



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:42 PM UTC

PDB ID : 1N1B / pdb_00001n1b
Title : Crystal Structure of (+)-Bornyl Diphosphate Synthase from Sage
Authors : Whittington, D.A.; Wise, M.L.; Urbansky, M.; Coates, R.M.; Croteau, R.B.;
Christianson, D.W.
Deposited on : 2002-10-17
Resolution : 2.00 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

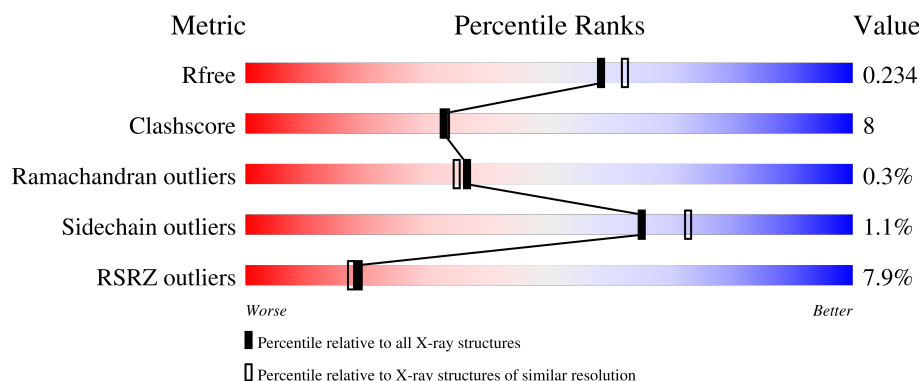
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 of 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	9% 76% 16% • 6%
1	B	549	6% 76% 17% • 5%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9156 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	516	Total	C	N	O	S	0	0	0
			4265	2768	702	777	18			
1	B	519	Total	C	N	O	S	0	1	0
			4290	2783	706	782	19			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		

- Molecule 3 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	Hg	0	0
			5	5		
3	B	6	Total	Hg	0	0
			6	6		

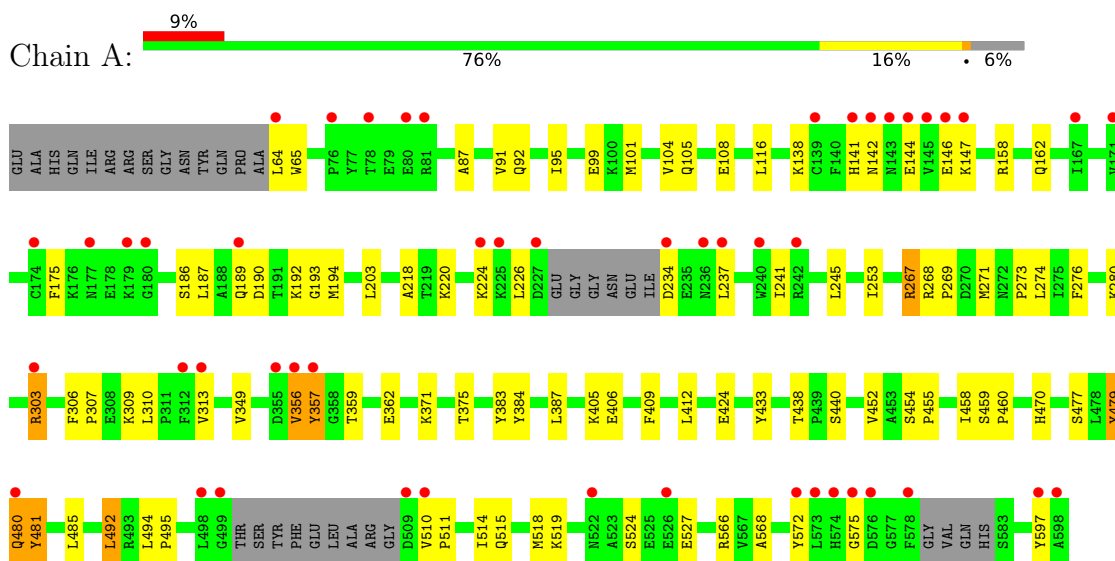
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	277	Total	O	0	0
			277	277		
4	B	311	Total	O	0	0
			311	311		

