



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2023 – 07:07 PM EDT

PDB ID : 7RUT
Title : Structure of Human ATP:Cobalamin Adenosyltransferase R190C bound to ATP
Authors : Mascarenhas, R.; Gouda, H.; Koutmos, M.; Banerjee, R.
Deposited on : 2021-08-18
Resolution : 1.50 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.35.1
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) : 2.35.1

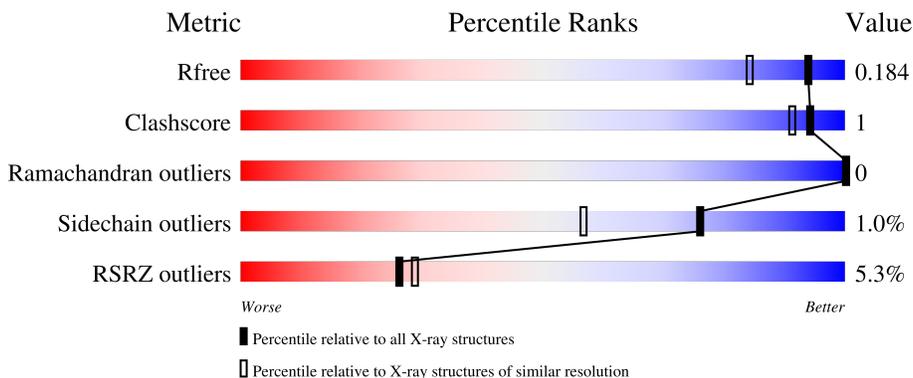
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

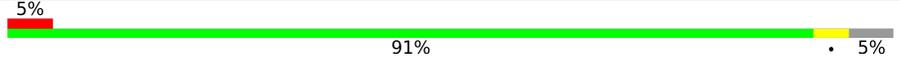
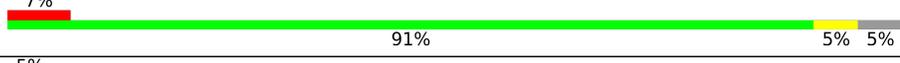
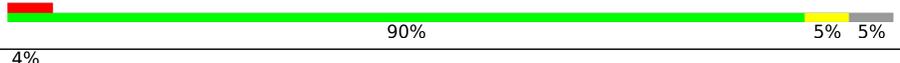
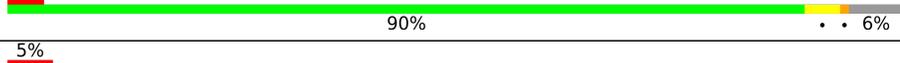
The reported resolution of this entry is 1.50 Å.

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.



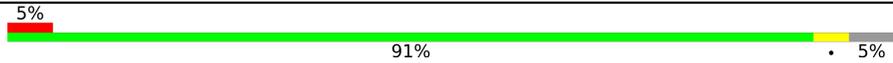
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2936 (1.50-1.50)
Clashscore	141614	3144 (1.50-1.50)
Ramachandran outliers	138981	3066 (1.50-1.50)
Sidechain outliers	138945	3064 (1.50-1.50)
RSRZ outliers	127900	2884 (1.50-1.50)

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	196	
1	B	196	
1	C	196	
1	D	196	
1	E	196	

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Mol	Chain	Length	Quality of chain
1	F	196	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a small red segment on the left labeled '5%', a large green segment in the middle labeled '91%', and a small yellow segment on the right labeled '5%'. The segments are separated by thin black lines.</p>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9572 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 Corrinoid adenosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	187	1427	900	238	281	8	0	1	0
1	C	186	1426	901	241	275	9	0	4	0
1	F	186	1436	906	243	279	8	0	3	0
1	B	187	1440	908	247	276	9	0	4	0
1	A	186	1458	925	246	279	8	0	5	0
1	D	185	1456	921	249	278	8	0	4	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	55	MET	-	initiating methionine	UNP Q96EY8
E	190	CYS	ARG	engineered mutation	UNP Q96EY8
C	55	MET	-	initiating methionine	UNP Q96EY8
C	190	CYS	ARG	engineered mutation	UNP Q96EY8
F	55	MET	-	initiating methionine	UNP Q96EY8
F	190	CYS	ARG	engineered mutation	UNP Q96EY8
B	55	MET	-	initiating methionine	UNP Q96EY8
B	190	CYS	ARG	engineered mutation	UNP Q96EY8
A	55	MET	-	initiating methionine	UNP Q96EY8
A	190	CYS	ARG	engineered mutation	UNP Q96EY8
D	55	MET	-	initiating methionine	UNP Q96EY8
D	190	CYS	ARG	engineered mutation	UNP Q96EY8

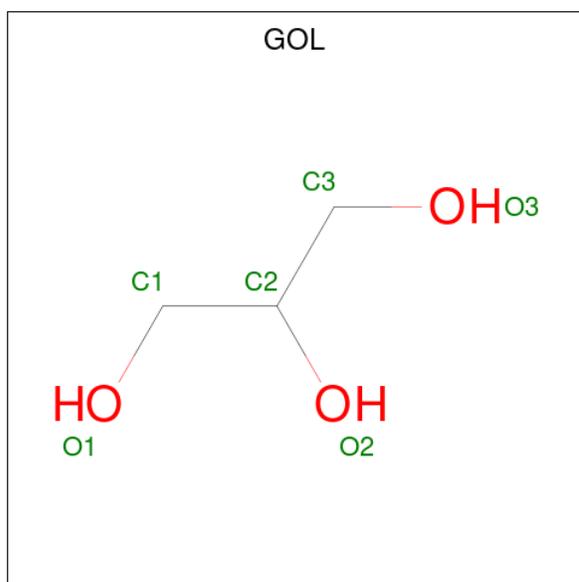
- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	E	1	Total Mg 1 1	0	0
2	C	2	Total Mg 2 2	0	0
2	F	2	Total Mg 2 2	0	0
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0
2	D	2	Total Mg 2 2	0	0

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total K 1 1	0	0
3	C	2	Total K 2 2	0	0
3	F	1	Total K 1 1	0	0
3	B	2	Total K 2 2	0	0
3	A	2	Total K 2 2	0	0
3	D	3	Total K 3 3	0	0

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	E	105	Total	O	0	0
			105	105		
6	C	109	Total	O	0	0
			109	109		
6	F	127	Total	O	0	0
			127	127		
6	B	128	Total	O	0	0
			128	128		
6	A	123	Total	O	0	0
			123	123		
6	D	125	Total	O	0	0
			125	125		

3 Residue-property plots [i](#)

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.

