



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:59 PM UTC

PDB ID : 5BUS / pdb_00005bus
Title : O-succinylbenzoate Coenzyme A Synthetase (MenE) from Bacillus Subtilis, in complex with AMP
Authors : Chen, Y.; Sun, Y.; Song, H.; Guo, Z.
Deposited on : 2015-06-04
Resolution : 2.60 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

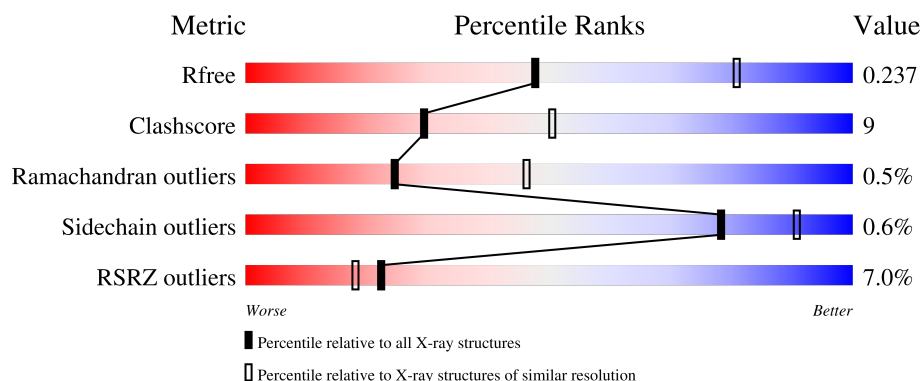
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	494	6% 84% 13% ••
1	B	494	6% 60% 16% • 23%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6493 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 2-succinylbenzoate--CoA ligase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	2	0
			3552	2256	600	673	23			
1	B	381	Total	C	N	O	S	0	2	0
			2829	1808	467	533	21			

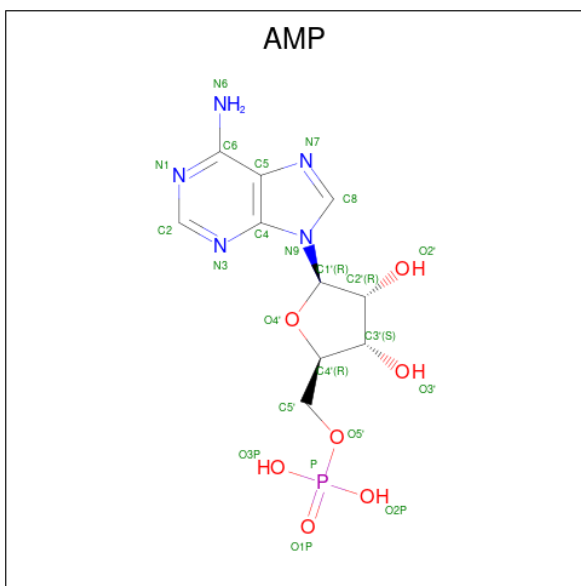
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP P23971
A	0	GLY	-	expression tag	UNP P23971
A	487	HIS	-	expression tag	UNP P23971
A	488	HIS	-	expression tag	UNP P23971
A	489	HIS	-	expression tag	UNP P23971
A	490	HIS	-	expression tag	UNP P23971
A	491	HIS	-	expression tag	UNP P23971
A	492	HIS	-	expression tag	UNP P23971
B	-1	MET	-	initiating methionine	UNP P23971
B	0	GLY	-	expression tag	UNP P23971
B	487	HIS	-	expression tag	UNP P23971
B	488	HIS	-	expression tag	UNP P23971
B	489	HIS	-	expression tag	UNP P23971
B	490	HIS	-	expression tag	UNP P23971
B	491	HIS	-	expression tag	UNP P23971
B	492	HIS	-	expression tag	UNP P23971

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 3 is ADENOSINE MONOPHOSPHATE (CCD ID: AMP) (formula: $C_{10}H_{14}N_5O_7P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		
3	B	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	44	Total	O	0	0
			44	44		
4	B	18	Total	O	0	0
			18	18		