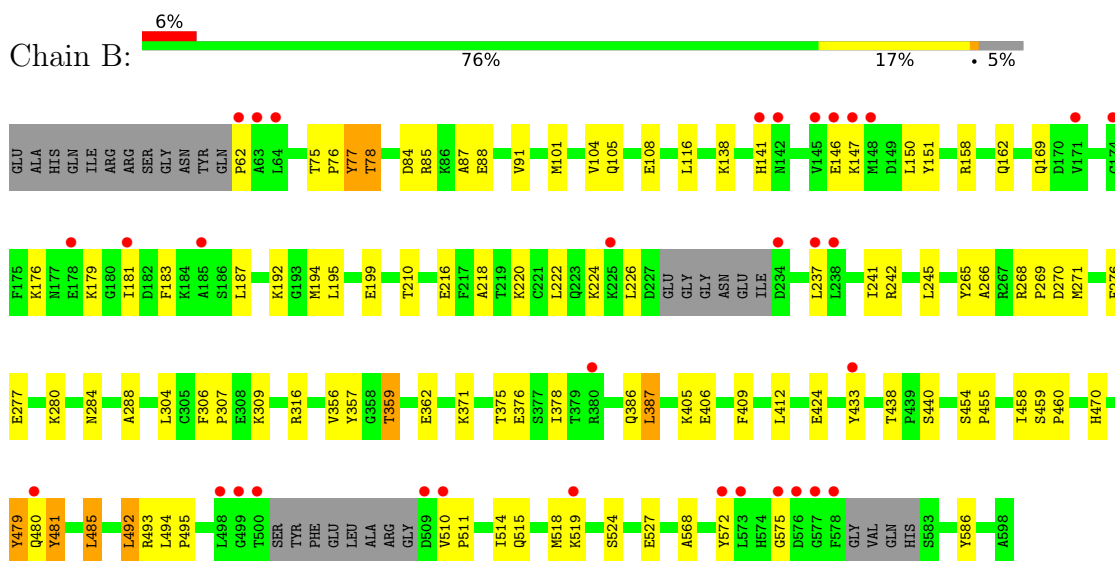
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: (+)-bornyl diphosphate synthase



- 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	102.06Å 116.77Å 120.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.83 – 2.00 19.83 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.9 (19.83-2.00) 95.9 (19.83-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.58 (at 2.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.203 , 0.234 0.203 , 0.234	Depositor DCC
R_{free} test set	3756 reflections (3.86%)	wwPDB-VP
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9156	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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, HG

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	0/4378	0.82	8/5931 (0.1%)
1	B	0.38	0/4404	0.86	12/5967 (0.2%)
All	All	0.37	0/8782	0.84	20/11898 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	480	GLN	N-CA-C	-14.08	96.12	113.01
1	A	480	GLN	N-CA-C	-13.56	96.00	112.54
1	B	356	VAL	N-CA-C	9.85	120.03	111.56
1	B	357	TYR	N-CA-C	9.45	122.92	111.40
1	B	359	THR	N-CA-C	-7.18	100.86	110.55
1	B	479	TYR	N-CA-C	7.10	121.77	113.18
1	B	440	SER	N-CA-C	-6.54	102.36	110.41
1	A	440	SER	N-CA-C	-6.49	102.42	110.41
1	B	316	ARG	N-CA-C	6.12	117.92	109.54
1	B	116	LEU	N-CA-C	-6.10	105.48	113.17
1	A	458	ILE	N-CA-C	5.98	116.15	110.53
1	B	458	ILE	N-CA-C	5.85	116.03	110.53
1	A	383	TYR	N-CA-C	5.72	117.97	111.11
1	A	479	TYR	N-CA-C	5.59	119.95	113.18
1	A	116	LEU	N-CA-C	-5.53	106.21	113.17
1	B	493	ARG	N-CA-C	5.51	116.96	111.07
1	A	357	TYR	N-CA-C	5.43	119.86	112.04
1	B	78	THR	N-CA-C	-5.41	106.61	114.12
1	B	356	VAL	CB-CA-C	-5.33	106.37	111.44
1	A	356	VAL	N-CA-C	5.19	120.13	109.34

There are no chirality outliers.

There are no planarity outliers.

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	4265	0	4152	68	0
1	B	4290	0	4174	67	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	5	0	0	0	0
3	B	6	0	0	0	0
4	A	277	0	0	1	0
4	B	311	0	0	1	0
All	All	9156	0	8326	133	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 8.

