



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:00 PM GMT

PDB ID : 1N20
Title : (+)-Bornyl Diphosphate Synthase: Complex with Mg and 3-aza-2,3-dihydrogeranyl diphosphate
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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

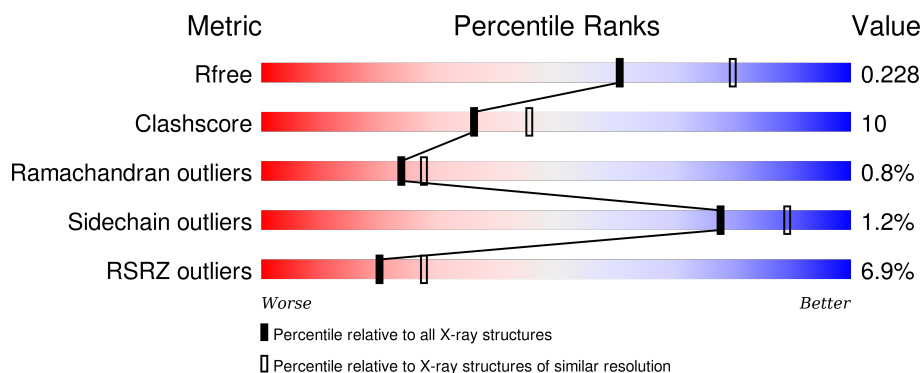
1 Overall quality at a glance

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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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. 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>8%</div> <div>79%</div> <div>17%</div> <div>• •</div> </div>
1	B	549	<div> <div>5%</div> <div>76%</div> <div>19%</div> <div>• •</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9030 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.

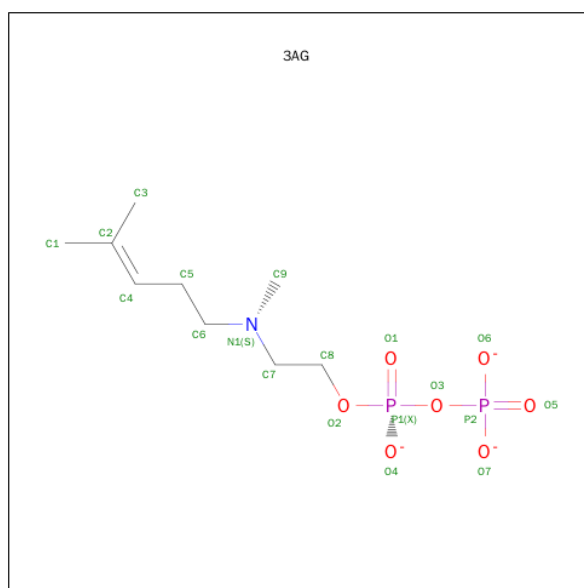
- Molecule 1 is a protein called (+)-bornyl diphosphate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	532	Total	C	N	O	S	0	0	0
			4379	2841	728	793	17			
1	B	527	Total	C	N	O	S	0	0	0
			4338	2817	716	788	17			

- 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 2-[METHYL-(4-METHYL-PENT-3-ENYL)-AMINO]-ETHYL-DIPHOSPHATE (three-letter code: 3AG) (formula: C₉H₁₈NO₇P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			19	9	1	7	2		
3	B	1	Total	C	N	O	P	0	0
			19	9	1	7	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total	O	0	0
			146	146		
4	B	123	Total	O	0	0
			123	123		

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.

- Chain A:
-
- 8% 79% 17%
- GLU
ALA
HIS
GLN
T54
R55
R56
G58
N59
Y60
L64
W65
T78
E79
E80
R81
Q92
M101
V104
Q106
L107
E108
K127
N135
E136
H137
K138
C139
F140
H141
ASN
ASN
GLU
VAL
GLU
LYS
MET
D149
L150
Y151
R158
L159
Q162
Q169
D170
V171
- F175
K176
E178
K179
G180
I181
D182
F183
K184
A185
S186
L187
A188
Q189
D190
T191
K192
L195
E199
E208
D227
GLY
GLY
ASN
GLU
ILE
D234
L237
L238
L239
W240
I241
I253
Q254
S255
A258
I262
L274
I275
F276
K280
- W300
K309
L310
V313
R314
D315
R316
E319
E329
T342
I343
I344
V345
L346
A347
V348
V349
I350
D355
V356
T359
L360
D361
E362
K371
R372
K373
C388
V392
K405
E406
L422
V423
E424
E429
H434
V452
A453
S454
P455
A456
I457
I458
V470

- Chain B:
-
- 5% 76% 19% . .

