



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 07:14 am BST

PDB ID : 1N24
Title : (+)-Bornyl diphosphate synthase: Complex with Mg and product
Authors : Whittington, D.A.; Wise, M.L.; Urbansky, M.; Coates, R.M.; Croteau, R.B.; Christianson, D.W.
Deposited on : 2002-10-21
Resolution : 2.30 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
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) : 2.11

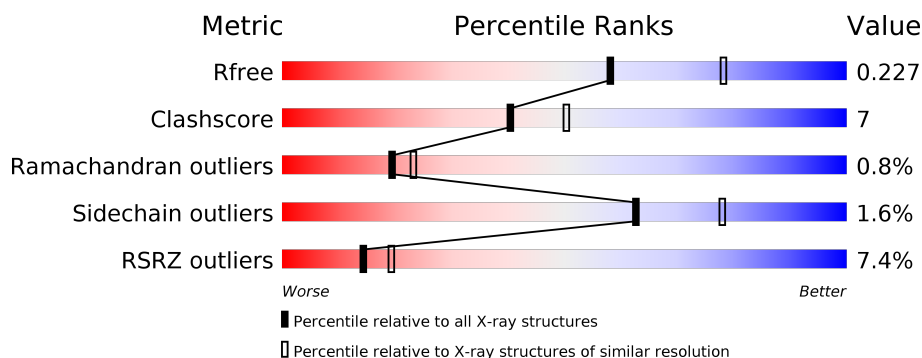
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	549	<div> <div>6%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	549	<div> <div>8%</div> <div> <div></div> <div>79%</div> <div>17%</div> <div>• •</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9146 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.

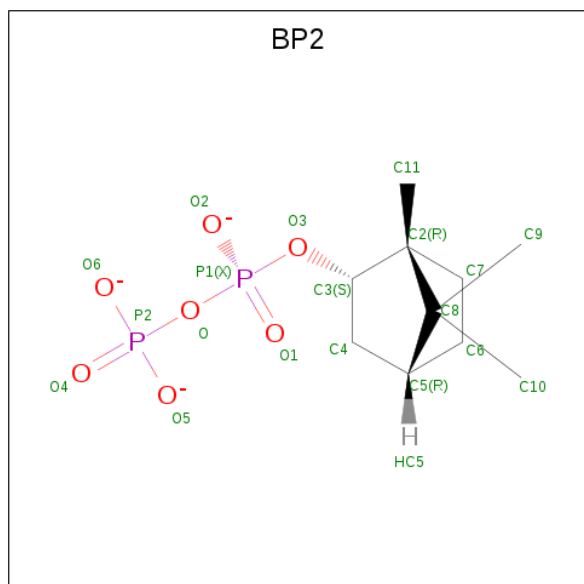
- Molecule 1 is a protein called (+)-bornyl diphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4408	2859	731	801	17			
1	B	527	Total	C	N	O	S	0	0	0
			4346	2817	717	794	18			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Mg	0	0
			3	3		
2	A	3	Total	Mg	0	0
			3	3		

- Molecule 3 is (+)-BORNYL DIPHOSPHATE (three-letter code: BP2) (formula: C₁₀H₁₇O₇P₂).

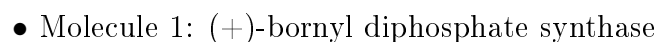


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			19	10	7	2		
3	B	1	Total	C	O	P	0	0
			19	10	7	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	191	Total	O	0	0
			191	191		
4	B	157	Total	O	0	0
			157	157		

- Molecule 1: (+)-bornyl diphosphate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.20Å 117.53Å 120.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.9 (30.00-2.30) 94.8 (30.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.19 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.210 , 0.238 0.204 , 0.227	Depositor DCC
R_{free} test set	2465 reflections (3.86%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.012 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9146	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, BP2

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	0.41	1/4527 (0.0%)	0.58	3/6135 (0.0%)
1	B	0.38	0/4460	0.54	1/6042 (0.0%)
All	All	0.40	1/8987 (0.0%)	0.56	4/12177 (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	A	0	1
1	B	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	76	PRO	CA-C	-8.17	1.36	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	76	PRO	CA-N-CD	-10.30	97.08	111.50
1	A	76	PRO	CB-CA-C	-5.25	98.88	112.00
1	B	480	GLN	N-CA-C	-5.17	97.05	111.00
1	A	76	PRO	O-C-N	-5.07	114.58	122.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	481	TYR	Sidechain
1	B	481	TYR	Sidechain

5.2 Too-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 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	4408	0	4293	62	0
1	B	4346	0	4236	65	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	19	0	17	0	0
3	B	19	0	17	0	0
4	A	191	0	0	5	0
4	B	157	0	0	3	0
All	All	9146	0	8563	125	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 7.

