



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:54 AM UTC

PDB ID : 1ATP / pdb_00001atp
Title : 2.2 angstrom refined crystal structure of the catalytic subunit of cAMP-dependent protein kinase complexed with MNATP and a peptide inhibitor
Authors : Zheng, J.; Trafny, E.A.; Knighton, D.R.; Xuong, N.-H.; Taylor, S.S.; Teneyck, L.F.; Sowadski, J.M.
Deposited on : 1993-01-08
Resolution : 2.20 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

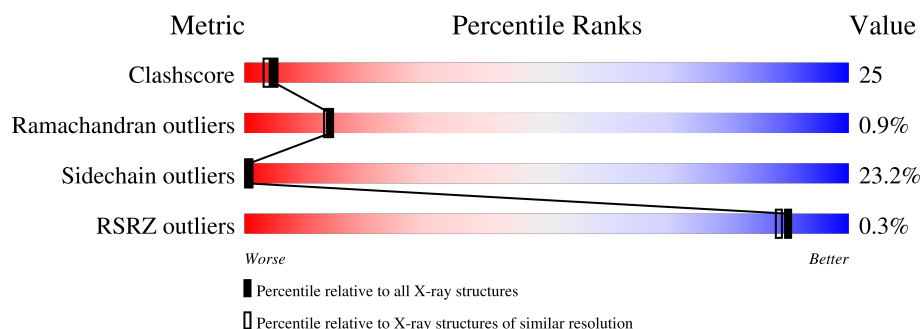
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	E	350	
2	I	20	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3070 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 cAMP-DEPENDENT PROTEIN KINASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	E	336	Total	C	N	O	P	S	0	0	0
			2777	1800	465	502	2	8			

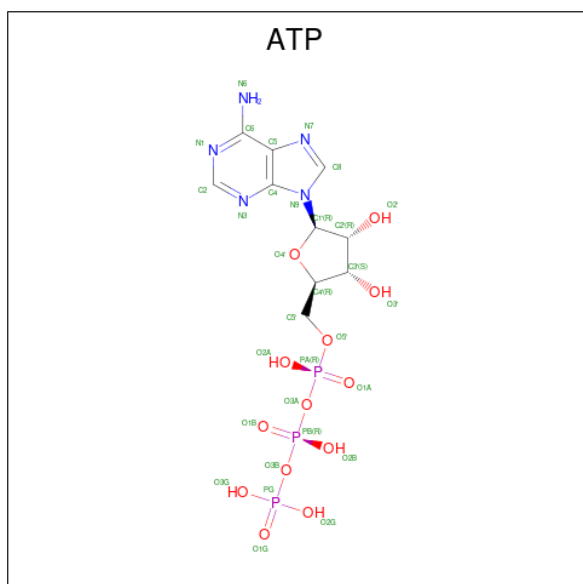
- Molecule 2 is a protein called PEPTIDE INHIBITOR PKI(5-24).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	20	Total	C	N	O	0	0	0
			157	94	32	31			

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total	Mn	0	0
			2	2		

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



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

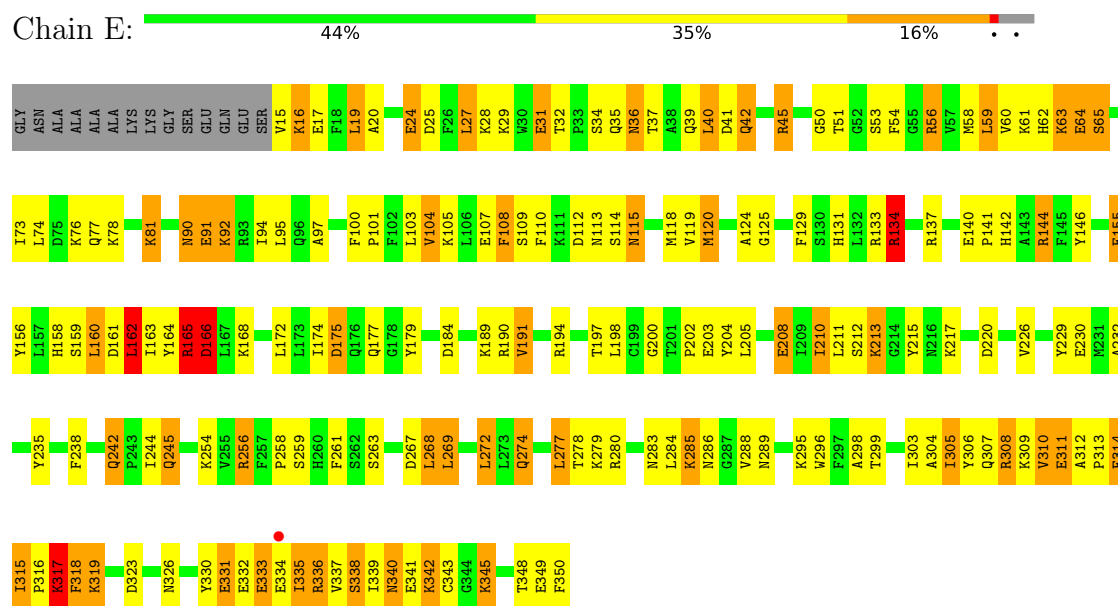
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	89	Total	O	0	0
			89	89		
5	I	14	Total	O	0	0
			14	14		

