



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:20 AM UTC

PDB ID : 1I7L / pdb_00001i7l
Title : CRYSTAL STRUCTURE ANALYSIS OF THE COMPLEX OF THE C DOMAIN OF SYNAPSIN II FROM RAT WITH ATP
Authors : Esser, L.; Palnitkar, M.; Deisenhofer, J.
Deposited on : 2001-03-09
Resolution : 2.35 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

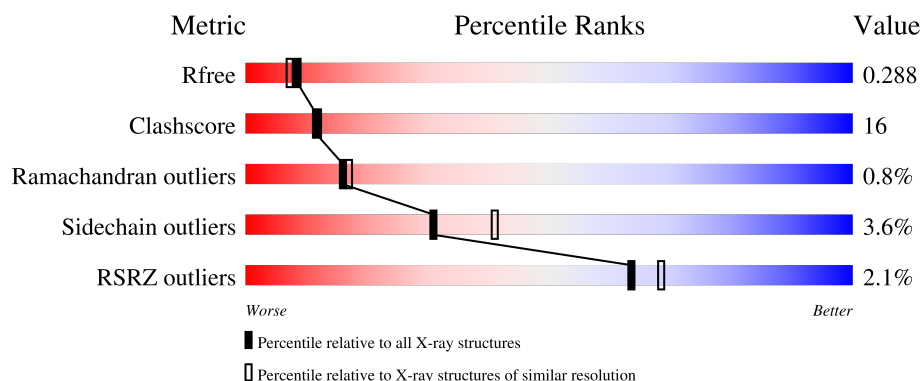
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

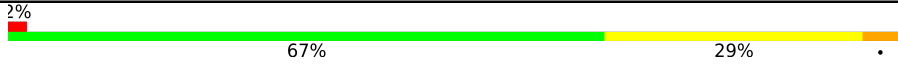
The reported resolution of this entry is 2.35 Å.

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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	309	
1	B	309	

2 Entry composition [i](#)

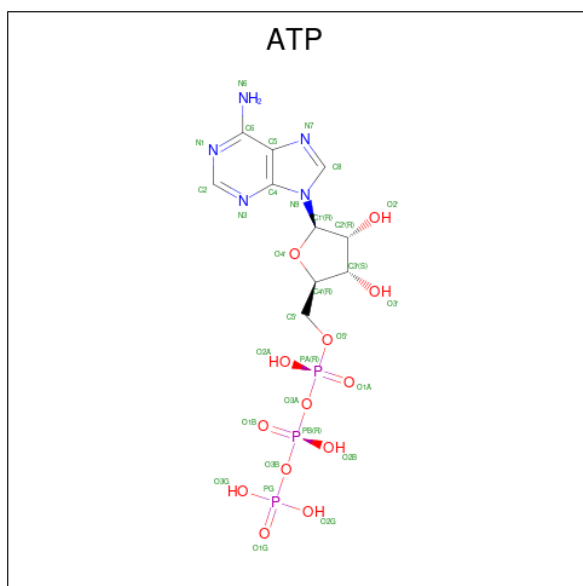
There are 3 unique types of molecules in this entry. The entry contains 5181 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SYNAPSIN II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	309	Total	C	N	O	S	55	0	0
			2462	1575	416	453	18			
1	B	309	Total	C	N	O	S	30	0	0
			2462	1575	416	453	18			

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



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

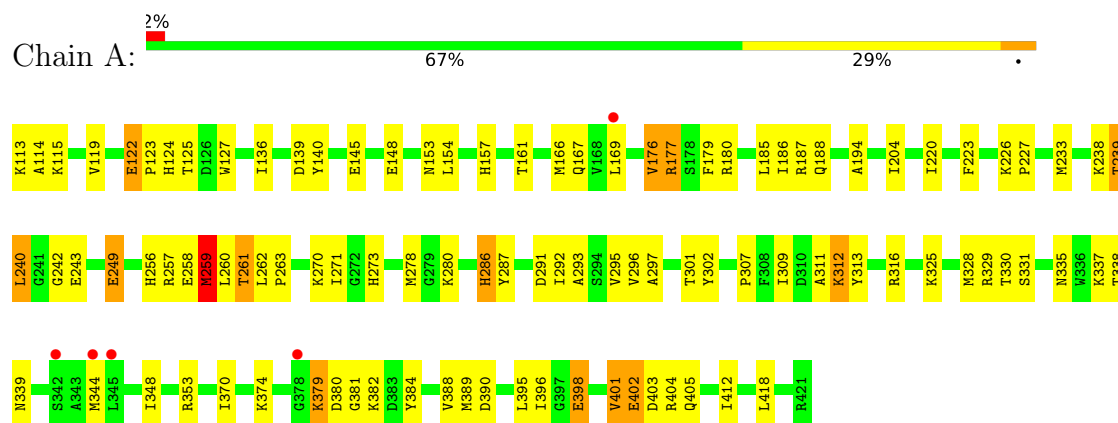
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	94	Total 94	O 94	0	0
3	B	101	Total 101	O 101	0	0

