



# Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 08:52 am BST

PDB ID : 6AVP  
Title : Staphylococcus aureus Type II pantothenate kinase in complex with ADP and pantothenate analog Phosphate-MeO-N5Pan  
Authors : Chen, Y.; Antoshchenko, T.; Strauss, E.; Barnard, L.; Huang, Y.H.  
Deposited on : 2017-09-04  
Resolution : 2.30 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

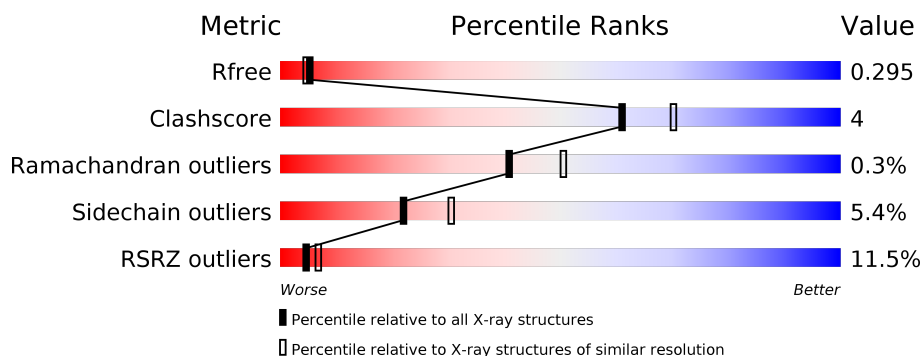
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	265	<div> <div>6%</div> <div>88%</div> <div>12%</div> <div>.</div> </div>
1	B	265	<div> <div>9%</div> <div>89%</div> <div>11%</div> </div>
1	C	265	<div> <div>11%</div> <div>86%</div> <div>11%</div> <div>..</div> </div>
1	D	265	<div> <div>19%</div> <div>92%</div> <div>5%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

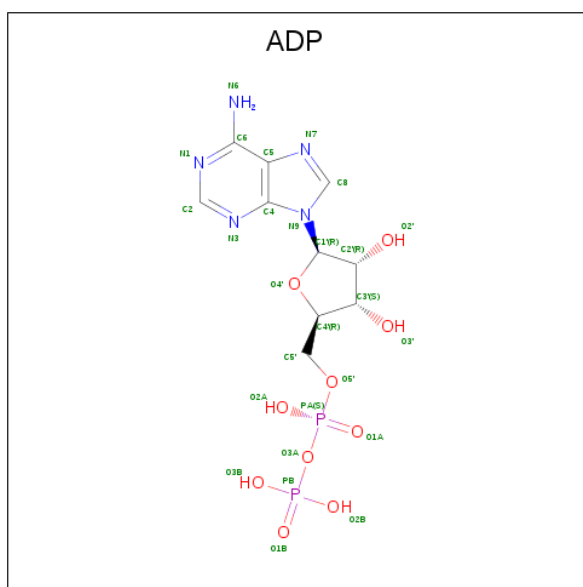
There are 5 unique types of molecules in this entry. The entry contains 8435 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Type II pantothenate kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	265	Total	C	N	O	S	0	0	0
			2033	1287	349	391	6			
1	B	265	Total	C	N	O	S	0	0	0
			2033	1287	349	391	6			
1	C	259	Total	C	N	O	S	0	0	0
			1982	1258	339	380	5			
1	D	259	Total	C	N	O	S	0	0	0
			1982	1258	339	380	5			

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula:  $C_{10}H_{15}N_5O_{10}P_2$ ).



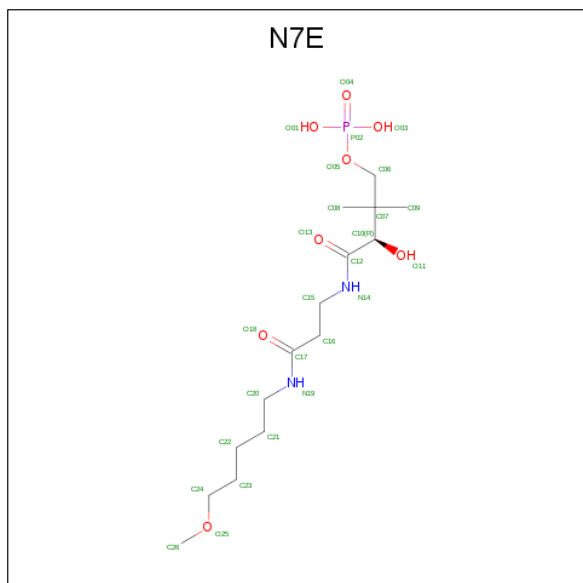
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0
			27	10	5	10	2	
2	B	1	Total	C	N	O	P	0
			27	10	5	10	2	

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is N 3 -[(2R)-2-hydroxy-3,3-dimethyl-4-(phosphonooxy)butanoyl]-N-(5-methoxy pentyl)-beta-alaninamide (three-letter code: N7E) (formula: C<sub>15</sub>H<sub>31</sub>N<sub>2</sub>O<sub>8</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			26	15	2	8	1		
3	B	1	Total	C	N	O	P	0	0
			26	15	2	8	1		
3	C	1	Total	C	N	O	P	0	0
			26	15	2	8	1		
3	D	1	Total	C	N	O	P	0	0
			26	15	2	8	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

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

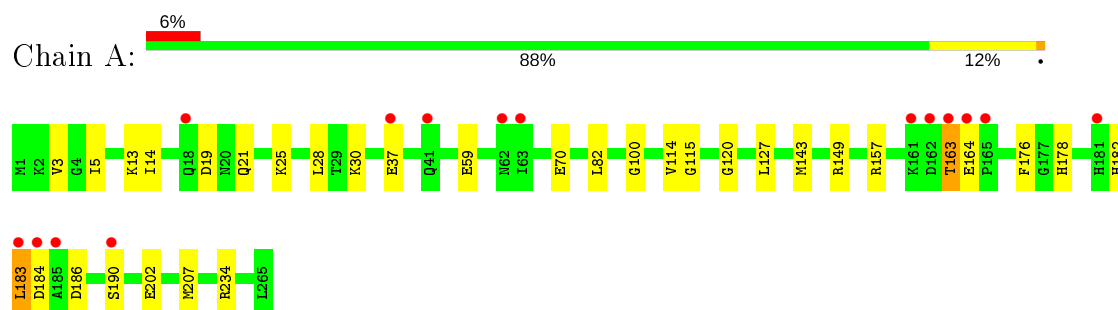
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	47	Total	O	0	0
			47	47		
5	B	46	Total	O	0	0
			46	46		
5	C	57	Total	O	0	0
			57	57		
5	D	39	Total	O	0	0
			39	39		

