.4
1	A	229	GLY	5.1
1	A	224	MET	4.8
1	A	43	PHE	4.4
1	B	90	ARG	4.0
1	B	43	PHE	3.7
1	B	4	ASN	3.6
1	B	56	ASN	3.5
1	A	90	ARG	3.4
1	B	49	VAL	3.2
1	B	306	ASN	3.2
1	B	229	GLY	3.2
1	B	224	MET	3.1
1	A	306	ASN	3.0
1	A	191	PRO	2.9
1	B	36	LEU	2.8
1	A	124	THR	2.8
1	A	252	THR	2.8
1	B	37	LEU	2.7
1	A	42	ASN	2.6
1	A	230	PHE	2.5
1	B	191	PRO	2.5
1	A	46	ASN	2.5
1	A	307	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	225	SER	2.4
1	B	230	PHE	2.4
1	B	178	LEU	2.3
1	A	45	LEU	2.2
1	A	212	ILE	2.1
1	B	91	GLY	2.1
1	A	302	ASN	2.1
1	A	92	GLN	2.1
1	A	171	VAL	2.1
1	B	92	GLN	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	1302	1/1	0.31	2.39	53,53,53,53	0
2	MG	A	1301	1/1	0.21	0.88	42,42,42,42	0
2	MG	B	1304	1/1	0.19	0.87	45,45,45,45	0
3	M0N	A	901	20/20	0.11	-1.00	38,48,60,61	0
3	M0N	B	902	20/20	0.10	-1.21	46,51,61,61	0
2	MG	B	1303	1/1	0.09	-1.94	44,44,44,44	0

6.5 Other polymers ⓘ

There are no such residues in this entry.