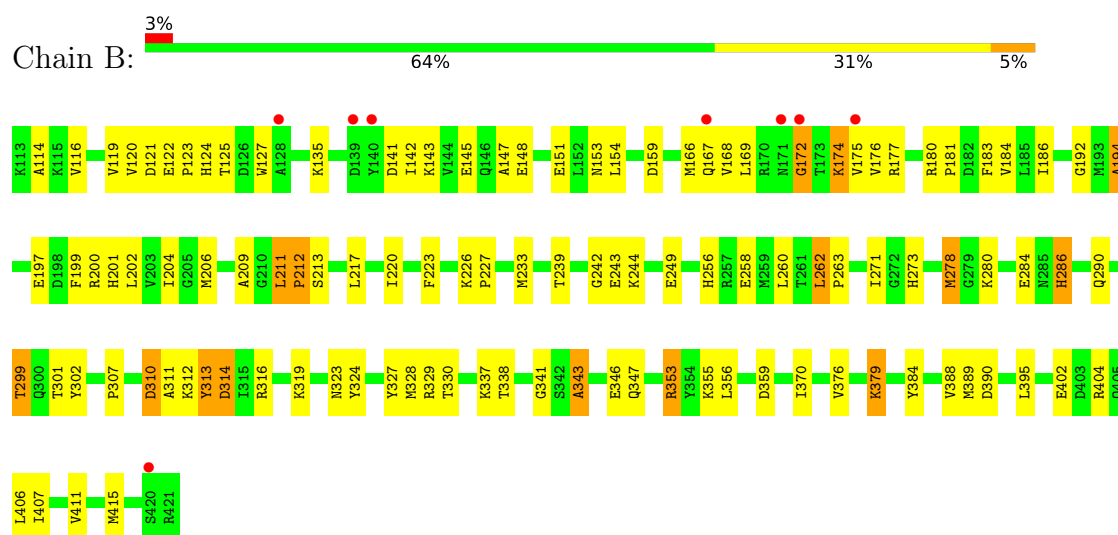
3 Residue-property plots

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: SYNAPSIN II



• Molecule 1: SYNAPSIN II



4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.28Å 121.28Å 165.89Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	12.00 – 2.35 12.00 – 2.35	Depositor EDS
% Data completeness (in resolution range)	92.9 (12.00-2.35) 90.2 (12.00-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.11Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.220 , 0.289 0.222 , 0.288	Depositor DCC
R_{free} test set	6028 reflections (7.94%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.232	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5181	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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

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	0.93	0/2517	1.26	24/3400 (0.7%)
1	B	0.92	0/2517	1.22	26/3400 (0.8%)
All	All	0.92	0/5034	1.24	50/6800 (0.7%)

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	311	ALA	N-CA-C	9.27	124.51	110.14
1	A	176	VAL	N-CA-C	9.02	120.03	106.42
1	B	311	ALA	N-CA-C	8.93	123.96	109.59
1	A	122	GLU	CA-C-N	8.84	128.48	119.82
1	A	122	GLU	C-N-CA	8.84	128.48	119.82
1	B	262	LEU	CA-C-N	8.73	129.17	120.52
1	B	262	LEU	C-N-CA	8.73	129.17	120.52
1	A	114	ALA	N-CA-C	7.37	121.15	109.50
1	B	284	GLU	N-CA-C	7.15	122.10	113.23
1	A	396	ILE	N-CA-C	7.09	120.41	108.86
1	A	287	TYR	N-CA-C	-6.99	104.39	113.12
1	A	240	LEU	N-CA-C	6.83	121.36	112.89
1	A	260	LEU	N-CA-C	-6.73	105.07	113.28
1	A	204	ILE	N-CA-C	-6.67	104.02	110.42
1	B	262	LEU	N-CA-C	6.60	117.96	109.65
1	B	260	LEU	N-CA-C	-6.52	104.81	112.89
1	B	142	ILE	N-CA-C	6.39	117.07	107.80
1	B	256	HIS	N-CA-C	6.27	118.92	111.71
1	B	194	ALA	N-CA-C	-6.23	101.76	110.35
1	B	212	PRO	N-CA-C	-6.09	102.32	111.57
1	B	343	ALA	N-CA-C	6.08	118.40	108.55
1	A	379	LYS	N-CA-C	-5.89	104.80	112.23

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	123	PRO	N-CA-C	5.87	121.55	113.98
1	B	213	SER	N-CA-C	5.84	117.01	108.14
1	A	312	LYS	N-CA-C	-5.78	104.41	112.45
1	A	176	VAL	CB-CA-C	-5.74	105.43	111.23
1	A	418	LEU	N-CA-C	5.73	120.55	113.50
1	B	370	ILE	CB-CA-C	-5.62	104.04	110.73
1	A	242	GLY	N-CA-C	5.61	120.88	113.37
1	B	313	TYR	N-CA-C	5.59	116.64	108.14
1	A	278	MET	N-CA-C	5.43	118.22	110.24
1	B	192	GLY	N-CA-C	-5.41	100.36	113.18
1	A	185	LEU	N-CA-C	-5.38	99.65	108.41
1	B	271	ILE	N-CA-C	5.35	115.95	108.89
1	B	278	MET	N-CA-C	5.35	117.78	110.24
1	B	211	LEU	N-CA-C	5.33	116.41	109.64
1	A	382	LYS	N-CA-C	5.22	118.52	110.42
1	A	271	ILE	N-CA-C	5.22	115.61	107.99
1	B	353	ARG	N-CA-C	-5.19	105.52	111.07
1	B	286	HIS	N-CA-C	5.17	117.32	111.11
1	B	239	THR	N-CA-C	5.16	118.03	111.69
1	A	194	ALA	N-CA-C	-5.10	103.19	110.59
1	B	204	ILE	N-CA-C	-5.09	105.54	110.42
1	A	139	ASP	N-CA-C	5.07	120.09	113.30
1	B	299	THR	N-CA-C	-5.06	107.05	113.18
1	A	412	ILE	N-CA-C	-5.05	105.78	110.53
1	B	379	LYS	N-CA-C	-5.05	106.38	112.54
1	A	370	ILE	CB-CA-C	-5.03	104.74	110.73
1	B	312	LYS	N-CA-C	-5.03	105.50	111.69
1	B	310	ASP	N-CA-C	-5.01	100.90	108.67

