



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 12:02 am BST

PDB ID : 1ZP9  
Title : Crystal Structure of full-length A.fulgidus Rio1 Serine Kinase bound to ATP and Mn<sup>2+</sup> ions.  
Authors : Wlodawer, A.; LaRonde-LeBlanc, N.  
Deposited on : 2005-05-16  
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

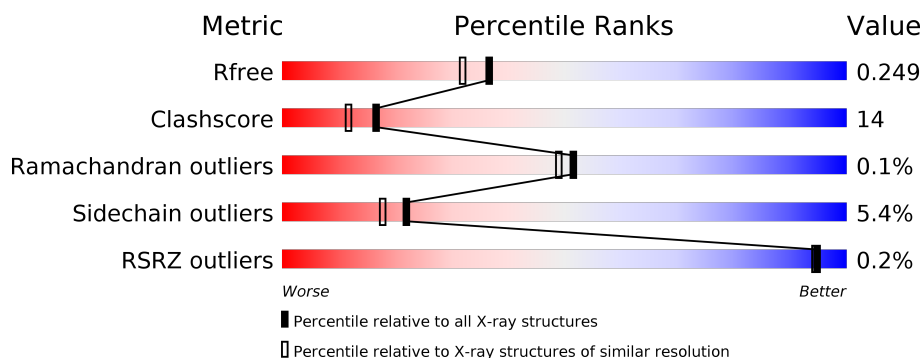
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	258	<div> <div>75%</div> <div>18%</div> <div>• 5%</div> </div>
1	B	258	<div> <div>76%</div> <div>15%</div> <div>• 6%</div> </div>
1	C	258	<div> <div>74%</div> <div>16%</div> <div>• 7%</div> </div>
1	D	258	<div> <div>69%</div> <div>22%</div> <div>• 5%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8879 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 Rio1 kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	245	Total	C	N	O	Se	0	0	0
			1971	1268	320	373	10			
1	B	242	Total	C	N	O	Se	0	0	0
			1963	1266	318	371	8			
1	C	241	Total	C	N	O	Se	0	0	0
			1927	1242	313	363	9			
1	D	244	Total	C	N	O	Se	0	0	0
			1974	1274	320	371	9			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	51	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	77	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	92	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	104	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	141	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	147	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	203	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	213	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	223	MSE	MET	MODIFIED RESIDUE	UNP O28471
A	251	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	1	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	51	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	77	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	92	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	104	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	141	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	147	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	203	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	213	MSE	MET	MODIFIED RESIDUE	UNP O28471
B	223	MSE	MET	MODIFIED RESIDUE	UNP O28471

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Chain	Residue	Modelled	Actual	Comment	Reference
B	251	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	1	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	51	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	77	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	92	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	104	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	141	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	147	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	203	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	213	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	223	MSE	MET	MODIFIED RESIDUE	UNP O28471
C	251	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	1	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	51	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	77	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	92	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	104	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	141	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	147	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	203	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	213	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	223	MSE	MET	MODIFIED RESIDUE	UNP O28471
D	251	MSE	MET	MODIFIED RESIDUE	UNP O28471

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mn 1 1	0	0
2	A	1	Total Mn 1 1	0	0
2	D	1	Total Mn 1 1	0	0
2	C	1	Total Mn 1 1	0	0

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>).



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

- Molecule 4 is water.

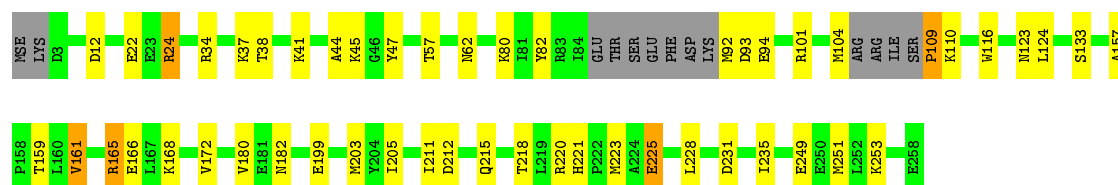
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	239	Total	O	0	0
			239	239		
4	B	257	Total	O	0	0
			257	257		
4	C	211	Total	O	0	0
			211	211		
4	D	209	Total	O	0	0
			209	209		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

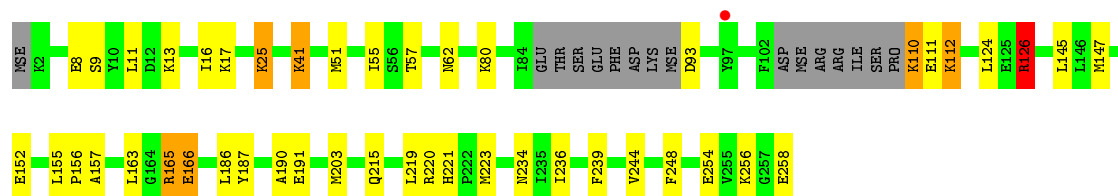
- Molecule 1: Rio1 kinase

Chain A: 



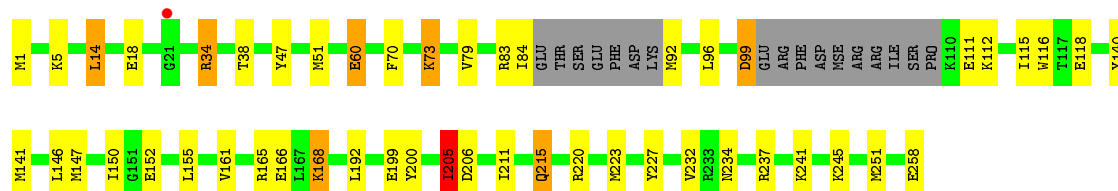
- Molecule 1: Rio1 kinase

Chain B: 



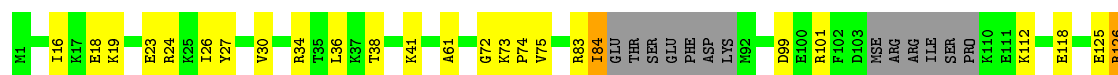
- Molecule 1: Rio1 kinase

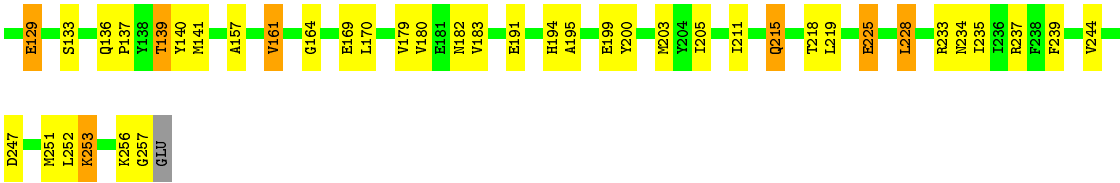
Chain C: 



- Molecule 1: Rio1 kinase

Chain D: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.31Å 80.37Å 121.32Å 90.00° 90.02° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 26.65 – 2.00	Depositor EDS
% Data completeness (in resolution range)	91.2 (30.00-2.00) 91.1 (26.65-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.11	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.48 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.177 , 0.249 0.179 , 0.249	Depositor DCC
$R_{free}$ test set	3196 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	21.1	Xtriage
Anisotropy	0.181	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 41.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.199 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	8879	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.39% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: MN, 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	0.85	0/1995	0.84	3/2665 (0.1%)
1	B	0.88	0/1989	0.86	2/2659 (0.1%)
1	C	0.80	0/1951	0.79	1/2610 (0.0%)
1	D	0.82	1/2000 (0.1%)	0.83	1/2673 (0.0%)
All	All	0.84	1/7935 (0.0%)	0.83	7/10607 (0.1%)