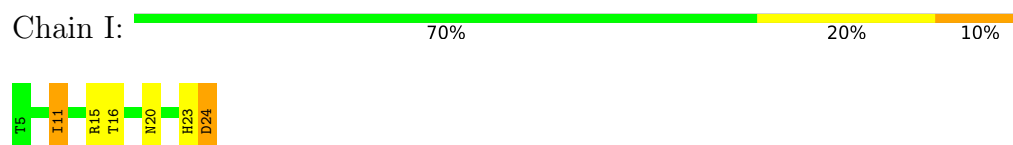
3 Residue-property plots

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: cAMP-DEPENDENT PROTEIN KINASE



• Molecule 2: PEPTIDE INHIBITOR PKI(5-24)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.58Å 76.28Å 80.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.20 10.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.20) 76.5 (10.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.21 (at 2.15Å)	Xtriage
Refinement program	TNT, X-PLOR	Depositor
R, R_{free}	0.177 , (Not available) 0.192 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 126.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3070	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% 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, MN, SEP, TPO

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	E	1.16	7/2825 (0.2%)	1.63	36/3804 (0.9%)
2	I	1.11	0/159	1.50	1/212 (0.5%)
All	All	1.16	7/2984 (0.2%)	1.63	37/4016 (0.9%)

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	E	1	2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	165	ARG	C-N	-13.94	1.14	1.33
1	E	166	ASP	C-N	6.70	1.43	1.33
1	E	25	ASP	CG-OD2	6.32	1.37	1.25
1	E	56	ARG	NE-CZ	-6.05	1.26	1.33
1	E	162	LEU	C-N	5.93	1.41	1.33
1	E	56	ARG	CZ-NH2	-5.03	1.26	1.33
1	E	323	ASP	C-O	-5.01	1.17	1.24

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	165	ARG	O-C-N	-13.34	96.90	122.11
1	E	104	VAL	CB-CA-C	-8.79	101.61	111.45
1	E	124	ALA	N-CA-C	7.84	120.73	111.71
1	E	317	LYS	N-CA-CB	7.20	121.88	110.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	90	ASN	CB-CA-C	-7.13	99.69	110.88
1	E	134	ARG	N-CA-CB	7.00	120.14	109.91
1	E	165	ARG	CA-C-N	6.91	134.73	121.54
1	E	165	ARG	C-N-CA	6.91	134.73	121.54
1	E	280	ARG	N-CA-C	6.88	119.77	109.25
1	E	134	ARG	N-CA-C	6.87	118.46	110.97
1	E	54	PHE	CA-CB-CG	-6.72	107.08	113.80
1	E	63	LYS	N-CA-C	6.60	118.47	111.28
1	E	333	GLU	N-CA-C	-6.59	100.73	110.48
1	E	134	ARG	CB-CA-C	6.56	121.06	110.96
1	E	274	GLN	CB-CA-C	6.51	121.08	110.22
1	E	210	ILE	CA-CB-CG2	6.25	121.12	110.50
1	E	115	ASN	N-CA-C	6.16	118.32	109.14
1	E	213	LYS	N-CA-C	6.02	121.36	113.30
1	E	210	ILE	N-CA-CB	5.92	122.51	110.78
1	E	317	LYS	CB-CA-C	5.92	119.96	109.72
1	E	280	ARG	CB-CA-C	-5.89	100.63	110.29
1	E	166	ASP	O-C-N	-5.86	114.80	122.59
1	E	298	ALA	N-CA-C	5.83	120.00	113.01
1	E	191	VAL	CB-CA-C	-5.80	100.44	110.65
2	I	24	ASP	CA-CB-CG	-5.62	106.98	112.60
1	E	56	ARG	NE-CZ-NH1	-5.45	116.05	121.50
1	E	175	ASP	N-CA-CB	5.45	117.90	110.38
1	E	162	LEU	CA-C-N	5.42	130.54	123.06
1	E	162	LEU	C-N-CA	5.42	130.54	123.06
1	E	323	ASP	CA-CB-CG	5.40	118.00	112.60
1	E	59	LEU	N-CA-C	-5.38	101.20	109.76
1	E	165	ARG	NE-CZ-NH2	5.32	123.99	119.20
1	E	310	VAL	N-CA-C	5.29	116.80	109.45
1	E	215	TYR	N-CA-C	5.20	115.92	108.74
1	E	24	GLU	N-CA-C	5.15	116.58	111.07
1	E	155	GLU	CA-C-O	5.13	125.90	120.10
1	E	16	LYS	CA-C-O	5.10	126.14	120.63