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	2462	0	2441	77	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2462	0	2441	80	0
2	A	31	0	12	1	0
2	B	31	0	12	1	0
3	A	94	0	0	5	0
3	B	101	0	0	5	0
All	All	5181	0	4906	155	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 16.

All (155) 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:177:ARG:HH11	1:A:177:ARG:CB	1.86	0.88
1:B:186:ILE:HD11	1:B:206:MET:HE1	1.60	0.84
1:A:335:ASN:HD21	1:A:337:LYS:HB2	1.43	0.83
1:A:226:LYS:H	1:A:273:HIS:HD2	1.24	0.83
1:A:233:MET:SD	1:A:388:VAL:HG23	2.22	0.79
1:A:331:SER:HB2	1:A:338:THR:HG23	1.64	0.79
1:A:177:ARG:HH11	1:A:177:ARG:HB2	1.47	0.78
1:B:159:ASP:HB3	3:B:1080:HOH:O	1.83	0.77
1:A:113:LYS:HG2	1:A:140:TYR:CE1	2.20	0.77
1:B:226:LYS:H	1:B:273:HIS:HD2	1.32	0.76
1:B:338:THR:HG22	1:B:338:THR:O	1.87	0.74
1:A:122:GLU:HB2	1:A:124:HIS:CE1	2.25	0.72
1:B:330:THR:O	1:B:343:ALA:HB1	1.91	0.71
1:A:270:LYS:HG2	1:A:280:LYS:HG2	1.73	0.71
1:B:169:LEU:HD13	1:B:174:LYS:HG3	1.73	0.70
1:A:166:MET:HE3	1:A:179:PHE:CE1	2.25	0.70
1:B:233:MET:SD	1:B:388:VAL:HG23	2.32	0.69
1:A:256:HIS:CD2	1:A:297:ALA:HB2	2.26	0.69
1:A:335:ASN:ND2	1:A:337:LYS:H	1.91	0.69
1:A:145:GLU:HB3	1:A:166:MET:HE1	1.75	0.68
1:B:151:GLU:HB3	1:B:167:GLN:O	1.94	0.67
1:B:280:LYS:HE2	2:B:801:ATP:O3A	1.95	0.67
1:A:177:ARG:HB2	1:A:177:ARG:NH1	2.09	0.67
1:A:115:LYS:HB2	1:A:136:ILE:HD11	1.76	0.66
1:B:154:LEU:C	1:B:154:LEU:HD12	2.20	0.66
1:B:226:LYS:HG3	1:B:273:HIS:CD2	2.32	0.65
1:B:166:MET:HG2	1:B:168:VAL:HG23	1.79	0.65
1:A:309:ILE:HG21	1:A:384:TYR:CD2	2.32	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:LYS:N	1:A:273:HIS:HD2	1.96	0.63
1:A:256:HIS:CE1	1:A:257:ARG:HG3	2.32	0.63
1:A:226:LYS:HG2	1:A:273:HIS:CD2	2.33	0.62
1:A:122:GLU:OE1	1:A:124:HIS:HE1	1.82	0.62
1:A:154:LEU:O	1:B:258:GLU:HG3	1.99	0.62
1:A:379:LYS:C	1:A:381:GLY:H	2.08	0.62
1:B:135:LYS:HD3	1:B:141:ASP:OD1	1.99	0.62
1:A:153:ASN:HD22	1:A:154:LEU:H	1.47	0.61
1:B:125:THR:CG2	1:B:404:ARG:HD3	2.31	0.61
1:B:125:THR:HG21	1:B:404:ARG:HD3	1.83	0.61
1:B:153:ASN:HD22	1:B:154:LEU:H	1.48	0.60
1:B:226:LYS:N	1:B:273:HIS:HD2	2.00	0.60
1:B:122:GLU:HB3	1:B:123:PRO:HD2	1.84	0.60
1:A:239:THR:HG22	1:A:240:LEU:HD23	1.84	0.58
1:A:259:MET:HE1	1:A:296:VAL:HG21	1.85	0.58
1:B:244:LYS:HE2	1:B:356:LEU:HD21	1.85	0.58
1:A:154:LEU:HD12	1:A:154:LEU:C	2.29	0.58
1:B:168:VAL:HG12	1:B:169:LEU:H	1.68	0.58
1:A:177:ARG:HH11	1:A:177:ARG:HB3	1.67	0.56
1:B:168:VAL:HG12	1:B:169:LEU:N	2.20	0.56
1:B:249:GLU:O	1:B:307:PRO:HD2	2.05	0.56
