



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:09 PM UTC

PDB ID : 5BUR / pdb_00005bur
Title : O-succinylbenzoate Coenzyme A Synthetase (MenE) from Bacillus Subtilis, in Complex with ATP and Magnesium Ion
Authors : Chen, Y.; Sun, Y.; Song, H.; Guo, Z.
Deposited on : 2015-06-04
Resolution : 2.82 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

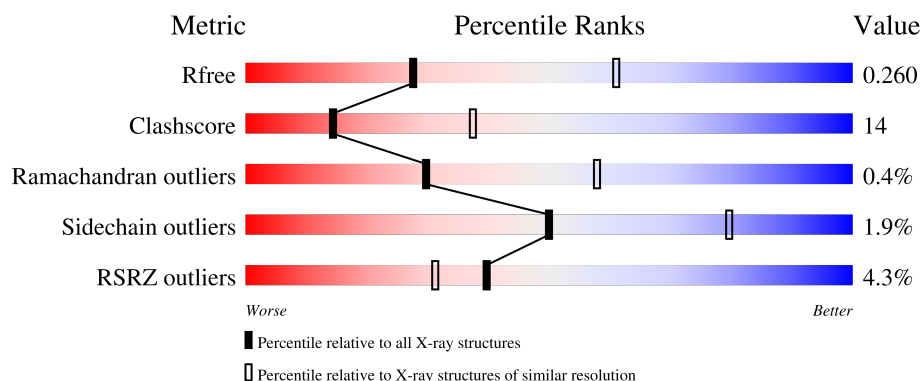
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

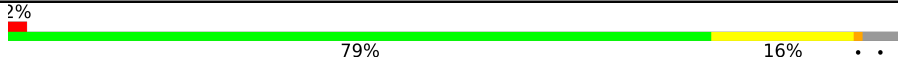
The reported resolution of this entry is 2.82 Å.

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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	
1	B	494	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6269 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	475	Total	C	N	O	S	0	1	0
			3514	2241	590	660	23			
1	B	365	Total	C	N	O	S	0	0	0
			2632	1680	439	493	20			

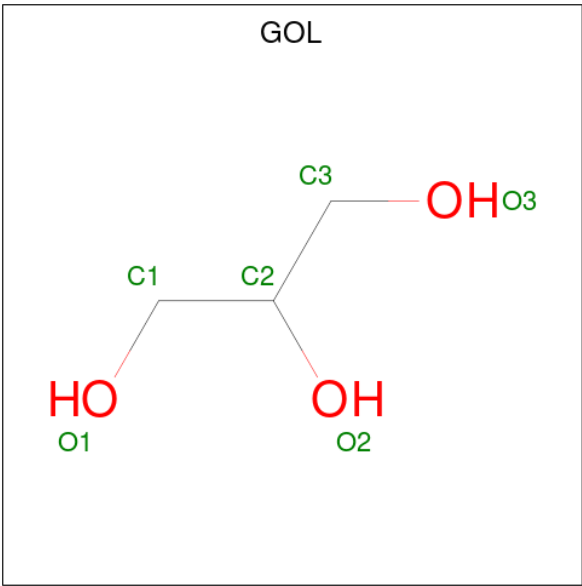
There are 16 discrepancies between the modelled and reference sequences:

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	3	Total	Cl	0	0
			3	3		
2	B	3	Total	Cl	0	0
			3	3		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).

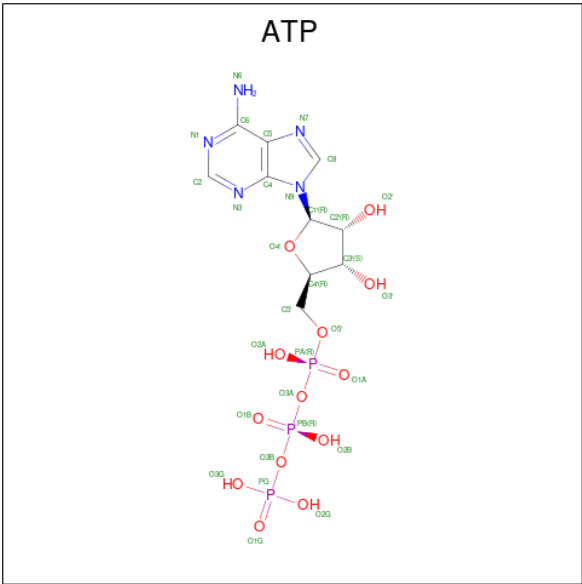


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

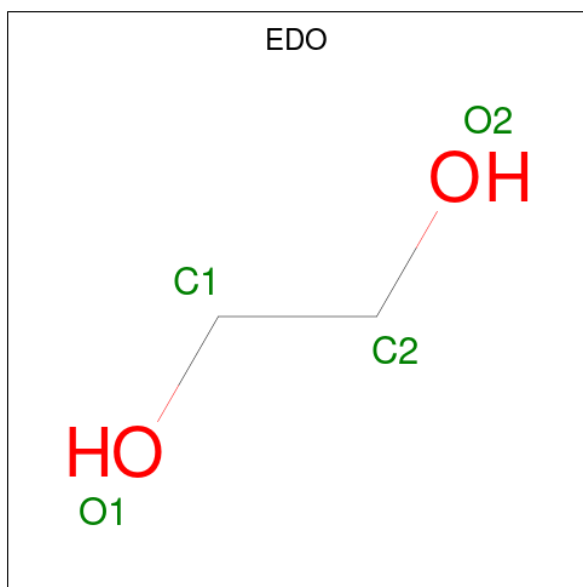
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
5	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).