4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	121.81Å 121.81Å 98.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.52 – 2.60 38.52 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.52-2.60) 99.5 (38.52-2.60)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.83 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.190 , 0.234 0.200 , 0.237	Depositor DCC
R_{free} test set	1997 reflections (4.56%)	wwPDB-VP
Wilson B-factor (Å ²)	60.6	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 74.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6493	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.56	0/3628	0.89	3/4936 (0.1%)
1	B	0.51	0/2885	0.87	2/3920 (0.1%)
All	All	0.54	0/6513	0.88	5/8856 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	380	LEU	N-CA-C	8.77	120.61	111.14
1	B	16	THR	CA-C-N	7.27	127.90	119.47
1	B	16	THR	C-N-CA	7.27	127.90	119.47
1	A	406	HIS	CA-C-N	5.75	125.88	119.32
1	A	406	HIS	C-N-CA	5.75	125.88	119.32

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3552	0	3307	41	0
1	B	2829	0	2633	67	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	23	0	12	0	0
3	B	23	0	12	4	0
4	A	44	0	0	1	0
4	B	18	0	0	0	0
All	All	6493	0	5964	107	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 9.

All (107) 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:329:GLU:HG2	1:B:330:PRO:HD2	1.48	0.96
1:A:400:GLU:OE1	1:A:474:ARG:NH2	2.09	0.86
1:B:329:GLU:HG2	1:B:330:PRO:CD	2.07	0.83
1:B:183:THR:HG22	1:B:185:GLN:H	1.43	0.83
1:B:365:THR:HG23	1:B:367:ASP:H	1.48	0.77
1:B:43:GLU:HG2	1:B:135:ILE:HG12	1.70	0.74
1:A:44:GLN:NE2	1:A:127:LEU:O	2.22	0.73
1:A:459:PHE:HA	1:A:483:GLY:HA3	1.69	0.72
1:B:65:GLU:CG	1:B:128:MET:HE1	2.21	0.70
1:B:336:ILE:HB	1:B:365:THR:HG21	1.75	0.68
1:A:63:ARG:NH1	1:A:107:ASP:OD2	2.28	0.66
1:B:19:ARG:NH2	1:B:184:GLU:HG3	2.13	0.64
1:B:65:GLU:HG3	1:B:128:MET:HE1	1.79	0.63
1:A:40:ARG:O	1:A:44:GLN:HG3	1.99	0.63
1:A:301:GLU:H	1:A:301:GLU:CD	2.06	0.61
1:B:23:ILE:HG12	1:B:28:THR:HG22	1.81	0.61
1:B:162:VAL:HG13	1:B:344:MET:HG3	1.82	0.61
1:A:62:ASN:HB2	1:A:217:GLN:HA	1.82	0.61
1:A:84:ASN:HB3	1:A:87:LEU:HG	1.81	0.61
1:B:41:MET:HE2	1:B:128:MET:HE2	1.84	0.60
1:B:290:GLU:HG3	1:B:344:MET:HE3	1.83	0.60
1:B:107:ASP:OD1	1:B:109:SER:HB3	2.03	0.59
1:B:41:MET:CE	1:B:128:MET:HE2	2.33	0.59
1:A:288:MET:HG3	1:A:290:GLU:HB2	1.84	0.59
1:B:19:ARG:HH22	1:B:184:GLU:HG3	1.69	0.58
1:A:337:MET:HE2	1:A:364:LYS:HG2	1.84	0.58
1:B:62:ASN:HB2	1:B:217:GLN:HA	1.84	0.58
1:B:93:LEU:HD22	1:B:118:ILE:HD12	1.86	0.56
1:A:464:LEU:O	1:A:466:ARG:HG3	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:LEU:O	1:B:92:ARG:NH2	2.41	0.55
1:A:477:LEU:O	1:A:479:ASP:HA	2.06	0.54
1:B:288:MET:HG3	1:B:290:GLU:HB2	1.90	0.54
1:A:162:VAL:HG13	1:A:344:MET:SD	2.49	0.53
1:A:117:HIS:H	1:A:117:HIS:CD2	2.26	0.53
1:B:206:SER:HB2	1:B:213:VAL:HG23	1.90	0.53
1:A:461:LEU:HD21	1:A:480:ALA:HB3	1.91	0.51
1:B:339:LYS:HD2	1:B:362:TRP:CE2	2.45	0.51
1:B:253:CYS:HB2	1:B:280:PHE:CE1	2.46	0.51
1:B:113:LYS:HA	1:B:116:GLU:HG2	1.93	0.51
1:B:290:GLU:HG3	1:B:344:MET:CE	2.41	0.51
1:B:225:LEU:HD21	1:B:246:LEU:HD22	1.93	0.50
1:B:16:THR:HB	1:B:19:ARG:HG3	1.92	0.50
1:A:429:TYR:CE2	1:A:459:PHE:HD2	2.30	0.50