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.08Å 117.54Å 120.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.77 – 2.30 28.77 – 2.29	Depositor EDS
% Data completeness (in resolution range)	92.8 (28.77-2.30) 92.3 (28.77-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.16 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.205 , 0.228 0.205 , 0.228	Depositor DCC
R_{free} test set	2410 reflections (4.05%)	DCC
Wilson B-factor (Å ²)	37.2	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.0	EDS
Estimated twinning fraction	0.011 for -h,l,k	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 62502 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9030	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, 3AG

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	0.37	0/4498	0.57	2/6098 (0.0%)
1	B	0.37	0/4455	0.56	3/6040 (0.0%)
All	All	0.37	0/8953	0.57	5/12138 (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	B	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	316	ARG	NE-CZ-NH1	-9.42	115.59	120.30
1	B	316	ARG	NE-CZ-NH2	-9.39	115.61	120.30
1	A	316	ARG	NE-CZ-NH2	8.96	124.78	120.30
1	B	316	ARG	NE-CZ-NH1	8.43	124.52	120.30
1	B	480	GLN	N-CA-C	-5.24	96.84	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
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	4379	0	4257	83	0
1	B	4338	0	4225	82	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	19	0	18	3	0
3	B	19	0	18	3	0
4	A	146	0	0	5	0
4	B	123	0	0	5	0
All	All	9030	0	8518	165	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 10.