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	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	225	GLU	CG-CD	6.85	1.62	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	126	ARG	NE-CZ-NH2	-8.39	116.11	120.30
1	B	126	ARG	NE-CZ-NH1	7.99	124.30	120.30
1	A	24	ARG	NE-CZ-NH2	-7.65	116.48	120.30
1	A	109	PRO	N-CA-CB	6.23	110.78	103.30
1	A	24	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	C	205	ILE	CB-CA-C	-5.54	100.53	111.60
1	D	126	ARG	NE-CZ-NH1	5.42	123.01	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	72	GLY	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	1971	0	1948	47	0
1	B	1963	0	1953	48	0
1	C	1927	0	1915	56	0
1	D	1974	0	1967	77	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	31	0	12	3	0
3	B	31	0	12	1	0
3	C	31	0	12	1	0
3	D	31	0	12	2	0
4	A	239	0	0	17	0
4	B	257	0	0	25	0
4	C	211	0	0	22	0
4	D	209	0	0	30	0
All	All	8879	0	7831	224	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 (224) 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:D:247:ASP:HB3	4:D:1570:HOH:O	1.51	1.10
1:A:47:TYR:HD2	4:A:1753:HOH:O	1.41	1.02
1:D:257:GLY:HA3	4:D:1835:HOH:O	1.61	0.99
1:D:38:THR:HG23	1:D:141:MSE:HE2	1.43	0.99