1:B:186:ASP:OD2	1:B:258:ARG:NH2	2.41	0.50
1:B:266:PRO:O	1:B:268:PRO:HD3	2.11	0.50
1:B:293:SER:OG	1:B:294:GLN:N	2.45	0.50
1:B:344:MET:HE1	1:B:347:TYR:CE1	2.47	0.49
1:B:295:ILE:HD13	1:B:338:VAL:HB	1.94	0.49
1:B:367:ASP:CG	3:B:503:AMP:HO2'	2.19	0.49
1:A:313:PRO:HG3	1:A:319:ILE:HG12	1.95	0.49
1:B:197:ILE:HG12	1:B:292:CYS:HA	1.94	0.49
1:A:400:GLU:CD	1:A:474:ARG:HH22	2.20	0.49
1:B:365:THR:CG2	1:B:367:ASP:H	2.22	0.49
1:B:275:CYS:HB3	1:B:280:PHE:HB2	1.95	0.49
1:B:189:LEU:HB3	1:B:236:ILE:HG13	1.93	0.49
1:A:335:GLU:OE1	1:A:383:ARG:NH1	2.46	0.48
1:A:328:CYS:HB3	1:A:332:GLU:HB3	1.96	0.48
1:B:187:ARG:HG2	1:B:212:THR:HB	1.96	0.48
1:A:386:LEU:HD11	1:A:393:ASN:HB3	1.96	0.48
1:B:301:GLU:H	1:B:301:GLU:CD	2.21	0.48
1:B:180:LEU:O	1:B:258:ARG:NH1	2.47	0.47
1:B:367:ASP:OD2	3:B:503:AMP:O2'	2.32	0.47
1:B:371:LEU:HD23	1:B:377:LEU:HA	1.96	0.47
1:A:197:ILE:HG12	1:A:292:CYS:HA	1.97	0.47
1:A:387:ILE:O	1:A:394:ILE:N	2.48	0.46
1:B:192:LEU:HD21	1:B:242:MET:SD	2.56	0.46
1:B:365:THR:HG23	1:B:367:ASP:N	2.25	0.46
1:B:304:MET:HA	1:B:305:GLU:HA	1.49	0.46
1:B:329:GLU:HG2	1:B:330:PRO:HD3	1.94	0.46
1:A:175:SER:OG	1:A:315:PHE:O	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:LEU:HD22	1:B:91:GLU:CD	2.41	0.45
1:B:162:VAL:HG13	1:B:344:MET:HE2	1.99	0.44
1:B:292:CYS:O	1:B:293:SER:CB	2.65	0.44
1:A:123:ASP:HB3	1:A:126:GLU:HB2	1.99	0.44
1:A:193:PRO:HG2	1:A:196[B]:HIS:ND1	2.33	0.44
1:A:140:TYR:OH	1:B:318:GLU:OE1	2.31	0.43
1:B:290:GLU:OE1	1:B:290:GLU:N	2.45	0.43
1:A:477:LEU:C	1:A:479:ASP:HA	2.43	0.43
1:B:75:LEU:HD22	1:B:137:ILE:HD12	2.00	0.43
1:B:367:ASP:HA	1:B:382:ARG:HA	2.01	0.43
1:B:253:CYS:HB2	1:B:280:PHE:CZ	2.54	0.43
1:A:406:HIS:HA	1:A:407:PRO:HD3	1.90	0.43
1:B:9:LEU:HA	1:B:207:VAL:CG2	2.49	0.43
1:B:321:ILE:HD12	1:B:371:LEU:HG	1.99	0.43
1:B:43:GLU:CD	1:B:134:GLU:HA	2.43	0.43
1:A:125:ASP:HA	1:A:128:MET:HE2	2.01	0.42
1:A:399:VAL:HB	1:A:414:VAL:HG21	2.00	0.42
1:B:336:ILE:HB	1:B:365:THR:CG2	2.47	0.42
1:A:73:CYS:HB3	1:A:78:VAL:O	2.19	0.42
1:A:460:VAL:C	1:A:461:LEU:HD12	2.44	0.42
1:B:319:ILE:HD11	1:B:338:VAL:HG12	2.02	0.42
1:A:44:GLN:HE22	1:A:128:MET:C	2.28	0.42
1:A:117:HIS:HD2	4:A:601:HOH:O	2.02	0.42
1:B:367:ASP:CG	3:B:503:AMP:O2'	2.62	0.42
1:A:211:MET:HE2	1:A:211:MET:HB3	1.99	0.42
1:A:474:ARG:HA	1:A:477:LEU:CD1	2.49	0.42
1:A:344:MET:HE1	1:A:347:TYR:CD1	2.55	0.41
1:B:336:ILE:O	1:B:365:THR:HG22	2.20	0.41
1:B:253:CYS:HA	1:B:254:PRO:HD3	1.89	0.41
1:A:380:LEU:O	1:A:381:ASP:HB3	2.21	0.41
1:B:239:VAL:HG22	1:B:242:MET:SD	2.61	0.41
1:A:168:ASN:HB3	1:A:292:CYS:SG	2.61	0.41
1:B:224:VAL:O	1:B:228:ILE:HG13	2.20	0.41
1:B:264:GLY:O	3:B:503:AMP:N6	2.54	0.41
1:B:162:VAL:HG22	1:B:344:MET:HE2	2.03	0.40
1:A:309:SER:HB2	1:A:378:TYR:CE2	2.56	0.40
1:B:300:PRO:C	1:B:302:PHE:H	2.30	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	479/494 (97%)	451 (94%)	26 (5%)	2 (0%)	30	51
1	B	377/494 (76%)	361 (96%)	14 (4%)	2 (0%)	24	46
All	All	856/988 (87%)	812 (95%)	40 (5%)	4 (0%)	24	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	293	SER
1	A	381	ASP
1	B	293	SER
1	B	221	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	347/421 (82%)	346 (100%)	1 (0%)	86	94
1	B	276/421 (66%)	273 (99%)	3 (1%)	65	84
All	All	623/842 (74%)	619 (99%)	4 (1%)	78	91