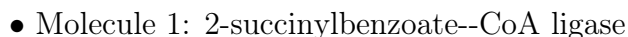


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	29	Total	O	0	0
			29	29		
7	B	15	Total	O	0	0
			15	15		

- Molecule 1: 2-succinylbenzoate--CoA ligase



LEU
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	121.83Å 121.83Å 97.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.39 – 2.82 36.39 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.8 (36.39-2.82) 99.7 (36.39-2.82)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.204 , 0.252 0.217 , 0.260	Depositor DCC
R_{free} test set	1999 reflections (4.05%)	wwPDB-VP
Wilson B-factor (Å ²)	69.2	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 52.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6269	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: ATP, EDO, MG, GOL, CL

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	0.65	1/3585 (0.0%)	0.96	13/4877 (0.3%)
1	B	0.71	1/2677 (0.0%)	0.99	7/3636 (0.2%)
All	All	0.67	2/6262 (0.0%)	0.97	20/8513 (0.2%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	454	ILE	C-N	5.41	1.40	1.33
1	B	254	PRO	N-CD	5.32	1.55	1.47

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	ALA	N-CA-C	-11.11	94.83	110.50
1	A	442	THR	N-CA-C	-8.43	103.51	113.88
1	B	118	ILE	CB-CA-C	-7.82	102.58	111.80
1	A	389	SER	N-CA-C	-7.37	96.51	108.52
1	A	158	LYS	CA-C-N	6.33	126.35	119.90
1	A	158	LYS	C-N-CA	6.33	126.35	119.90
1	A	454	ILE	CA-C-N	-6.28	113.48	119.76
1	A	454	ILE	C-N-CA	-6.28	113.48	119.76
1	A	269	LEU	CA-C-N	-6.06	112.77	119.19

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	LEU	C-N-CA	-6.06	112.77	119.19
1	A	434	LYS	CA-C-N	-5.51	112.96	119.84
1	A	434	LYS	C-N-CA	-5.51	112.96	119.84
1	B	253	CYS	CA-C-N	-5.44	113.03	119.84
1	B	253	CYS	C-N-CA	-5.44	113.03	119.84
1	A	25	GLU	CB-CA-C	-5.38	109.92	117.23
1	B	259	CYS	N-CA-C	5.27	116.63	107.93
1	A	346	SER	N-CA-C	5.23	117.02	108.55
1	B	246	LEU	N-CA-C	-5.22	104.40	111.55
1	B	238	ALA	N-CA-C	5.11	116.06	108.60
1	A	432	LEU	N-CA-C	5.09	117.27	109.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	389	SER	Peptide

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	3514	0	3314	60	0
1	B	2632	0	2362	109	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	6	0	8	0	0
4	A	1	0	0	0	0
5	A	31	0	12	5	0
5	B	31	0	12	7	0
6	A	4	0	6	0	0
7	A	29	0	0	1	0
7	B	15	0	0	1	0
All	All	6269	0	5714	168	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 14.

All (168) 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:344:MET:CE	1:B:346:SER:C	2.15	1.19
1:B:23:ILE:HD11	1:B:214:VAL:HG22	1.21	1.16
1:B:23:ILE:HD11	1:B:214:VAL:CG2	1.76	1.15
1:B:23:ILE:CD1	1:B:214:VAL:HG13	1.75	1.15
1:B:152:THR:CB	5:B:504:ATP:O3G	1.95	1.15
1:A:392:GLU:CB	1:A:393:ASN:HA	1.75	1.13
1:B:152:THR:HB	5:B:504:ATP:O3G	1.53	1.08
1:B:306:LYS:CB	1:B:309:SER:CB	2.34	1.06
1:B:274:GLU:HA	1:B:275:CYS:C	1.78	1.05
1:B:152:THR:OG1	5:B:504:ATP:O3G	1.77	1.02
1:B:23:ILE:HD11	1:B:214:VAL:HG13	1.45	0.98
1:B:23:ILE:HD11	1:B:214:VAL:CG1	1.93	0.98
1:B:274:GLU:HA	1:B:275:CYS:O	1.62	0.97
1:B:23:ILE:CD1	1:B:214:VAL:HG22	1.93	0.97
1:B:274:GLU:CA	1:B:275:CYS:C	2.37	0.97
1:B:344:MET:HE1	1:B:347:TYR:N	1.80	0.94
1:A:455:PRO:HB2	1:A:458:PHE:HE1	1.32	0.93
1:B:344:MET:HE1	1:B:346:SER:C	1.94	0.93
1:A:440:GLU:HA	1:A:442:THR:H	1.36	0.91
1:B:344:MET:HE3	1:B:346:SER:C	2.01	0.85
1:A:386:LEU:HD11	1:A:393:ASN:HB3	1.58	0.85
1:B:23:ILE:HD11	1:B:214:VAL:CB	2.07	0.84
1:A:455:PRO:HB2	1:A:458:PHE:CE1	2.13	0.82
1:B:87:LEU:O	1:B:92:ARG:NH2	2.13	0.81
1:A:392:GLU:CB	1:A:393:ASN:CA	2.57	0.81
1:B:23:ILE:HD12	1:B:214:VAL:HG13	1.63	0.80
1:B:353:ALA:O	1:B:357:SER:OG	1.98	0.79
1:B:45:LEU:O	1:B:47:ALA:O	2.01	0.78
1:B:93:LEU:HD22	1:B:118:ILE:HD11	1.66	0.78
1:A:410:ALA:HB2	1:A:433:HIS:HD2	1.46	0.78
1:B:274:GLU:CB	1:B:275:CYS:C	2.57	0.77
1:B:23:ILE:CG1	1:B:214:VAL:HA	2.16	0.75
1:A:84:ASN:HB3	1:A:87:LEU:HG	1.68	0.75
1:A:440:GLU:HA	1:A:442:THR:N	2.03	0.74
1:B:92:ARG:NH1	1:B:115:TYR:OH	2.21	0.74
1:B:45:LEU:C	1:B:47:ALA:O	2.30	0.74
1:A:152:THR:OG1	5:A:506:ATP:O3G	2.06	0.73
1:B:23:ILE:HG13	1:B:214:VAL:HA	1.69	0.72
1:B:344:MET:CE	1:B:347:TYR:N	2.45	0.72
1:B:8:TRP:CE2	1:B:141:MET:HE1	2.25	0.71