1:A:233:MET:CE	1:A:388:VAL:HG23	2.34	0.56
1:B:244:LYS:HE2	1:B:356:LEU:CD2	2.36	0.55
1:B:411:VAL:O	1:B:415:MET:HG3	2.06	0.55
1:A:166:MET:HE3	1:A:179:PHE:CZ	2.42	0.55
1:A:177:ARG:CB	1:A:177:ARG:NH1	2.64	0.55
1:A:125:THR:HG23	1:A:404:ARG:HD3	1.88	0.54
1:A:243:GLU:HB3	3:A:1074:HOH:O	2.07	0.54
1:B:299:THR:OG1	1:B:301:THR:HG22	2.07	0.54
1:A:353:ARG:HD3	3:A:1097:HOH:O	2.07	0.54
1:B:120:VAL:HA	1:B:147:ALA:O	2.08	0.54
1:A:316:ARG:HG3	1:A:374:LYS:HG2	1.90	0.54
1:B:151:GLU:OE2	1:B:168:VAL:HG13	2.08	0.54
1:A:119:VAL:HG11	1:A:127:TRP:CD1	2.43	0.53
1:A:301:THR:HG23	1:A:302:TYR:O	2.09	0.53
1:B:206:MET:HE3	1:B:217:LEU:CD2	2.39	0.52
1:A:339:ASN:HB2	2:A:800:ATP:O2G	2.09	0.52
1:A:398:GLU:CD	1:A:398:GLU:H	2.17	0.52
1:B:122:GLU:HA	1:B:148:GLU:OE2	2.10	0.52
1:B:338:THR:O	1:B:338:THR:CG2	2.57	0.52
1:A:226:LYS:HB2	1:A:227:PRO:HD3	1.90	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:ARG:HD2	1:A:389:MET:CE	2.39	0.52
1:B:310:ASP:OD1	1:B:379:LYS:HD2	2.08	0.52
1:A:115:LYS:CG	1:A:136:ILE:HD11	2.40	0.52
1:B:199:PHE:HA	1:B:201:HIS:CE1	2.45	0.52
1:B:286:HIS:CE1	1:B:290:GLN:NE2	2.78	0.51
1:B:143:LYS:NZ	1:B:145:GLU:OE2	2.42	0.51
1:B:194:ALA:HB2	1:B:278:MET:SD	2.51	0.50
1:A:115:LYS:CB	1:A:136:ILE:HD11	2.41	0.50
1:B:233:MET:CE	1:B:388:VAL:HG23	2.42	0.49
1:A:325:LYS:HE2	1:A:403:ASP:OD1	2.11	0.49
1:B:319:LYS:HD2	1:B:324:TYR:HE1	1.76	0.49
1:A:313:TYR:HB2	1:A:329:ARG:O	2.12	0.49
1:B:206:MET:HE3	1:B:217:LEU:HD21	1.95	0.49
1:B:184:VAL:CG2	1:B:211:LEU:HD23	2.43	0.48
1:B:121:ASP:OD1	1:B:122:GLU:N	2.46	0.48
1:A:249:GLU:O	1:A:307:PRO:HD2	2.14	0.48
1:A:113:LYS:HG2	1:A:140:TYR:CD1	2.48	0.48
1:B:327:TYR:CE1	1:B:347:GLN:HB2	2.48	0.48
1:A:312:LYS:O	1:A:330:THR:HG23	2.14	0.47
1:B:119:VAL:HG11	1:B:127:TRP:CD1	2.49	0.47
1:B:314:ASP:HB2	1:B:329:ARG:HB3	1.97	0.47
1:A:176:VAL:O	1:A:176:VAL:HG12	2.14	0.47
1:B:184:VAL:HG23	1:B:211:LEU:HD23	1.96	0.47
1:B:262:LEU:HD11	1:B:286:HIS:CD2	2.50	0.47
1:A:243:GLU:N	3:A:1074:HOH:O	2.48	0.47
1:B:186:ILE:HD12	1:B:220:ILE:HD13	1.97	0.46
1:A:259:MET:CE	1:A:293:ALA:HA	2.45	0.46
1:B:355:LYS:HE2	1:B:359:ASP:OD2	2.16	0.46
1:A:258:GLU:HG3	1:B:154:LEU:O	2.14	0.46
1:A:329:ARG:HA	1:A:344:MET:O	2.15	0.46
1:A:402:GLU:O	1:A:405:GLN:HB3	2.14	0.46
1:B:176:VAL:O	1:B:176:VAL:HG12	2.15	0.46
1:A:301:THR:HG23	1:A:302:TYR:N	2.30	0.46
1:A:316:ARG:HD2	1:A:389:MET:HE1	1.97	0.46
1:B:200:ARG:HD3	3:B:1118:HOH:O	2.16	0.46
1:B:116:VAL:CG1	1:B:181:PRO:HA	2.46	0.46
1:A:331:SER:CB	1:A:338:THR:HG23	2.39	0.46