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	E	134	ARG	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	165	ARG	Mainchain
1	E	166	ASP	Mainchain

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	E	2777	0	2762	142	0
2	I	157	0	146	8	0
3	E	2	0	0	0	0
4	E	31	0	12	0	0
5	E	89	0	0	9	0
5	I	14	0	0	0	0
All	All	3070	0	2920	148	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 25.

All (148) 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:E:305:ILE:HG13	1:E:310:VAL:HG21	1.23	1.11
1:E:310:VAL:HG12	1:E:311:GLU:HG3	1.50	0.90
1:E:285:LYS:H	1:E:285:LYS:HD2	1.40	0.86
1:E:284:LEU:HB3	1:E:285:LYS:HD2	1.60	0.82
1:E:345:LYS:H	1:E:345:LYS:HD2	1.47	0.79
1:E:303:ILE:O	1:E:307:GLN:HG3	1.85	0.76
1:E:303:ILE:HD12	1:E:303:ILE:H	1.50	0.75
1:E:305:ILE:HG13	1:E:310:VAL:CG2	2.10	0.74
1:E:319:LYS:HD2	1:E:319:LYS:H	1.52	0.73
1:E:163:ILE:HG12	1:E:217:LYS:HD3	1.71	0.72
1:E:335:ILE:HD12	5:E:596:HOH:O	1.89	0.72
1:E:285:LYS:HD2	1:E:285:LYS:N	2.04	0.71
1:E:64:GLU:H	1:E:64:GLU:CD	1.98	0.71
1:E:62:HIS:HD2	1:E:65:SER:HB3	1.54	0.71
1:E:285:LYS:HG2	1:E:286:ASN:N	2.07	0.70
1:E:203:GLU:OE2	2:I:15:ARG:HD3	1.91	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:334:GLU:HB2	1:E:336:ARG:HH12	1.58	0.68
1:E:133:ARG:NH2	1:E:230:GLU:OE2	2.26	0.68
1:E:277:LEU:O	1:E:283:ASN:ND2	2.27	0.68
1:E:36:ASN:ND2	1:E:110:PHE:HB2	2.10	0.67
1:E:56:ARG:HH12	1:E:333:GLU:CD	2.03	0.66
1:E:164:TYR:CE2	1:E:166:ASP:C	2.73	0.66
1:E:200:GLY:HA3	1:E:205:LEU:HD21	1.76	0.66
1:E:303:ILE:HD12	1:E:303:ILE:N	2.10	0.65
1:E:242:GLN:HG3	1:E:245:GLN:OE1	1.96	0.65
1:E:285:LYS:HG2	1:E:286:ASN:H	1.60	0.65
1:E:15:VAL:N	1:E:17:GLU:OE1	2.30	0.65
1:E:274:GLN:HG2	1:E:279:LYS:HB3	1.78	0.64
1:E:53:SER:OG	2:I:20:ASN:HB3	1.98	0.64
1:E:58:MET:CE	1:E:335:ILE:HD11	2.27	0.64
1:E:335:ILE:HD13	1:E:335:ILE:N	2.12	0.64
1:E:45:ARG:NH2	1:E:333:GLU:O	2.31	0.63
1:E:90:ASN:O	1:E:94:ILE:HG13	1.99	0.63
1:E:77:GLN:NE2	1:E:342:LYS:HG3	2.13	0.63
1:E:39:GLN:O	1:E:42:GLN:HG3	2.00	0.61
1:E:338:SEP:HB2	5:E:586:HOH:O	2.00	0.61
1:E:338:SEP:O2P	1:E:340:ASN:ND2	2.33	0.61
1:E:100:PHE:CG	1:E:101:PRO:HD2	2.35	0.60
1:E:189:LYS:HD3	1:E:191:VAL:CG2	2.31	0.60
1:E:62:HIS:CD2	1:E:65:SER:HB3	2.37	0.59
1:E:125:GLY:HA3	1:E:174:ILE:O	2.02	0.59
1:E:163:ILE:HG22	1:E:165:ARG:HG3	1.83	0.59
1:E:204:TYR:CE2	1:E:226:VAL:HG12	2.37	0.59