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	221	VAL
1	B	182	ILE

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Mol	Chain	Res	Type
1	B	224	VAL
1	B	327	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	240	GLN
1	B	61	GLN
1	B	163	GLN
1	B	226	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	AMP	B	503	-	25,25,25	1.55	6 (24%)	37,38,38	1.92	9 (24%)
3	AMP	A	503	-	25,25,25	1.84	7 (28%)	37,38,38	1.98	9 (24%)

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	AMP	B	503	-	-	3/10/26/26	0/3/3/3
3	AMP	A	503	-	-	0/10/26/26	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	503	AMP	C5-C4	3.98	1.46	1.39
3	A	503	AMP	C5-N7	-3.88	1.32	1.39
3	A	503	AMP	C8-N9	-3.38	1.31	1.37
3	A	503	AMP	C5-C4	3.19	1.44	1.39
3	B	503	AMP	C5-N7	-3.10	1.33	1.39
3	A	503	AMP	C4-N9	-3.06	1.31	1.37
3	B	503	AMP	C4-N9	-2.73	1.32	1.37
3	A	503	AMP	P-O2P	-2.73	1.44	1.54
3	A	503	AMP	P-O3P	-2.42	1.45	1.54
3	B	503	AMP	C5-C6	2.23	1.47	1.41
3	B	503	AMP	C8-N9	-2.21	1.33	1.37
3	A	503	AMP	C5-C6	2.10	1.46	1.41
3	B	503	AMP	C8-N7	2.01	1.35	1.31