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:ALA:HB2	1:D:205:ILE:HD11	1.45	0.99
1:C:241:LYS:HD2	4:C:1406:HOH:O	1.63	0.96
1:A:205:ILE:HB	4:A:1561:HOH:O	1.67	0.93
1:A:22:GLU:HG2	4:A:1568:HOH:O	1.67	0.93
1:D:141:MSE:HE3	4:D:1439:HOH:O	1.70	0.89
1:D:161:VAL:HG13	1:D:199:GLU:HG3	1.55	0.88
1:A:223:MSE:HE2	4:A:1661:HOH:O	1.73	0.88
1:B:145:LEU:HD11	4:B:1850:HOH:O	1.74	0.88
1:B:110:LYS:HG2	1:B:111:GLU:N	1.89	0.86
1:D:233:ARG:O	1:D:237:ARG:HG3	1.76	0.85
1:B:221:HIS:HD2	1:B:223:MSE:H	1.21	0.85
1:C:232:VAL:HG22	1:C:251:MSE:HE2	1.60	0.84
1:B:80:LYS:HD2	4:B:1722:HOH:O	1.77	0.83
1:D:180:VAL:HG11	1:D:251:MSE:HE2	1.59	0.83
1:B:110:LYS:HG2	1:B:111:GLU:H	1.47	0.80
1:A:221:HIS:HD2	1:A:223:MSE:H	1.30	0.80
1:D:161:VAL:HG13	1:D:199:GLU:CG	2.12	0.80
1:C:234:ASN:ND2	1:C:237:ARG:HH22	1.80	0.79
1:A:218:THR:HG22	1:A:220:ARG:H	1.45	0.79
1:D:157:ALA:CB	1:D:205:ILE:HD11	2.12	0.78
1:C:234:ASN:HD22	1:C:237:ARG:HH22	1.30	0.78
1:D:253:LYS:HG3	4:D:1190:HOH:O	1.84	0.77
1:C:168:LYS:HE3	4:C:1556:HOH:O	1.84	0.76
1:D:141:MSE:CE	4:D:1439:HOH:O	2.30	0.76
1:D:180:VAL:HG11	1:D:251:MSE:CE	2.17	0.75
1:D:251:MSE:HE3	1:D:251:MSE:HA	1.69	0.73
1:C:1:MSE:HA	4:C:1737:HOH:O	1.88	0.73
1:B:220:ARG:NH2	4:B:1351:HOH:O	2.23	0.72
1:A:41:LYS:HE3	4:A:1323:HOH:O	1.90	0.72
1:C:70:PHE:HD1	4:C:1761:HOH:O	1.73	0.72
1:A:94:GLU:HB2	4:A:1661:HOH:O	1.91	0.71
1:D:164:GLY:HA2	4:D:1130:HOH:O	1.91	0.71
1:C:79:VAL:HG22	1:C:146:LEU:CD2	2.20	0.70
1:C:111:GLU:O	1:C:115:ILE:HD12	1.91	0.70
1:D:180:VAL:CG1	1:D:251:MSE:HE2	2.21	0.70
1:A:37:LYS:HD2	4:A:1908:HOH:O	1.92	0.69
1:A:166:GLU:HG2	4:A:1717:HOH:O	1.90	0.69
1:A:221:HIS:CD2	1:A:223:MSE:H	2.10	0.69
1:A:38:THR:HG22	4:A:1653:HOH:O	1.92	0.69
1:D:99:ASP:OD2	1:D:101:ARG:HD3	1.93	0.68
1:A:101:ARG:HH22	1:A:123:ASN:HD21	1.42	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:111:GLU:HB3	4:B:1867:HOH:O	1.93	0.68
1:D:84:ILE:HG21	4:D:1169:HOH:O	1.94	0.68
1:C:92:MSE:HE2	1:C:112:LYS:HG2	1.75	0.67
1:A:161:VAL:HG13	1:A:199:GLU:HG3	1.77	0.67
1:C:92:MSE:CE	1:C:112:LYS:HG2	2.25	0.66
1:B:165:ARG:NH1	4:B:1696:HOH:O	2.27	0.66
1:C:92:MSE:CE	1:C:96:LEU:HD11	2.26	0.66
1:D:133:SER:H	1:D:182:ASN:ND2	1.92	0.66
1:D:218:THR:HG22	1:D:219:LEU:N	2.11	0.66
1:C:1:MSE:CA	4:C:1737:HOH:O	2.42	0.65
1:B:147:MSE:HB3	4:B:1850:HOH:O	1.96	0.65
1:B:221:HIS:CD2	1:B:223:MSE:H	2.10	0.64
1:B:93:ASP:N	4:B:1724:HOH:O	2.30	0.64
1:D:112:LYS:HE2	4:D:1874:HOH:O	1.97	0.64
1:D:118:GLU:HG3	1:D:140:TYR:CZ	2.31	0.64
1:B:165:ARG:HD2	4:B:1538:HOH:O	1.97	0.64
1:C:14:LEU:HG	1:C:51:MSE:HG3	1.79	0.64
1:D:233:ARG:NH1	4:D:1713:HOH:O	2.27	0.64
1:D:191:GLU:C	1:D:219:LEU:HD13	2.17	0.64
1:A:109:PRO:N	4:A:1669:HOH:O	2.31	0.63
1:D:126:ARG:O	1:D:129:GLU:HG3	1.98	0.63
1:D:84:ILE:CG2	4:D:1169:HOH:O	2.48	0.62
1:C:84:ILE:HD11	4:C:1596:HOH:O	2.00	0.62
1:C:118:GLU:HG3	1:C:140:TYR:CZ	2.34	0.62
1:B:111:GLU:HG2	4:B:1607:HOH:O	1.99	0.61
1:D:234:ASN:ND2	4:D:1500:HOH:O	2.32	0.61
1:D:161:VAL:CG1	1:D:199:GLU:HG3	2.28	0.60
1:C:234:ASN:HD22	1:C:237:ARG:NH2	1.98	0.59
1:D:126:ARG:HD2	4:D:1041:HOH:O	2.01	0.59
1:B:157:ALA:HB3	1:B:203:MSE:HE2	1.83	0.59
1:B:223:MSE:CE	4:B:1644:HOH:O	2.51	0.59
1:C:241:LYS:HG2	1:D:26:ILE:HG12	1.84	0.59
1:C:79:VAL:HG22	1:C:146:LEU:HD22	1.84	0.59
1:C:92:MSE:HE2	1:C:96:LEU:HD11	1.84	0.58
1:A:218:THR:HB	4:A:1170:HOH:O	2.02	0.58
1:C:92:MSE:HE3	1:C:116:TRP:CE3	2.39	0.58
1:D:225:GLU:HG3	4:D:1252:HOH:O	2.03	0.58
1:D:126:ARG:CZ	4:D:1846:HOH:O	2.50	0.58
1:D:157:ALA:HB2	1:D:205:ILE:CD1	2.28	0.58
1:A:218:THR:HG22	1:A:220:ARG:N	2.17	0.57
1:A:133:SER:H	1:A:182:ASN:ND2	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:232:VAL:HG13	1:C:251:MSE:SE	2.55	0.57
1:A:161:VAL:HG13	1:A:199:GLU:CG	2.35	0.57
1:C:211:ILE:HD11	4:C:1334:HOH:O	2.03	0.57
1:B:234:ASN:ND2	4:B:1165:HOH:O	2.33	0.57
1:A:211:ILE:HD11	4:A:1741:HOH:O	2.04	0.56
1:A:249:GLU:OE2	1:A:253:LYS:NZ	2.35	0.56
1:D:237:ARG:NH2	4:D:1221:HOH:O	2.38	0.56
1:A:34:ARG:NH1	1:A:110:LYS:HD3	2.21	0.56
1:B:110:LYS:HG3	4:B:1607:HOH:O	2.05	0.56
1:D:179:VAL:HG11	1:D:235:ILE:HD12	1.88	0.56
1:D:83:ARG:HD3	4:D:1330:HOH:O	2.05	0.56
1:A:218:THR:CG2	1:A:220:ARG:H	2.16	0.56
1:B:110:LYS:CG	1:B:111:GLU:H	2.17	0.55
1:B:165:ARG:NH1	4:B:1751:HOH:O	2.38	0.55
1:B:41:LYS:HD2	4:B:1579:HOH:O	2.04	0.55
1:C:223:MSE:HG2	1:C:227:TYR:CE1	2.42	0.55
1:D:252:LEU:O	1:D:256:LYS:HG3	2.06	0.55
1:C:92:MSE:HB2	4:C:1582:HOH:O	2.06	0.55
1:D:157:ALA:CB	1:D:205:ILE:CD1	2.85	0.55
1:B:17:LYS:HE3	4:B:1308:HOH:O	2.06	0.55
1:C:18:GLU:CB	4:C:1347:HOH:O	2.55	0.55
1:C:215:GLN:NE2	1:C:215:GLN:H	2.04	0.55
1:D:73:LYS:HB2	1:D:74:PRO:HD2	1.88	0.55
1:A:104:MSE:HG3	4:A:1781:HOH:O	2.07	0.54
1:A:101:ARG:HH12	1:A:123:ASN:ND2	2.04	0.54
1:A:41:LYS:O	1:A:45:LYS:HG3	2.08	0.54
1:B:163:LEU:O	1:B:166:GLU:HB2	2.07	0.54
1:C:220:ARG:HD2	4:C:1425:HOH:O	2.08	0.53
1:D:73:LYS:HE2	4:D:1913:HOH:O	2.07	0.53
1:D:129:GLU:HB3	4:D:1020:HOH:O	2.08	0.53
1:D:203:MSE:HE2	1:D:211:ILE:HG21	1.91	0.53
1:D:83:ARG:CD	4:D:1330:HOH:O	2.56	0.53