1:E:56:ARG:NH2	1:E:333:GLU:OE1	2.25	0.59
1:E:285:LYS:H	1:E:285:LYS:CD	2.15	0.59
1:E:155:GLU:HG3	1:E:288:VAL:HG11	1.84	0.59
1:E:304:ALA:HA	1:E:309:LYS:HD3	1.83	0.59
1:E:158:HIS:HE1	1:E:220:ASP:OD2	1.86	0.58
1:E:304:ALA:HB1	1:E:309:LYS:HG2	1.86	0.58
1:E:58:MET:HE3	1:E:73:ILE:CD1	2.35	0.57
1:E:286:ASN:O	1:E:289:ASN:HB2	2.04	0.57
1:E:189:LYS:HD3	1:E:191:VAL:HG21	1.86	0.57
1:E:144:ARG:HB2	1:E:296:TRP:CH2	2.40	0.56
1:E:160:LEU:O	1:E:162:LEU:HD13	2.04	0.56
1:E:331:GLU:O	1:E:331:GLU:HG3	2.04	0.56
1:E:91:GLU:CD	1:E:118:MET:HE1	2.30	0.56
1:E:175:ASP:OD2	1:E:308:ARG:NH1	2.29	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:310:VAL:HG12	1:E:311:GLU:CG	2.33	0.55
1:E:29:LYS:HD3	1:E:97:ALA:HA	1.88	0.54
1:E:35:GLN:HE21	1:E:350:PHE:C	2.14	0.54
1:E:303:ILE:H	1:E:303:ILE:CD1	2.20	0.54
1:E:310:VAL:HG12	1:E:311:GLU:N	2.23	0.54
1:E:334:GLU:C	1:E:335:ILE:HD13	2.33	0.53
1:E:162:LEU:HD12	1:E:190:ARG:HA	1.91	0.53
1:E:284:LEU:CB	1:E:285:LYS:HD2	2.36	0.53
1:E:56:ARG:NH1	1:E:333:GLU:HB2	2.23	0.52
2:I:23:HIS:CD2	2:I:24:ASP:N	2.77	0.52
1:E:131:HIS:HA	1:E:134:ARG:NH1	2.23	0.52
1:E:345:LYS:HD2	1:E:345:LYS:N	2.20	0.52
1:E:168:LYS:O	1:E:172:LEU:HG	2.11	0.51
1:E:58:MET:HE1	1:E:335:ILE:HD11	1.93	0.51
1:E:76:LYS:HG3	1:E:114:SER:O	2.10	0.51
1:E:36:ASN:HD21	1:E:110:PHE:HB2	1.75	0.51
1:E:345:LYS:H	1:E:345:LYS:CD	2.06	0.51
1:E:56:ARG:HH12	1:E:333:GLU:HB2	1.76	0.51
1:E:163:ILE:HG22	1:E:165:ARG:CG	2.41	0.51
1:E:140:GLU:HB2	1:E:141:PRO:HD3	1.93	0.50
1:E:41:ASP:OD1	1:E:42:GLN:HG2	2.11	0.50
1:E:129:PHE:CE1	1:E:133:ARG:NH1	2.80	0.50
1:E:133:ARG:NH1	5:E:480:HOH:O	2.43	0.49
1:E:108:PHE:HD1	1:E:119:VAL:HB	1.78	0.49
1:E:334:GLU:HB2	1:E:336:ARG:NH1	2.24	0.49
1:E:34:SER:O	1:E:92:LYS:NZ	2.43	0.49
1:E:35:GLN:HG2	1:E:350:PHE:OXT	2.13	0.48
2:I:23:HIS:CD2	2:I:24:ASP:H	2.31	0.48
1:E:100:PHE:HB2	1:E:156:TYR:CE2	2.48	0.48
1:E:258:PRO:HG2	1:E:261:PHE:CE1	2.49	0.48
2:I:11:ILE:CD1	2:I:16:THR:HG21	2.44	0.48
1:E:256:ARG:HB3	1:E:256:ARG:CZ	2.40	0.47
1:E:76:LYS:HE2	1:E:112:ASP:O	2.13	0.47
1:E:155:GLU:O	1:E:159:SER:HB3	2.14	0.47
1:E:17:GLU:O	1:E:20:ALA:HB3	2.14	0.47
2:I:23:HIS:O	2:I:24:ASP:O	2.33	0.47
1:E:77:GLN:CD	1:E:342:LYS:HG3	2.40	0.46
1:E:29:LYS:HB2	1:E:97:ALA:CB	2.46	0.46
1:E:332:GLU:O	1:E:334:GLU:OE2	2.33	0.46
1:E:316:PRO:HA	5:E:535:HOH:O	2.16	0.46