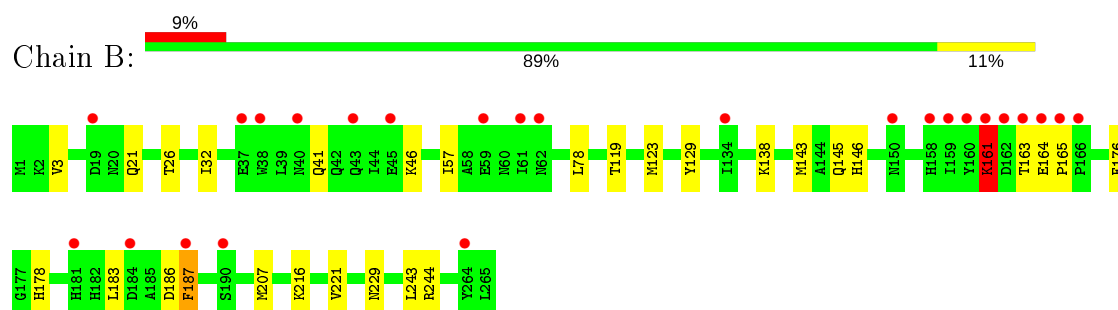
### 3 Residue-property plots [i](#)

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.

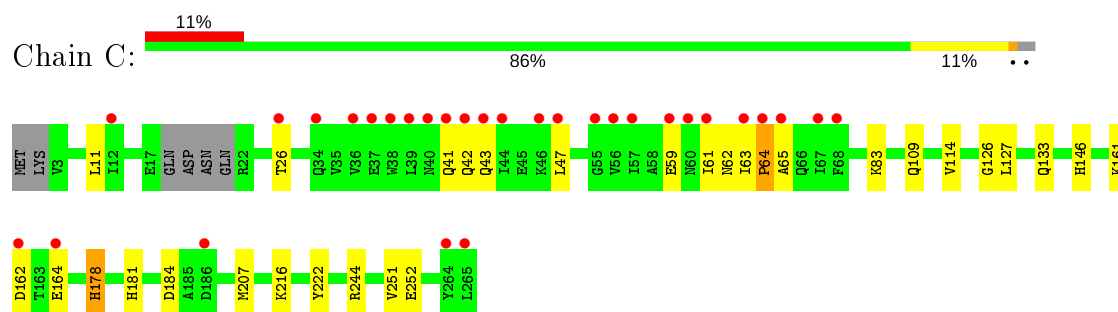
- Molecule 1: Type II pantothenate kinase



- Molecule 1: Type II pantothenate kinase

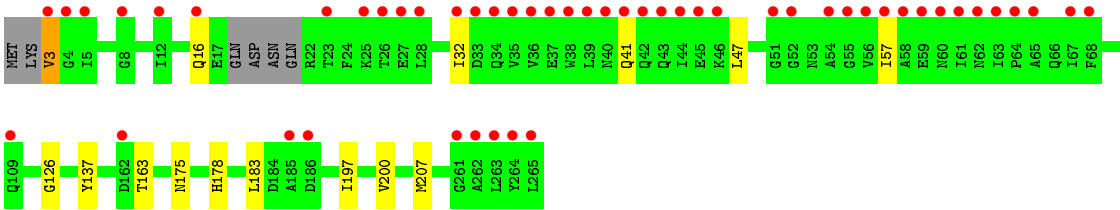


- Molecule 1: Type II pantothenate kinase



- Molecule 1: Type II pantothenate kinase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.57Å 56.70Å 133.07Å 90.00° 99.24° 90.00°	Depositor
Resolution (Å)	131.35 – 2.30 44.95 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.4 (131.35-2.30) 99.4 (44.95-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.245 , 0.295 0.247 , 0.295	Depositor DCC
$R_{free}$ test set	2410 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.8	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 38.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8435	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.71% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: N7E, MG, ADP

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	0.48	0/2068	0.65	0/2801
1	B	0.56	0/2068	0.69	0/2801
1	C	0.56	0/2016	0.69	0/2731
1	D	0.46	0/2016	0.64	0/2731
All	All	0.52	0/8168	0.67	0/11064

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	2033	0	2021	23	0
1	B	2033	0	2021	25	0
1	C	1982	0	1969	23	0
1	D	1982	0	1969	8	0
2	A	27	0	12	0	0
2	B	27	0	12	0	0
2	C	27	0	12	0	0
2	D	27	0	12	1	0
3	A	26	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	26	0	0	0	0
3	C	26	0	0	0	0
3	D	26	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	47	0	0	18	0
5	B	46	0	0	11	0
5	C	57	0	0	12	0
5	D	39	0	0	0	0
All	All	8435	0	8028	73	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 4.