1:A:157:ALA:HB3	1:A:203:MSE:HE2	1.89	0.53
1:A:225:GLU:OE2	1:A:228:LEU:HD23	2.09	0.53
1:B:165:ARG:NH2	4:B:1591:HOH:O	2.29	0.53
1:C:168:LYS:HG3	4:C:1683:HOH:O	2.08	0.53
1:D:16:ILE:HG22	1:D:24:ARG:HG3	1.90	0.52
1:A:212:ASP:OD1	3:A:260:ATP:O3G	2.26	0.52
1:A:57:THR:HA	1:A:62:ASN:HD22	1.75	0.52
1:C:112:LYS:NZ	4:C:1678:HOH:O	2.43	0.52
1:D:218:THR:HG22	1:D:219:LEU:H	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ALA:HB2	4:A:1731:HOH:O	2.10	0.51
1:B:110:LYS:CG	1:B:111:GLU:N	2.68	0.51
1:D:61:ALA:CB	3:D:260:ATP:O1G	2.59	0.51
1:A:80:LYS:NZ	3:A:260:ATP:O3G	2.38	0.51
1:A:47:TYR:CD2	4:A:1753:HOH:O	2.31	0.51
1:D:253:LYS:CG	4:D:1190:HOH:O	2.53	0.51
1:C:38:THR:HG23	1:C:141:MSE:HG2	1.91	0.51
1:D:253:LYS:CE	4:D:1732:HOH:O	2.59	0.51
1:B:41:LYS:HE3	4:C:1584:HOH:O	2.10	0.51
1:D:41:LYS:HD3	1:D:141:MSE:HE1	1.93	0.51
1:D:126:ARG:NH1	4:D:1846:HOH:O	2.44	0.50
1:A:218:THR:HG22	1:A:220:ARG:HB2	1.94	0.50
1:D:141:MSE:HE3	4:D:1686:HOH:O	2.11	0.50
1:D:125:GLU:O	1:D:129:GLU:HG2	2.11	0.50
1:A:12:ASP:OD1	1:A:24:ARG:NH2	2.42	0.50
1:A:165:ARG:HD3	1:A:165:ARG:N	2.25	0.50
1:B:16:ILE:HD11	1:B:51:MSE:HB2	1.93	0.49
1:C:223:MSE:HG2	1:C:227:TYR:CZ	2.47	0.49
1:D:139:THR:HG21	4:D:1473:HOH:O	2.11	0.49
1:B:165:ARG:NH1	4:B:1538:HOH:O	2.44	0.49
1:B:186:LEU:O	1:B:190:ALA:HB3	2.13	0.49
1:C:150:ILE:CG2	1:C:205:ILE:HD11	2.43	0.49
1:C:223:MSE:HB3	1:C:227:TYR:CE2	2.48	0.48
1:B:17:LYS:HE2	4:B:1484:HOH:O	2.13	0.48
1:C:150:ILE:HG22	1:C:205:ILE:HD11	1.96	0.48
1:A:34:ARG:NH2	4:A:1652:HOH:O	2.46	0.47
1:B:126:ARG:HD3	1:B:190:ALA:O	2.14	0.47
1:B:220:ARG:NE	4:B:1270:HOH:O	2.43	0.47
1:B:9:SER:O	1:B:13:LYS:HG3	2.15	0.47
1:C:70:PHE:O	1:C:73:LYS:HD2	2.15	0.47
1:D:30:VAL:HA	1:D:83:ARG:HH22	1.79	0.47
1:B:239:PHE:HB3	1:B:244:VAL:HB	1.96	0.47
1:C:258:GLU:HB3	4:C:1401:HOH:O	2.15	0.46
1:A:218:THR:CG2	1:A:220:ARG:HB2	2.45	0.46
1:C:161:VAL:HG21	1:C:200:TYR:CE1	2.50	0.46
1:C:99:ASP:HB2	4:C:1403:HOH:O	2.14	0.46
1:A:124:LEU:HD21	4:A:1888:HOH:O	2.14	0.46
1:D:218:THR:HG21	4:D:1380:HOH:O	2.15	0.46
1:D:61:ALA:HB2	3:D:260:ATP:O1G	2.16	0.46
1:D:239:PHE:HB3	1:D:244:VAL:HB	1.98	0.45
1:D:194:HIS:O	1:D:195:ALA:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:218:THR:CG2	4:D:1380:HOH:O	2.64	0.45
1:D:253:LYS:HD3	4:D:1777:HOH:O	2.16	0.45
1:B:110:LYS:CG	4:B:1607:HOH:O	2.63	0.45
1:B:187:TYR:O	1:B:191:GLU:HA	2.17	0.45
1:D:136:GLN:NE2	1:D:137:PRO:HD2	2.32	0.45
1:B:112:LYS:HD3	4:B:1398:HOH:O	2.16	0.45
1:A:92:MSE:HG2	1:A:116:TRP:CZ3	2.53	0.44
1:D:215:GLN:NE2	1:D:215:GLN:H	2.15	0.44
1:D:118:GLU:HG3	1:D:140:TYR:OH	2.16	0.44
1:D:161:VAL:HG21	1:D:200:TYR:CD1	2.53	0.44
1:D:253:LYS:HE2	4:D:1732:HOH:O	2.18	0.44
1:D:218:THR:CG2	1:D:219:LEU:N	2.80	0.44
1:B:256:LYS:HE2	1:B:258:GLU:OE1	2.17	0.44
1:D:251:MSE:CE	1:D:251:MSE:HA	2.42	0.44
1:B:157:ALA:O	1:B:203:MSE:HE3	2.18	0.44
1:B:8:GLU:HG2	4:B:1467:HOH:O	2.18	0.44
1:A:82:TYR:HE1	3:A:260:ATP:O1G	2.01	0.43
1:B:155:LEU:HA	1:B:156:PRO:HD3	1.79	0.43
1:C:245:LYS:HD2	1:D:83:ARG:CZ	2.49	0.43
1:C:70:PHE:CD1	4:C:1761:HOH:O	2.56	0.43
1:A:159:THR:OG1	1:A:161:VAL:HG23	2.18	0.43
1:B:112:LYS:CD	4:B:1398:HOH:O	2.66	0.43
1:A:180:VAL:HG22	1:A:251:MSE:HE3	2.01	0.43
1:C:161:VAL:HG23	1:C:199:GLU:HG3	2.00	0.43
1:B:157:ALA:O	1:B:203:MSE:CE	2.66	0.43
1:B:25:LYS:HA	1:B:25:LYS:HD2	1.50	0.43
1:C:166:GLU:HG3	4:C:1466:HOH:O	2.18	0.42
1:D:139:THR:CG2	4:D:1473:HOH:O	2.68	0.42
1:D:228:LEU:HD13	1:D:252:LEU:CD1	2.49	0.42
1:D:233:ARG:NH2	4:D:1277:HOH:O	2.52	0.42
1:B:57:THR:HA	1:B:62:ASN:HD22	1.85	0.42
1:A:34:ARG:HH12	1:A:110:LYS:HD3	1.83	0.42
1:A:231:ASP:O	1:A:235:ILE:HG12	2.19	0.42
1:B:223:MSE:HE1	4:B:1644:HOH:O	2.18	0.42
1:A:225:GLU:HA	1:A:225:GLU:OE2	2.20	0.42
1:D:34:ARG:HG2	1:D:34:ARG:H	1.60	0.42
1:D:179:VAL:O	1:D:183:VAL:HG23	2.20	0.42
1:C:92:MSE:CE	1:C:116:TRP:HB2	2.50	0.41
1:B:55:ILE:HG21	3:B:260:ATP:H1'	2.02	0.41
1:C:79:VAL:HG22	1:C:146:LEU:HD23	1.99	0.41
1:D:18:GLU:HA	1:D:24:ARG:HD3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:245:LYS:HZ2	1:D:83:ARG:NH2	2.19	0.41
1:C:241:LYS:HE2	1:D:26:ILE:HD11	2.02	0.41
1:C:60:GLU:HG3	3:C:260:ATP:O2G	2.20	0.41
1:A:157:ALA:O	1:A:203:MSE:HE3	2.21	0.41
1:C:92:MSE:HE1	1:C:112:LYS:HG2	2.00	0.41
1:C:223:MSE:HB2	1:C:223:MSE:HE2	1.96	0.41
1:C:237:ARG:HD3	4:C:1420:HOH:O	2.21	0.41
1:B:254:GLU:HG3	4:B:1894:HOH:O	2.20	0.41
1:C:47:TYR:HE1	4:C:1761:HOH:O	2.03	0.41
1:C:147:MSE:HE1	4:C:1013:HOH:O	2.21	0.40
1:B:236:ILE:CD1	1:B:248:PHE:HA	2.51	0.40
1:C:83:ARG:NH2	4:C:1241:HOH:O	2.53	0.40
1:D:23:GLU:O	1:D:27:TYR:HD1	2.04	0.40
1:C:34:ARG:H	1:C:34:ARG:HG2	1.69	0.40
1:C:60:GLU:HG2	4:C:1011:HOH:O	2.21	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	239/258 (93%)	233 (98%)	6 (2%)	0	100	100
1	B	236/258 (92%)	225 (95%)	11 (5%)	0	100	100
1	C	235/258 (91%)	226 (96%)	8 (3%)	1 (0%)	34	30
1	D	238/258 (92%)	232 (98%)	6 (2%)	0	100	100
All	All	948/1032 (92%)	916 (97%)	31 (3%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	C	206	ASP