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:MET:HE3	1:B:346:SER:O	1.91	0.71
1:B:248:GLU:HA	1:B:249:GLU:CB	2.21	0.70
1:A:187:ARG:HG2	1:A:212:THR:HB	1.74	0.70
5:A:506:ATP:O2B	7:A:601:HOH:O	2.09	0.69
1:B:344:MET:HE1	1:B:347:TYR:CA	2.23	0.68
1:B:93:LEU:HD23	1:B:118:ILE:HD12	1.77	0.67
1:B:328:CYS:HB3	1:B:332:GLU:HG3	1.76	0.66
1:A:439:GLY:HA3	1:A:442:THR:N	2.10	0.66
1:A:247:LEU:O	1:A:278:LYS:NZ	2.30	0.65
1:B:93:LEU:HD22	1:B:118:ILE:CD1	2.26	0.65
1:A:437:SER:OG	1:A:440:GLU:O	2.11	0.64
1:B:183:THR:HG22	1:B:185:GLN:H	1.62	0.64
1:B:367:ASP:OD2	5:B:504:ATP:O2'	2.13	0.63
1:A:410:ALA:HB2	1:A:433:HIS:CD2	2.31	0.63
1:B:344:MET:HE2	1:B:346:SER:C	2.19	0.63
1:A:4:GLU:OE1	1:B:339:LYS:NZ	2.31	0.63
1:B:47:ALA:O	1:B:48:HIS:HB2	1.97	0.63
1:A:445:CYS:HB3	1:A:454:ILE:HG23	1.81	0.63
1:B:330:PRO:O	1:B:331:TYR:HB2	1.99	0.61
1:B:274:GLU:CA	1:B:275:CYS:O	2.41	0.61
1:A:439:GLY:HA2	1:A:440:GLU:C	2.26	0.61
1:B:344:MET:CE	1:B:346:SER:O	2.45	0.61
1:B:23:ILE:HG13	1:B:23:ILE:O	2.01	0.61
1:B:93:LEU:CD2	1:B:118:ILE:CD1	2.79	0.61
1:B:58:ILE:HB	1:B:82:LEU:HD23	1.84	0.59
1:B:23:ILE:CD1	1:B:214:VAL:CG1	2.55	0.59
1:B:253:CYS:SG	1:B:257:ILE:HB	2.42	0.59
1:B:349:ASN:O	1:B:350:ARG:HD3	2.02	0.59
1:B:93:LEU:CD2	1:B:118:ILE:HD12	2.32	0.59
1:B:315:PHE:O	1:B:316:SER:CB	2.51	0.58
1:B:344:MET:CE	1:B:346:SER:CA	2.81	0.58
1:A:121:THR:O	1:A:122:ILE:HG13	2.05	0.57
1:B:7:ASN:HA	1:B:141:MET:HG3	1.87	0.57
1:A:388:ILE:HA	1:A:392:GLU:CB	2.35	0.57
1:B:306:LYS:CB	1:B:309:SER:CA	2.81	0.56
1:B:336:ILE:O	1:B:365:THR:HG22	2.05	0.56
1:B:326:GLN:HG3	1:B:327:VAL:H	1.71	0.56
1:B:344:MET:HE2	1:B:346:SER:CA	2.35	0.56
1:B:23:ILE:CG1	1:B:214:VAL:HG13	2.34	0.55
1:A:406:HIS:NE2	1:A:441:LEU:CB	2.69	0.55
1:B:274:GLU:HA	1:B:275:CYS:CB	2.35	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:TRP:CZ2	1:B:141:MET:HE1	2.41	0.54
1:B:248:GLU:CA	1:B:249:GLU:CB	2.85	0.54
1:B:344:MET:HE1	1:B:347:TYR:HA	1.90	0.54
1:B:275:CYS:N	1:B:276:ARG:HA	2.23	0.54
1:B:288:MET:HB2	5:B:504:ATP:H5'1	1.89	0.53
1:A:152:THR:CB	5:A:506:ATP:O3G	2.57	0.53
1:B:344:MET:O	1:B:344:MET:HG3	2.07	0.53
1:B:79:LYS:HD3	1:B:146:THR:HB	1.91	0.52
1:B:74:PHE:HE2	1:B:204:PHE:HE1	1.57	0.52
1:A:19:ARG:NH2	1:A:184:GLU:HG2	2.24	0.52
1:B:306:LYS:CB	1:B:309:SER:HA	2.38	0.52
1:A:106:THR:OG1	1:A:107:ASP:N	2.41	0.52
1:A:140:TYR:OH	1:B:318:GLU:OE1	2.14	0.52
1:A:136:GLU:CD	1:A:137:ILE:HA	2.35	0.51
1:B:328:CYS:HB3	1:B:332:GLU:CG	2.39	0.51
1:A:344:MET:HE1	1:A:347:TYR:CD1	2.46	0.51
1:B:325:GLY:N	1:B:326:GLN:OE1	2.44	0.51
1:A:25:GLU:OE2	1:A:25:GLU:HA	2.10	0.51
1:A:466:ARG:HD3	1:A:470:ASN:O	2.11	0.51
1:B:274:GLU:CB	1:B:275:CYS:O	2.59	0.50
1:B:81:VAL:HG22	1:B:149:LEU:HB3	1.94	0.49
1:B:274:GLU:CB	1:B:276:ARG:N	2.75	0.49
1:A:455:PRO:CB	1:A:458:PHE:HE1	2.16	0.49
1:B:297:THR:O	1:B:311:GLY:HA3	2.13	0.49
1:A:439:GLY:CA	1:A:440:GLU:C	2.85	0.49
1:B:59:LEU:HD12	1:B:85:THR:HG22	1.95	0.49