All (73) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:LEU:HD23	5:A:433:HOH:O	1.47	1.12
1:A:182:HIS:CD2	5:A:405:HOH:O	2.00	1.11
1:C:222:TYR:O	5:C:401:HOH:O	1.77	1.02
1:B:146:HIS:N	5:B:401:HOH:O	1.68	0.95
1:B:145:GLN:CA	5:B:401:HOH:O	2.15	0.94
1:C:63:ILE:CG1	1:C:64:PRO:HD2	1.98	0.94
1:C:63:ILE:HG13	1:C:64:PRO:HD2	1.49	0.93
1:C:164:GLU:OE1	5:C:402:HOH:O	1.94	0.85
1:B:123:MET:HE2	1:D:175:ASN:HB2	1.57	0.85
1:B:145:GLN:HB2	5:B:401:HOH:O	1.77	0.84
1:B:145:GLN:N	5:B:401:HOH:O	2.13	0.81
1:C:252:GLU:OE2	5:C:403:HOH:O	1.98	0.80
1:A:120:GLY:HA3	5:A:415:HOH:O	1.84	0.75
1:C:63:ILE:HG13	1:C:64:PRO:CD	2.16	0.75
1:C:164:GLU:CD	5:C:404:HOH:O	2.29	0.72
1:B:163:THR:HG22	1:B:164:GLU:N	2.07	0.69
1:C:63:ILE:HG12	1:C:64:PRO:HD2	1.75	0.67
1:A:100:GLY:HA2	5:A:415:HOH:O	1.94	0.66
1:A:114:VAL:C	5:A:403:HOH:O	2.33	0.66
1:A:115:GLY:N	5:A:403:HOH:O	2.28	0.66
1:C:164:GLU:OE1	5:C:404:HOH:O	2.12	0.66
1:A:182:HIS:HD2	5:A:405:HOH:O	1.55	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:GLU:OE2	5:A:402:HOH:O	2.17	0.60
1:C:244:ARG:NH2	5:C:405:HOH:O	2.18	0.60
1:B:163:THR:CG2	1:B:164:GLU:N	2.65	0.59
1:B:145:GLN:CB	5:B:401:HOH:O	2.33	0.58
5:A:422:HOH:O	1:C:114:VAL:HA	2.03	0.58
1:C:164:GLU:HB3	5:C:404:HOH:O	2.04	0.56
1:B:119:THR:HA	1:B:123:MET:HE3	1.88	0.56
1:C:178:HIS:HA	5:C:406:HOH:O	2.05	0.55
1:C:244:ARG:NH1	5:C:405:HOH:O	2.33	0.54
1:C:181:HIS:CD2	5:C:406:HOH:O	2.61	0.53
1:B:163:THR:O	1:B:165:PRO:HD3	2.09	0.53
1:A:100:GLY:N	5:A:401:HOH:O	1.88	0.52
1:A:207:MET:SD	1:C:207:MET:SD	3.08	0.51
1:B:145:GLN:HG2	5:B:430:HOH:O	2.11	0.51
1:B:207:MET:CE	5:B:440:HOH:O	2.60	0.50
1:C:146:HIS:HD2	5:C:449:HOH:O	1.94	0.50
1:A:5:ILE:HG12	1:A:14:ILE:HG12	1.93	0.49
1:A:157:ARG:NH1	5:A:404:HOH:O	2.42	0.49
1:A:30:LYS:HD3	5:A:425:HOH:O	2.12	0.48
1:A:183:LEU:O	1:A:184:ASP:HB2	2.14	0.48
1:A:176:PHE:O	1:C:126:GLY:HA3	2.13	0.48
1:C:63:ILE:CG1	1:C:64:PRO:CD	2.78	0.47
1:D:3:VAL:HG23	1:D:16:GLN:HG2	1.96	0.47
1:B:138:LYS:HG3	5:B:429:HOH:O	2.13	0.47
1:C:181:HIS:CG	5:C:406:HOH:O	2.68	0.47
1:B:207:MET:HE2	5:B:440:HOH:O	2.15	0.46
1:D:137:TYR:CZ	2:D:301:ADP:C8	3.04	0.46
1:B:186:ASP:O	1:B:187:PHE:HB2	2.16	0.46
1:C:47:LEU:HD23	1:C:65:ALA:HB2	1.96	0.46
1:B:129:TYR:C	1:B:129:TYR:CD1	2.87	0.46
1:D:32:ILE:HD11	1:D:57:ILE:HG13	1.98	0.45
1:B:207:MET:SD	1:D:207:MET:SD	3.15	0.45
1:A:100:GLY:CA	5:A:415:HOH:O	2.60	0.44
1:D:3:VAL:CG1	1:D:47:LEU:HD13	2.48	0.44
1:B:78:LEU:HD11	1:B:221:VAL:HG11	2.00	0.43
1:B:32:ILE:HD11	1:B:57:ILE:HG13	2.00	0.43
1:B:229:ASN:CG	5:B:402:HOH:O	2.57	0.43
1:B:145:GLN:CG	5:B:430:HOH:O	2.67	0.42
1:A:190:SER:HB2	5:A:434:HOH:O	2.18	0.42
1:C:61:ILE:CG2	1:C:62:ASN:N	2.82	0.42
1:B:161:LYS:HD2	1:B:161:LYS:HA	1.72	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:THR:HG21	5:A:442:HOH:O	2.20	0.42
1:D:197:ILE:HA	1:D:200:VAL:HG12	2.02	0.42
1:B:186:ASP:OD1	1:B:186:ASP:N	2.50	0.41
1:B:176:PHE:O	1:D:126:GLY:HA3	2.20	0.41
1:B:164:GLU:HA	1:B:165:PRO:HD2	1.75	0.41
1:A:120:GLY:CA	5:A:415:HOH:O	2.58	0.41
1:A:30:LYS:HE3	5:A:414:HOH:O	2.21	0.41
1:A:127:LEU:HD11	1:C:127:LEU:HD11	2.03	0.40
1:A:143:MET:HG2	5:A:434:HOH:O	2.21	0.40
1:A:149:ARG:NH2	1:A:202:GLU:OE2	2.51	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	263/265 (99%)	250 (95%)	13 (5%)	0	100	100
1	B	263/265 (99%)	248 (94%)	13 (5%)	2 (1%)	19	23
1	C	255/265 (96%)	248 (97%)	6 (2%)	1 (0%)	34	42
1	D	255/265 (96%)	245 (96%)	10 (4%)	0	100	100
All	All	1036/1060 (98%)	991 (96%)	42 (4%)	3 (0%)	41	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	187	PHE
1	B	161	LYS
1	C	64	PRO