1:B:194:ALA:HB3	1:B:197:GLU:CD	2.42	0.45
1:B:346:GLU:HG2	3:B:1181:HOH:O	2.16	0.45
1:A:261:THR:HG22	1:A:261:THR:O	2.17	0.44
1:B:227:PRO:HG2	1:B:302:TYR:CE1	2.52	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:ILE:HG21	1:A:384:TYR:CE2	2.53	0.44
1:A:125:THR:CG2	1:A:404:ARG:HD3	2.48	0.44
1:B:114:ALA:HA	1:B:141:ASP:O	2.16	0.44
1:B:169:LEU:HD13	1:B:174:LYS:CG	2.46	0.44
1:B:175:VAL:HG11	1:B:177:ARG:HD2	1.99	0.44
1:A:145:GLU:HB3	1:A:166:MET:CE	2.46	0.44
1:B:202:LEU:O	1:B:206:MET:HG3	2.18	0.43
1:A:291:ASP:O	1:A:295:VAL:HG23	2.18	0.43
1:A:148:GLU:CG	1:A:188:GLN:HE22	2.32	0.43
1:A:167:GLN:HG2	1:A:176:VAL:HG22	2.01	0.43
1:B:242:GLY:O	1:B:243:GLU:C	2.61	0.43
1:B:323:ASN:CB	1:B:406:LEU:HD13	2.49	0.43
1:B:407:ILE:O	1:B:411:VAL:HG23	2.19	0.43
1:B:116:VAL:HG12	1:B:181:PRO:HA	2.01	0.42
1:A:223:PHE:CE2	1:A:390:ASP:HB3	2.54	0.42
1:A:233:MET:HE3	1:A:388:VAL:HG23	2.00	0.42
1:B:153:ASN:HD22	1:B:154:LEU:N	2.14	0.42
1:B:395:LEU:HD23	1:B:395:LEU:HA	1.79	0.42
1:B:223:PHE:CE2	1:B:390:ASP:HB3	2.54	0.42
1:B:313:TYR:CG	1:B:328:MET:HE2	2.54	0.42
1:B:353:ARG:HH11	1:B:353:ARG:HG3	1.84	0.42
1:A:187:ARG:NH2	1:A:395:LEU:HG	2.34	0.42
1:A:186:ILE:HD12	1:A:220:ILE:HD13	2.02	0.42
1:A:292:ILE:O	1:A:296:VAL:HG23	2.20	0.42
1:B:356:LEU:HD23	3:B:1137:HOH:O	2.20	0.42
1:B:376:VAL:HG23	1:B:384:TYR:HB2	2.00	0.42
1:B:169:LEU:HD11	1:B:172:GLY:C	2.44	0.41
1:B:183:PHE:CD1	1:B:184:VAL:N	2.88	0.41
1:B:316:ARG:HD2	1:B:389:MET:CE	2.50	0.41
1:A:153:ASN:HD22	1:A:154:LEU:N	2.17	0.41
1:B:211:LEU:HA	1:B:212:PRO:HD3	1.87	0.41
1:A:157:HIS:HD2	1:A:161:THR:O	2.03	0.41
1:A:238:LYS:HE3	1:A:238:LYS:HB2	1.70	0.41
1:A:403:ASP:O	1:A:404:ARG:C	2.63	0.41
1:B:226:LYS:CG	1:B:273:HIS:CD2	3.03	0.41
1:A:353:ARG:NE	3:A:1097:HOH:O	2.52	0.41
1:B:249:GLU:HG3	3:B:1140:HOH:O	2.21	0.41
1:A:295:VAL:C	1:A:297:ALA:N	2.78	0.41
1:A:353:ARG:CD	3:A:1097:HOH:O	2.67	0.40
1:B:186:ILE:HD12	1:B:220:ILE:CD1	2.51	0.40
1:B:209:ALA:HB3	1:B:211:LEU:CD1	2.51	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:LYS:HB2	1:B:337:LYS:HE2	2.02	0.40
1:A:401:VAL:HG23	1:A:404:ARG:NH2	2.37	0.40
1:B:233:MET:HE3	1:B:388:VAL:HG23	2.02	0.40
1:A:262:LEU:CD1	1:A:286:HIS:HB2	2.52	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	307/309 (99%)	286 (93%)	19 (6%)	2 (1%)	18	20
1	B	307/309 (99%)	289 (94%)	15 (5%)	3 (1%)	12	12
All	All	614/618 (99%)	575 (94%)	34 (6%)	5 (1%)	16	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	172	GLY
1	B	341	GLY
1	A	259	MET
1	A	380	ASP
1	B	124	HIS