1:B:294:GLN:H	1:B:294:GLN:CD	2.21	0.48
1:A:309:SER:HB2	1:A:378:TYR:CE2	2.48	0.48
1:B:133:GLU:O	1:B:135:ILE:N	2.46	0.48
1:B:189:LEU:HD22	1:B:236:ILE:HD13	1.95	0.48
1:A:187:ARG:NH1	1:A:231:HIS:O	2.39	0.48
1:B:277:GLU:O	1:B:279:GLY:HA2	2.14	0.48
1:B:332:GLU:HG3	1:B:332:GLU:O	2.12	0.48
1:B:367:ASP:CG	5:B:504:ATP:O2'	2.57	0.48
1:B:192:LEU:HD21	1:B:242:MET:SD	2.54	0.48
1:A:197:ILE:HG12	1:A:292:CYS:HA	1.96	0.47
1:A:285:SER:O	5:A:506:ATP:N6	2.45	0.47
1:B:350:ARG:HB3	1:B:353:ALA:HB3	1.95	0.47
1:B:20:ILE:O	1:B:212:THR:OG1	2.29	0.46
1:A:440:GLU:CA	1:A:442:THR:H	2.19	0.46
1:B:253:CYS:HB3	1:B:279:GLY:O	2.15	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:PRO:HG2	1:A:196:HIS:ND1	2.32	0.45
1:A:62:ASN:O	1:A:217:GLN:HG3	2.16	0.45
5:B:504:ATP:H5'2	5:B:504:ATP:H8	1.82	0.45
1:A:182:ILE:HG13	1:B:181:GLY:HA2	1.99	0.45
1:B:336:ILE:HB	1:B:365:THR:HG21	1.98	0.45
1:A:62:ASN:HB2	1:A:217:GLN:HA	1.99	0.44
1:B:74:PHE:CE2	1:B:204:PHE:HE1	2.34	0.44
1:A:406:HIS:HB3	1:A:408:ALA:N	2.32	0.44
1:A:253:CYS:HB2	1:A:280:PHE:CD1	2.53	0.44
1:B:318:GLU:HB2	1:B:339:LYS:HB3	2.00	0.44
1:B:4:GLU:N	7:B:603:HOH:O	2.35	0.44
1:B:358:PHE:CE1	1:B:363:LEU:HD13	2.53	0.44
1:B:62:ASN:HB2	1:B:217:GLN:HA	1.98	0.44
1:B:248:GLU:HA	1:B:249:GLU:C	2.43	0.43
1:B:62:ASN:O	1:B:217:GLN:HG2	2.17	0.43
1:B:149:LEU:HD12	1:B:162:VAL:O	2.18	0.43
1:B:162:VAL:HG13	1:B:344:MET:SD	2.59	0.43
1:A:295:ILE:HG23	1:A:296:VAL:HG23	2.00	0.43
1:A:24:TYR:O	1:A:25:GLU:HB2	2.17	0.43
1:A:292:CYS:O	1:A:293:SER:OG	2.30	0.43
1:A:406:HIS:N	1:A:407:PRO:HA	2.33	0.43
1:B:365:THR:HG23	1:B:367:ASP:N	2.34	0.43
1:A:464:LEU:O	1:A:466:ARG:HG3	2.19	0.42
1:A:463:ARG:C	1:A:464:LEU:HD23	2.44	0.42
1:B:220:SER:O	1:B:224:VAL:HB	2.19	0.42
1:B:9:LEU:HD21	1:B:35:PHE:HA	2.01	0.42
1:A:136:GLU:HA	1:A:137:ILE:HA	1.68	0.42
1:A:74:PHE:HE2	1:A:204:PHE:HE1	1.68	0.42
1:A:59:LEU:HD12	1:A:85:THR:HG22	2.02	0.42
1:A:114:GLU:O	1:A:117:HIS:HE1	2.03	0.42
1:B:211:MET:HE2	1:B:211:MET:HB3	1.94	0.41
1:B:315:PHE:O	1:B:316:SER:HB2	2.20	0.41
1:A:400:GLU:O	1:A:404:LEU:HD13	2.20	0.41
1:B:118:ILE:HG21	1:B:118:ILE:HD13	1.85	0.41
1:A:152:THR:HB	5:A:506:ATP:O3G	2.20	0.41
1:B:59:LEU:HD13	1:B:92:ARG:HD3	2.02	0.41
1:B:318:GLU:OE2	1:B:339:LYS:HD3	2.21	0.41
1:A:79:LYS:HE3	1:A:146:THR:HB	2.03	0.41
1:A:344:MET:HE1	1:A:347:TYR:CE1	2.56	0.41
1:B:288:MET:HG3	1:B:290:GLU:H	1.86	0.41
1:A:337:MET:HE2	1:A:364:LYS:HG2	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:ILE:HD12	1:A:370:TYR:HA	2.02	0.40
1:B:158:LYS:HA	1:B:159:PRO:HD2	1.88	0.40
1:B:347:TYR:HB2	1:B:350:ARG:HB2	2.03	0.40
1:A:269:LEU:O	1:A:273:GLU:HG3	2.21	0.40
1:A:472:LEU:HD22	1:A:474:ARG:HG3	2.02	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	472/494 (96%)	438 (93%)	32 (7%)	2 (0%)	30	58
1	B	351/494 (71%)	324 (92%)	26 (7%)	1 (0%)	36	64
All	All	823/988 (83%)	762 (93%)	58 (7%)	3 (0%)	30	58