### 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	215/215 (100%)	201 (94%)	14 (6%)	17	23
1	B	215/215 (100%)	203 (94%)	12 (6%)	21	29
1	C	209/215 (97%)	194 (93%)	15 (7%)	14	18
1	D	209/215 (97%)	204 (98%)	5 (2%)	49	66
All	All	848/860 (99%)	802 (95%)	46 (5%)	22	30

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	VAL
1	A	13	LYS
1	A	19	ASP
1	A	21	GLN
1	A	25	LYS
1	A	28	LEU
1	A	37	GLU
1	A	59	GLU
1	A	163	THR
1	A	164	GLU
1	A	178	HIS
1	A	183	LEU
1	A	186	ASP
1	A	234	ARG
1	B	3	VAL
1	B	21	GLN
1	B	26	THR
1	B	41	GLN
1	B	46	LYS
1	B	143	MET
1	B	161	LYS
1	B	178	HIS
1	B	183	LEU
1	B	216	LYS

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Mol	Chain	Res	Type
1	B	243	LEU
1	B	244	ARG
1	C	11	LEU
1	C	26	THR
1	C	41	GLN
1	C	42	GLN
1	C	43	GLN
1	C	59	GLU
1	C	83	LYS
1	C	109	GLN
1	C	133	GLN
1	C	161	LYS
1	C	162	ASP
1	C	178	HIS
1	C	184	ASP
1	C	216	LYS
1	C	251	VAL
1	D	3	VAL
1	D	41	GLN
1	D	163	THR
1	D	178	HIS
1	D	183	LEU

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

Mol	Chain	Res	Type
1	A	178	HIS
1	A	182	HIS
1	A	253	ASN
1	B	178	HIS
1	B	182	HIS
1	C	43	GLN
1	C	133	GLN
1	C	146	HIS
1	C	178	HIS
1	C	181	HIS
1	D	146	HIS

### 5.3.3 RNA ⓘ

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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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	N7E	C	302	4	22,25,25	0.87	0	31,33,33	1.65	4 (12%)
3	N7E	A	302	4	22,25,25	0.73	0	31,33,33	1.65	7 (22%)
2	ADP	D	301	4	24,29,29	1.08	2 (8%)	29,45,45	1.20	3 (10%)
2	ADP	A	301	4	24,29,29	1.02	2 (8%)	29,45,45	1.44	4 (13%)
3	N7E	D	302	4	22,25,25	0.87	0	31,33,33	1.53	4 (12%)
3	N7E	B	302	4	22,25,25	0.87	0	31,33,33	1.62	7 (22%)
2	ADP	C	301	4	24,29,29	1.15	2 (8%)	29,45,45	1.39	5 (17%)
2	ADP	B	301	4	24,29,29	1.06	2 (8%)	29,45,45	1.35	4 (13%)

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	N7E	C	302	4	-	3/31/31/31	-
3	N7E	A	302	4	-	1/31/31/31	-
2	ADP	D	301	4	-	0/12/32/32	0/3/3/3
2	ADP	A	301	4	-	6/12/32/32	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	N7E	D	302	4	-	2/31/31/31	-
3	N7E	B	302	4	-	2/31/31/31	-
2	ADP	C	301	4	-	0/12/32/32	0/3/3/3
2	ADP	B	301	4	-	1/12/32/32	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	ADP	O4'-C1'	3.00	1.45	1.41
2	D	301	ADP	C5-C4	2.75	1.48	1.40
2	C	301	ADP	C5-C4	2.63	1.47	1.40
2	A	301	ADP	O4'-C1'	2.59	1.44	1.41
2	B	301	ADP	C5-C4	2.30	1.47	1.40
2	A	301	ADP	C5-C4	2.23	1.46	1.40
2	D	301	ADP	C2-N3	2.22	1.35	1.32
2	C	301	ADP	C2-N3	2.14	1.35	1.32

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	302	N7E	O03-P02-O05	-5.37	92.45	106.73
3	C	302	N7E	O03-P02-O05	-4.66	94.35	106.73
3	C	302	N7E	O01-P02-O05	-4.28	95.35	106.73
2	A	301	ADP	N3-C2-N1	-4.14	122.21	128.68
3	A	302	N7E	O03-P02-O05	-3.87	96.43	106.73
3	A	302	N7E	C15-C16-C17	-3.68	106.22	112.36
2	B	301	ADP	N3-C2-N1	-3.57	123.10	128.68
3	B	302	N7E	O05-P02-O04	-3.55	96.52	106.47
2	C	301	ADP	N3-C2-N1	-3.28	123.55	128.68
3	B	302	N7E	C15-C16-C17	-3.27	106.92	112.36
3	B	302	N7E	O01-P02-O05	-3.20	98.22	106.73
3	B	302	N7E	O03-P02-O05	-3.10	98.48	106.73
3	C	302	N7E	O03-P02-O01	3.10	119.48	107.64
3	A	302	N7E	O03-P02-O01	3.09	119.46	107.64
3	D	302	N7E	O03-P02-O01	3.05	119.29	107.64
2	D	301	ADP	N3-C2-N1	-3.02	123.96	128.68
2	D	301	ADP	C4-C5-N7	-2.88	106.40	109.40
3	A	302	N7E	O05-P02-O04	-2.81	98.60	106.47
2	C	301	ADP	C3'-C2'-C1'	2.78	105.17	100.98
3	A	302	N7E	O01-P02-O05	-2.77	99.37	106.73
3	B	302	N7E	O03-P02-O01	2.74	118.10	107.64
2	C	301	ADP	O4'-C1'-C2'	-2.72	102.95	106.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	ADP	C4-C5-N7	-2.67	106.61	109.40
3	D	302	N7E	O01-P02-O05	-2.64	99.71	106.73
2	C	301	ADP	C4-C5-N7	-2.59	106.70	109.40
3	B	302	N7E	O11-C10-C07	-2.51	104.34	110.25
3	C	302	N7E	C23-C22-C21	-2.48	101.82	114.42
2	A	301	ADP	C4-C5-N7	-2.43	106.87	109.40
2	D	301	ADP	O3B-PB-O2B	2.42	116.90	107.64
2	A	301	ADP	PA-O3A-PB	-2.40	124.58	132.83
2	B	301	ADP	O3B-PB-O2B	2.36	116.67	107.64
3	B	302	N7E	O01-P02-O04	2.36	119.92	110.68
2	B	301	ADP	PA-O3A-PB	-2.26	125.07	132.83
2	A	301	ADP	C2-N1-C6	2.17	122.47	118.75
2	C	301	ADP	C1'-N9-C4	2.16	130.44	126.64
3	A	302	N7E	C20-N19-C17	-2.13	118.88	122.84
3	A	302	N7E	O01-P02-O04	2.04	118.68	110.68
3	D	302	N7E	O05-P02-O04	-2.03	100.77	106.47

There are no chirality outliers.