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	266/266 (100%)	252 (95%)	14 (5%)	20	25
1	B	266/266 (100%)	261 (98%)	5 (2%)	50	65
All	All	532/532 (100%)	513 (96%)	19 (4%)	31	41

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	LEU
1	A	177	ARG
1	A	180	ARG
1	A	239	THR
1	A	249	GLU
1	A	259	MET
1	A	261	THR
1	A	263	PRO
1	A	286	HIS
1	A	328	MET
1	A	348	ILE
1	A	398	GLU
1	A	401	VAL
1	A	402	GLU
1	B	174	LYS
1	B	180	ARG
1	B	263	PRO
1	B	314	ASP
1	B	402	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	HIS
1	A	153	ASN
1	A	188	GLN
1	A	232	GLN
1	A	250	GLN
1	A	256	HIS
1	A	273	HIS
1	A	275	HIS
1	A	286	HIS
1	A	290	GLN
1	A	323	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	335	ASN
1	B	153	ASN
1	B	201	HIS
1	B	232	GLN
1	B	250	GLN
1	B	256	HIS
1	B	273	HIS
1	B	286	HIS
1	B	290	GLN
1	B	322	ASN
1	B	323	ASN
1	B	399	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](#)

2 ligands are modelled in this entry.

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	$\# Z > 2$
2	ATP	B	801	-	32,33,33	1.21	3 (9%)	48,52,52	0.99	1 (2%)
2	ATP	A	800	-	32,33,33	1.18	2 (6%)	48,52,52	0.94	2 (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
2	ATP	B	801	-	-	2/22/38/38	0/3/3/3
2	ATP	A	800	-	-	3/22/38/38	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	ATP	PB-O3A	-4.11	1.55	1.59
2	A	800	ATP	PB-O3B	3.79	1.63	1.59
2	B	801	ATP	PB-O3B	2.72	1.62	1.59
2	A	800	ATP	C4-N3	2.40	1.38	1.34
2	B	801	ATP	C3'-C4'	-2.10	1.47	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	ATP	O2B-PB-O3A	2.82	114.91	107.27
2	A	800	ATP	O2B-PB-O3A	2.17	113.13	107.27
2	A	800	ATP	O3G-PG-O2G	2.03	115.42	107.80

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	ATP	C5'-O5'-PA-O1A
2	A	800	ATP	PB-O3A-PA-O2A
2	A	800	ATP	PB-O3A-PA-O1A
2	B	801	ATP	PB-O3A-PA-O1A
2	B	801	ATP	PB-O3A-PA-O2A

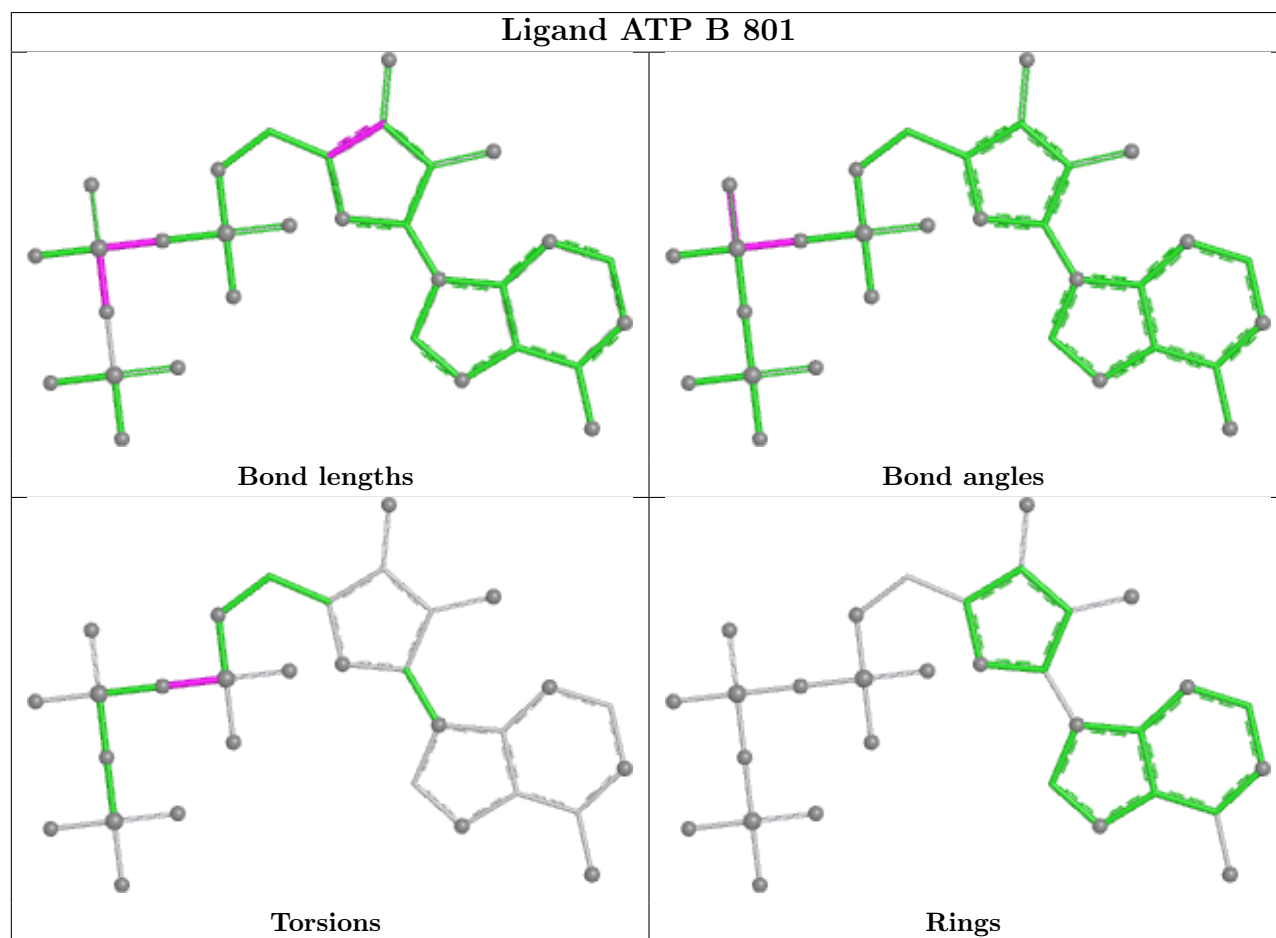
There are no ring outliers.