All (133) 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:138:LYS:HA	1:A:141:HIS:CE1	2.20	0.76
1:A:303:ARG:HH11	1:A:303:ARG:HB3	1.50	0.76
1:A:514:ILE:HG22	1:A:518:MET:HE2	1.69	0.75
1:B:492:LEU:HD13	1:B:568:ALA:HB2	1.70	0.74
1:A:492:LEU:HD13	1:A:568:ALA:HB2	1.71	0.72
1:B:276:PHE:CZ	1:B:280:LYS:HD2	2.25	0.72
1:A:276:PHE:CZ	1:A:280:LYS:HD2	2.25	0.71
1:B:514:ILE:HG22	1:B:518:MET:HE2	1.73	0.70
1:B:479:TYR:C	1:B:481:TYR:N	2.45	0.68
1:B:479:TYR:C	1:B:481:TYR:H	1.95	0.66
1:A:310:LEU:O	1:A:313:VAL:HG12	1.96	0.65
1:A:479:TYR:C	1:A:481:TYR:H	2.01	0.65
1:A:494:LEU:HB2	1:A:495:PRO:HD3	1.78	0.65
1:A:479:TYR:C	1:A:481:TYR:N	2.45	0.63
1:B:146:GLU:HG3	1:B:147:LYS:H	1.64	0.62
1:A:146:GLU:HG3	1:A:147:LYS:H	1.65	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:494:LEU:HB2	1:B:495:PRO:HD3	1.80	0.61
1:A:313:VAL:HG21	1:A:349:VAL:HG13	1.80	0.61
1:B:187:LEU:HB3	1:B:194:MET:HE3	1.81	0.61
1:A:101:MET:HG2	1:A:105:GLN:HB2	1.82	0.61
1:A:64:LEU:HG	1:A:65:TRP:H	1.64	0.61
1:B:62:PRO:HG3	1:B:586:TYR:CD2	2.36	0.60
1:A:267:ARG:HH11	1:A:267:ARG:HG2	1.67	0.59
1:B:378:ILE:HG12	1:B:386:GLN:HG2	1.84	0.58
1:A:158:ARG:O	1:A:162:GLN:HG3	2.04	0.58
1:B:138:LYS:HA	1:B:141:HIS:CE1	2.39	0.58
1:B:158:ARG:O	1:B:162:GLN:HG3	2.04	0.58
1:B:359:THR:OG1	1:B:362:GLU:HG3	2.04	0.57
1:A:280:LYS:HE3	1:A:597:TYR:CD2	2.39	0.57
1:B:306:PHE:HB2	1:B:307:PRO:HD3	1.87	0.57
1:A:313:VAL:HG21	1:A:349:VAL:HG22	1.86	0.57
1:B:524:SER:OG	1:B:527:GLU:HG3	2.05	0.57
1:B:218:ALA:O	1:B:222:LEU:HD13	2.05	0.56
1:B:309:LYS:HG3	1:B:387:LEU:HD23	1.87	0.56
1:A:175:PHE:HB3	1:A:187:LEU:HD11	1.88	0.56
1:A:303:ARG:HB3	1:A:303:ARG:NH1	2.19	0.56
1:B:226:LEU:HD22	1:B:245:LEU:HD12	1.87	0.55
1:B:179:LYS:HB3	1:B:181:ILE:HD13	1.87	0.55
1:B:265:TYR:CE2	1:B:271:MET:HG2	2.41	0.55
1:B:409:PHE:HB2	1:B:470:HIS:CD2	2.42	0.55
1:B:309:LYS:HG3	1:B:387:LEU:CD2	2.36	0.55
1:B:181:ILE:HD12	1:B:181:ILE:N	2.22	0.55
1:B:237:LEU:O	1:B:241:ILE:HG13	2.07	0.55
1:A:524:SER:OG	1:A:527:GLU:HG3	2.07	0.54
1:B:101:MET:HG2	1:B:105:GLN:HB2	1.89	0.54
1:A:237:LEU:O	1:A:241:ILE:HG13	2.08	0.53
1:A:220:LYS:O	1:A:224:LYS:HG3	2.08	0.52
1:B:376:GLU:HG3	4:B:1135:HOH:O	2.09	0.52
1:A:142:ASN:HB3	1:A:144:GLU:HG3	1.92	0.51
1:A:104:VAL:O	1:A:108:GLU:HG3	2.09	0.51
1:A:357:TYR:HD2	1:A:384:TYR:CZ	2.29	0.50
1:B:459:SER:OG	1:B:460:PRO:HD3	2.12	0.50
1:B:150:LEU:HD21	1:B:194:MET:HE2	1.93	0.50