All (15) torsion outliers are listed below:

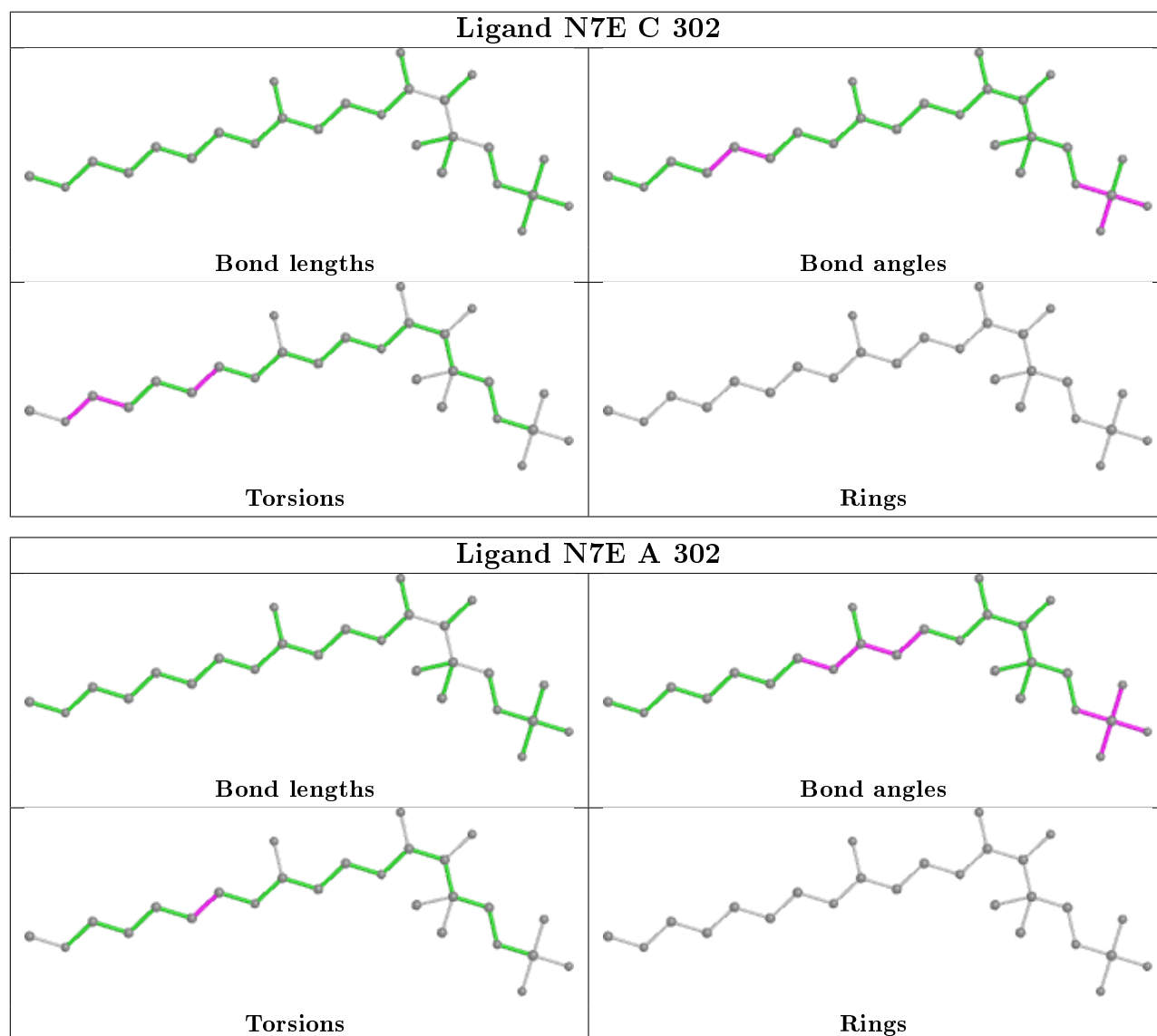
Mol	Chain	Res	Type	Atoms
2	A	301	ADP	C5'-O5'-PA-O1A
2	A	301	ADP	C5'-O5'-PA-O2A
2	A	301	ADP	C5'-O5'-PA-O3A
3	A	302	N7E	N19-C20-C21-C22
3	B	302	N7E	C22-C23-C24-O25
3	C	302	N7E	N19-C20-C21-C22
3	C	302	N7E	C22-C23-C24-O25
2	A	301	ADP	O4'-C4'-C5'-O5'
2	A	301	ADP	C3'-C4'-C5'-O5'
3	B	302	N7E	C20-C21-C22-C23
3	D	302	N7E	C20-C21-C22-C23
2	A	301	ADP	PB-O3A-PA-O2A
2	B	301	ADP	PB-O3A-PA-O1A
3	D	302	N7E	C22-C23-C24-O25
3	C	302	N7E	C23-C24-O25-C26