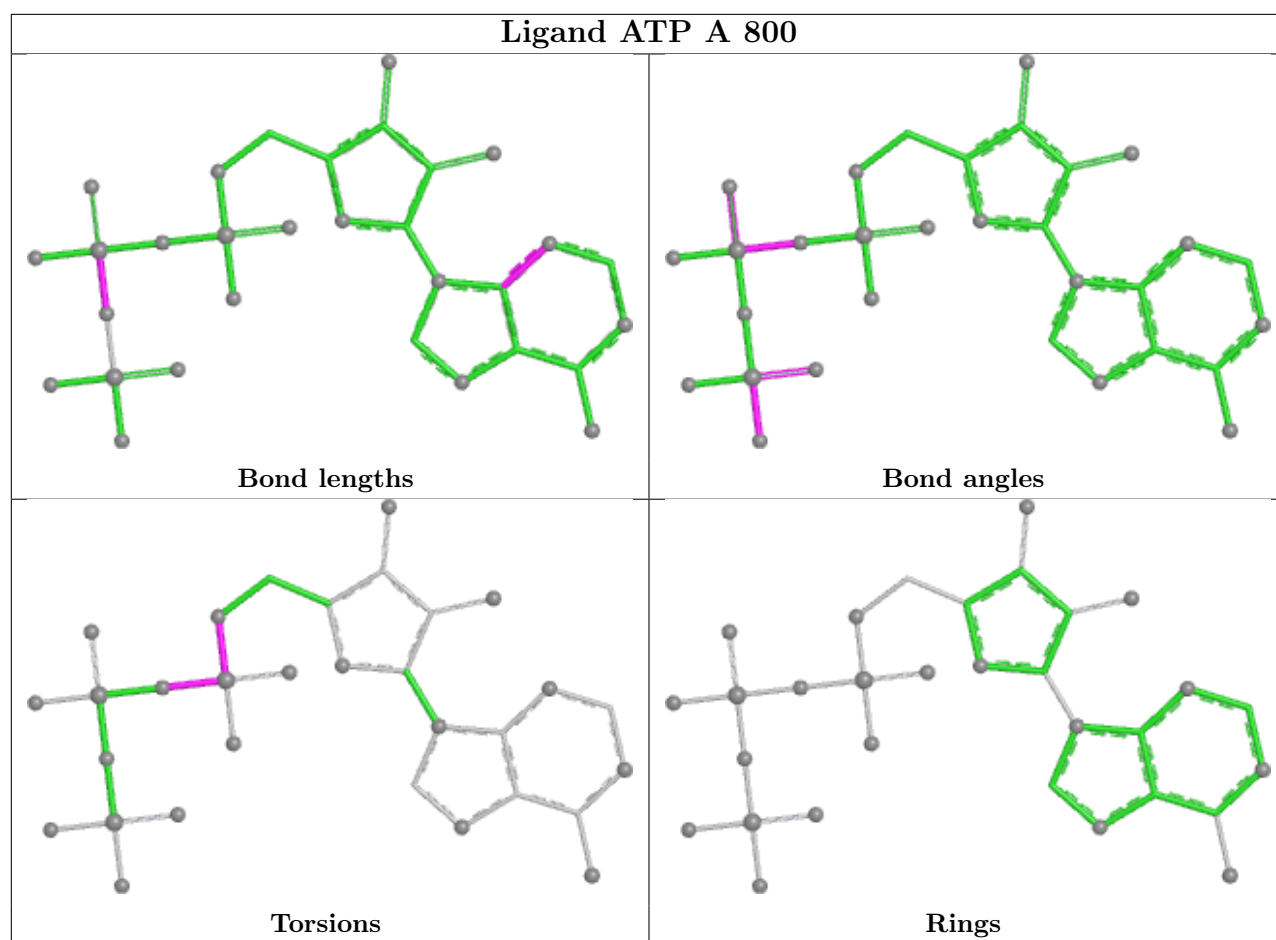
2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	801	ATP	1	0
2	A	800	ATP	1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	303/309 (98%)	-0.05	5 (1%) 69 74	15, 37, 64, 92	5 (1%)
1	B	305/309 (98%)	-0.08	8 (2%) 57 63	15, 37, 80, 98	0
All	All	608/618 (98%)	-0.06	13 (2%) 63 68	15, 37, 76, 98	5 (0%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	344	MET	3.4
1	B	128	ALA	2.9
1	A	342	SER	2.9
1	B	420	SER	2.7
1	B	171	ASN	2.3
1	A	378	GLY	2.3
1	B	172	GLY	2.3
1	A	169	LEU	2.2
1	B	175	VAL	2.3
1	B	139	ASP	2.2
1	B	167	GLN	2.1
1	B	140	TYR	2.1
1	A	345	LEU	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](#)