All (125) 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:429:GLU:HB3	1:A:510:VAL:HG21	1.28	1.13
1:B:429:GLU:HB3	1:B:510:VAL:HG21	1.28	1.13
1:B:146:GLU:HG2	1:B:147:LYS:H	1.22	1.05
1:B:141:HIS:O	1:B:142:ASN:OD1	1.86	0.93
1:B:143:ASN:OD1	1:B:171:VAL:HG22	1.73	0.88
1:B:492:LEU:HD13	1:B:568:ALA:HB2	1.55	0.88
1:A:310:LEU:O	1:A:313:VAL:HG22	1.81	0.80
1:A:492:LEU:HD13	1:A:568:ALA:HB2	1.62	0.80
1:B:143:ASN:HD21	1:B:168:SER:HB3	1.50	0.76
1:B:143:ASN:OD1	1:B:171:VAL:CG2	2.34	0.76
1:A:429:GLU:CB	1:A:510:VAL:HG21	2.14	0.75
1:B:146:GLU:HG2	1:B:147:LYS:N	2.02	0.72
1:B:81:ARG:HH12	1:B:82:HIS:CE1	2.12	0.66
1:B:143:ASN:ND2	1:B:168:SER:HB3	2.10	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:HIS:O	1:B:142:ASN:CG	2.33	0.66
1:A:494:LEU:HB2	1:A:495:PRO:HD3	1.77	0.66
1:B:276:PHE:CZ	1:B:280:LYS:HD2	2.30	0.66
1:B:494:LEU:HB2	1:B:495:PRO:HD3	1.77	0.65
1:A:225:LYS:HE3	1:A:233:ILE:HD11	1.80	0.64
1:B:479:TYR:C	1:B:481:TYR:H	1.96	0.63
1:B:479:TYR:C	1:B:481:TYR:N	2.43	0.63
1:B:498:LEU:HD13	1:B:532:VAL:HG11	1.80	0.63
1:B:238:LEU:O	1:B:242:ARG:HG3	1.99	0.62
1:A:123:GLN:O	1:A:127:LYS:HG3	2.00	0.62
1:A:429:GLU:HB3	1:A:510:VAL:CG2	2.17	0.62
1:B:492:LEU:O	1:B:492:LEU:HD12	1.99	0.61
1:A:276:PHE:CZ	1:A:280:LYS:HD2	2.35	0.60
1:A:92:GLN:HB3	1:A:274:LEU:HD13	1.83	0.60
1:A:224:LYS:C	1:A:226:LEU:H	2.05	0.60
1:B:266:ALA:HB2	1:B:276:PHE:CE1	2.37	0.59
1:A:189:GLN:HG3	1:A:190:ASP:N	2.17	0.59
1:A:452:VAL:HG22	1:A:492:LEU:HG	1.84	0.59
1:B:169:GLN:NE2	1:B:210:THR:HB	2.18	0.59
1:A:585:THR:HG22	1:A:589:ILE:HG13	1.84	0.58
1:B:452:VAL:HG22	1:B:492:LEU:HG	1.86	0.56
1:B:81:ARG:NH1	1:B:82:HIS:CE1	2.72	0.56
1:B:175:PHE:HB3	1:B:187:LEU:HD11	1.88	0.56
1:B:276:PHE:CE2	1:B:280:LYS:HD2	2.41	0.56
1:B:500:THR:HG22	1:B:576:ASP:HB2	1.88	0.55
1:A:101:MET:CE	1:A:105:GLN:HB3	2.37	0.55
1:B:104:VAL:O	1:B:108:GLU:HG3	2.07	0.54
1:A:510:VAL:HG23	4:A:815:HOH:O	2.06	0.54
1:B:492:LEU:HD13	1:B:568:ALA:CB	2.35	0.54
1:A:104:VAL:O	1:A:108:GLU:HG3	2.07	0.54
1:B:253:ILE:HD11	1:B:566:ARG:HB2	1.90	0.53
1:B:452:VAL:CG2	1:B:492:LEU:HG	2.37	0.53
1:A:355:ASP:OD1	1:A:356:VAL:HG23	2.08	0.53
1:A:303:ARG:NH2	4:A:990:HOH:O	2.28	0.53
1:A:373:TRP:HB2	4:A:912:HOH:O	2.08	0.53
1:A:452:VAL:CG2	1:A:492:LEU:HG	2.39	0.53
1:A:500:THR:O	1:A:504:GLU:HG3	2.09	0.53
1:A:233:ILE:HG22	1:A:234:ASP:N	2.23	0.53
1:A:454:SER:HB2	1:A:455:PRO:HD3	1.91	0.52
1:B:417:LYS:HG3	4:B:839:HOH:O	2.10	0.52
1:A:189:GLN:HG3	1:A:190:ASP:H	1.75	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:SER:O	1:A:189:GLN:HG2	2.10	0.51
1:B:89:LEU:O	1:B:93:VAL:HG23	2.10	0.51
1:A:280:LYS:HG2	1:A:597:TYR:CD2	2.46	0.51
1:A:479:TYR:C	1:A:481:TYR:N	2.54	0.51
1:B:148:MET:CE	1:B:174:CYS:HB3	2.41	0.50
1:A:253:ILE:HD11	1:A:566:ARG:HB2	1.92	0.50
1:A:480:GLN:O	1:A:481:TYR:HB2	2.11	0.50
1:A:181:ILE:HG22	1:A:182:ASP:OD1	2.11	0.50
1:A:375:THR:HB	1:B:412:LEU:HD23	1.93	0.50