### 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	205/216 (95%)	198 (97%)	7 (3%)	37	36
1	B	206/216 (95%)	194 (94%)	12 (6%)	20	15
1	C	200/216 (93%)	187 (94%)	13 (6%)	17	12
1	D	206/216 (95%)	194 (94%)	12 (6%)	20	15
All	All	817/864 (95%)	773 (95%)	44 (5%)	22	18

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	ASP
1	A	161	VAL
1	A	165	ARG
1	A	168	LYS
1	A	172	VAL
1	A	215	GLN
1	A	225	GLU
1	B	11	LEU
1	B	25	LYS
1	B	41	LYS
1	B	110	LYS
1	B	112	LYS
1	B	124	LEU
1	B	126	ARG
1	B	152	GLU
1	B	165	ARG
1	B	166	GLU
1	B	215	GLN
1	B	219	LEU
1	C	5	LYS
1	C	14	LEU

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Mol	Chain	Res	Type
1	C	34	ARG
1	C	60	GLU
1	C	73	LYS
1	C	99	ASP
1	C	152	GLU
1	C	155	LEU
1	C	165	ARG
1	C	168	LYS
1	C	192	LEU
1	C	205	ILE
1	C	215	GLN
1	D	19	LYS
1	D	36	LEU
1	D	75	VAL
1	D	84	ILE
1	D	129	GLU
1	D	139	THR
1	D	161	VAL
1	D	169	GLU
1	D	170	LEU
1	D	215	GLN
1	D	228	LEU
1	D	253	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	123	ASN
1	A	182	ASN
1	A	188	GLN
1	A	215	GLN
1	A	221	HIS
1	B	62	ASN
1	B	136	GLN
1	B	177	ASN
1	B	188	GLN
1	B	215	GLN
1	B	221	HIS
1	C	177	ASN
1	C	188	GLN
1	C	215	GLN

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Mol	Chain	Res	Type
1	C	234	ASN
1	D	136	GLN
1	D	182	ASN
1	D	188	GLN
1	D	215	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 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	$\# Z  > 2$
3	ATP	D	260	2	26,33,33	1.34	2 (7%)	31,52,52	1.36	4 (12%)
3	ATP	C	260	2	26,33,33	1.15	2 (7%)	31,52,52	1.54	2 (6%)
3	ATP	B	260	2	26,33,33	1.24	2 (7%)	31,52,52	1.68	5 (16%)
3	ATP	A	260	2	26,33,33	0.82	0	31,52,52	1.74	5 (16%)

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	ATP	D	260	2	-	2/18/38/38	0/3/3/3
3	ATP	C	260	2	-	0/18/38/38	0/3/3/3
3	ATP	B	260	2	-	2/18/38/38	0/3/3/3
3	ATP	A	260	2	-	1/18/38/38	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	260	ATP	C2-N3	5.00	1.40	1.32
3	B	260	ATP	C2-N3	4.24	1.38	1.32
3	C	260	ATP	C2-N3	4.04	1.38	1.32
3	D	260	ATP	C2-N1	3.24	1.39	1.33
3	B	260	ATP	C2-N1	3.01	1.39	1.33
3	C	260	ATP	O4'-C1'	2.22	1.44	1.41

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	260	ATP	N3-C2-N1	-6.84	117.99	128.68
3	C	260	ATP	N3-C2-N1	-6.58	118.39	128.68
3	B	260	ATP	N3-C2-N1	-4.84	121.12	128.68
3	B	260	ATP	O3G-PG-O3B	4.55	119.91	104.64
3	D	260	ATP	N3-C2-N1	-4.32	121.92	128.68
3	D	260	ATP	O3G-PG-O3B	3.28	115.62	104.64
3	A	260	ATP	PB-O3B-PG	-3.11	122.15	132.83
3	A	260	ATP	O3G-PG-O1G	-2.49	100.95	110.68
3	B	260	ATP	O2G-PG-O3B	2.44	112.81	104.64
3	C	260	ATP	C4-C5-N7	-2.28	107.02	109.40
3	A	260	ATP	O3G-PG-O3B	2.28	112.27	104.64
3	B	260	ATP	C4-C5-N7	-2.26	107.05	109.40
3	D	260	ATP	C4-C5-N7	-2.12	107.19	109.40
3	A	260	ATP	C2-N1-C6	2.10	122.35	118.75
3	D	260	ATP	PB-O3B-PG	-2.06	125.75	132.83
3	B	260	ATP	C1'-N9-C4	-2.04	123.05	126.64

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	260	ATP	PB-O3B-PG-O1G

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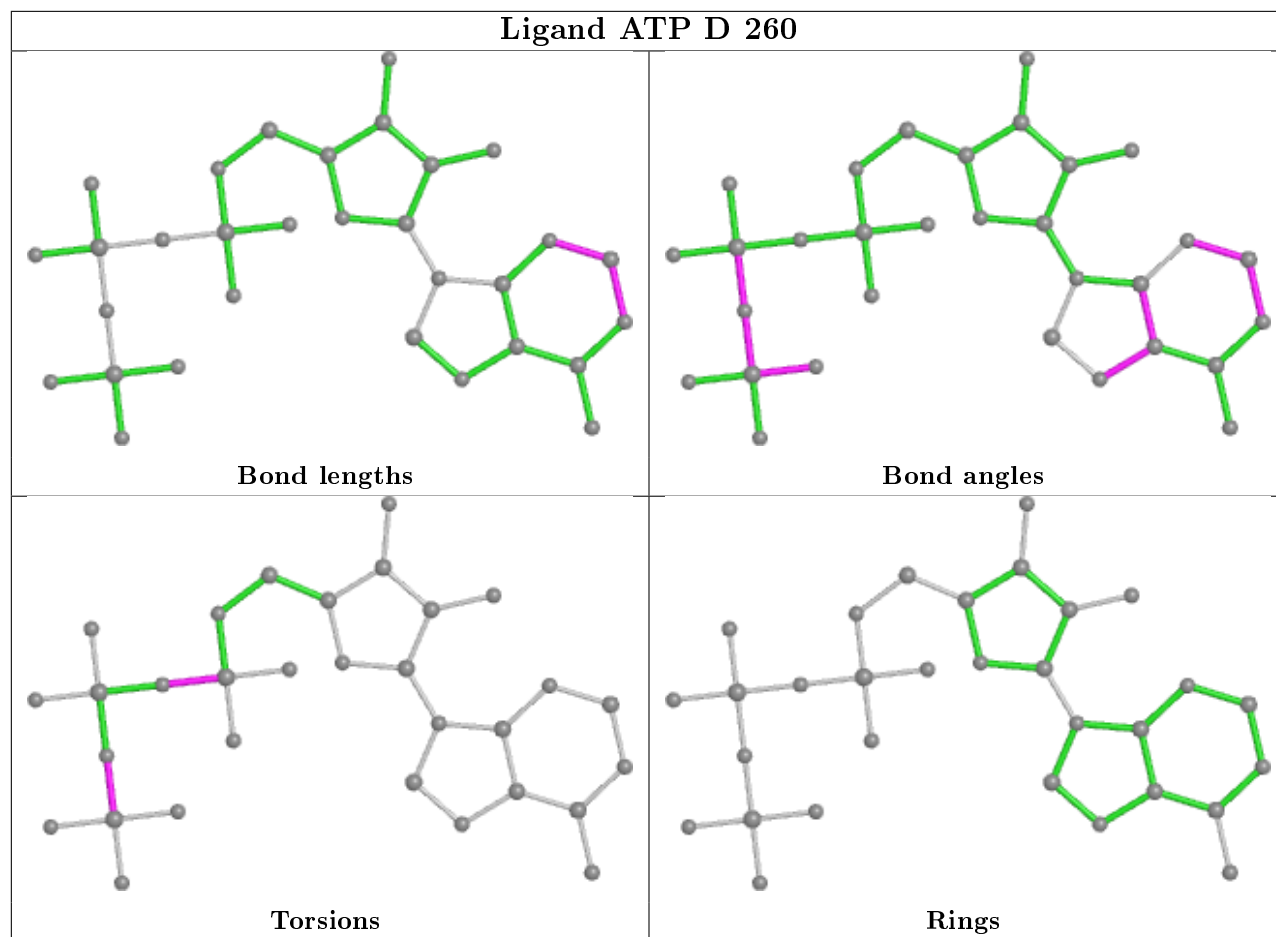
Mol	Chain	Res	Type	Atoms
3	B	260	ATP	PB-O3A-PA-O1A
3	D	260	ATP	PB-O3A-PA-O1A
3	D	260	ATP	PB-O3B-PG-O1G
3	B	260	ATP	PB-O3A-PA-O2A