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	503	AMP	C5-C4-N3	-5.86	118.65	126.72
3	B	503	AMP	C5-C4-N3	-5.47	119.18	126.72
3	B	503	AMP	N3-C4-N9	4.22	134.34	127.17
3	A	503	AMP	C4-C5-N7	-4.07	105.93	110.58
3	A	503	AMP	N3-C4-N9	3.97	133.91	127.17
3	A	503	AMP	C2-N3-C4	3.78	121.06	111.83
3	B	503	AMP	C2'-C1'-N9	-3.72	104.06	113.30
3	B	503	AMP	C2-N3-C4	3.61	120.65	111.83
3	A	503	AMP	C2'-C1'-N9	-3.37	104.94	113.30
3	A	503	AMP	N3-C2-N1	-3.23	123.69	128.58
3	B	503	AMP	C4-C5-N7	-3.18	106.94	110.58
3	B	503	AMP	N3-C2-N1	-3.08	123.92	128.58
3	A	503	AMP	C6-C5-N7	2.73	137.35	132.09
3	B	503	AMP	C4-N9-C8	2.48	108.34	105.74
3	A	503	AMP	C5-N7-C8	2.38	107.19	103.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	503	AMP	O3P-P-O5'	-2.16	101.05	106.67
3	B	503	AMP	C5-N7-C8	2.15	106.82	103.45
3	B	503	AMP	C6-C5-N7	2.05	136.04	132.09

There are no chirality outliers.

All (3) torsion outliers are listed below:

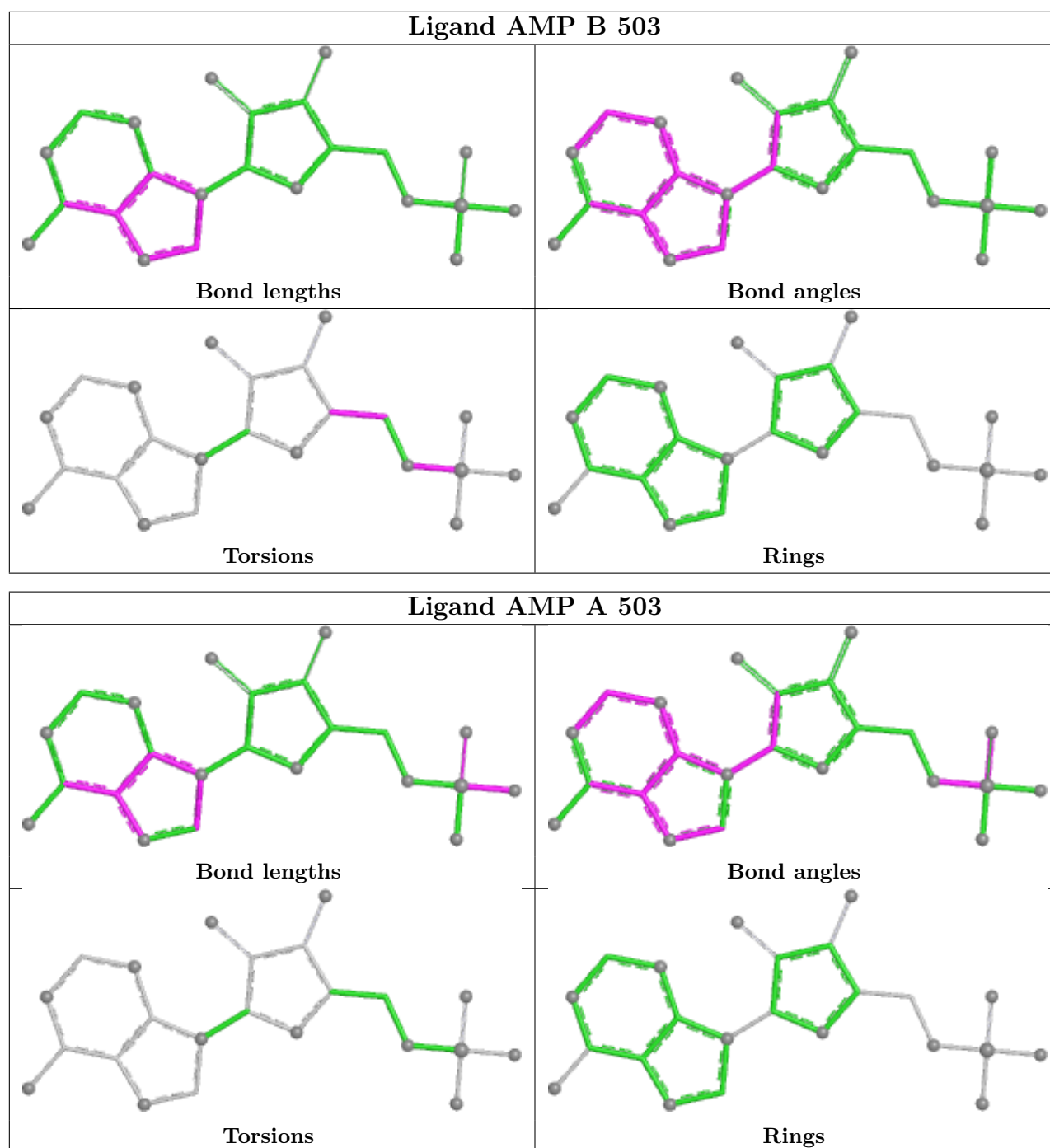
Mol	Chain	Res	Type	Atoms
3	B	503	AMP	C5'-O5'-P-O2P
3	B	503	AMP	C3'-C4'-C5'-O5'
3	B	503	AMP	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	503	AMP	4	0