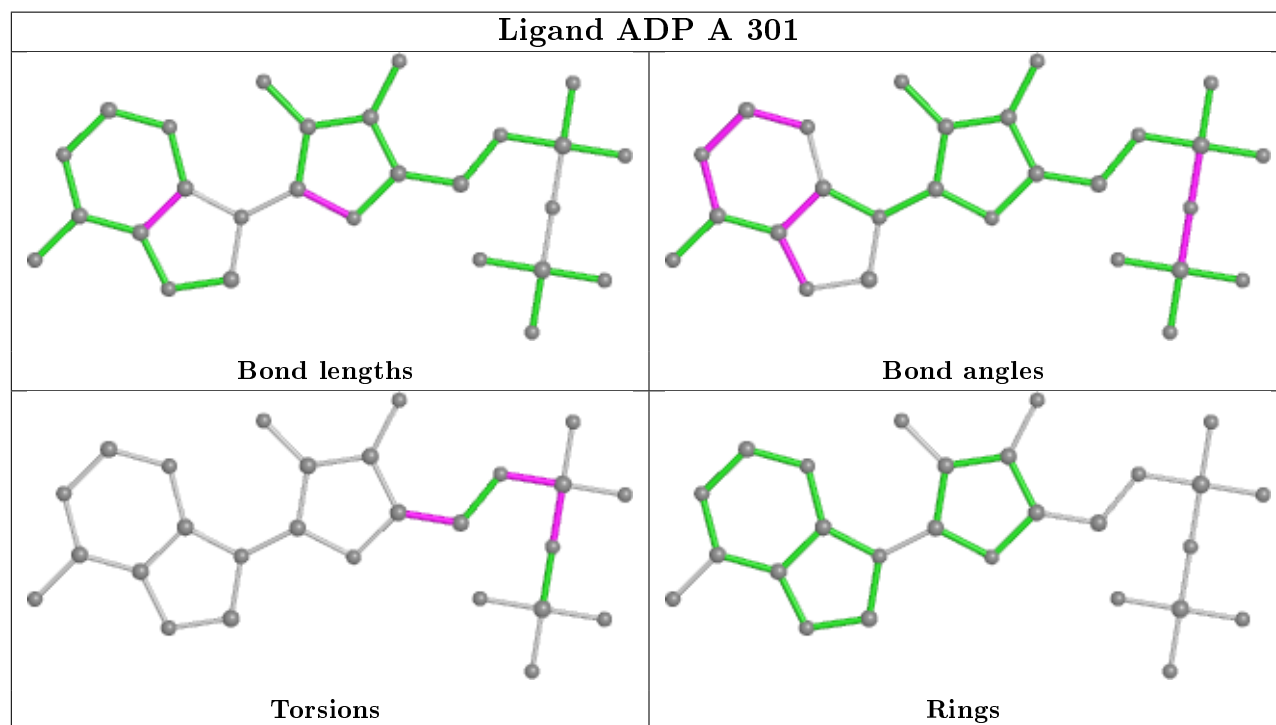
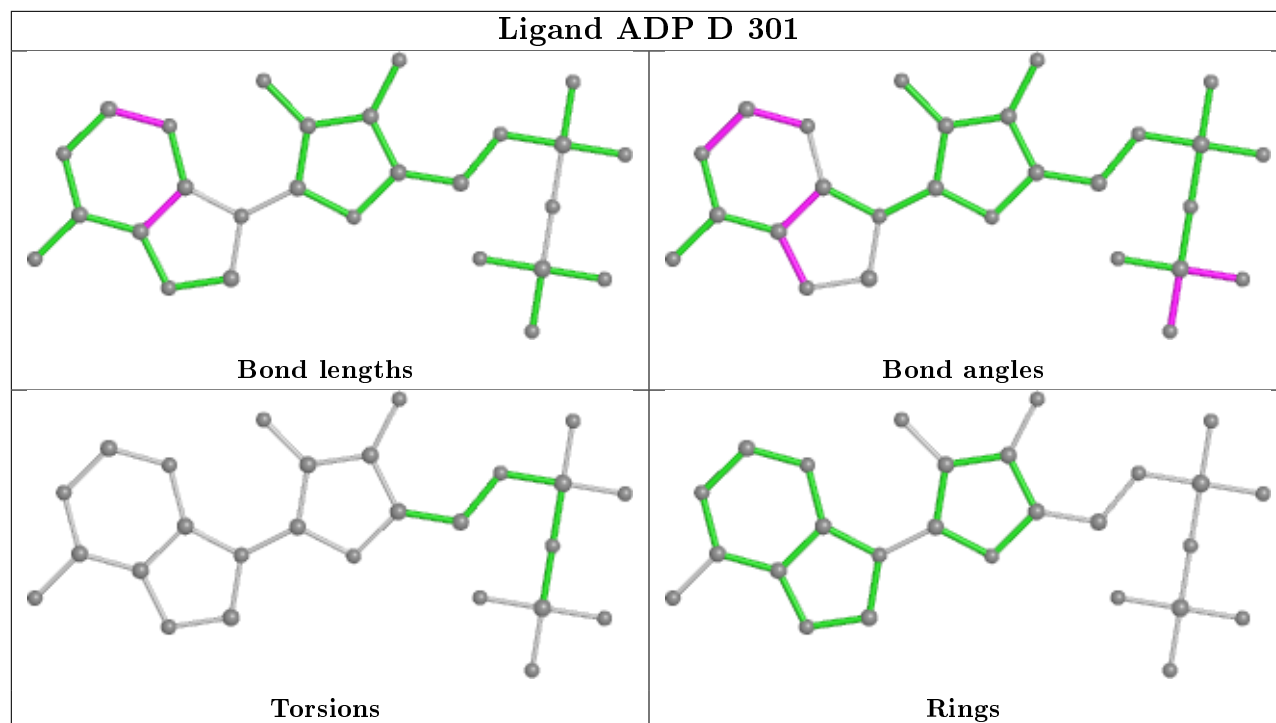
There are no ring outliers.