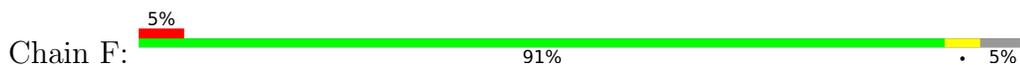
- Molecule 1: Corrinoic adenosyltransferase



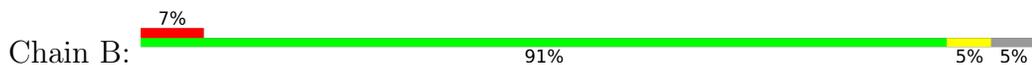
- Molecule 1: Corrinoic adenosyltransferase



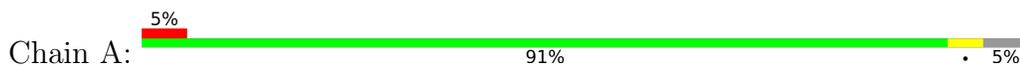
- Molecule 1: Corrinoic adenosyltransferase



- Molecule 1: Corrinoic adenosyltransferase



- Molecule 1: Corrinoic adenosyltransferase



- Molecule 1: Corrinoic adenosyltransferase

Chain D:  4% 90% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.15Å 75.77Å 95.84Å 90.00° 91.30° 90.00°	Depositor
Resolution (Å)	40.49 – 1.50 47.91 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.49-1.50) 99.9 (47.91-1.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 1.50Å)	Xtrriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.150 , 0.184 0.150 , 0.184	Depositor DCC
R_{free} test set	8845 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	20.4	Xtrriage
Anisotropy	0.314	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.000 for k,h,-l 0.001 for -k,-h,-l 0.037 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	9572	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.56% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ATP, K, MG

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	0.34	0/1486	0.48	0/2011
1	B	0.35	0/1467	0.52	0/1983
1	C	0.33	0/1459	0.47	0/1975
1	D	0.33	0/1484	0.49	0/2005
1	E	0.34	0/1453	0.48	0/1965
1	F	0.34	0/1463	0.49	0/1981
All	All	0.34	0/8812	0.49	0/11920

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1458	0	1430	5	0
1	B	1440	0	1404	4	0
1	C	1426	0	1396	5	0
1	D	1456	0	1440	4	0
1	E	1427	0	1393	2	0
1	F	1436	0	1399	3	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	1	0	0	0	0
2	F	2	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	3	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	A	31	0	12	0	0
4	B	31	0	12	0	0
4	C	31	0	12	0	0
4	D	31	0	12	0	0
4	E	31	0	12	0	0
4	F	31	0	12	0	0
5	A	6	0	8	1	0
6	A	123	0	0	0	0
6	B	128	0	0	1	0
6	C	109	0	0	0	0
6	D	125	0	0	0	0
6	E	105	0	0	0	0
6	F	127	0	0	0	0
All	All	9572	0	8542	23	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 1.

All (23) 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:C:103:VAL:HG13	1:C:108:HIS:HB2	1.72	0.71
1:A:140:LEU:HD23	1:A:143:THR:HG21	1.77	0.67
1:A:103:VAL:HG13	1:A:108:HIS:HB2	1.78	0.65
5:A:501:GOL:O3	5:A:501:GOL:O1	2.22	0.58
1:E:103:VAL:HG13	1:E:108:HIS:HB2	1.86	0.56
1:A:122:GLN:HA	1:A:122:GLN:OE1	2.07	0.55
1:A:80:ASP:OD2	1:A:82[B]:VAL:HG22	2.08	0.53
1:C:151:LEU:HG	1:C:155:GLN:HE21	1.72	0.53
1:C:72:THR:HA	1:C:139:HIS:HD2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:72:THR:HA	1:C:139:HIS:CD2	2.45	0.51
1:D:197:VAL:HG23	1:D:198:PRO:HD3	1.95	0.48
1:B:174:SER:HB2	1:B:236:LYS:NZ	2.31	0.45
1:D:103:VAL:HG13	1:D:108:HIS:HB2	1.99	0.44
1:F:103:VAL:HG13	1:F:108:HIS:HB2	2.00	0.43
1:A:140:LEU:HA	1:A:143:THR:HG23	2.00	0.43
1:B:186:ARG:HD3	1:B:221:PHE:HB2	2.01	0.42
1:F:167:LEU:HD21	1:F:225:ARG:CZ	2.50	0.42
1:C:159:LYS:O	1:C:162:SER:HB2	2.20	0.42
1:B:217[A]:SER:OG	6:B:401:HOH:O	2.20	0.42
1:D:239:MET:SD	1:D:240:LYS:N	2.93	0.41
1:F:160:TYR:O	1:F:164:LEU:HD12	2.20	0.40
1:D:199:LEU:HD23	1:D:202:MET:CE	2.51	0.40
1:E:109:THR:HA	1:B:162[A]:SER:O	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	A	189/196 (96%)	187 (99%)	2 (1%)	0	100	100
1	B	189/196 (96%)	187 (99%)	2 (1%)	0	100	100
1	C	188/196 (96%)	186 (99%)	2 (1%)	0	100	100
1	D	187/196 (95%)	185 (99%)	2 (1%)	0	100	100
1	E	184/196 (94%)	182 (99%)	2 (1%)	0	100	100
1	F	187/196 (95%)	185 (99%)	2 (1%)	0	100	100
All	All	1124/1176 (96%)	1112 (99%)	12 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	154/164 (94%)	153 (99%)	1 (1%)	86	74
1	B	150/164 (92%)	147 (98%)	3 (2%)	55	25
1	C	151/164 (92%)	150 (99%)	1 (1%)	84	69
1	D	155/164 (94%)	153 (99%)	2 (1%)	69	44
1	E	152/164 (93%)	151 (99%)	1 (1%)	84	69
1	F	151/164 (92%)	150 (99%)	1 (1%)	84	69
All	All	913/984 (93%)	904 (99%)	9 (1%)	76	57

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	215	ARG
1	C	215	ARG
1	F	215	ARG
1	B	108	HIS
1	B	215	ARG
1	B	241	ASN
1	A	215	ARG
1	D	215	ARG
1	D	239	MET

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

Mol	Chain	Res	Type
1	C	155	GLN
1	F	122	GLN
1	F	201	GLN
1	F	233	ASN
1	B	108	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 20 are monoatomic - leaving 7 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	B	304	3,2	26,33,33	1.11	3 (11%)	31,52,52	1.35	3 (9%)
4	ATP	D	306	3,2	26,33,33	1.02	2 (7%)	31,52,52	1.16	2 (6%)
4	ATP	C	305	3,2	26,33,33	1.08	3 (11%)	31,52,52	1.34	3 (9%)
4	ATP	E	303	3,2	26,33,33	1.02	2 (7%)	31,52,52	1.22	2 (6%)
5	GOL	A	501	-	5,5,5	0.91	0	5,5,5	0.97	0
4	ATP	F	304	3,2	26,33,33	1.05	2 (7%)	31,52,52	1.19	2 (6%)
4	ATP	A	505	3,2	26,33,33	1.12	3 (11%)	31,52,52	1.29	2 (6%)

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ATP	B	304	3,2	-	2/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ATP	D	306	3,2	-	3/18/38/38	0/3/3/3
4	ATP	C	305	3,2	-	4/18/38/38	0/3/3/3
4	ATP	E	303	3,2	-	4/18/38/38	0/3/3/3
5	GOL	A	501	-	-	0/4/4/4	-
4	ATP	F	304	3,2	-	3/18/38/38	0/3/3/3
4	ATP	A	505	3,2	-	4/18/38/38	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	505	ATP	O4'-C1'	3.12	1.45	1.41
4	B	304	ATP	O4'-C1'	2.73	1.44	1.41
4	F	304	ATP	O4'-C1'	2.65	1.44	1.41
4	F	304	ATP	C5-C4	2.64	1.47	1.40
4	B	304	ATP	C2-N3	2.63	1.36	1.32
4	C	305	ATP	O4'-C1'	2.61	1.44	1.41
4	D	306	ATP	C5-C4	2.58	1.47	1.40
4	E	303	ATP	O4'-C1'	2.54	1.44	1.41
4	A	505	ATP	C5-C4	2.54	1.47	1.40
4	C	305	ATP	C5-C4	2.51	1.47	1.40
4	B	304	ATP	C5-C4	2.49	1.47	1.40
4	E	303	ATP	C5-C4	2.48	1.47	1.40
4	A	505	ATP	C2-N3	2.12	1.35	1.32
4	D	306	ATP	C2-N3	2.08	1.35	1.32
4	C	305	ATP	C2-N3	2.06	1.35	1.32