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	293	SER
1	B	293	SER
1	A	465	PRO

5.3.2 Protein sidechains [i](#)

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%)	342 (99%)	5 (1%)	59	84
1	B	240/421 (57%)	234 (98%)	6 (2%)	42	74
All	All	587/842 (70%)	576 (98%)	11 (2%)	50	79

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

Mol	Chain	Res	Type
1	A	221	VAL
1	A	321	ILE
1	A	327	VAL
1	A	402	VAL
1	A	471	LYS
1	B	148	THR
1	B	185	GLN
1	B	213	VAL
1	B	224	VAL
1	B	314	LEU
1	B	332	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	163	GLN
1	A	179	ASN
1	A	240	GLN
1	A	284	GLN
1	A	373	ASN
1	A	433	HIS
1	B	61	GLN

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 11 ligands modelled in this entry, 7 are monoatomic - leaving 4 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
5	ATP	A	506	4	32,33,33	1.60	5 (15%)	48,52,52	1.97	10 (20%)
5	ATP	B	504	-	32,33,33	1.35	4 (12%)	48,52,52	1.85	7 (14%)
6	EDO	A	507	-	3,3,3	0.33	0	2,2,2	0.25	0
3	GOL	A	504	-	5,5,5	0.28	0	5,5,5	0.34	0

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
5	ATP	A	506	4	-	11/22/38/38	0/3/3/3
5	ATP	B	504	-	-	9/22/38/38	0/3/3/3
6	EDO	A	507	-	-	1/1/1/1	-
3	GOL	A	504	-	-	0/4/4/4	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	506	ATP	C5-N7	-4.06	1.31	1.39
5	B	504	ATP	C5-C4	3.94	1.46	1.39
5	A	506	ATP	C8-N9	-3.11	1.32	1.37
5	B	504	ATP	C5-N7	-3.00	1.33	1.39
5	A	506	ATP	C5-C4	2.77	1.44	1.39
5	A	506	ATP	C4-N9	-2.75	1.32	1.37

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	504	ATP	C5-C6	2.22	1.47	1.41
5	B	504	ATP	C4-N9	-2.20	1.33	1.37
5	A	506	ATP	PA-O2A	-2.09	1.45	1.55

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	506	ATP	C5-C4-N3	-6.75	117.42	126.72
5	B	504	ATP	C5-C4-N3	-6.53	117.72	126.72
5	A	506	ATP	N3-C4-N9	5.08	135.81	127.17
5	B	504	ATP	N3-C4-N9	4.91	135.51	127.17
5	A	506	ATP	C2-N3-C4	3.98	121.56	111.83
5	B	504	ATP	C2-N3-C4	3.93	121.42	111.83
5	A	506	ATP	C4-C5-N7	-3.53	106.55	110.58
5	B	504	ATP	C4-C5-N7	-3.39	106.71	110.58
5	A	506	ATP	N3-C2-N1	-3.38	123.47	128.58
5	B	504	ATP	N3-C2-N1	-3.15	123.82	128.58
5	A	506	ATP	C5-N7-C8	2.61	107.56	103.45
5	B	504	ATP	C3'-C2'-C1'	2.50	106.20	101.46
5	B	504	ATP	C5-N7-C8	2.41	107.24	103.45
5	A	506	ATP	O3B-PB-O1B	-2.24	103.95	110.70
5	A	506	ATP	O3G-PG-O2G	2.18	115.99	107.80
5	A	506	ATP	C3'-C2'-C1'	2.05	105.33	101.46
5	A	506	ATP	C2'-C1'-N9	-2.04	108.23	113.30

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	506	ATP	PB-O3B-PG-O2G
5	A	506	ATP	PB-O3B-PG-O3G
5	A	506	ATP	C5'-O5'-PA-O1A
5	A	506	ATP	C5'-O5'-PA-O2A
5	A	506	ATP	C5'-O5'-PA-O3A
5	B	504	ATP	PB-O3B-PG-O2G
5	B	504	ATP	C5'-O5'-PA-O1A
5	B	504	ATP	C5'-O5'-PA-O2A
5	B	504	ATP	C5'-O5'-PA-O3A
6	A	507	EDO	O1-C1-C2-O2
5	A	506	ATP	C3'-C4'-C5'-O5'
5	A	506	ATP	PB-O3A-PA-O5'
5	B	504	ATP	PG-O3B-PB-O1B

Continued on next page...