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 [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	481/494 (97%)	0.21	28 (5%) 29 23	24, 67, 107, 129	2 (0%)
1	B	381/494 (77%)	0.36	32 (8%) 17 13	37, 73, 124, 152	2 (0%)
All	All	862/988 (87%)	0.27	60 (6%) 22 18	24, 70, 118, 152	4 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	373	ASN	5.0
1	B	381	ASP	5.0
1	B	304	MET	4.4
1	B	324	ASP	4.2
1	B	268	PRO	4.1
1	B	153	SER	3.9
1	A	390	GLY	3.9
1	A	388	ILE	3.8
1	B	196[A]	HIS	3.6
1	B	305	GLU	3.5
1	A	430	LEU	3.4
1	A	251	ASN	3.2
1	B	26	ASP	3.1
1	A	391	GLY	3.0
1	B	270	PRO	3.0
1	B	250	THR	2.9
1	A	26	ASP	2.9
1	A	421	LYS	2.9
1	B	386	LEU	2.9
1	A	475	ASN	2.9
1	B	266	PRO	2.8
1	B	307	LEU	2.8
1	B	3	THR	2.8
1	A	485	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	484	GLU	2.8
1	A	126	GLU	2.7
1	A	135	ILE	2.7
1	B	247	LEU	2.7
1	A	458	PHE	2.7
1	A	424	LYS	2.7
1	B	241	THR	2.6
1	A	419	ASP	2.5
1	B	23	ILE	2.5
1	A	479	ASP	2.5
1	A	459	PHE	2.5
1	A	420	LYS	2.4
1	B	382	ARG	2.4
1	B	383	ARG	2.4
1	B	375	GLY	2.4
1	A	229	ASN	2.4
1	A	461	LEU	2.3
1	B	269	LEU	2.3
1	B	271	LEU	2.3
1	B	145	ALA	2.2
1	A	217	GLN	2.2
1	A	478	LYS	2.2
1	B	274	GLU	2.2
1	A	437	SER	2.2
1	B	154	GLY	2.2
1	A	116	GLU	2.1
1	B	155	THR	2.1
1	A	425	VAL	2.1
1	A	440	GLU	2.1
1	B	27	GLN	2.1
1	B	287	GLY	2.1
1	B	384	SER	2.0
1	B	379	VAL	2.0
1	A	481	ARG	2.0
1	A	473	LEU	2.0
1	B	229	ASN	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 [i](#)