All (165) 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:B:429:GLU:HB3	1:B:510:VAL:HG21	1.25	1.12
1:A:429:GLU:HB3	1:A:510:VAL:HG21	1.28	1.09
1:B:492:LEU:HD13	1:B:568:ALA:HB2	1.51	0.93
1:A:492:LEU:HD13	1:A:568:ALA:HB2	1.60	0.84
1:A:313:VAL:HG11	1:A:349:VAL:HG22	1.64	0.78
1:B:160:LEU:HD13	1:B:167:ILE:HD11	1.67	0.77
1:A:186:SER:O	1:A:189:GLN:HG2	1.88	0.73
1:B:429:GLU:CB	1:B:510:VAL:HG21	2.14	0.72
1:B:89:LEU:O	1:B:93:VAL:HG23	1.91	0.71
1:A:429:GLU:CB	1:A:510:VAL:HG21	2.15	0.71
1:B:313:VAL:HG21	1:B:349:VAL:HA	1.72	0.71
1:A:457:ILE:CD1	3:A:900:3AG:HC32	2.21	0.70
1:A:104:VAL:O	1:A:108:GLU:HG3	1.90	0.70
1:B:277:GLU:HA	1:B:280:LYS:HE2	1.76	0.66
1:B:316:ARG:HH11	1:B:579:GLY:HA2	1.60	0.66
1:B:492:LEU:HD13	1:B:568:ALA:CB	2.25	0.66
1:A:189:GLN:HG3	1:A:190:ASP:N	2.10	0.65
1:B:61:GLN:HB2	1:B:316:ARG:HH22	1.62	0.65
1:A:276:PHE:CZ	1:A:280:LYS:HD2	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:429:GLU:HB3	1:B:510:VAL:CG2	2.16	0.64
1:A:316:ARG:HH11	1:A:579:GLY:HA2	1.63	0.63
1:B:104:VAL:O	1:B:108:GLU:HG3	1.99	0.63
1:A:494:LEU:HB2	1:A:495:PRO:HD3	1.80	0.63
1:B:310:LEU:O	1:B:313:VAL:HG12	1.99	0.63
1:B:494:LEU:HB2	1:B:495:PRO:HD3	1.79	0.63
1:A:56:ARG:HH21	1:A:56:ARG:HG2	1.63	0.62
1:B:158:ARG:O	1:B:162:GLN:HG3	2.00	0.61
1:A:134:TYR:CE1	1:A:141:HIS:HB3	2.36	0.60
1:A:429:GLU:HB3	1:A:510:VAL:CG2	2.17	0.60
1:A:189:GLN:HG3	1:A:190:ASP:H	1.65	0.60
1:A:158:ARG:O	1:A:162:GLN:HG3	2.02	0.60
1:A:280:LYS:HE2	4:A:974:HOH:O	2.03	0.59
1:B:218:ALA:O	1:B:222:LEU:HB2	2.01	0.59
1:A:360:LEU:HD13	1:A:434:HIS:HD2	1.67	0.59
1:B:316:ARG:HD2	1:B:578:PHE:O	2.03	0.59
1:A:280:LYS:HE3	1:A:597:TYR:CE2	2.38	0.59
1:A:92:GLN:HB3	1:A:274:LEU:HD13	1.83	0.58
1:B:492:LEU:HD12	1:B:492:LEU:O	2.02	0.58
1:A:313:VAL:CG1	1:A:349:VAL:HG22	2.32	0.58
1:B:479:TYR:C	1:B:481:TYR:N	2.48	0.57
1:A:585:THR:HA	4:A:1043:HOH:O	2.04	0.57
1:A:280:LYS:HE3	1:A:597:TYR:CD2	2.39	0.57
1:A:316:ARG:HD2	1:A:319:GLU:OE2	2.03	0.57
1:A:253:ILE:HD11	1:A:566:ARG:HB2	1.86	0.57
1:B:505:LEU:HD13	1:B:515:GLN:OE1	2.05	0.56
1:A:355:ASP:OD1	1:A:356:VAL:HG23	2.06	0.56
1:A:524:SER:OG	1:A:527:GLU:HG3	2.05	0.56
1:B:225:LYS:HG3	1:B:233:ILE:HD11	1.87	0.55
1:B:360:LEU:HD13	1:B:434:HIS:HD2	1.72	0.55
1:B:524:SER:OG	1:B:527:GLU:HG3	2.06	0.55
1:A:577:GLY:HA2	1:A:581:GLN:O	2.07	0.55
1:B:492:LEU:HD21	3:B:901:3AG:HC31	1.87	0.55
1:A:457:ILE:HD12	3:A:900:3AG:HC32	1.89	0.55
1:B:112:ASP:CG	1:B:268:ARG:HH12	2.10	0.54
1:B:101:MET:HG2	1:B:105:GLN:HB2	1.90	0.54