1:E:268:LEU:HD22	1:E:272:LEU:HD22	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:312:ALA:HA	1:E:313:PRO:HD2	1.84	0.45
1:E:314:PHE:C	1:E:315:ILE:HD13	2.42	0.45
1:E:317:LYS:HA	1:E:317:LYS:HD2	1.44	0.45
2:I:23:HIS:CG	2:I:24:ASP:N	2.82	0.45
1:E:319:LYS:H	1:E:319:LYS:CD	2.26	0.45
1:E:45:ARG:O	1:E:45:ARG:HG2	1.98	0.45
1:E:208:GLU:HB3	1:E:213:LYS:HG2	1.99	0.45
1:E:285:LYS:CG	1:E:286:ASN:N	2.79	0.45
1:E:338:SEP:OG	1:E:339:ILE:N	2.51	0.44
1:E:50:GLY:HA2	1:E:330:TYR:CE1	2.52	0.44
1:E:164:TYR:HE2	1:E:166:ASP:C	2.25	0.44
1:E:100:PHE:CD2	1:E:101:PRO:HD2	2.52	0.44
1:E:244:ILE:HG13	1:E:245:GLN:N	2.33	0.44
1:E:345:LYS:N	1:E:345:LYS:CD	2.78	0.44
1:E:211:LEU:HD23	1:E:211:LEU:HA	1.75	0.43
1:E:142:HIS:CD2	1:E:146:TYR:CE2	3.06	0.43
1:E:177:GLN:HG2	1:E:179:TYR:CD1	2.53	0.43
1:E:230:GLU:HA	1:E:235:TYR:O	2.18	0.43
1:E:204:TYR:CE2	1:E:226:VAL:CG1	3.01	0.43
1:E:142:HIS:NE2	1:E:146:TYR:CE2	2.87	0.43
1:E:189:LYS:NZ	5:E:428:HOH:O	2.49	0.43
1:E:58:MET:HE1	1:E:335:ILE:CD1	2.49	0.43
1:E:19:LEU:HD12	1:E:19:LEU:HA	1.89	0.42
1:E:310:VAL:HG12	1:E:311:GLU:H	1.84	0.42
1:E:334:GLU:CB	1:E:336:ARG:HH12	2.30	0.42
1:E:105:LYS:O	1:E:120:MET:HB3	2.20	0.42
1:E:314:PHE:C	1:E:314:PHE:CD1	2.97	0.42
1:E:35:GLN:NE2	1:E:350:PHE:O	2.46	0.42
1:E:40:LEU:HD23	1:E:40:LEU:HA	1.80	0.42
1:E:36:ASN:HA	1:E:109:SER:O	2.20	0.41
1:E:107:GLU:N	1:E:119:VAL:O	2.51	0.41
1:E:76:LYS:HG3	1:E:115:ASN:HA	2.02	0.41
1:E:310:VAL:CG1	1:E:311:GLU:HG3	2.35	0.41
1:E:229:TYR:C	1:E:229:TYR:CD1	2.98	0.41
1:E:269:LEU:HD12	1:E:269:LEU:HA	1.85	0.41
1:E:36:ASN:ND2	1:E:110:PHE:CB	2.81	0.41
1:E:212:SER:HA	5:E:415:HOH:O	2.20	0.41
1:E:137:ARG:NH1	1:E:232:ALA:O	2.53	0.41
1:E:305:ILE:O	1:E:306:TYR:C	2.59	0.41
1:E:41:ASP:OD1	1:E:42:GLN:N	2.54	0.41
1:E:100:PHE:HB3	1:E:103:LEU:HG	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:194:ARG:HE	1:E:194:ARG:HB2	1.80	0.41
1:E:318:PHE:HB3	5:E:684:HOH:O	2.19	0.41
1:E:319:LYS:HD2	1:E:319:LYS:N	2.28	0.41
1:E:343:CYS:N	5:E:659:HOH:O	2.24	0.41
2:I:11:ILE:HD12	2:I:16:THR:HG21	2.03	0.41
1:E:56:ARG:HH11	1:E:56:ARG:HD2	1.41	0.41
1:E:272:LEU:HD12	1:E:272:LEU:HA	1.87	0.40
1:E:81:LYS:NZ	5:E:464:HOH:O	2.36	0.40
1:E:27:LEU:O	1:E:31:GLU:HG2	2.21	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	E	332/350 (95%)	309 (93%)	20 (6%)	3 (1%)	14	14
2	I	18/20 (90%)	16 (89%)	2 (11%)	0	100	100
All	All	350/370 (95%)	325 (93%)	22 (6%)	3 (1%)	14	14