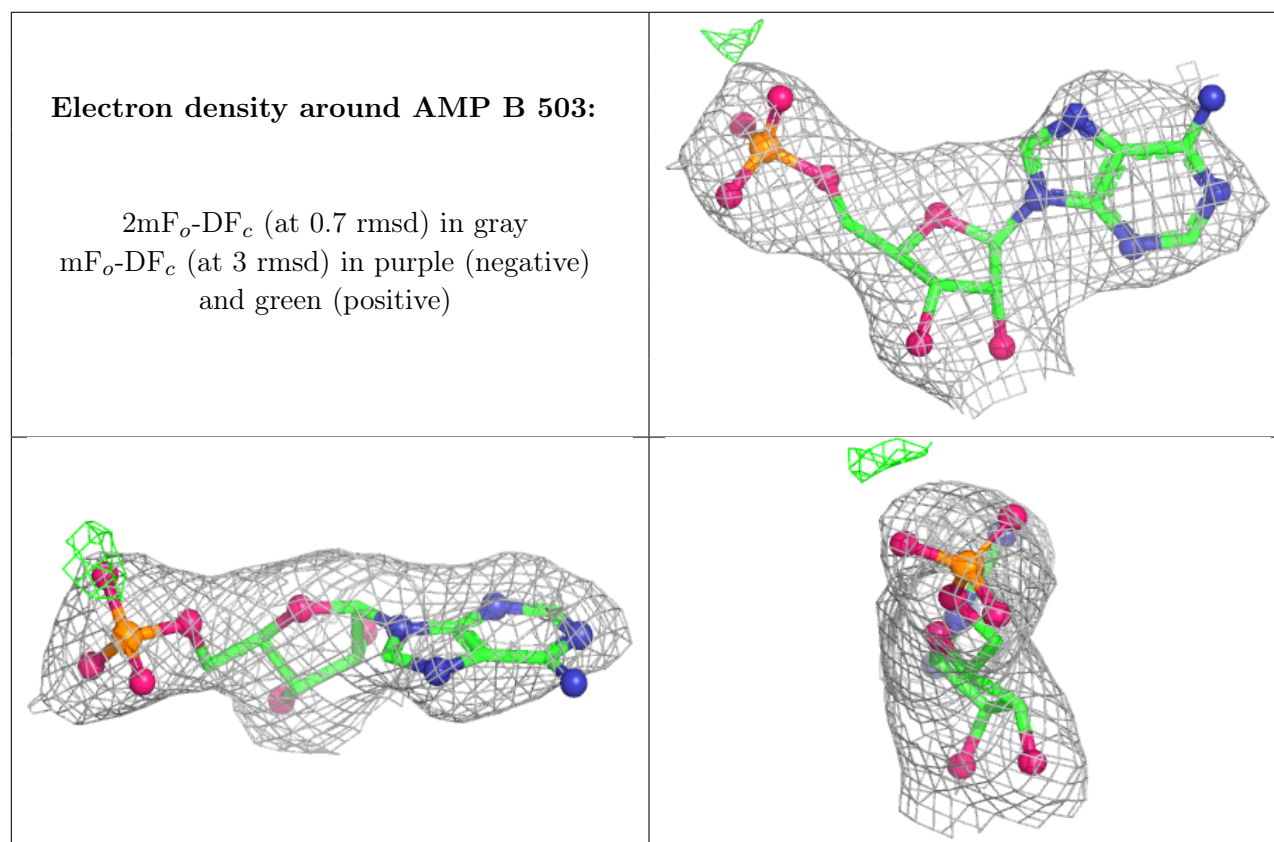
There are no oligosaccharides in this entry.