1:A:479:TYR:C	1:A:481:TYR:N	2.51	0.54
1:B:454:SER:HB2	1:B:455:PRO:HD3	1.89	0.54
1:A:101:MET:HG2	1:A:105:GLN:HB2	1.90	0.53
1:B:380:ARG:HH21	1:B:380:ARG:HG2	1.71	0.53
1:A:582:HIS:CD2	1:A:582:HIS:H	2.27	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:454:SER:HB2	1:A:455:PRO:HD3	1.91	0.52
1:B:253:ILE:HD11	1:B:566:ARG:HB2	1.91	0.52
1:B:528:ALA:O	1:B:532:VAL:HG23	2.09	0.52
1:B:313:VAL:HG23	1:B:352:ASP:HB2	1.92	0.52
1:A:458:ILE:HG21	1:A:485:LEU:HD22	1.91	0.52
1:A:528:ALA:O	1:A:532:VAL:HG23	2.09	0.52
1:A:429:GLU:O	1:A:510:VAL:HG22	2.10	0.51
1:B:266:ALA:HB2	1:B:276:PHE:CE1	2.45	0.51
1:B:313:VAL:CG2	1:B:349:VAL:HA	2.39	0.51
1:A:373:TRP:HB2	4:A:909:HOH:O	2.11	0.51
1:B:277:GLU:O	1:B:280:LYS:HG2	2.10	0.51
1:B:226:LEU:HD21	1:B:242:ARG:HG2	1.93	0.51
1:B:429:GLU:O	1:B:510:VAL:HG22	2.11	0.50
1:A:510:VAL:HG23	4:A:945:HOH:O	2.10	0.50
1:B:585:THR:HG22	1:B:589:ILE:HG13	1.94	0.50
1:A:480:GLN:O	1:A:481:TYR:HB2	2.12	0.50
1:B:309:LYS:C	1:B:310:LEU:HD12	2.32	0.50
1:A:359:THR:OG1	1:A:362:GLU:HG3	2.13	0.49
1:B:382:PRO:HD3	4:B:972:HOH:O	2.13	0.49
1:B:498:LEU:CD1	1:B:532:VAL:HG11	2.42	0.49
1:A:276:PHE:CE1	1:A:280:LYS:HD2	2.48	0.49
1:A:316:ARG:HH11	1:A:579:GLY:CA	2.26	0.48
1:A:360:LEU:HD13	1:A:434:HIS:CD2	2.48	0.48
1:B:226:LEU:HD22	1:B:245:LEU:HD12	1.95	0.48
1:B:160:LEU:CD1	1:B:167:ILE:HD11	2.41	0.48
1:A:316:ARG:HD2	1:A:319:GLU:CD	2.33	0.48
1:A:171:VAL:O	1:A:171:VAL:HG12	2.13	0.48
1:B:215:ARG:HD3	4:B:1016:HOH:O	2.13	0.48
1:A:254:GLN:NE2	4:A:928:HOH:O	2.45	0.48
1:A:452:VAL:HG22	1:A:492:LEU:HG	1.96	0.47
1:B:360:LEU:HD13	1:B:434:HIS:CD2	2.49	0.47
1:A:405:LYS:NZ	1:A:406:GLU:HG2	2.29	0.47
1:B:237:LEU:O	1:B:241:ILE:HG13	2.15	0.47
1:A:182:ASP:HA	1:A:213:LEU:HD21	1.96	0.47
1:B:61:GLN:CB	1:B:316:ARG:HH22	2.27	0.47
1:B:194:MET:HG3	1:B:221:CYS:SG	2.54	0.47
1:B:359:THR:OG1	1:B:362:GLU:HG3	2.14	0.47
1:A:64:LEU:HD13	1:A:586:TYR:CD1	2.50	0.47
1:A:576:ASP:HB3	1:A:581:GLN:HB2	1.97	0.47
1:A:177:ASN:HA	1:A:184:LYS:HG2	1.96	0.47
1:A:500:THR:O	1:A:504:GLU:HG3	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:LEU:O	1:B:113:LEU:HG	2.15	0.46
1:B:510:VAL:HG23	4:B:1005:HOH:O	2.15	0.46
1:B:479:TYR:C	1:B:481:TYR:H	2.07	0.45
1:A:237:LEU:O	1:A:241:ILE:HG13	2.16	0.45
1:A:60:TYR:CE1	1:A:314:ARG:HG3	2.52	0.45
1:B:373:TRP:HB2	4:B:959:HOH:O	2.16	0.45
1:A:344:ILE:HG12	1:A:457:ILE:HD13	1.99	0.45
1:B:168:SER:O	1:B:171:VAL:HG23	2.17	0.45
1:A:503:PHE:CD2	1:A:580:VAL:HG11	2.51	0.45
1:B:88:GLU:O	1:B:92:GLN:HG2	2.17	0.45