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	184	ASP
1	E	36	ASN
1	E	202	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	E	295/303 (97%)	224 (76%)	71 (24%)	1	0
2	I	15/15 (100%)	14 (93%)	1 (7%)	15	17
All	All	310/318 (98%)	238 (77%)	72 (23%)	1	0

All (72) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	16	LYS
1	E	19	LEU
1	E	24	GLU
1	E	27	LEU
1	E	28	LYS
1	E	31	GLU
1	E	32	THR
1	E	37	THR
1	E	40	LEU
1	E	42	GLN
1	E	45	ARG
1	E	51	THR
1	E	59	LEU
1	E	60	VAL
1	E	61	LYS
1	E	63	LYS
1	E	64	GLU
1	E	65	SER
1	E	74	LEU
1	E	78	LYS
1	E	81	LYS
1	E	91	GLU
1	E	92	LYS
1	E	95	LEU
1	E	104	VAL
1	E	108	PHE
1	E	113	ASN
1	E	120	MET
1	E	134	ARG
1	E	144	ARG
1	E	160	LEU
1	E	161	ASP
1	E	162	LEU

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Mol	Chain	Res	Type
1	E	198	LEU
1	E	208	GLU
1	E	210	ILE
1	E	238	PHE
1	E	242	GLN
1	E	245	GLN
1	E	254	LYS
1	E	256	ARG
1	E	259	SER
1	E	263	SER
1	E	267	ASP
1	E	268	LEU
1	E	269	LEU
1	E	272	LEU
1	E	277	LEU
1	E	278	THR
1	E	285	LYS
1	E	295	LYS
1	E	299	THR
1	E	305	ILE
1	E	308	ARG
1	E	311	GLU
1	E	314	PHE
1	E	315	ILE
1	E	317	LYS
1	E	318	PHE
1	E	319	LYS
1	E	326	ASN
1	E	331	GLU
1	E	335	ILE
1	E	336	ARG
1	E	337	VAL
1	E	340	ASN
1	E	341	GLU
1	E	342	LYS
1	E	345	LYS
1	E	348	THR
1	E	349	GLU
2	I	11	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	35	GLN
1	E	36	ASN
1	E	62	HIS
1	E	77	GLN
1	E	99	ASN
1	E	113	ASN
1	E	158	HIS
1	E	271	ASN
1	E	326	ASN
2	I	20	ASN
2	I	23	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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$
1	SEP	E	338	1	8,9,10	1.07	0	7,12,14	4.88	2 (28%)
1	TPO	E	197	1	8,10,11	0.96	0	10,14,16	1.28	2 (20%)

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
1	SEP	E	338	1	-	6/6/8/10	-
1	TPO	E	197	1	-	1/9/11/13	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	338	SEP	O3P-P-O1P	-9.61	73.39	110.83
1	E	338	SEP	O2P-P-O1P	8.51	144.00	110.83
1	E	197	TPO	P-OG1-CB	-2.22	117.29	123.33
1	E	197	TPO	CG2-CB-CA	-2.12	109.12	113.26

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	E	197	TPO	O-C-CA-CB
1	E	338	SEP	C-CA-CB-OG
1	E	338	SEP	CB-OG-P-O1P
1	E	338	SEP	CB-OG-P-O2P
1	E	338	SEP	CB-OG-P-O3P
1	E	338	SEP	CA-CB-OG-P
1	E	338	SEP	N-CA-CB-OG

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	338	SEP	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
4	ATP	E	355	3	32,33,33	1.05	2 (6%)	48,52,52	1.05	4 (8%)

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
4	ATP	E	355	3	-	3/22/38/38	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	355	ATP	PG-O3G	-3.06	1.43	1.54
4	E	355	ATP	PB-O3A	-2.69	1.56	1.59

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	355	ATP	C4-N9-C1'	3.19	134.08	126.63
4	E	355	ATP	C1'-N9-C8	-2.93	120.60	127.09
4	E	355	ATP	O2G-PG-O3B	2.11	111.70	104.64
4	E	355	ATP	O2B-PB-O3B	2.01	112.71	107.27

There are no chirality outliers.