1:A:515:GLN:H	1:A:515:GLN:CD	2.15	0.50
1:A:570:PHE:HB2	4:A:858:HOH:O	2.11	0.50
1:B:169:GLN:HE22	1:B:210:THR:H	1.60	0.49
1:A:225:LYS:CE	1:A:233:ILE:HD11	2.41	0.49
1:B:232:GLU:O	1:B:233:ILE:HG13	2.12	0.49
1:A:280:LYS:HE3	1:A:597:TYR:HD2	1.78	0.48
1:B:310:LEU:O	1:B:313:VAL:HG22	2.13	0.48
1:A:175:PHE:HB3	1:A:187:LEU:HD11	1.95	0.48
1:B:355:ASP:OD1	1:B:356:VAL:HG23	2.13	0.48
1:A:101:MET:HE1	1:A:105:GLN:HB3	1.96	0.48
1:A:429:GLU:O	1:A:510:VAL:HG22	2.14	0.48
1:B:524:SER:OG	1:B:527:GLU:HG3	2.14	0.48
1:B:454:SER:HB2	1:B:455:PRO:HD3	1.95	0.47
1:A:224:LYS:C	1:A:226:LEU:N	2.68	0.47
1:B:125:GLU:OE1	1:B:125:GLU:N	2.42	0.47
1:A:89:LEU:O	1:A:93:VAL:HG23	2.15	0.46
1:B:459:SER:OG	1:B:460:PRO:HD3	2.16	0.46
1:B:303:ARG:NH2	4:B:894:HOH:O	2.48	0.46
1:B:237:LEU:O	1:B:241:ILE:HG13	2.15	0.46
1:A:101:MET:HE2	1:A:105:GLN:HB3	1.98	0.46
1:A:87:ALA:O	1:A:91:VAL:HG23	2.15	0.46
1:A:524:SER:OG	1:A:527:GLU:HG3	2.15	0.45
1:B:233:ILE:O	1:B:234:ASP:C	2.53	0.45
1:A:412:LEU:HD23	1:B:375:THR:HB	1.97	0.45
1:A:374:ASP:HB2	4:A:887:HOH:O	2.17	0.45
1:A:459:SER:OG	1:A:460:PRO:HD3	2.16	0.45
1:A:162:GLN:HG3	1:A:203:LEU:HD21	1.99	0.44
1:A:276:PHE:CE1	1:A:280:LYS:HD2	2.52	0.44
1:A:280:LYS:HE3	1:A:597:TYR:CD2	2.53	0.44
1:B:510:VAL:HG23	4:B:959:HOH:O	2.17	0.44
1:A:359:THR:OG1	1:A:362:GLU:HG3	2.17	0.44
1:A:561:ALA:O	1:A:564:ILE:HG22	2.17	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:237:LEU:O	1:A:241:ILE:HG13	2.19	0.43
1:B:458:ILE:HG21	1:B:485:LEU:HD22	2.00	0.43
1:A:169:GLN:OE1	1:A:210:THR:HB	2.18	0.43
1:B:148:MET:HE2	1:B:174:CYS:HB3	2.00	0.43
1:A:477:SER:O	1:A:480:GLN:HG2	2.19	0.43
1:A:577:GLY:O	1:A:582:HIS:HA	2.18	0.43
1:B:162:GLN:HG3	1:B:203:LEU:HD21	2.00	0.43
1:B:565:GLY:O	1:B:569:GLN:HG3	2.18	0.43
1:B:139:CYS:HB2	1:B:146:GLU:HB2	2.01	0.42
1:B:561:ALA:O	1:B:564:ILE:HG22	2.19	0.42
1:A:171:VAL:O	1:A:171:VAL:HG12	2.20	0.42
1:A:233:ILE:HG22	1:A:234:ASP:H	1.84	0.42
1:A:565:GLY:O	1:A:569:GLN:HG3	2.20	0.42
1:B:482:HIS:CE1	1:B:484:ILE:HG13	2.54	0.42
1:B:271:MET:HG2	1:B:272:ASN:N	2.34	0.42
1:B:480:GLN:O	1:B:481:TYR:HB2	2.20	0.42
1:B:101:MET:HG2	1:B:105:GLN:HB2	2.01	0.41
1:B:245:LEU:HD23	1:B:245:LEU:HA	1.89	0.41
1:A:175:PHE:N	1:A:175:PHE:CD1	2.88	0.41
1:A:151:TYR:CD1	1:A:192:LYS:HG2	2.56	0.41
1:B:171:VAL:HG12	1:B:171:VAL:O	2.21	0.41
1:B:175:PHE:CD1	1:B:175:PHE:N	2.88	0.41
1:B:151:TYR:CD1	1:B:192:LYS:HG2	2.56	0.41
1:A:162:GLN:CG	1:A:203:LEU:HD21	2.51	0.41
1:B:85:ARG:HH21	1:B:277:GLU:HG2	1.86	0.41
1:B:122:PHE:O	1:B:126:ILE:HG13	2.21	0.40
1:B:309:LYS:C	1:B:310:LEU:HD12	2.42	0.40
1:B:367:THR:O	1:B:371:LYS:HG3	2.22	0.40
1:B:359:THR:OG1	1:B:362:GLU:HG3	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	528/549 (96%)	507 (96%)	17 (3%)	4 (1%)	19	23
1	B	519/549 (94%)	489 (94%)	26 (5%)	4 (1%)	19	23
All	All	1047/1098 (95%)	996 (95%)	43 (4%)	8 (1%)	19	23