1:A:452:VAL:CG2	1:A:492:LEU:HG	2.47	0.45
1:A:175:PHE:HB3	1:A:187:LEU:HD11	1.98	0.45
1:B:313:VAL:HG22	1:B:314:ARG:N	2.32	0.45
1:A:309:LYS:C	1:A:310:LEU:HD12	2.37	0.45
1:A:64:LEU:HD13	1:A:586:TYR:HD1	1.82	0.44
1:A:151:TYR:CE1	1:A:192:LYS:HE2	2.52	0.44
1:A:212:GLU:OE2	1:A:215:ARG:NH1	2.46	0.44
1:A:492:LEU:HD21	3:A:900:3AG:HC31	1.99	0.44
1:A:195:LEU:O	1:A:199:GLU:HG2	2.18	0.43
1:A:258:ALA:O	1:A:262:ILE:HG13	2.18	0.43
1:A:169:GLN:HE22	1:A:208:GLU:HA	1.83	0.43
1:B:480:GLN:O	1:B:481:TYR:HB2	2.18	0.43
1:A:65:TRP:CZ2	1:A:586:TYR:CD1	3.06	0.43
1:B:111:HIS:CD2	1:B:158:ARG:HH22	2.36	0.43
1:A:342:THR:O	1:A:345:VAL:HG12	2.18	0.43
1:A:92:GLN:CB	1:A:274:LEU:HD13	2.48	0.43
1:B:175:PHE:HB3	1:B:187:LEU:HD11	2.01	0.43
1:B:151:TYR:CD1	1:B:192:LYS:HG2	2.53	0.43
1:B:300:TRP:CH2	1:B:342:THR:HG23	2.54	0.43
1:B:233:ILE:HG22	1:B:234:ASP:N	2.33	0.42
1:A:151:TYR:CD1	1:A:192:LYS:HG2	2.53	0.42
1:B:63:ALA:HB2	4:B:1023:HOH:O	2.19	0.42
1:B:433:TYR:CD2	1:B:434:HIS:ND1	2.88	0.42
1:A:255:SER:OG	1:A:329:GLU:HG2	2.19	0.42
1:B:457:ILE:CD1	3:B:901:3AG:HC32	2.49	0.42
1:B:405:LYS:NZ	1:B:406:GLU:HG2	2.35	0.42
1:B:498:LEU:HD11	1:B:532:VAL:HG11	2.02	0.42
1:A:481:TYR:HD2	1:A:481:TYR:HA	1.65	0.42
1:A:60:TYR:CD1	1:A:314:ARG:HG3	2.54	0.42
1:A:492:LEU:HD12	1:A:492:LEU:O	2.20	0.41
1:A:316:ARG:HH11	1:A:579:GLY:C	2.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:355:ASP:OD1	1:B:356:VAL:HG23	2.19	0.41
1:A:388:CYS:O	1:A:392:VAL:HG23	2.20	0.41
1:A:169:GLN:NE2	1:A:208:GLU:HA	2.36	0.41
1:A:371:LYS:HD3	1:A:424:GLU:OE2	2.20	0.41
1:B:360:LEU:HB2	1:B:434:HIS:HD2	1.86	0.41
1:B:195:LEU:O	1:B:199:GLU:HG2	2.20	0.41
1:A:56:ARG:NH2	1:A:56:ARG:HG2	2.34	0.41
1:B:175:PHE:N	1:B:175:PHE:CD1	2.88	0.41
1:B:316:ARG:HD3	1:B:319:GLU:OE2	2.21	0.41
1:B:167:ILE:CG2	1:B:171:VAL:HG21	2.51	0.41
1:B:263:ASP:OD1	1:B:597:TYR:OH	2.34	0.41
1:B:342:THR:O	1:B:345:VAL:HG12	2.21	0.41
1:B:388:CYS:O	1:B:392:VAL:HG23	2.21	0.41
1:B:500:THR:O	1:B:504:GLU:HG3	2.21	0.40
1:B:371:LYS:HD3	1:B:424:GLU:OE2	2.20	0.40
1:B:458:ILE:HG21	1:B:485:LEU:HD22	2.03	0.40
1:B:352:ASP:O	1:B:356:VAL:HB	2.21	0.40
1:B:360:LEU:HB2	1:B:434:HIS:CD2	2.56	0.40
1:B:224:LYS:C	1:B:226:LEU:H	2.24	0.40
1:A:300:TRP:CH2	1:A:342:THR:HG23	2.57	0.40
1:B:578:PHE:CE2	3:B:901:3AG:HC61	2.56	0.40
1:A:107:LEU:HD22	1:A:159:LEU:HD12	2.02	0.40
1:A:479:TYR:C	1:A:481:TYR:H	2.14	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	526/549 (96%)	505 (96%)	17 (3%)	4 (1%)	24 27
1	B	521/549 (95%)	498 (96%)	19 (4%)	4 (1%)	24 27