1:A:64:LEU:HG	1:A:65:TRP:N	2.27	0.50
1:B:85:ARG:NH2	1:B:277:GLU:HG2	2.28	0.49
1:A:438:THR:HG21	1:A:519:LYS:HD3	1.94	0.49
1:B:104:VAL:O	1:B:108:GLU:HG3	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:PHE:CE1	1:B:280:LYS:HD2	2.47	0.48
1:B:515:GLN:CD	1:B:515:GLN:H	2.22	0.48
1:B:77:TYR:CD1	1:B:284:ASN:HB3	2.49	0.48
1:A:309:LYS:C	1:A:310:LEU:HD12	2.38	0.48
1:A:515:GLN:H	1:A:515:GLN:CD	2.21	0.48
1:B:438:THR:HG21	1:B:519:LYS:HD3	1.95	0.48
1:B:433:TYR:CD1	1:B:510:VAL:HG22	2.49	0.48
1:B:492:LEU:CD1	1:B:568:ALA:HB2	2.41	0.47
1:A:433:TYR:CD1	1:A:510:VAL:HG22	2.49	0.47
1:A:271:MET:HE3	1:A:273:PRO:HA	1.95	0.47
1:B:226:LEU:HD21	1:B:242:ARG:HG2	1.95	0.47
1:A:146:GLU:HG3	1:A:147:LYS:N	2.30	0.46
1:A:192:LYS:HG3	4:A:1147:HOH:O	2.15	0.46
1:A:253:ILE:HD11	1:A:566:ARG:HB2	1.97	0.46
1:A:492:LEU:HD12	1:A:492:LEU:O	2.16	0.46
1:B:85:ARG:NH2	1:B:277:GLU:CG	2.79	0.46
1:B:146:GLU:HG3	1:B:147:LYS:N	2.30	0.46
1:B:220:LYS:O	1:B:224:LYS:HG3	2.17	0.45
1:B:169:GLN:OE1	1:B:210:THR:HB	2.17	0.45
1:B:78:THR:HG23	1:B:288:ALA:HB1	1.97	0.45
1:A:375:THR:HB	1:B:412:LEU:HD23	1.98	0.45
1:A:405:LYS:NZ	1:A:406:GLU:HG2	2.32	0.45
1:B:176:LYS:HA	1:B:183:PHE:HA	1.99	0.45
1:A:371:LYS:HD3	1:A:424:GLU:OE2	2.16	0.45
1:B:405:LYS:NZ	1:B:406:GLU:HG2	2.32	0.45
1:B:194:MET:HB3	1:B:222:LEU:HD11	1.98	0.45
1:B:454:SER:HB2	1:B:455:PRO:HD3	1.98	0.44
1:A:452:VAL:CG2	1:A:492:LEU:HG	2.47	0.44
1:A:459:SER:OG	1:A:460:PRO:HD3	2.17	0.44
1:B:276:PHE:CE2	1:B:280:LYS:HD2	2.53	0.44
1:B:151:TYR:CD1	1:B:192:LYS:HG2	2.53	0.44
1:A:280:LYS:HE3	1:A:597:TYR:CE2	2.53	0.43
1:A:409:PHE:CG	1:A:470:HIS:CD2	3.07	0.43
1:A:481:TYR:HD2	1:A:481:TYR:HA	1.70	0.43
1:B:409:PHE:HB2	1:B:470:HIS:NE2	2.34	0.43
1:A:186:SER:O	1:A:189:GLN:HG2	2.19	0.43
1:A:268:ARG:HA	1:A:269:PRO:HD3	1.90	0.43
1:A:306:PHE:HB2	1:A:307:PRO:HD3	2.01	0.42
1:B:266:ALA:HB2	1:B:276:PHE:CE1	2.54	0.42
1:B:485:LEU:HD23	1:B:485:LEU:N	2.34	0.42
1:A:454:SER:HB2	1:A:455:PRO:HD3	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:GLN:HB3	1:A:274:LEU:HD13	2.01	0.42
1:A:480:GLN:O	1:A:481:TYR:HB2	2.19	0.42
1:A:572:TYR:HA	1:A:575:GLY:O	2.20	0.42
1:A:158:ARG:HG3	1:A:203:LEU:HD11	2.01	0.42
1:B:405:LYS:HD2	1:B:405:LYS:C	2.44	0.42
1:B:481:TYR:HD2	1:B:481:TYR:HA	1.80	0.42
1:A:510:VAL:HA	1:A:511:PRO:HD3	1.88	0.42
1:B:75:THR:HA	1:B:76:PRO:HD3	1.88	0.42