Continued from previous page...

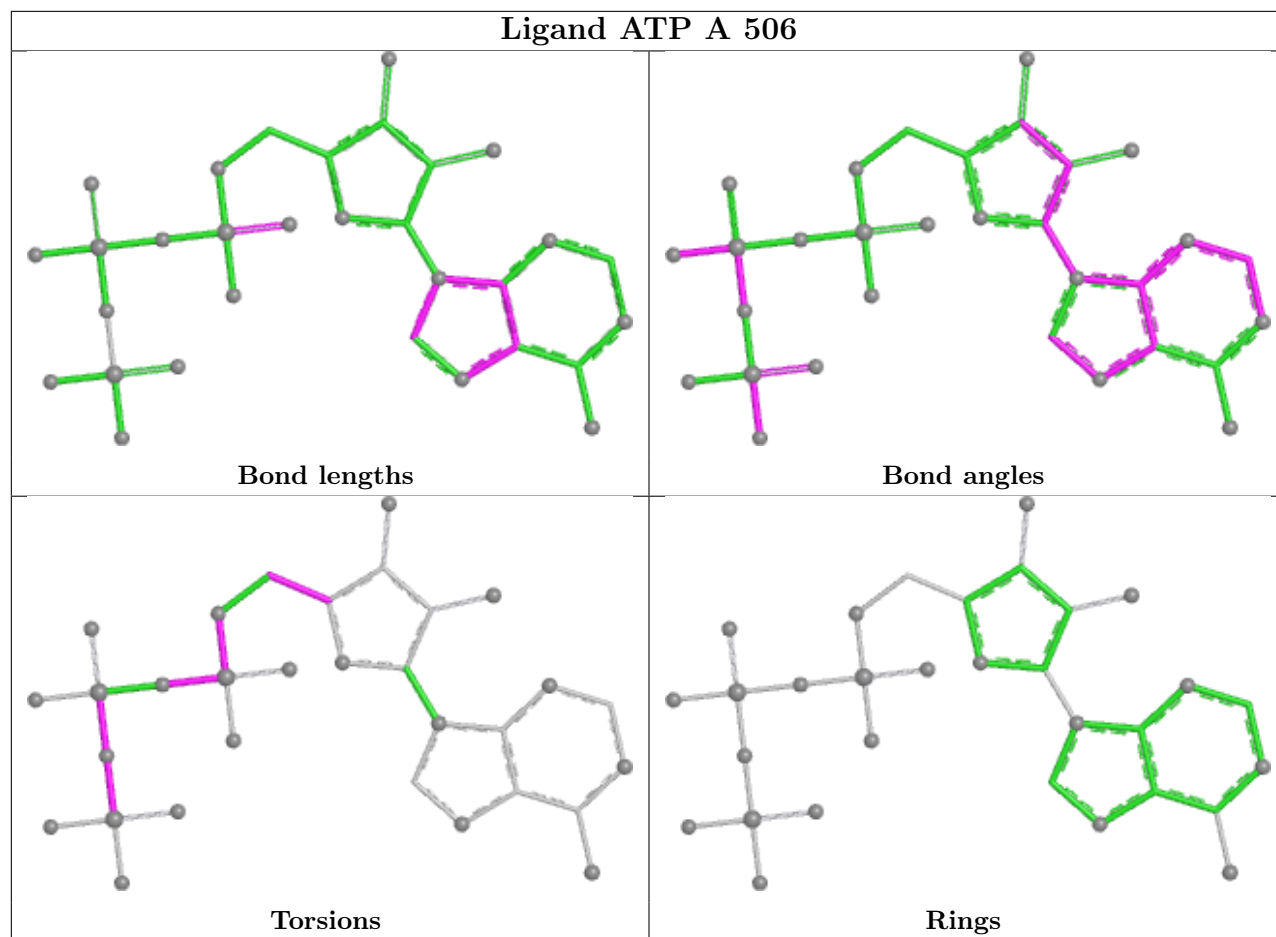
Mol	Chain	Res	Type	Atoms
5	A	506	ATP	O4'-C4'-C5'-O5'
5	B	504	ATP	C4'-C5'-O5'-PA
5	A	506	ATP	PB-O3B-PG-O1G
5	B	504	ATP	PB-O3B-PG-O1G
5	A	506	ATP	PG-O3B-PB-O1B
5	B	504	ATP	PG-O3B-PB-O2B
5	A	506	ATP	PG-O3B-PB-O2B
5	B	504	ATP	C3'-C4'-C5'-O5'