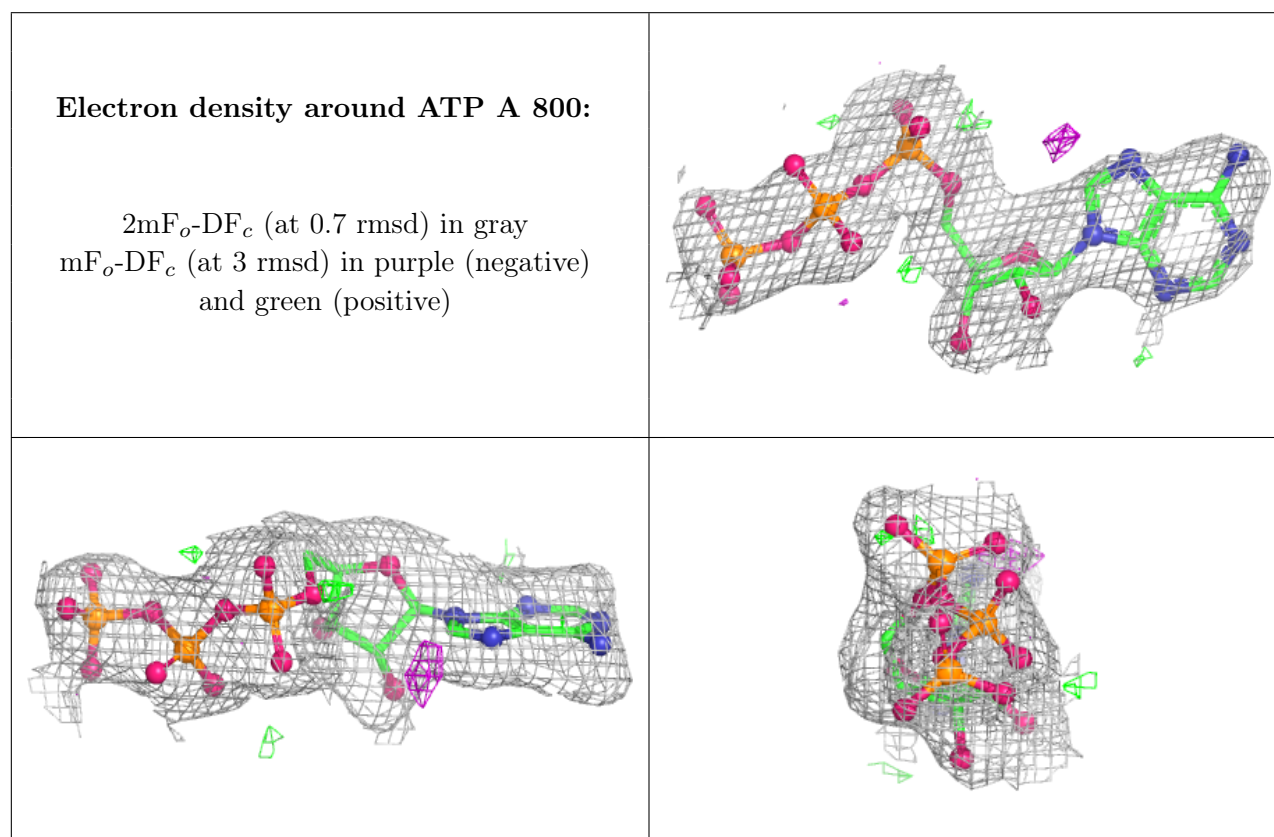
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

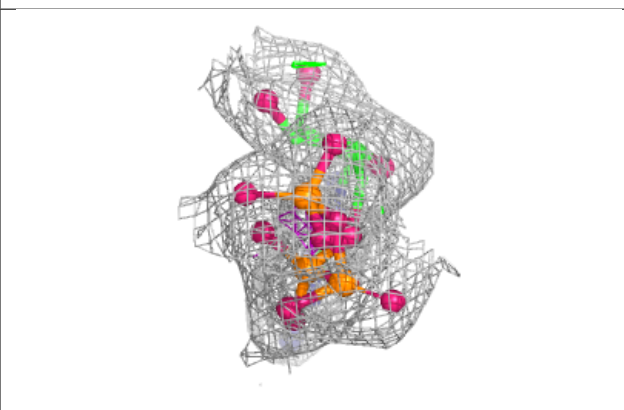
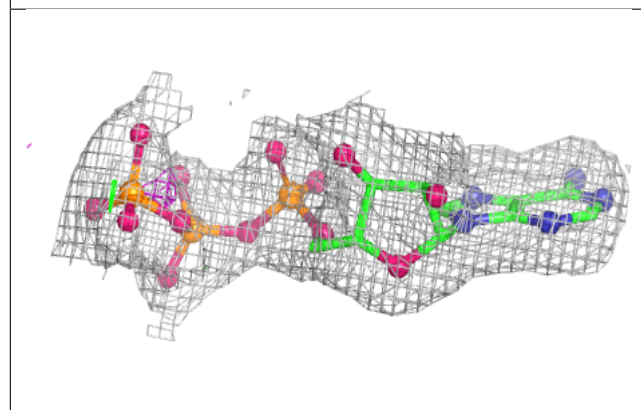
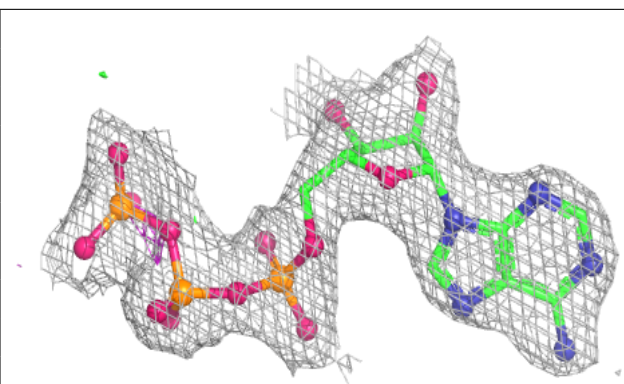
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ATP	A	800	31/31	0.90	0.10	50,57,69,70	0
2	ATP	B	801	31/31	0.94	0.07	25,32,45,50	0

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 801:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.