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	304	ATP	N3-C2-N1	-3.32	123.49	128.68
4	C	305	ATP	N3-C2-N1	-3.25	123.60	128.68
4	A	505	ATP	N3-C2-N1	-3.13	123.78	128.68
4	D	306	ATP	N3-C2-N1	-2.96	124.06	128.68
4	E	303	ATP	N3-C2-N1	-2.88	124.17	128.68
4	F	304	ATP	N3-C2-N1	-2.75	124.38	128.68
4	F	304	ATP	N6-C6-N1	2.46	123.68	118.57
4	B	304	ATP	C4-C5-N7	-2.43	106.86	109.40
4	A	505	ATP	C4-C5-N7	-2.43	106.87	109.40
4	C	305	ATP	C2-N1-C6	2.41	122.87	118.75
4	E	303	ATP	C4-C5-N7	-2.33	106.97	109.40
4	B	304	ATP	C1'-N9-C4	-2.24	122.70	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	305	ATP	C3'-C2'-C1'	2.16	104.22	100.98
4	D	306	ATP	C4-C5-N7	-2.04	107.27	109.40

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	D	306	ATP	C5'-O5'-PA-O2A
4	C	305	ATP	PB-O3B-PG-O1G
4	B	304	ATP	PB-O3B-PG-O1G
4	A	505	ATP	PB-O3B-PG-O1G
4	E	303	ATP	PG-O3B-PB-O1B
4	C	305	ATP	PG-O3B-PB-O1B
4	F	304	ATP	PG-O3B-PB-O1B
4	E	303	ATP	C5'-O5'-PA-O2A
4	B	304	ATP	PG-O3B-PB-O1B
4	A	505	ATP	PG-O3B-PB-O1B
4	A	505	ATP	PB-O3A-PA-O2A
4	F	304	ATP	PB-O3B-PG-O2G
4	D	306	ATP	C5'-O5'-PA-O3A
4	E	303	ATP	PG-O3B-PB-O2B
4	C	305	ATP	PG-O3B-PB-O2B
4	F	304	ATP	PG-O3B-PB-O2B
4	A	505	ATP	PG-O3B-PB-O2B
4	E	303	ATP	C5'-O5'-PA-O1A
4	C	305	ATP	C5'-O5'-PA-O2A
4	D	306	ATP	C5'-O5'-PA-O1A

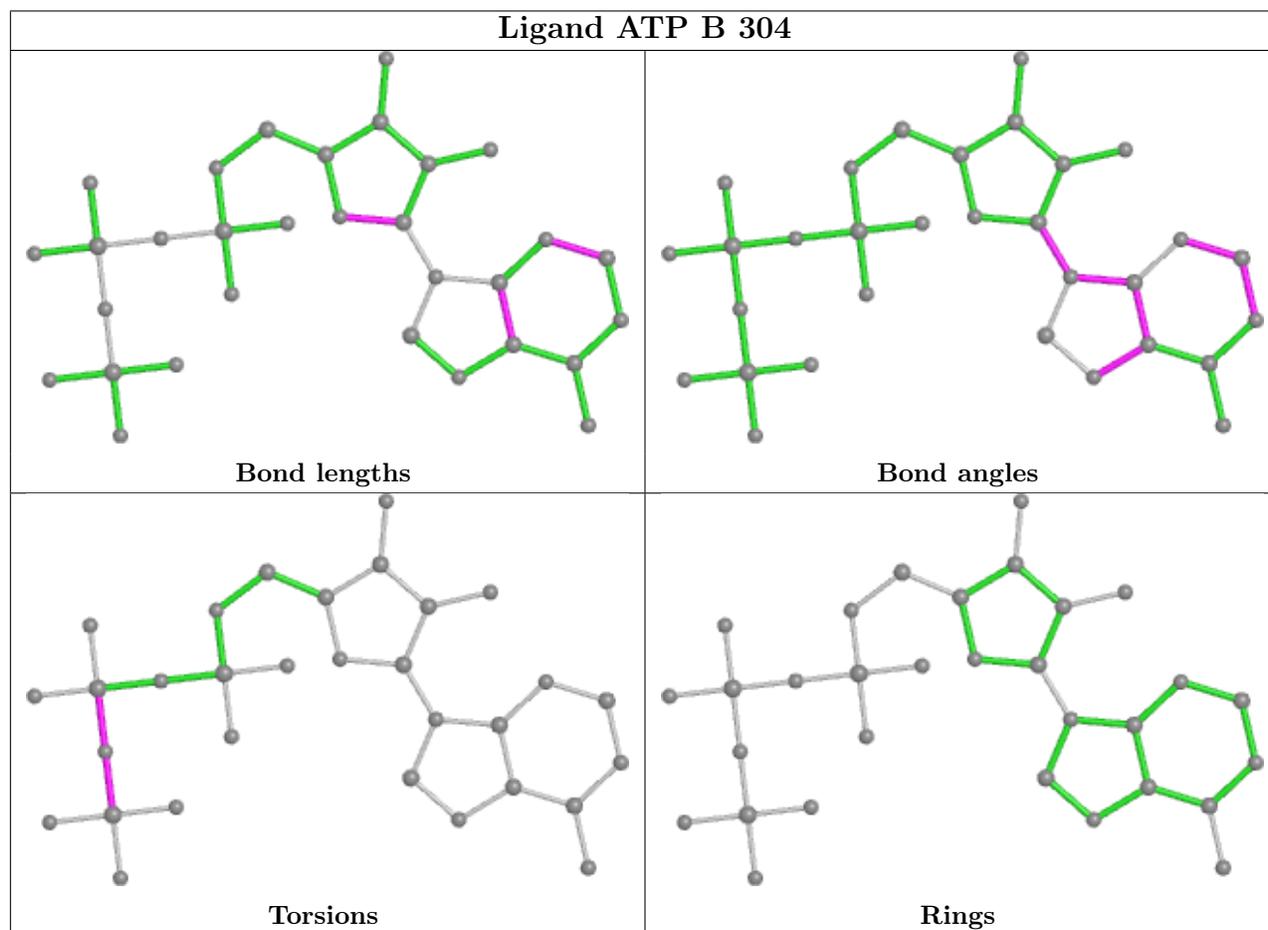
There are no ring outliers.