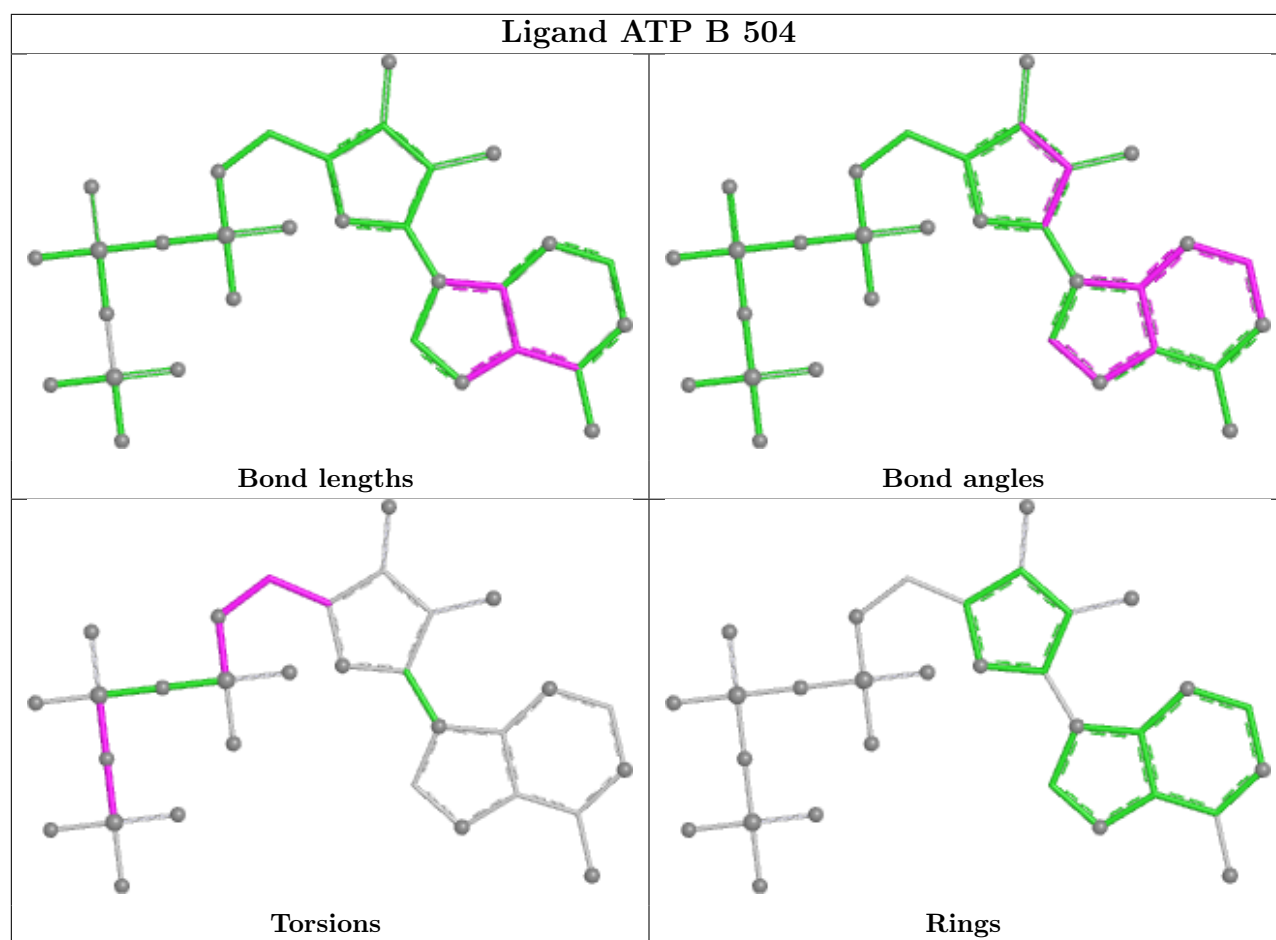
There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	506	ATP	5	0
5	B	504	ATP	7	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	475/494 (96%)	0.11	10 (2%) 63 54	27, 68, 99, 115	1 (0%)
1	B	365/494 (73%)	0.46	26 (7%) 22 16	46, 72, 108, 132	0
All	All	840/988 (85%)	0.26	36 (4%) 40 31	27, 69, 102, 132	1 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	253	CYS	5.5
1	A	418	GLU	4.2
1	A	393	ASN	3.7
1	B	220	SER	3.5
1	B	23	ILE	3.4
1	B	247	LEU	3.4
1	A	464	LEU	3.1
1	B	309	SER	3.0
1	B	327	VAL	2.9
1	A	456	ALA	2.8
1	B	303	SER	2.7
1	B	3	THR	2.7
1	B	381	ASP	2.7
1	B	320	LYS	2.6
1	B	304	MET	2.6
1	B	321	ILE	2.5
1	B	198	SER	2.4
1	A	438	ALA	2.4
1	B	280	PHE	2.4
1	B	223	ASP	2.4
1	B	275	CYS	2.3
1	A	27	GLN	2.3
1	B	22	LEU	2.3
1	B	252	ARG	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	268	PRO	2.3
1	B	246	LEU	2.2
1	B	27	GLN	2.2
1	B	277	GLU	2.2
1	A	185	GLN	2.1
1	A	49	SER	2.1
1	B	221	VAL	2.1
1	B	156	THR	2.1
1	A	388	ILE	2.0
1	B	47	ALA	2.0
1	A	114	GLU	2.0
1	B	357	SER	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](#)