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	481	TYR
1	A	77	TYR
1	A	177	ASN
1	B	136	GLU
1	B	142	ASN
1	B	481	TYR
1	A	225	LYS
1	B	77	TYR

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	468/480 (98%)	460 (98%)	8 (2%)	60	76
1	B	463/480 (96%)	456 (98%)	7 (2%)	65	79
All	All	931/960 (97%)	916 (98%)	15 (2%)	62	78

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	PRO
1	A	182	ASP
1	A	209	ASP
1	A	254	GLN
1	A	485	LEU
1	A	492	LEU
1	A	571	ILE
1	A	582	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	182	ASP
1	B	209	ASP
1	B	227	ASP
1	B	254	GLN
1	B	323	TRP
1	B	492	LEU
1	B	571	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	A	254	GLN
1	A	282	ASN
1	A	531	HIS
1	B	82	HIS
1	B	169	GLN
1	B	173	ASN
1	B	254	GLN
1	B	531	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](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 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	BP2	B	802	2	18,20,20	1.37	2 (11%)	32,35,35	1.98	7 (21%)
3	BP2	A	801	2	18,20,20	1.32	2 (11%)	32,35,35	1.94	6 (18%)

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	BP2	B	802	2	-	2/11/42/42	0/3/2/2
3	BP2	A	801	2	-	2/11/42/42	0/3/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	BP2	P2-O4	3.18	1.60	1.50
3	B	802	BP2	P1-O3	-3.03	1.52	1.60
3	B	802	BP2	P2-O4	2.90	1.59	1.50
3	A	801	BP2	P1-O3	-2.76	1.53	1.60