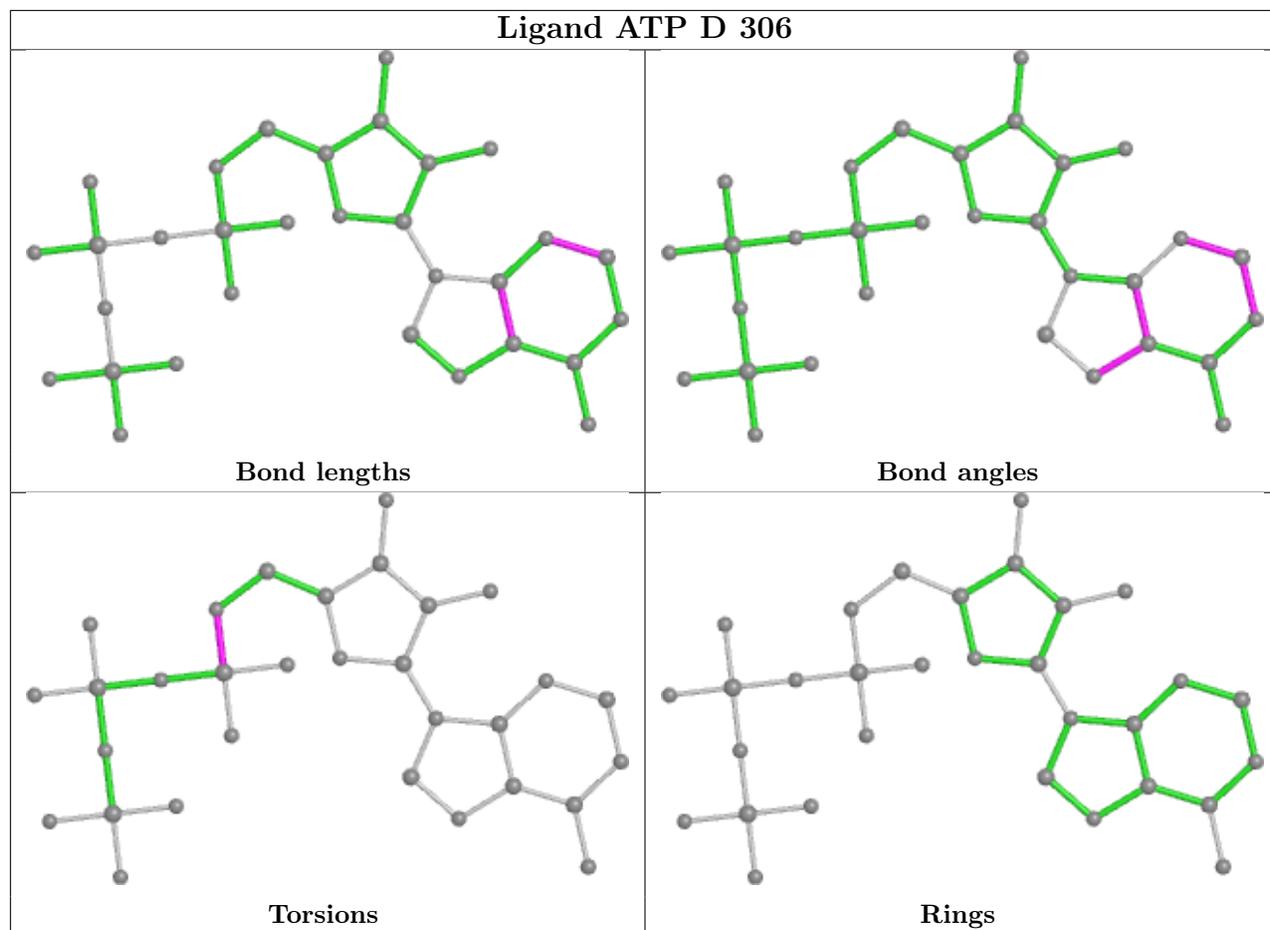
1 monomer is involved in 1 short contact:

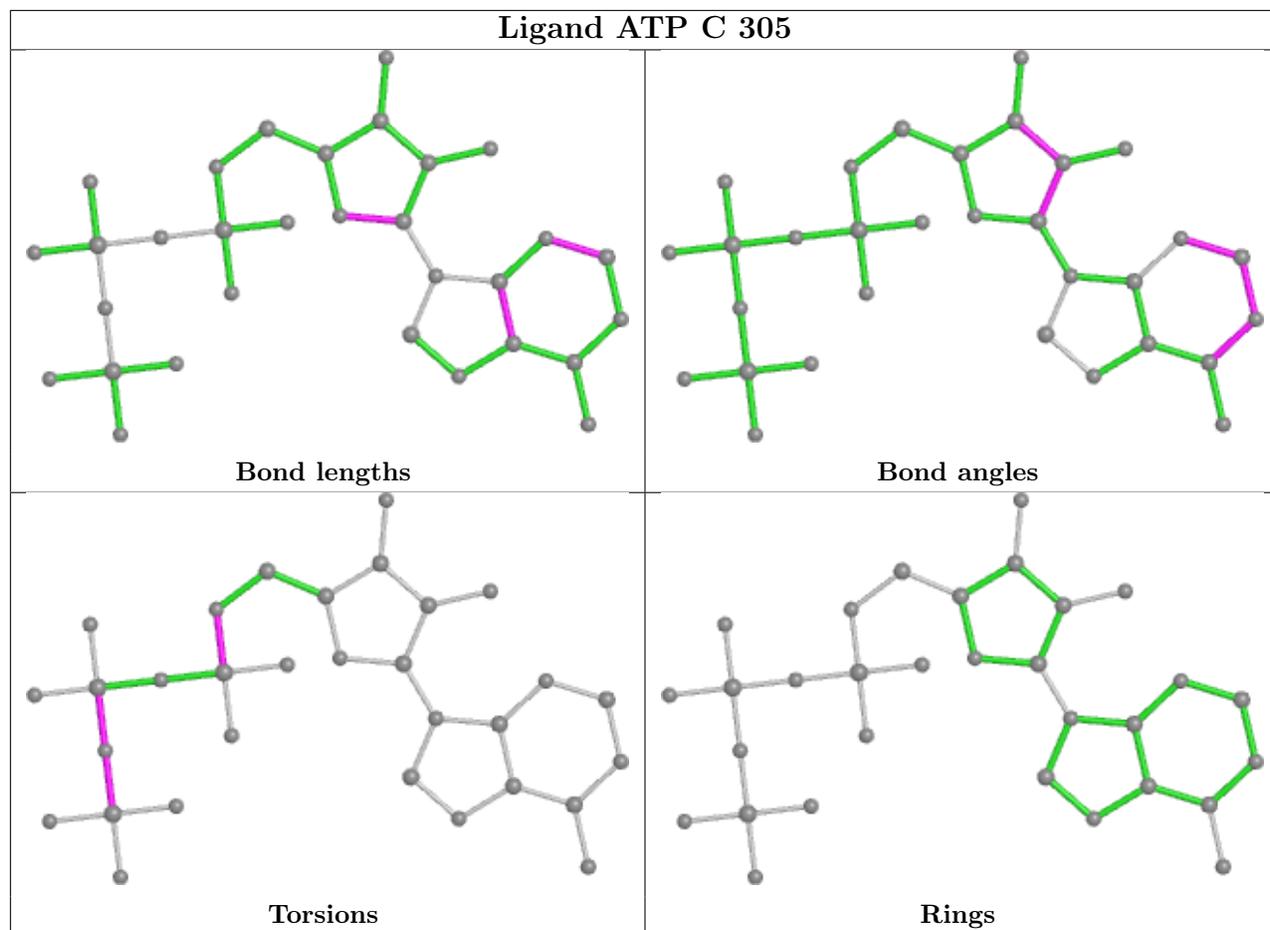
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	501	GOL	1	0