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1047/1098 (95%)	1003 (96%)	36 (3%)	8 (1%)	24 27

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	481	TYR
1	A	576	ASP
1	B	136	GLU
1	A	177	ASN
1	B	77	TYR
1	B	481	TYR
1	A	577	GLY
1	B	580	VAL

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	463/480 (96%)	458 (99%)	5 (1%)	80 90
1	B	460/480 (96%)	454 (99%)	6 (1%)	76 87
All	All	923/960 (96%)	912 (99%)	11 (1%)	78 89

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	ARG
1	A	226	LEU
1	A	254	GLN
1	A	492	LEU
1	A	582	HIS
1	B	231	ASN
1	B	254	GLN
1	B	323	TRP
1	B	379	THR
1	B	492	LEU

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Mol	Chain	Res	Type
1	B	582	HIS

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

Mol	Chain	Res	Type
1	A	173	ASN
1	A	254	GLN
1	A	434	HIS
1	B	82	HIS
1	B	173	ASN
1	B	231	ASN
1	B	254	GLN

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	3AG	A	900	2	17,18,18	2.19	4 (23%)	21,25,25	0.95	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	3AG	B	901	2	17,18,18	2.14	4 (23%)	21,25,25	1.00	1 (4%)

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	3AG	A	900	2	-	0/19/19/19	0/0/0/0
3	3AG	B	901	2	-	0/19/19/19	0/0/0/0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	900	3AG	P1-O1	2.29	1.59	1.51
3	B	901	3AG	P1-O1	2.53	1.60	1.51
3	B	901	3AG	P2-O5	3.97	1.59	1.51
3	A	900	3AG	P2-O5	4.07	1.59	1.51
3	B	901	3AG	P2-O6	4.81	1.60	1.51
3	A	900	3AG	P2-O7	4.95	1.60	1.51
3	B	901	3AG	P2-O7	5.01	1.60	1.51
3	A	900	3AG	P2-O6	5.10	1.60	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	900	3AG	C8-C7-N1	2.08	118.27	113.38
3	B	901	3AG	C8-C7-N1	2.13	118.39	113.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	900	3AG	3	0
3	B	901	3AG	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	532/549 (96%)	0.22	43 (8%) 15 21	23, 40, 78, 102	0
1	B	527/549 (95%)	0.21	30 (5%) 27 36	22, 43, 86, 105	0
All	All	1059/1098 (96%)	0.21	73 (6%) 20 27	22, 42, 82, 105	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	181	ILE	6.4
1	B	233	ILE	6.2
1	B	138	LYS	5.1
1	A	138	LYS	4.6
1	A	137	HIS	4.2
1	B	238	LEU	4.2
1	A	506	ALA	4.1
1	A	226	LEU	3.9
1	B	231	ASN	3.8
1	B	583	SER	3.7
1	B	100	LYS	3.7
1	A	80	GLU	3.6
1	B	226	LEU	3.6
1	B	137	HIS	3.4
1	B	181	ILE	3.4
1	A	140	PHE	3.3
1	B	180	GLY	3.3
1	B	136	GLU	3.3
1	B	580	VAL	3.3
1	B	139	CYS	3.3
1	B	140	PHE	3.3
1	A	54	ILE	3.3
1	A	503	PHE	3.3
1	A	505	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	59	ASN	3.2
1	A	177	ASN	3.0
1	A	580	VAL	3.0
1	A	180	GLY	2.9
1	A	238	LEU	2.9
1	A	141	HIS	2.8
1	A	582	HIS	2.8
1	A	586	TYR	2.8
1	B	223	GLN	2.7
1	B	178	GLU	2.7
1	B	189	GLN	2.7
1	A	136	GLU	2.7
1	A	350	ILE	2.6
1	A	453	ALA	2.6
1	A	134	TYR	2.6
1	A	343	ILE	2.6
1	A	347	ALA	2.6
1	A	457	ILE	2.5
1	A	189	GLN	2.5
1	A	598	ALA	2.5
1	B	224	LYS	2.5
1	A	346	LEU	2.5
1	A	344	ILE	2.4
1	B	574	HIS	2.4
1	B	586	TYR	2.4
1	A	57	SER	2.3
1	A	187	LEU	2.3
1	B	505	LEU	2.3
1	B	239	LEU	2.3
1	A	78	THR	2.3
1	B	242	ARG	2.3
1	B	95	ILE	2.2
1	A	422	LEU	2.2
1	B	598	ALA	2.2
1	A	348	THR	2.2
1	A	349	VAL	2.2
1	A	186	SER	2.2
1	A	179	LYS	2.2
1	B	177	ASN	2.2
1	B	236	ASN	2.1
1	B	524	SER	2.1
1	A	139	CYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	81	ARG	2.1
1	B	183	PHE	2.1
1	A	239	LEU	2.1
1	A	452	VAL	2.1
1	B	213	LEU	2.0
1	A	178	GLU	2.0
1	A	127	LYS	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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	3AG	B	901	19/19	0.96	0.17	0.43	29,42,53,53	0
3	3AG	A	900	19/19	0.97	0.19	0.15	25,37,43,43	0
2	MG	A	701	1/1	0.97	0.10	-1.09	29,29,29,29	0
2	MG	A	702	1/1	0.96	0.10	-1.54	41,41,41,41	0
2	MG	A	703	1/1	0.95	0.08	-1.56	38,38,38,38	0
2	MG	B	705	1/1	0.98	0.08	-2.81	33,33,33,33	0
2	MG	B	704	1/1	0.98	0.08	-3.38	44,44,44,44	0
2	MG	B	706	1/1	0.88	0.07	-4.39	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.