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	BP2	O-P1-O3	5.87	114.33	102.48
3	A	801	BP2	O-P1-O3	5.17	112.91	102.48
3	A	801	BP2	C4-C5-C8	4.82	108.57	102.52
3	B	802	BP2	C4-C5-C8	4.74	108.47	102.52
3	A	801	BP2	C8-C2-C3	4.28	106.43	100.63
3	B	802	BP2	C8-C2-C3	4.09	106.18	100.63
3	B	802	BP2	C2-C8-C5	-2.59	88.06	93.71
3	A	801	BP2	C2-C8-C5	-2.55	88.17	93.71
3	A	801	BP2	O3-C3-C2	2.34	113.22	110.03
3	B	802	BP2	O3-C3-C2	2.34	113.21	110.03
3	A	801	BP2	C6-C7-C2	-2.10	100.44	104.19
3	B	802	BP2	C7-C2-C8	2.06	104.84	101.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	802	BP2	C6-C7-C2	-2.03	100.57	104.19

There are no chirality outliers.

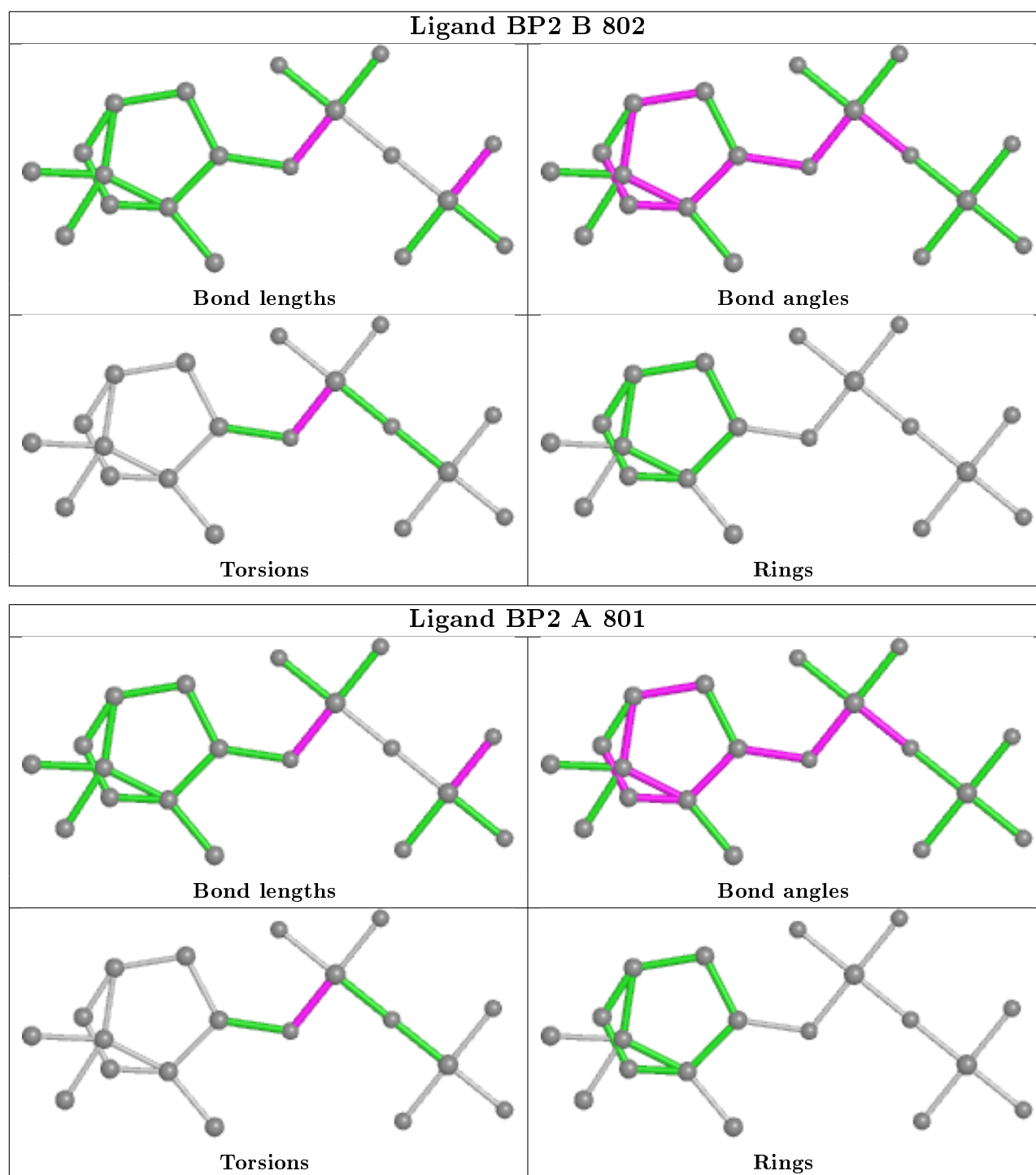
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	802	BP2	C3-O3-P1-O
3	A	801	BP2	C3-O3-P1-O
3	B	802	BP2	C3-O3-P1-O1
3	A	801	BP2	C3-O3-P1-O1