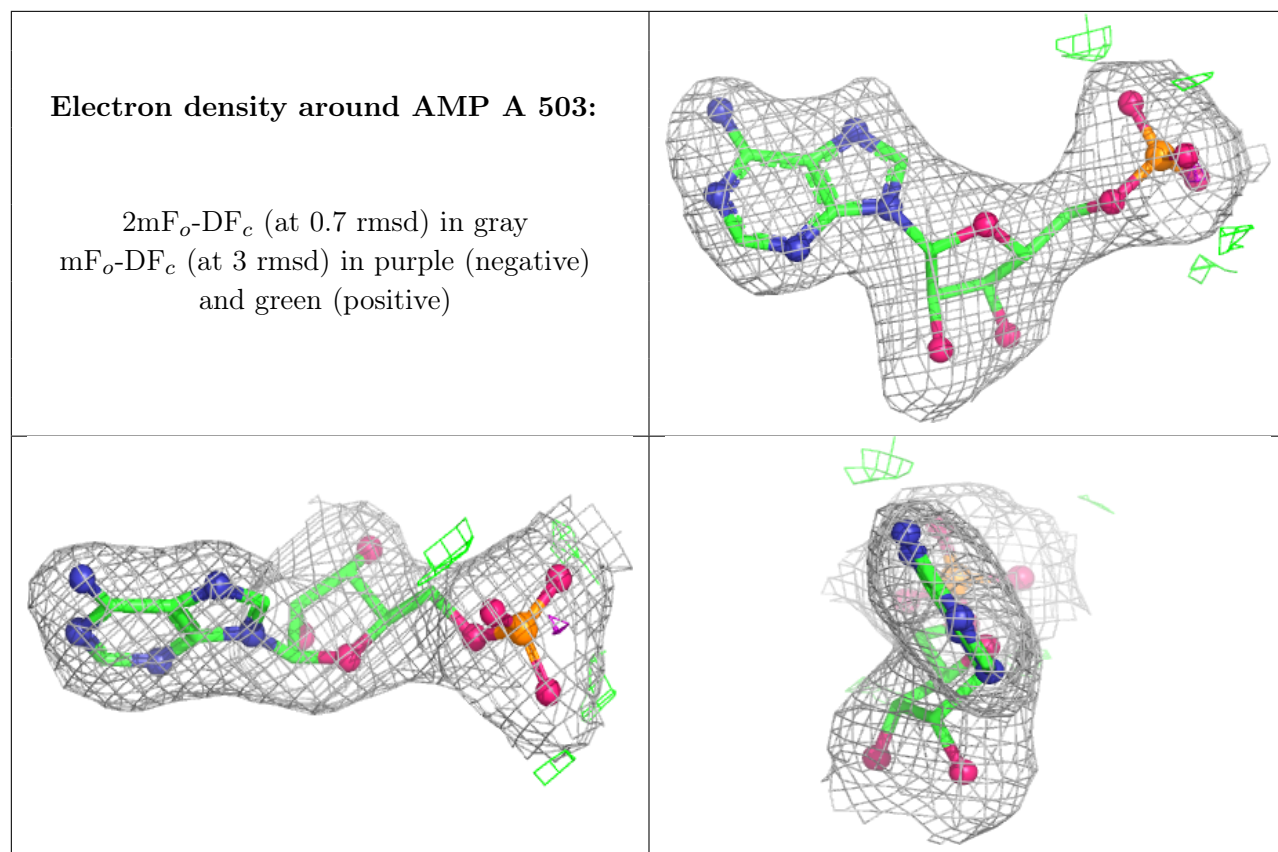
6.4 Ligands [i](#)

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	AMP	B	503	23/23	0.90	0.12	83,102,115,118	0
2	CL	B	501	1/1	0.91	0.19	113,113,113,113	0
2	CL	A	501	1/1	0.91	0.24	106,106,106,106	0
2	CL	B	502	1/1	0.94	0.12	108,108,108,108	0
2	CL	A	502	1/1	0.94	0.11	108,108,108,108	0
3	AMP	A	503	23/23	0.97	0.07	41,49,56,66	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.