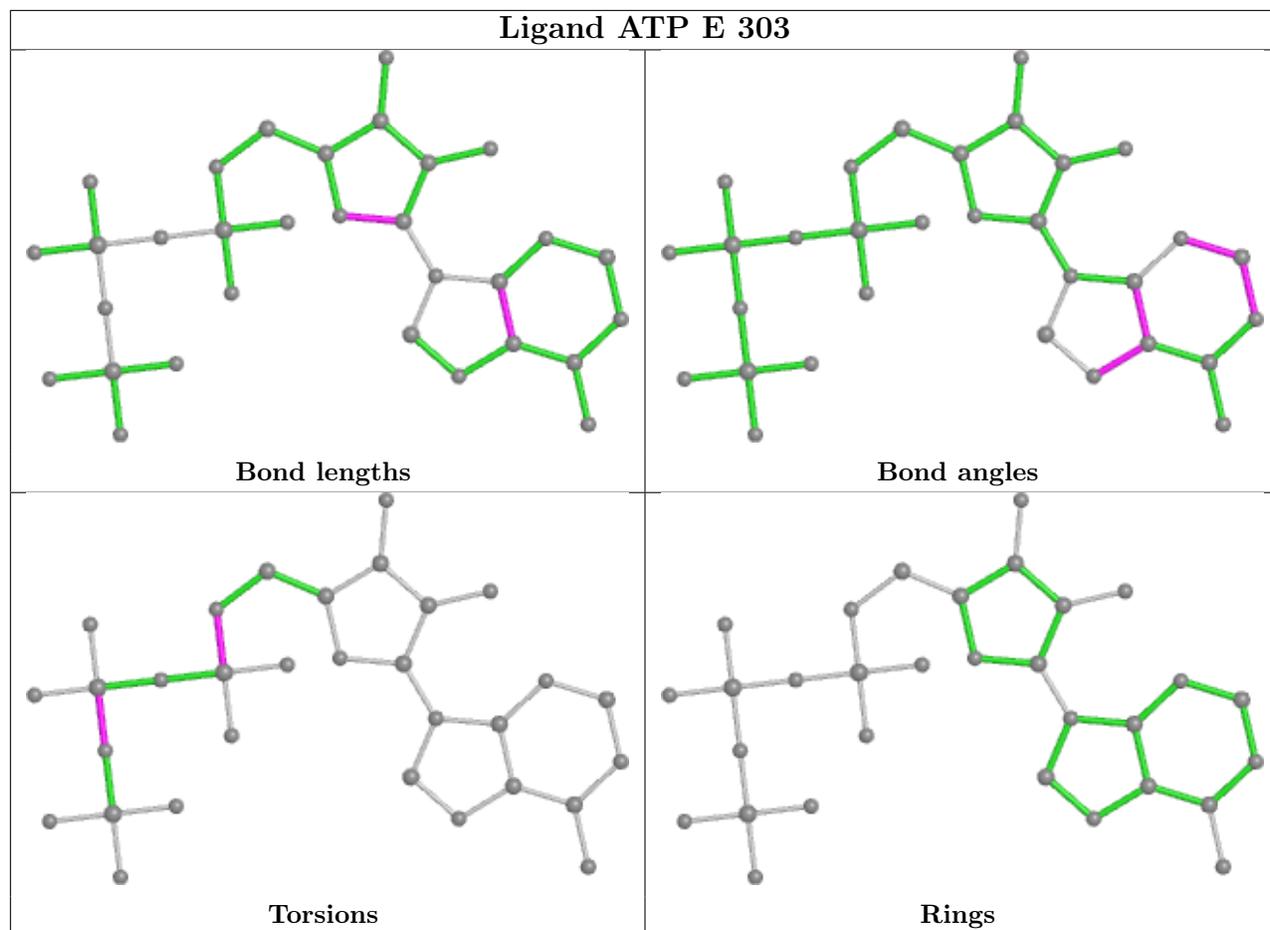
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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

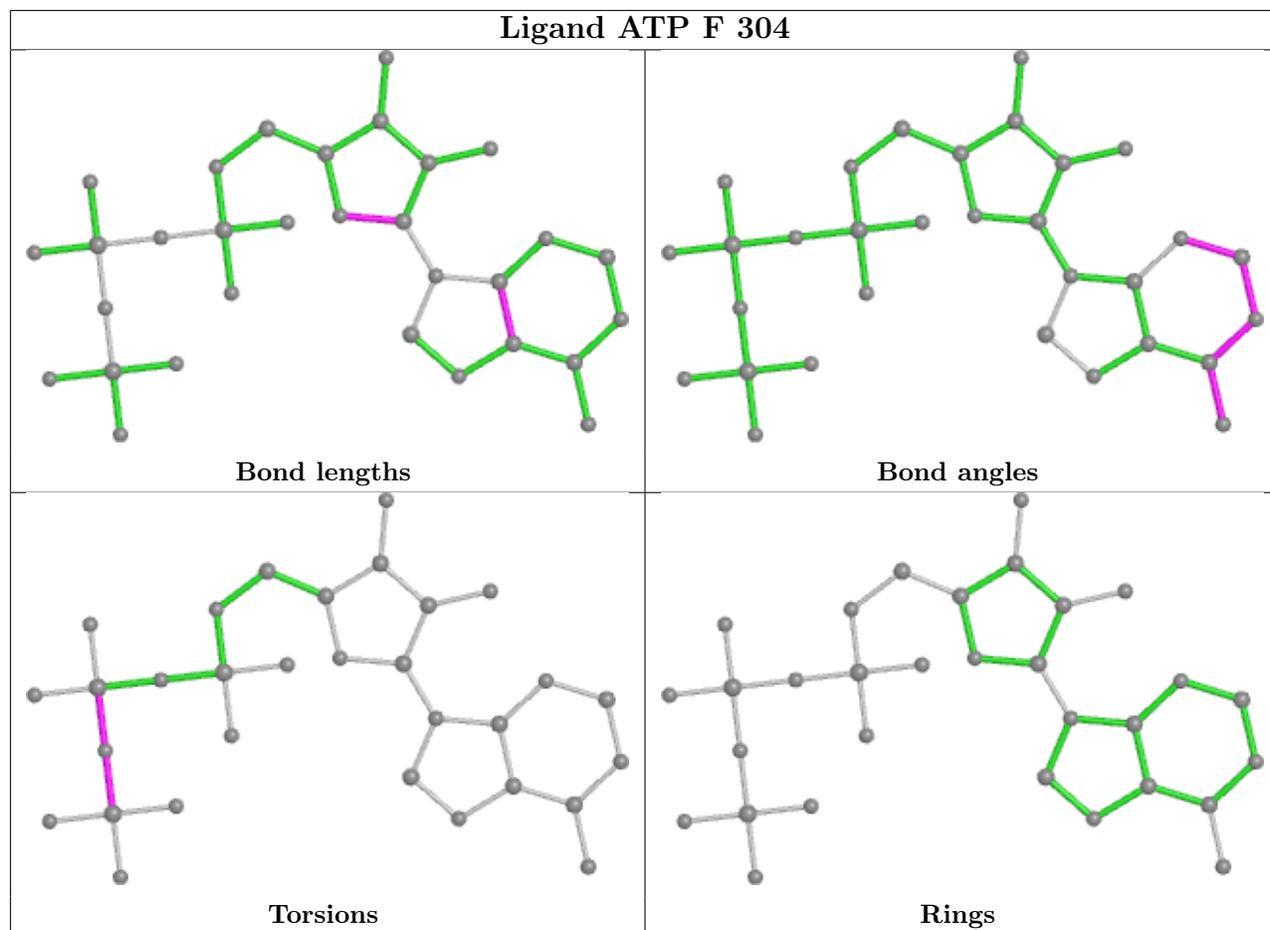
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