1:A:87:ALA:O	1:A:91:VAL:HG23	2.19	0.42
1:B:150:LEU:CD2	1:B:194:MET:HE2	2.50	0.42
1:B:268:ARG:HA	1:B:269:PRO:HD3	1.93	0.41
1:A:95:ILE:O	1:A:99:GLU:HG3	2.20	0.41
1:B:510:VAL:HA	1:B:511:PRO:HD3	1.90	0.41
1:A:280:LYS:HE3	1:A:597:TYR:HD2	1.81	0.41
1:B:84:ASP:O	1:B:88:GLU:HG3	2.21	0.41
1:A:359:THR:OG1	1:A:362:GLU:HG3	2.20	0.41
1:B:268:ARG:HB3	1:B:270:ASP:OD1	2.20	0.41
1:A:313:VAL:CG2	1:A:349:VAL:HG22	2.49	0.41
1:A:492:LEU:CD1	1:A:568:ALA:HB2	2.45	0.41
1:A:190:ASP:OD2	1:A:193:GLY:HA3	2.21	0.41
1:A:194:MET:HE3	1:A:218:ALA:HA	2.02	0.41
1:A:190:ASP:O	1:A:194:MET:HG2	2.20	0.41
1:A:226:LEU:HD22	1:A:245:LEU:HD12	2.03	0.41
1:A:234:ASP:HB3	1:A:237:LEU:HB3	2.03	0.41
1:A:405:LYS:HD2	1:A:405:LYS:C	2.46	0.41
1:A:452:VAL:HG22	1:A:492:LEU:HG	2.03	0.41
1:A:477:SER:O	1:A:480:GLN:HG2	2.20	0.41
1:B:216:GLU:O	1:B:220:LYS:HG3	2.21	0.41
1:B:371:LYS:HD3	1:B:424:GLU:OE2	2.20	0.41
1:A:412:LEU:HD23	1:B:375:THR:HB	2.03	0.41
1:B:87:ALA:O	1:B:91:VAL:HG23	2.20	0.41
1:B:245:LEU:HD23	1:B:245:LEU:HA	1.90	0.41
1:A:276:PHE:CE2	1:A:280:LYS:HD2	2.56	0.40
1:B:195:LEU:O	1:B:199:GLU:HG2	2.21	0.40
1:B:572:TYR:HA	1:B:575:GLY:O	2.22	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	508/549 (92%)	495 (97%)	12 (2%)	1 (0%)	43	42
1	B	512/549 (93%)	499 (98%)	11 (2%)	2 (0%)	30	27
All	All	1020/1098 (93%)	994 (98%)	23 (2%)	3 (0%)	36	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	481	TYR
1	B	77	TYR
1	B	481	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	455/480 (95%)	449 (99%)	6 (1%)	61	68
1	B	458/480 (95%)	454 (99%)	4 (1%)	70	78
All	All	913/960 (95%)	903 (99%)	10 (1%)	65	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	267	ARG
1	A	303	ARG
1	A	356	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	387	LEU
1	A	485	LEU
1	A	492	LEU
1	B	304	LEU
1	B	387	LEU
1	B	485	LEU
1	B	492	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	A	123	GLN
1	A	135	ASN
1	A	142	ASN
1	A	254	GLN
1	A	470	HIS
1	A	531	HIS
1	A	574	HIS
1	B	68	ASN
1	B	92	GLN
1	B	137	HIS
1	B	141	HIS
1	B	142	ASN
1	B	143	ASN
1	B	254	GLN
1	B	282	ASN
1	B	470	HIS
1	B	574	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 13 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	516/549 (93%)	0.31	49 (9%)	14 12	17, 33, 66, 77	0
1	B	519/549 (94%)	0.18	33 (6%)	25 24	12, 30, 64, 76	1 (0%)
All	All	1035/1098 (94%)	0.24	82 (7%)	18 17	12, 32, 65, 77	1 (0%)