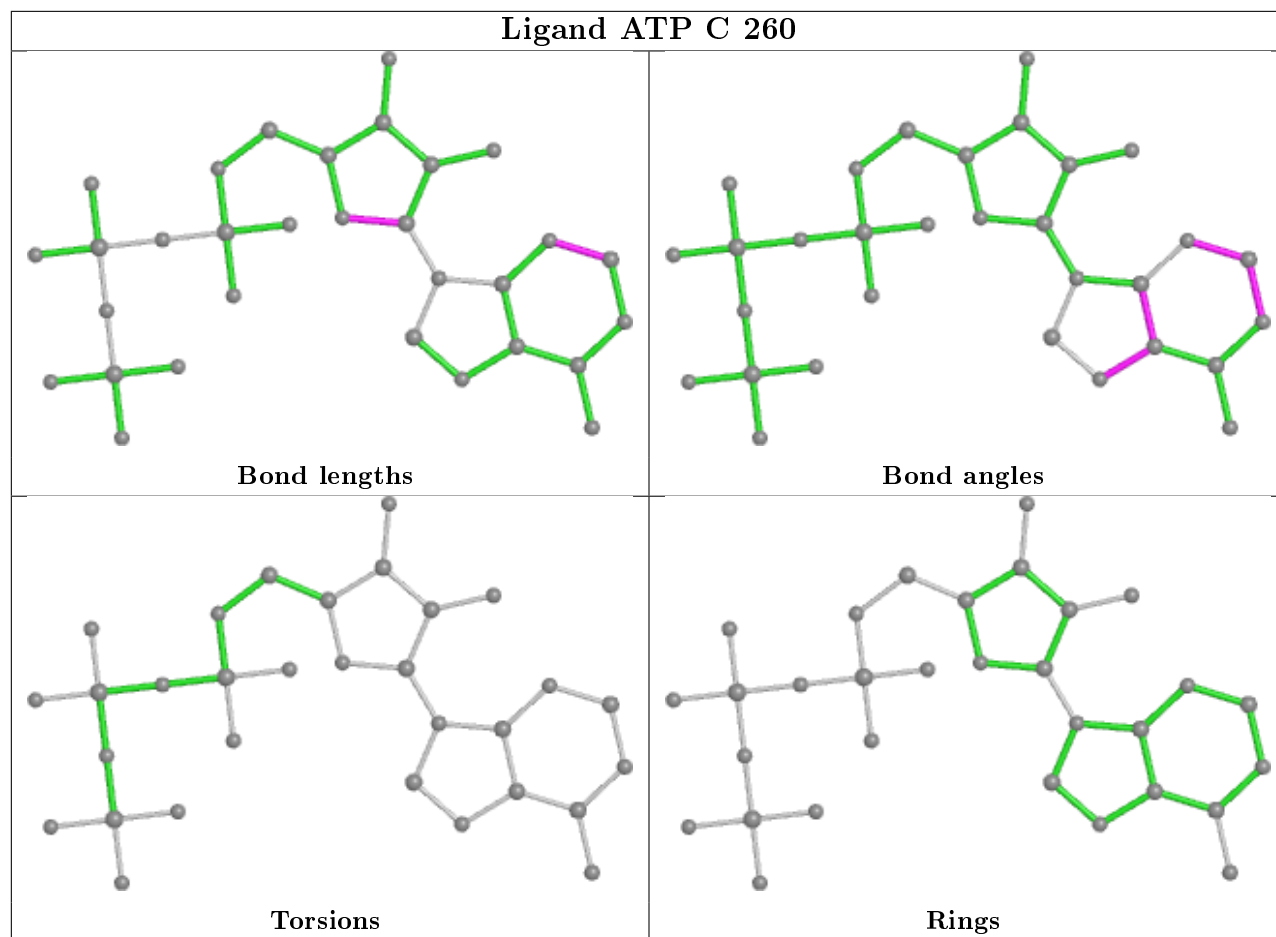
There are no ring outliers.

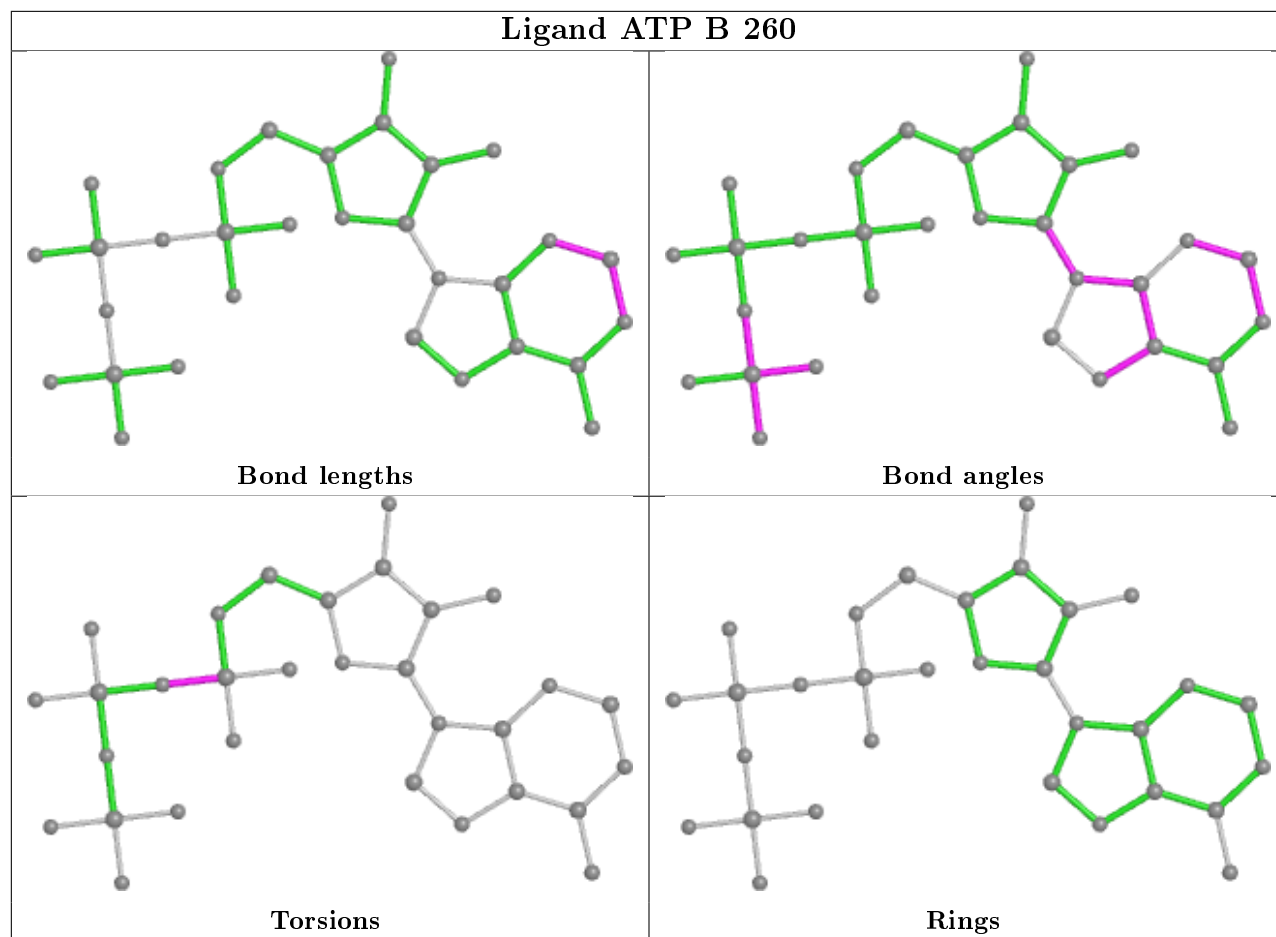
4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	260	ATP	2	0
3	C	260	ATP	1	0
3	B	260	ATP	1	0
3	A	260	ATP	3	0