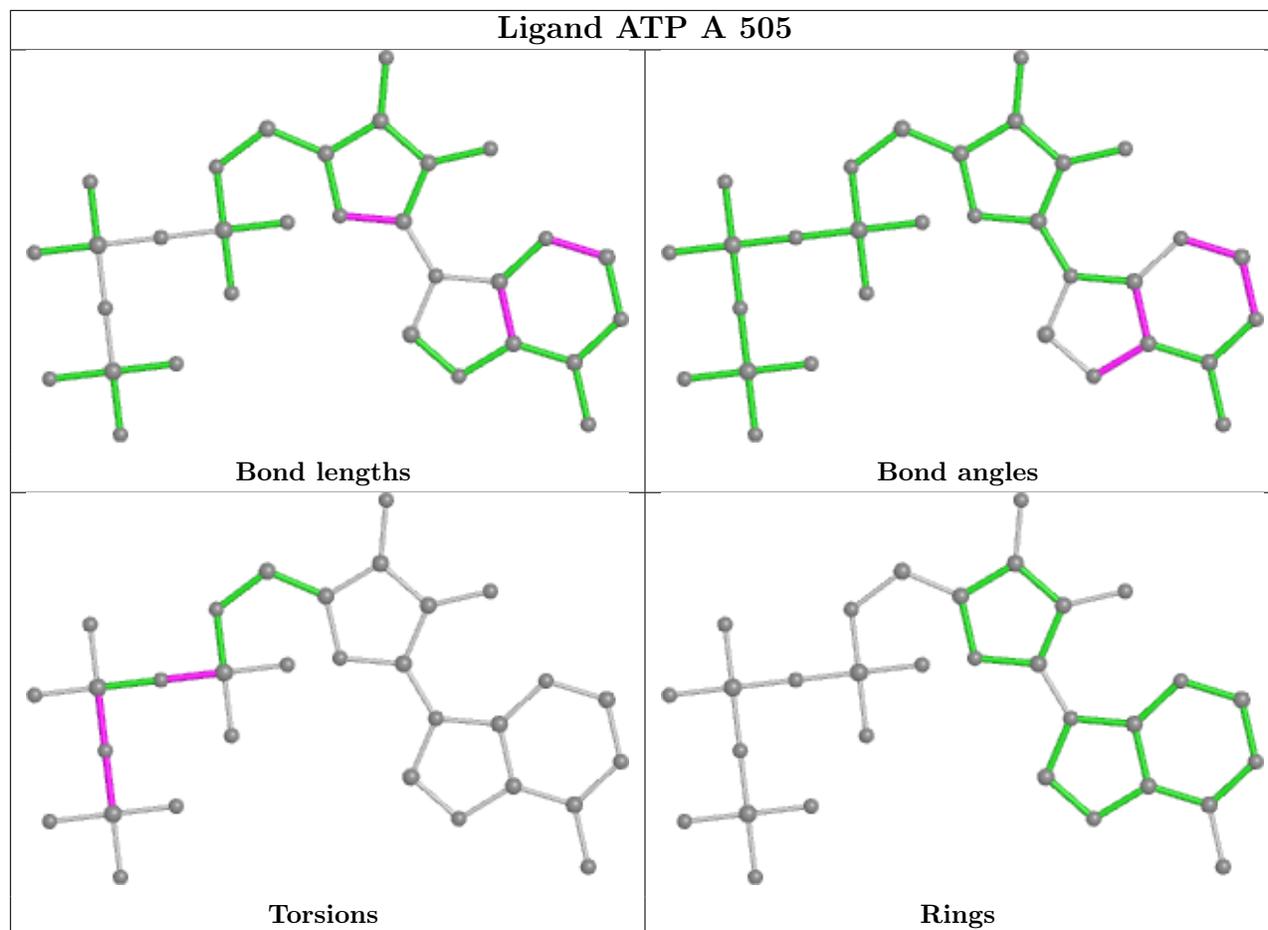












5.7 Other polymers [i](#)

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	186/196 (94%)	0.09	9 (4%) 30 33	16, 25, 61, 83	0
1	B	187/196 (95%)	0.12	14 (7%) 14 14	16, 25, 77, 142	0
1	C	186/196 (94%)	0.14	10 (5%) 25 28	16, 27, 67, 92	0
1	D	185/196 (94%)	0.10	7 (3%) 40 44	16, 25, 52, 79	0
1	E	187/196 (95%)	0.17	10 (5%) 26 29	17, 29, 61, 91	0
1	F	186/196 (94%)	0.10	9 (4%) 30 33	16, 25, 59, 87	0
All	All	1117/1176 (94%)	0.12	59 (5%) 26 29	16, 26, 64, 142	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	249	GLY	13.6
1	D	169	ALA	7.9
1	C	168	THR	6.7
1	C	107	GLY	6.5
1	E	168	THR	6.5
1	B	169	ALA	5.3
1	D	168	THR	5.1
1	B	168	THR	5.0
1	C	139	HIS	4.9
1	D	170	PHE	4.8
1	B	239	MET	4.8
1	E	132	CYS	4.8
1	E	248	GLU	4.8
1	A	168	THR	4.6
1	E	56	PRO	4.4
1	C	138	ALA	4.4
1	D	56	PRO	4.3
1	E	71	PHE	4.3
1	F	169	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
1	C	56	PRO	4.1
1	E	169	ALA	3.8
1	B	108	HIS	3.8
1	F	142	TYR	3.7
1	A	241	ASN	3.7
1	A	104	THR	3.7
1	C	241	ASN	3.6
1	F	168	THR	3.6
1	C	142	TYR	3.5
1	A	239	MET	3.4
1	C	71	PHE	3.3
1	C	169	ALA	3.3
1	B	241	ASN	3.2
1	D	239	MET	3.1
1	A	142	TYR	3.1
1	F	164	LEU	3.0
1	B	104	THR	2.8
1	B	242	ASP	2.8
1	F	237	ILE	2.8
1	B	238	TYR	2.8
1	B	234	GLN	2.7
1	B	142	TYR	2.7
1	A	56	PRO	2.6
1	E	167	LEU	2.6
1	E	245	ALA	2.5
1	A	109	THR	2.5
1	B	107	GLY	2.5
1	F	166	PRO	2.5
1	B	56	PRO	2.5
1	D	167	LEU	2.3
1	A	178	ILE	2.2
1	D	58	ILE	2.2
1	F	165	PRO	2.2
1	B	170	PHE	2.1
1	F	239	MET	2.1
1	C	170	PHE	2.1
1	F	241	ASN	2.1
1	A	240	LYS	2.1
1	E	144	THR	2.0
1	B	171	ILE	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 monosaccharides in this entry.