All (3) torsion outliers are listed below:

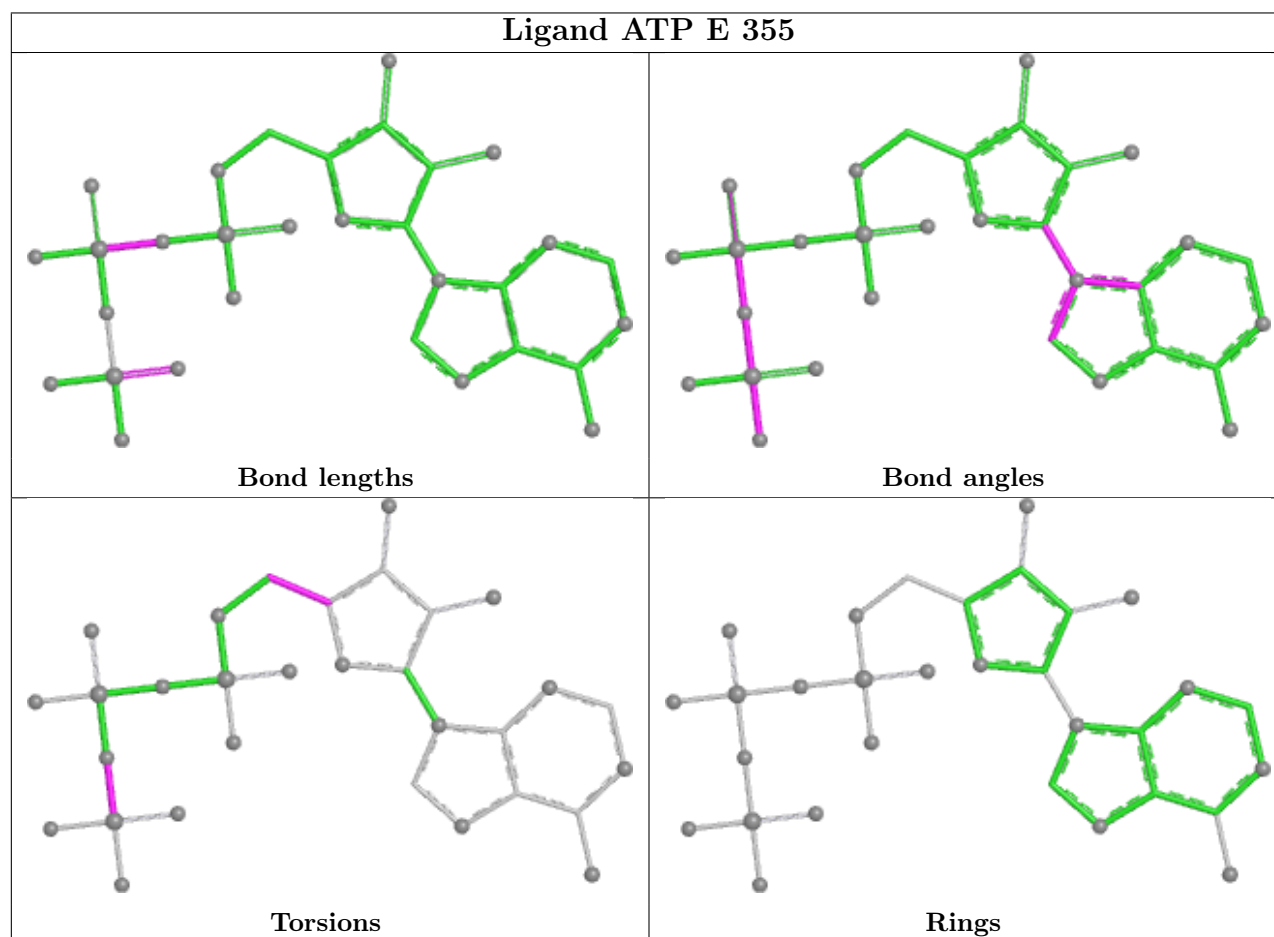
Mol	Chain	Res	Type	Atoms
4	E	355	ATP	PB-O3B-PG-O3G
4	E	355	ATP	PB-O3B-PG-O2G
4	E	355	ATP	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	165:ARG	C	166:ASP	N	1.14