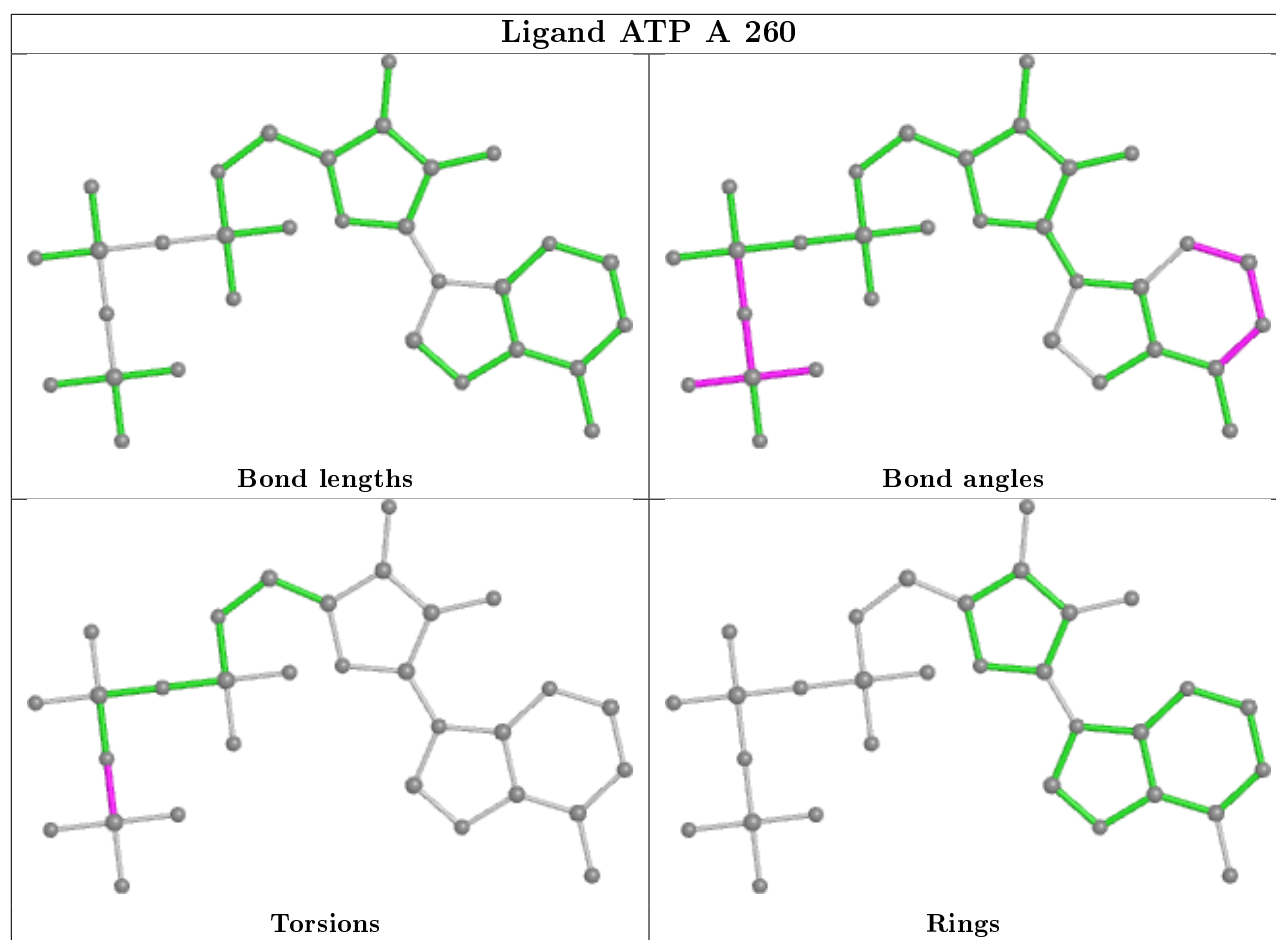
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	235/258 (91%)	-0.39	0 <b>100</b> <b>100</b>	11, 18, 38, 49	0
1	B	234/258 (90%)	-0.36	1 (0%) <b>92</b> <b>92</b>	10, 18, 34, 45	0
1	C	231/258 (89%)	-0.30	1 (0%) <b>92</b> <b>92</b>	12, 21, 39, 45	0
1	D	234/258 (90%)	-0.25	0 <b>100</b> <b>100</b>	13, 22, 38, 43	0
All	All	934/1032 (90%)	-0.33	2 (0%) <b>95</b> <b>94</b>	10, 20, 38, 49	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	21	GLY	3.4
1	B	97	TYR	2.4

### 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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

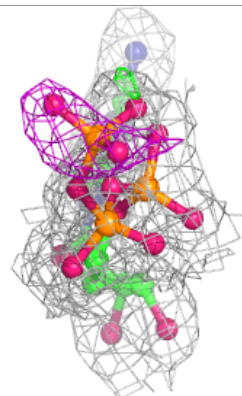
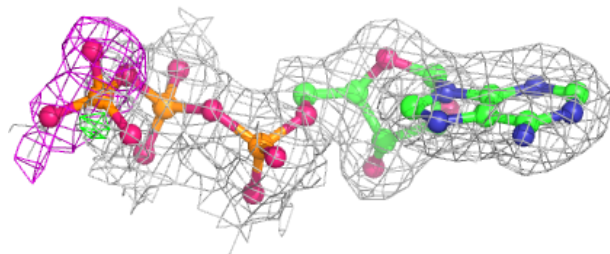
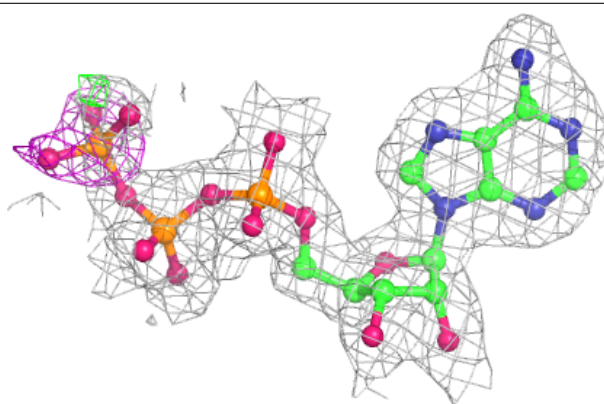
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ATP	D	260	31/31	0.96	0.12	10,15,25,27	4
3	ATP	B	260	31/31	0.97	0.11	8,13,23,26	4
3	ATP	C	260	31/31	0.97	0.12	8,13,27,29	4
3	ATP	A	260	31/31	0.98	0.10	9,14,22,25	4
2	MN	A	259	1/1	1.00	0.04	16,16,16,16	0
2	MN	B	259	1/1	1.00	0.05	14,14,14,14	0
2	MN	C	259	1/1	1.00	0.09	19,19,19,19	0
2	MN	D	259	1/1	1.00	0.06	19,19,19,19	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.