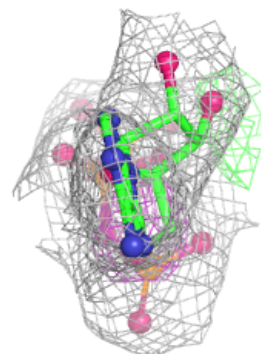
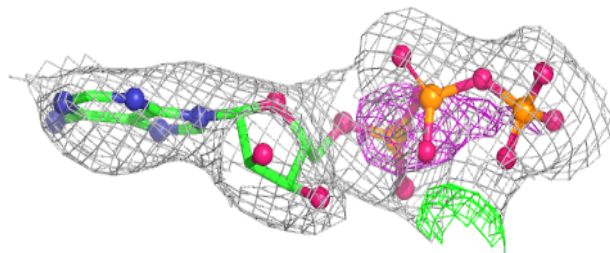
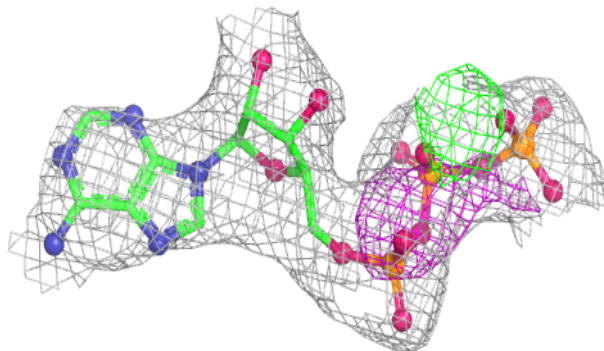
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

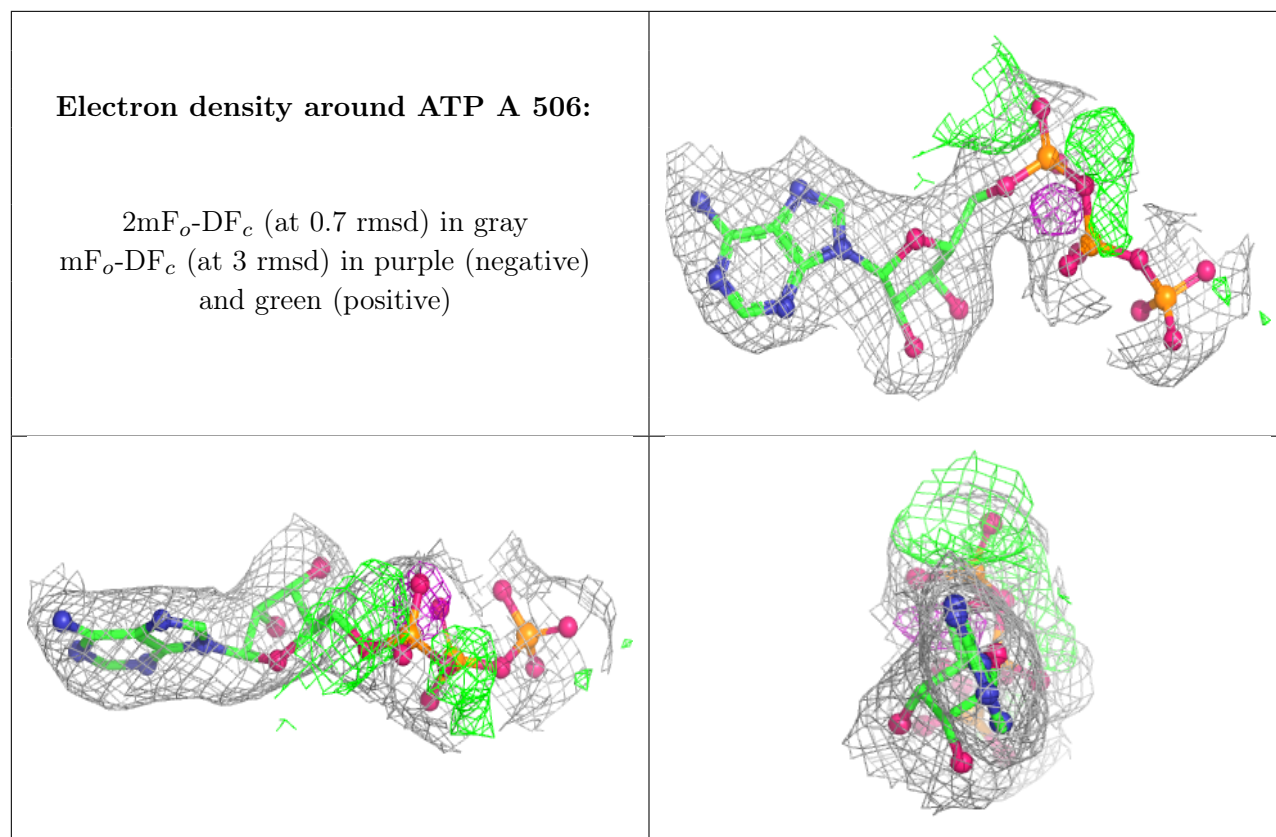
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	ATP	B	504	31/31	0.74	0.12	90,108,118,122	0
2	CL	A	503	1/1	0.78	0.17	100,100,100,100	0
3	GOL	A	504	6/6	0.79	0.14	86,91,92,94	0
6	EDO	A	507	4/4	0.84	0.24	58,70,70,79	0
2	CL	A	501	1/1	0.91	0.13	88,88,88,88	0
2	CL	B	502	1/1	0.91	0.18	93,93,93,93	0
2	CL	B	503	1/1	0.91	0.15	96,96,96,96	0
2	CL	A	502	1/1	0.92	0.28	80,80,80,80	0
5	ATP	A	506	31/31	0.92	0.10	46,58,73,75	0
4	MG	A	505	1/1	0.95	0.11	64,64,64,64	0
2	CL	B	501	1/1	0.95	0.10	74,74,74,74	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 ATP B 504:

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.