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	E	334/350 (95%)	-0.24	1 (0%) 90 88	15, 32, 67, 88	0
2	I	20/20 (100%)	-0.30	0 100 100	14, 26, 43, 86	0
All	All	354/370 (95%)	-0.25	1 (0%) 90 88	14, 32, 67, 88	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	334	GLU	2.6

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SEP	E	338	10/11	0.96	0.08	35,45,61,100	0
1	TPO	E	197	11/12	0.98	0.07	13,18,45,78	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

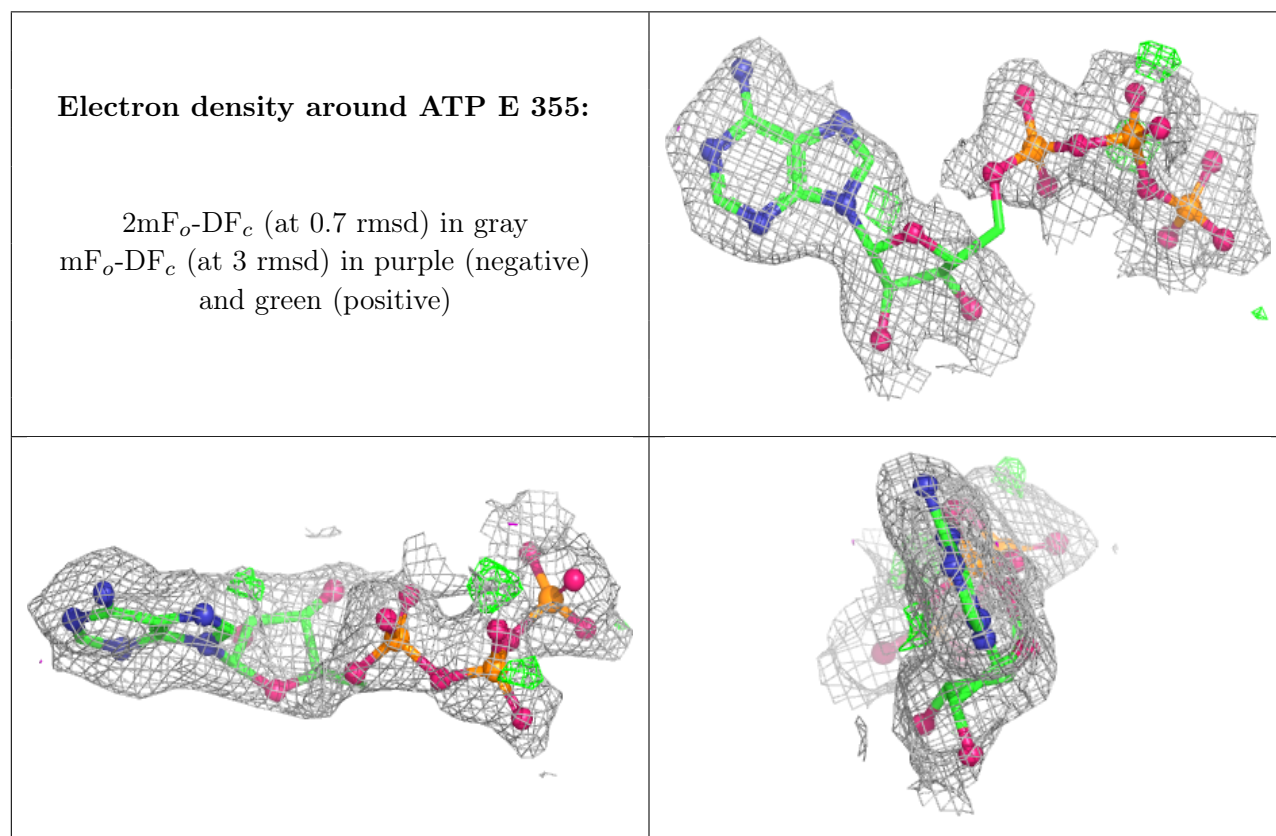
6.4 Ligands [i](#)

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	MN	E	351	1/1	0.98	0.03	23,23,23,23	0
4	ATP	E	355	31/31	0.98	0.06	1,19,45,58	0
3	MN	E	352	1/1	0.99	0.03	21,21,21,21	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 ⓘ

There are no such residues in this entry.