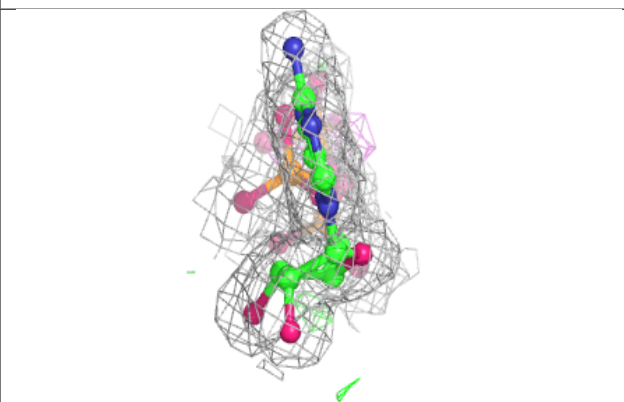
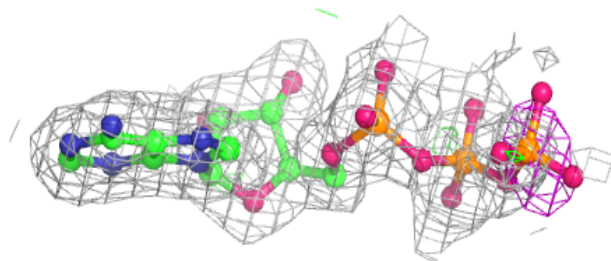
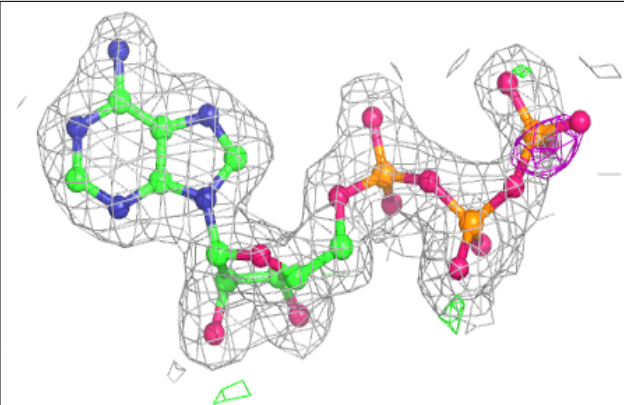
**Electron density around ATP D 260:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

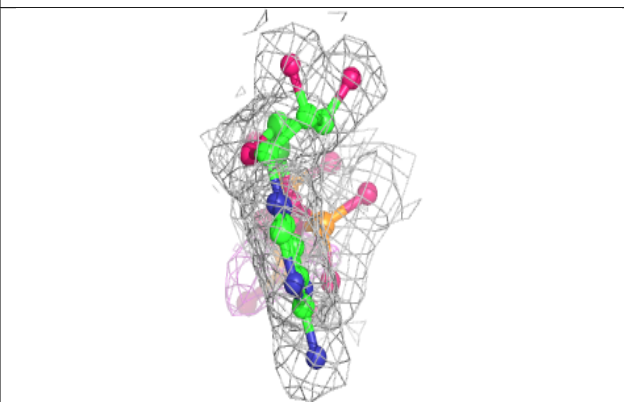
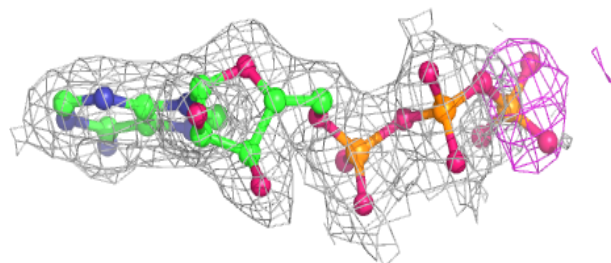
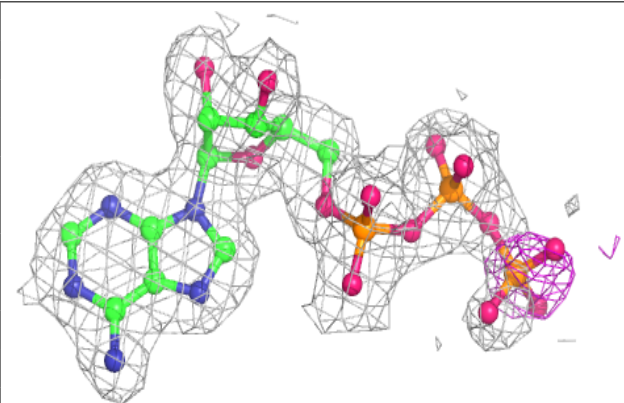


**Electron density around ATP B 260:**

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

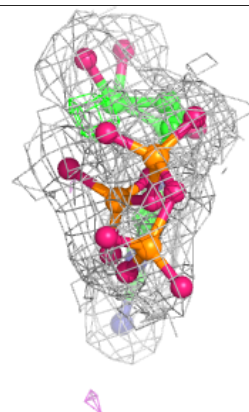
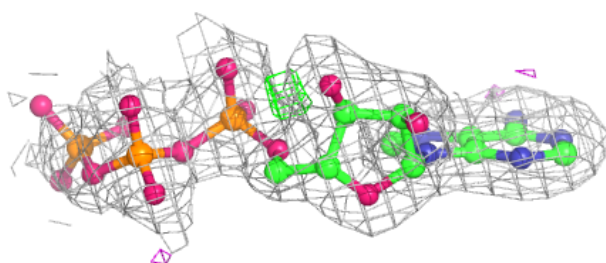
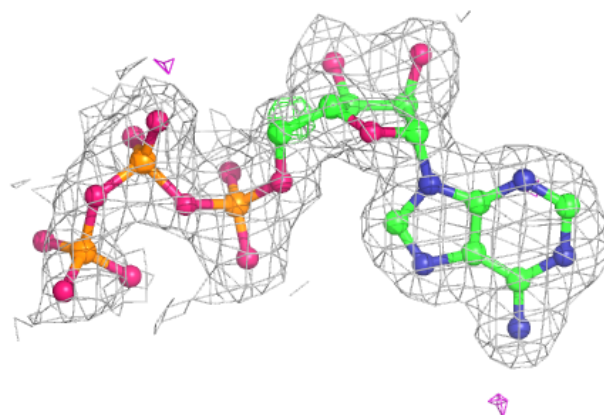
**Electron density around ATP C 260:**

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



**Electron density around ATP A 260:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.