6.4 Ligands [i](#)

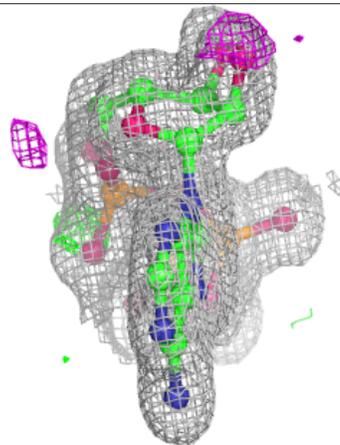
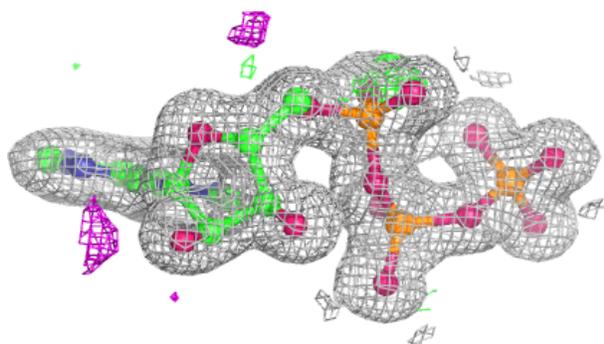
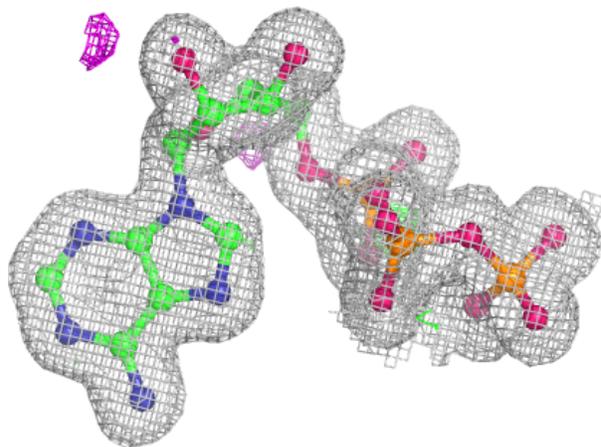
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MG	C	302	1/1	0.62	0.10	57,57,57,57	0
2	MG	F	302	1/1	0.78	0.07	50,50,50,50	0
5	GOL	A	501	6/6	0.83	0.09	66,67,67,67	0
2	MG	E	301	1/1	0.91	0.20	71,71,71,71	0
3	K	C	304	1/1	0.98	0.09	54,54,54,54	0
4	ATP	E	303	31/31	0.98	0.06	17,20,24,25	0
4	ATP	F	304	31/31	0.98	0.07	16,20,24,26	0
4	ATP	B	304	31/31	0.98	0.07	16,19,23,24	0
4	ATP	A	505	31/31	0.98	0.07	17,21,23,25	0
3	K	E	302	1/1	0.98	0.15	40,40,40,40	0
2	MG	D	302	1/1	0.99	0.05	19,19,19,19	0
4	ATP	C	305	31/31	0.99	0.06	16,19,22,23	0
4	ATP	D	306	31/31	0.99	0.06	17,21,25,27	0
3	K	D	305	1/1	0.99	0.07	48,48,48,48	0
3	K	B	303	1/1	1.00	0.04	24,24,24,24	0
3	K	A	503	1/1	1.00	0.07	26,26,26,26	0
3	K	A	504	1/1	1.00	0.10	24,24,24,24	0
3	K	D	303	1/1	1.00	0.08	25,25,25,25	0
3	K	D	304	1/1	1.00	0.05	24,24,24,24	0
2	MG	A	502	1/1	1.00	0.06	20,20,20,20	0
2	MG	D	301	1/1	1.00	0.06	18,18,18,18	0
2	MG	F	301	1/1	1.00	0.05	18,18,18,18	0
2	MG	C	301	1/1	1.00	0.05	18,18,18,18	0
3	K	C	303	1/1	1.00	0.08	25,25,25,25	0
2	MG	B	301	1/1	1.00	0.07	18,18,18,18	0
3	K	F	303	1/1	1.00	0.07	24,24,24,24	0
3	K	B	302	1/1	1.00	0.10	23,23,23,23	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.

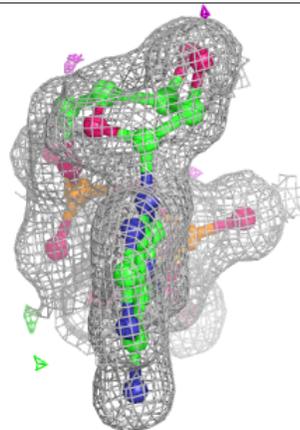
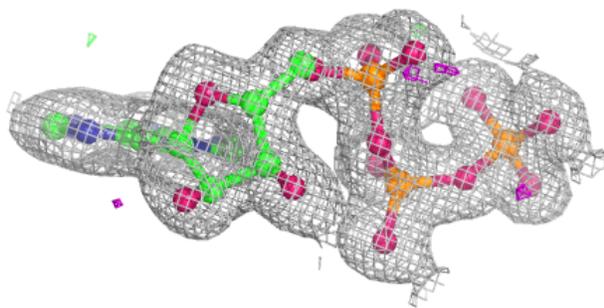
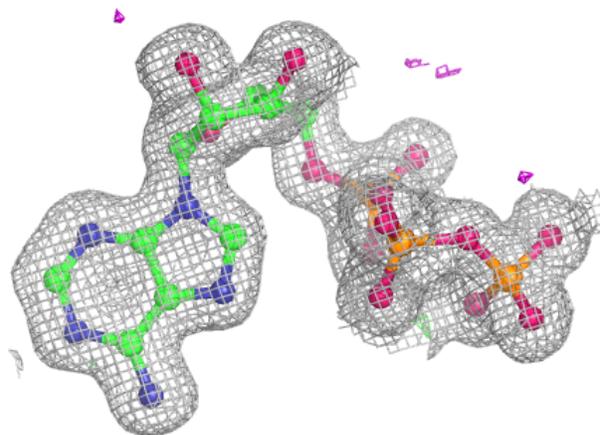
Electron density around ATP E 303:

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



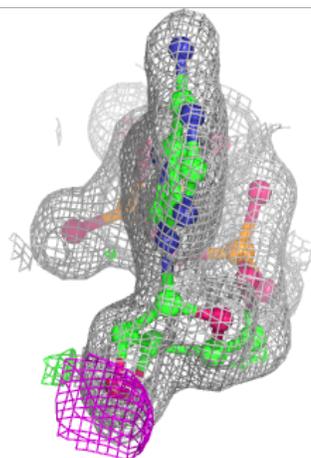
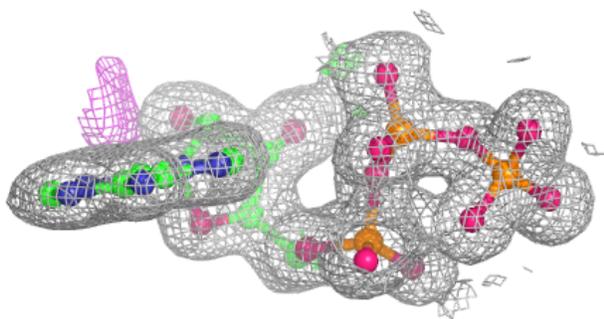
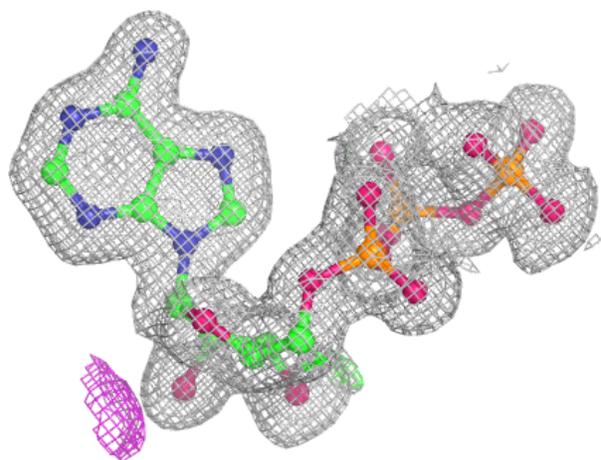
Electron density around ATP F 304:

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



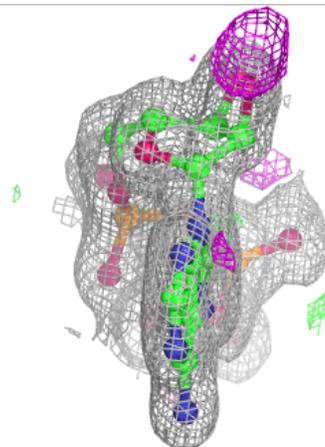
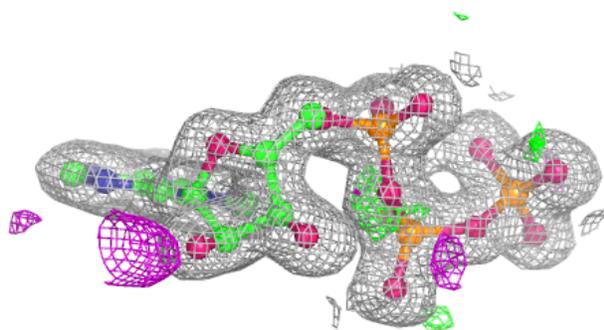
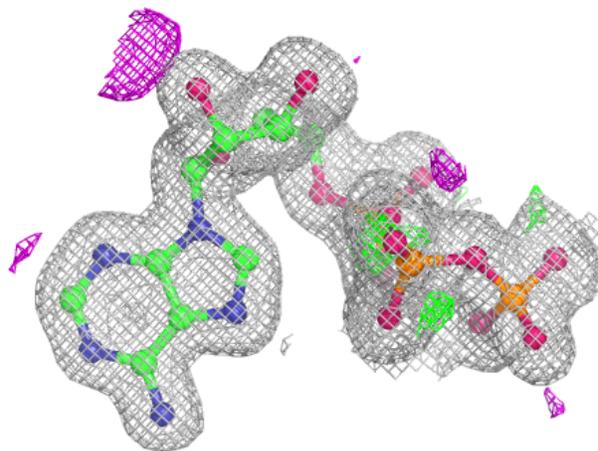
Electron density around ATP B 304:

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



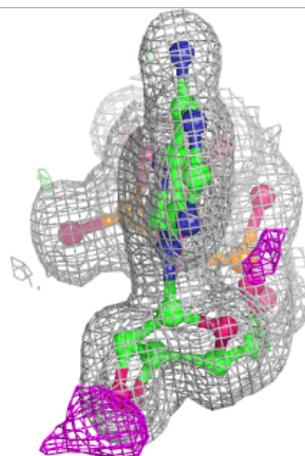
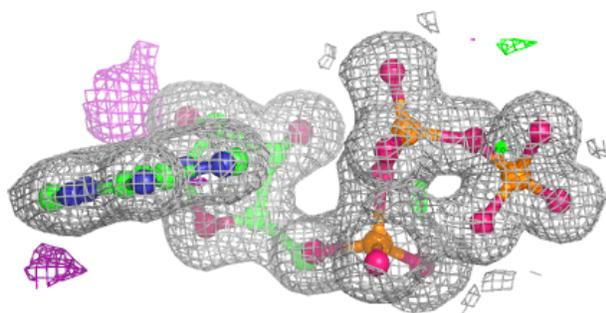
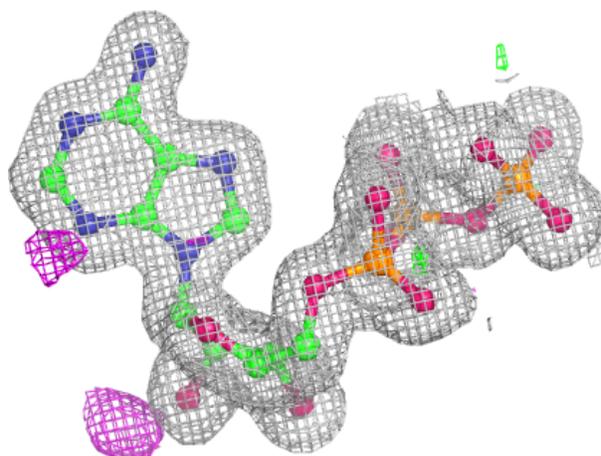
Electron density around ATP A 505:

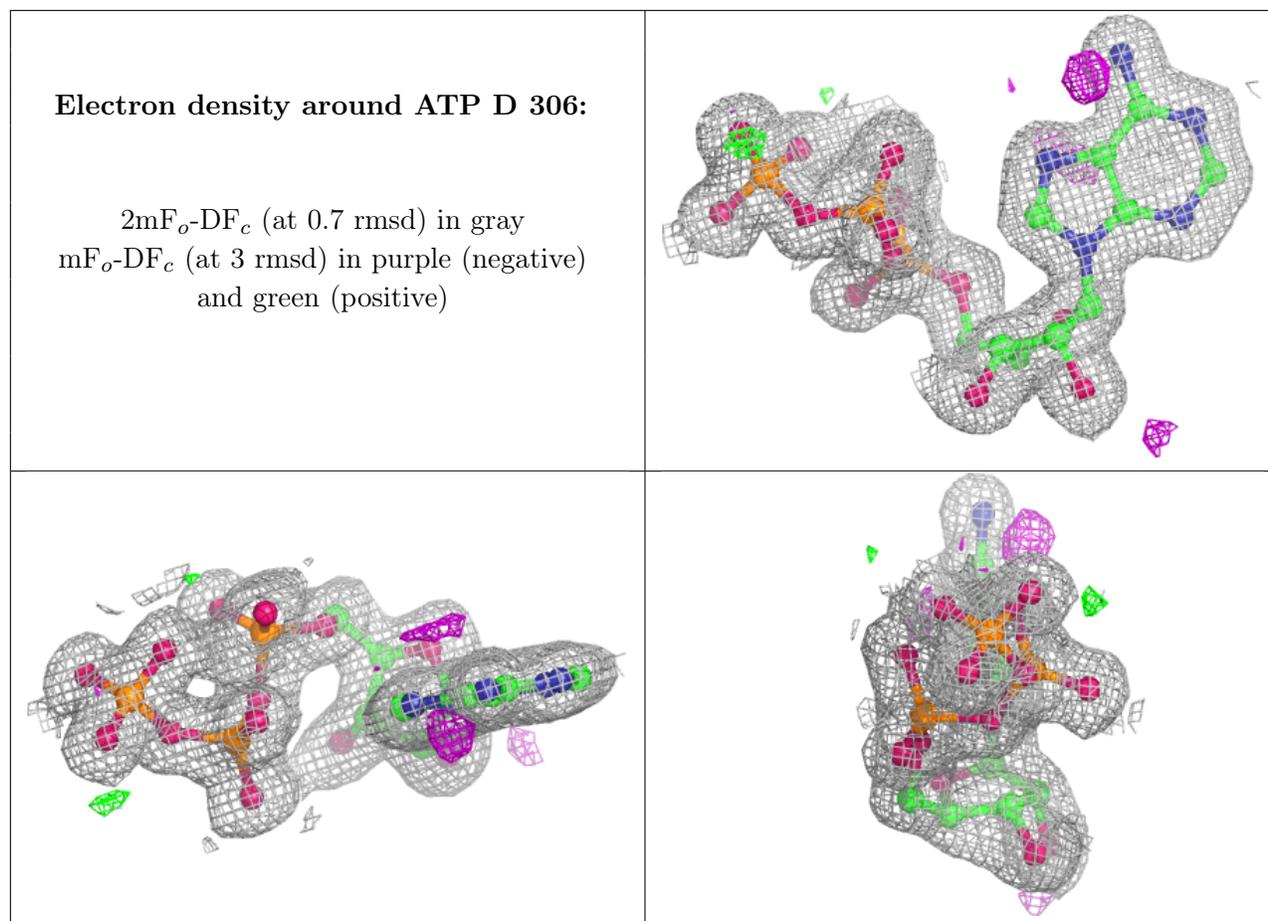
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ATP C 305:

$2mF_o-DF_c$ (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.