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	576	ASP	5.5
1	A	64	LEU	5.2
1	B	575	GLY	4.9
1	A	313	VAL	4.9
1	A	575	GLY	4.7
1	A	145	VAL	4.7
1	A	357	TYR	4.5
1	A	510	VAL	4.4
1	A	78	THR	4.1
1	A	499	GLY	4.1
1	B	578	PHE	4.0
1	A	171	VAL	3.8
1	B	148	MET	3.8
1	A	509	ASP	3.6
1	A	146	GLU	3.5
1	B	142	ASN	3.4
1	A	576	ASP	3.4
1	B	509	ASP	3.4
1	A	597	TYR	3.3
1	B	480	GLN	3.3
1	A	573	LEU	3.3
1	B	577	GLY	3.3
1	B	181	ILE	3.3
1	A	480	GLN	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	578	PHE	3.3
1	A	81	ARG	3.2
1	A	312	PHE	3.2
1	B	510	VAL	3.1
1	B	573	LEU	2.9
1	A	574	HIS	2.9
1	B	62	PRO	2.8
1	B	64	LEU	2.8
1	B	141	HIS	2.7
1	A	80	GLU	2.7
1	B	146	GLU	2.7
1	B	185	ALA	2.7
1	A	225	LYS	2.7
1	A	234	ASP	2.6
1	B	225	LYS	2.6
1	B	519	LYS	2.6
1	B	238	LEU	2.5
1	A	179	LYS	2.5
1	A	142	ASN	2.5
1	B	237	LEU	2.5
1	A	147	LYS	2.5
1	A	177	ASN	2.4
1	A	522	ASN	2.4
1	B	499	GLY	2.4
1	B	178	GLU	2.4
1	B	147	LYS	2.4
1	A	180	GLY	2.3
1	A	139	CYS	2.3
1	B	174	CYS	2.3
1	A	240	TRP	2.3
1	B	63	ALA	2.3
1	A	189	GLN	2.3
1	A	236	ASN	2.3
1	A	237	LEU	2.3
1	B	145	VAL	2.3
1	A	143	ASN	2.3
1	B	171	VAL	2.2
1	A	227	ASP	2.2
1	A	242	ARG	2.2
1	A	224	LYS	2.2
1	B	234	ASP	2.2
1	B	433	TYR	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	167	ILE	2.1
1	A	526	GLU	2.1
1	A	356	VAL	2.1
1	A	498	LEU	2.1
1	A	572	TYR	2.1
1	B	500	THR	2.1
1	A	598	ALA	2.1
1	A	141	HIS	2.1
1	B	380	ARG	2.1
1	A	144	GLU	2.0
1	B	498	LEU	2.0
1	A	174	CYS	2.0
1	A	355	ASP	2.0
1	A	303	ARG	2.0
1	B	572	TYR	2.0
1	A	76	PRO	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 oligosaccharides 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(Å ²)	Q<0.9
3	HG	B	913	1/1	0.87	0.13	53,53,53,53	1
3	HG	B	914	1/1	0.90	0.14	58,58,58,58	1
3	HG	A	906	1/1	0.91	0.13	59,59,59,59	1
3	HG	A	907	1/1	0.91	0.20	57,57,57,57	1
3	HG	B	909	1/1	0.92	0.31	34,34,34,34	1
3	HG	A	908	1/1	0.93	0.11	52,52,52,52	1
2	MG	B	901	1/1	0.93	0.09	39,39,39,39	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HG	B	915	1/1	0.94	0.09	57,57,57,57	1
2	MG	A	902	1/1	0.95	0.12	53,53,53,53	0
3	HG	A	903	1/1	0.96	0.28	33,33,33,33	1
3	HG	A	904	1/1	0.99	0.11	23,23,23,23	1
3	HG	B	910	1/1	0.99	0.05	26,26,26,26	1
3	HG	B	911	1/1	0.99	0.04	24,24,24,24	1

6.5 Other polymers [i](#)

There are no such residues in this entry.