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	534/549 (97%)	0.19	33 (6%)	20 26	21, 38, 71, 93	0
1	B	527/549 (95%)	0.30	46 (8%)	10 14	20, 41, 89, 114	0
All	All	1061/1098 (96%)	0.25	79 (7%)	14 19	20, 40, 81, 114	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	146	GLU	7.7
1	B	142	ASN	6.4
1	B	143	ASN	6.3
1	B	147	LYS	6.3
1	B	506	ALA	6.0
1	B	505	LEU	5.1
1	B	141	HIS	5.1
1	B	583	SER	4.9
1	B	148	MET	4.7
1	B	238	LEU	4.7
1	B	145	VAL	4.4
1	A	233	ILE	4.3
1	A	138	LYS	4.2
1	A	598	ALA	4.2
1	B	233	ILE	4.1
1	A	181	ILE	4.0
1	B	226	LEU	4.0
1	A	178	GLU	4.0
1	A	137	HIS	3.9
1	A	54	ILE	3.8
1	A	78	THR	3.8
1	B	584	LYS	3.8
1	B	235	GLU	3.7
1	A	344	ILE	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	80	GLU	3.5
1	A	343	ILE	3.4
1	A	235	GLU	3.4
1	B	574	HIS	3.3
1	B	586	TYR	3.3
1	B	138	LYS	3.3
1	B	178	GLU	3.2
1	B	578	PHE	3.2
1	A	179	LYS	3.1
1	B	232	GLU	3.1
1	B	189	GLN	3.1
1	B	225	LYS	3.0
1	B	231	ASN	2.9
1	A	180	GLY	2.9
1	A	189	GLN	2.8
1	B	577	GLY	2.8
1	B	100	LYS	2.8
1	A	224	LYS	2.8
1	B	224	LYS	2.8
1	A	346	LEU	2.7
1	B	507	ARG	2.7
1	A	453	ALA	2.7
1	B	274	LEU	2.6
1	A	234	ASP	2.6
1	B	228	GLU	2.6
1	B	180	GLY	2.5
1	B	242	ARG	2.5
1	B	62	PRO	2.5
1	B	227	ASP	2.4
1	B	177	ASN	2.4
1	A	347	ALA	2.4
1	A	350	ILE	2.4
1	B	587	GLU	2.4
1	A	134	TYR	2.3
1	A	506	ALA	2.3
1	B	80	GLU	2.3
1	A	141	HIS	2.3
1	A	457	ILE	2.3
1	B	140	PHE	2.3
1	A	345	VAL	2.2
1	B	510	VAL	2.2
1	A	61	GLN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	564	ILE	2.1
1	A	522	ASN	2.1
1	B	200	ALA	2.1
1	A	340	ALA	2.1
1	B	95	ILE	2.1
1	B	213	LEU	2.1
1	B	234	ASP	2.0
1	B	64	LEU	2.0
1	B	522	ASN	2.0
1	A	348	THR	2.0
1	A	127	LYS	2.0
1	A	140	PHE	2.0
1	B	181	ILE	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

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. 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.

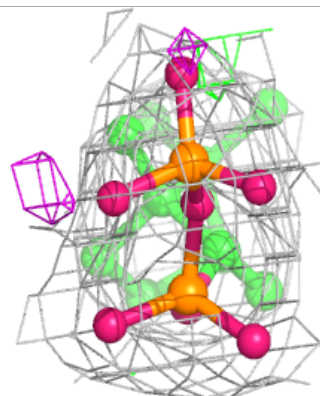
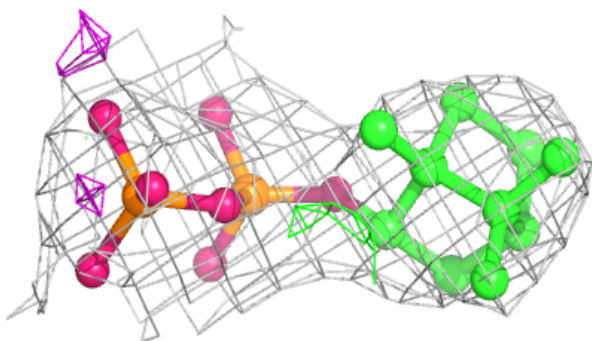
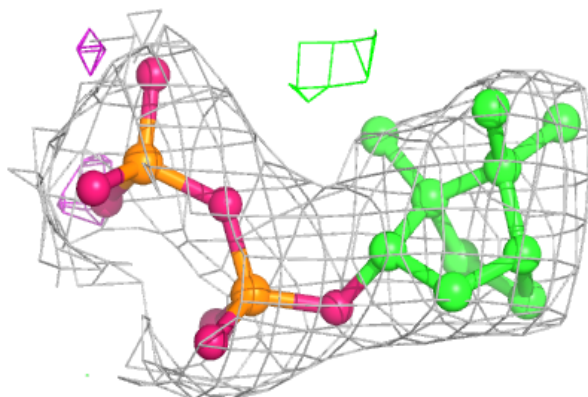
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MG	B	706	1/1	0.75	0.08	45,45,45,45	0
2	MG	B	704	1/1	0.96	0.04	25,25,25,25	0
2	MG	B	705	1/1	0.97	0.07	38,38,38,38	0
2	MG	A	703	1/1	0.97	0.11	37,37,37,37	0
3	BP2	B	802	19/19	0.97	0.13	36,44,49,49	0
2	MG	A	702	1/1	0.97	0.10	28,28,28,28	0
2	MG	A	701	1/1	0.98	0.11	29,29,29,29	0
3	BP2	A	801	19/19	0.98	0.17	26,33,35,37	0

The following is a graphical depiction of the model fit to experimental electron density of all

instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

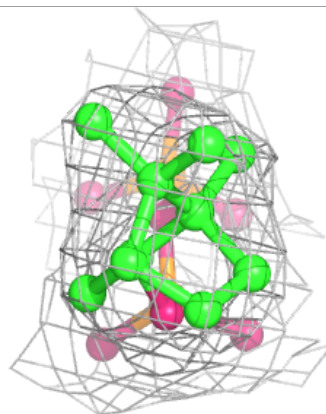
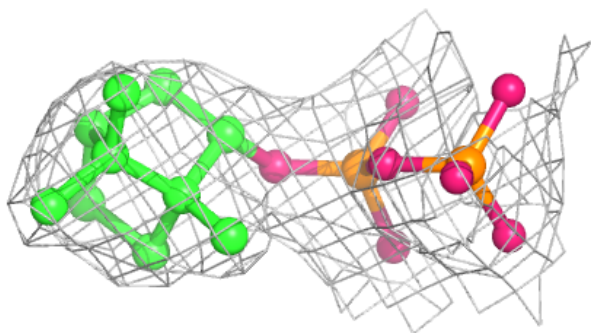
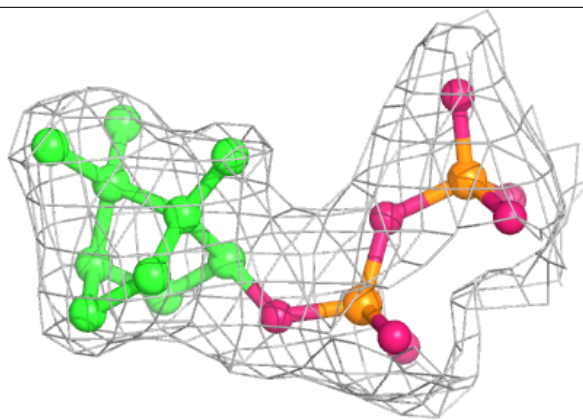
Electron density around BP2 B 802:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BP2 A 801:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.