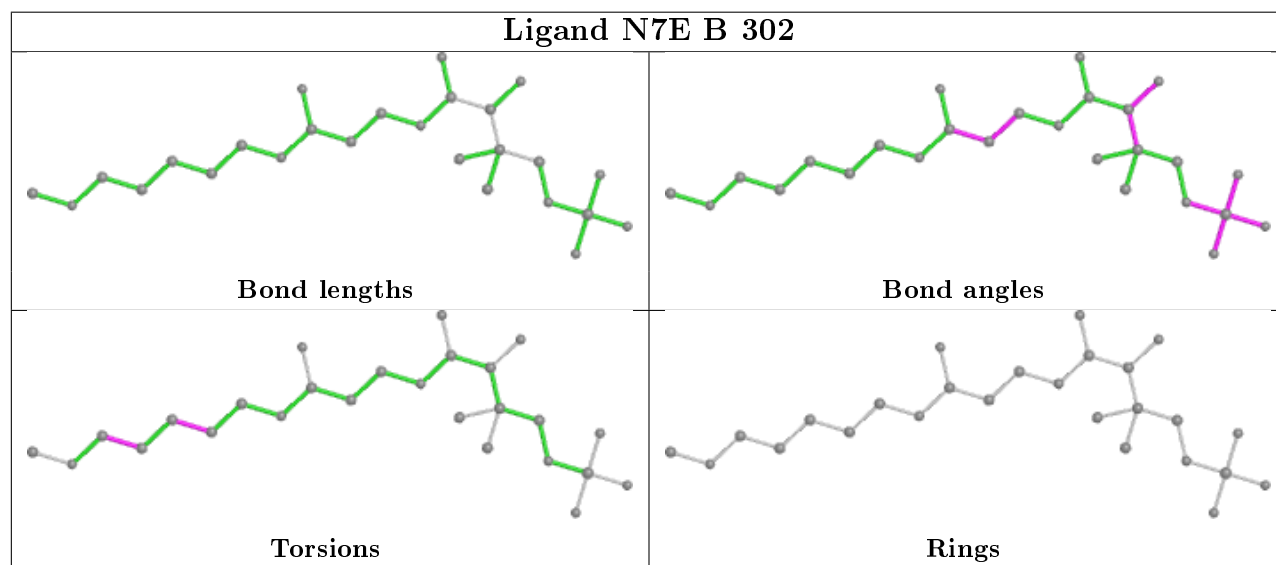
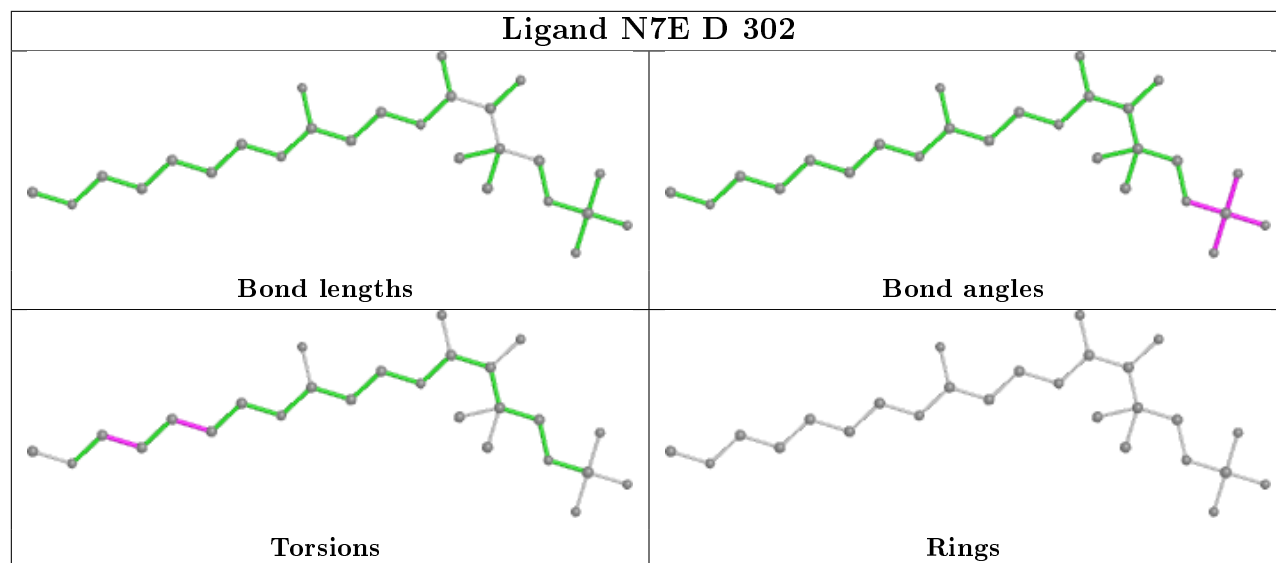
1 monomer is involved in 1 short contact:

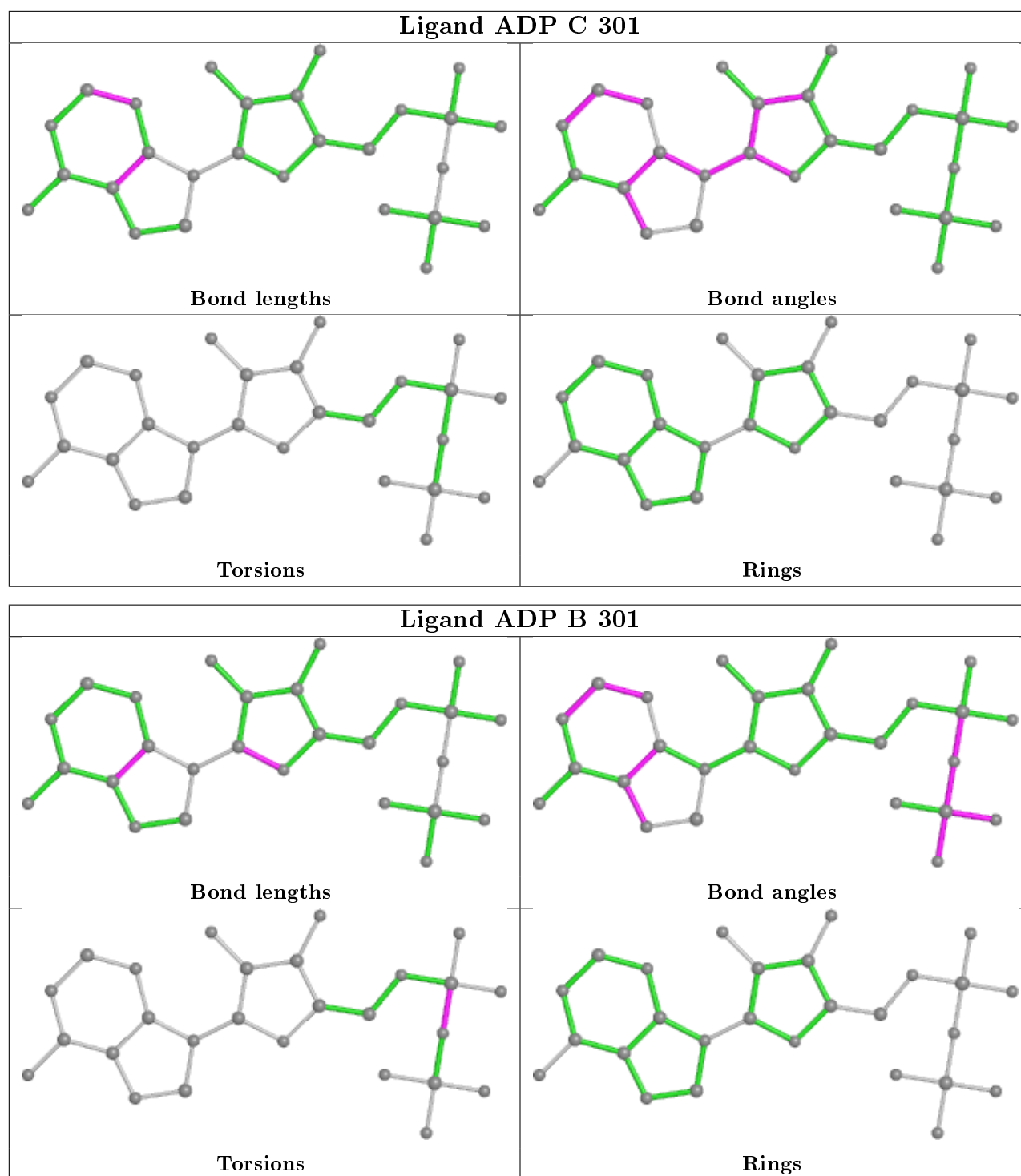
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	301	ADP	1	0

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 ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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, 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	265/265 (100%)	0.57	15 (5%) 23 30	14, 34, 59, 89	0
1	B	265/265 (100%)	0.64	25 (9%) 8 11	20, 38, 66, 92	0
1	C	259/265 (97%)	0.81	30 (11%) 4 6	13, 32, 84, 111	0
1	D	259/265 (97%)	1.04	51 (19%) 1 1	18, 38, 88, 100	0
All	All	1048/1060 (98%)	0.76	121 (11%) 4 7	13, 35, 77, 111	0

All (121) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	61	ILE	8.4
1	D	39	LEU	6.8
1	A	184	ASP	6.4
1	C	41	GLN	6.1
1	D	67	ILE	5.9
1	D	264	TYR	5.8
1	D	61	ILE	5.8
1	C	264	TYR	5.7
1	D	35	VAL	5.5
1	B	264	TYR	5.2
1	D	38	TRP	5.1
1	C	57	ILE	4.7
1	D	43	GLN	4.7
1	D	46	LYS	4.7
1	C	65	ALA	4.7
1	D	16	GLN	4.6
1	D	41	GLN	4.5
1	D	36	VAL	4.4
1	C	46	LYS	4.4
1	D	55	GLY	4.1
1	D	26	THR	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	161	LYS	4.1
1	C	55	GLY	4.0
1	C	265	LEU	4.0
1	B	161	LYS	4.0
1	D	186	ASP	4.0
1	C	64	PRO	4.0
1	D	37	GLU	3.8
1	B	162	ASP	3.7
1	D	42	GLN	3.7
1	B	59	GLU	3.7
1	C	40	ASN	3.7
1	B	37	GLU	3.6
1	B	62	ASN	3.6
1	C	42	GLN	3.6
1	D	60	ASN	3.5
1	D	62	ASN	3.5
1	D	56	VAL	3.5
1	C	44	ILE	3.4
1	D	58	ALA	3.4
1	A	162	ASP	3.3
1	D	3	VAL	3.3
1	D	44	ILE	3.2
1	D	64	PRO	3.2
1	C	68	PHE	3.2
1	D	59	GLU	3.2
1	D	265	LEU	3.2
1	D	65	ALA	3.2
1	C	59	GLU	3.2
1	C	37	GLU	3.2
1	C	60	ASN	3.2
1	D	54	ALA	3.1
1	D	40	ASN	3.1
1	D	34	GLN	3.0
1	C	38	TRP	2.9
1	B	43	GLN	2.9
1	B	61	ILE	2.9
1	A	165	PRO	2.9
1	D	162	ASP	2.9
1	D	33	ASP	2.8
1	C	34	GLN	2.8
1	B	158	HIS	2.8
1	C	47	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	43	GLN	2.7
1	A	163	THR	2.7
1	C	36	VAL	2.6
1	C	186	ASP	2.6
1	D	27	GLU	2.6
1	A	190	SER	2.6
1	A	185	ALA	2.6
1	D	63	ILE	2.6
1	B	160	TYR	2.6
1	A	181	HIS	2.6
1	C	162	ASP	2.6
1	D	109	GLN	2.6
1	D	68	PHE	2.5
1	B	181	HIS	2.5
1	D	5	ILE	2.5
1	D	32	ILE	2.5
1	D	25	LYS	2.5
1	B	163	THR	2.5
1	A	164	GLU	2.5
1	D	185	ALA	2.5
1	C	67	ILE	2.5
1	B	166	PRO	2.4
1	B	40	ASN	2.4
1	D	262	ALA	2.4
1	D	57	ILE	2.4
1	C	39	LEU	2.4
1	C	63	ILE	2.4
1	B	164	GLU	2.4
1	D	263	LEU	2.4
1	B	38	TRP	2.4
1	B	187	PHE	2.3
1	C	26	THR	2.3
1	A	63	ILE	2.3
1	C	12	ILE	2.3
1	B	45	GLU	2.3
1	D	45	GLU	2.3
1	B	19	ASP	2.3
1	A	37	GLU	2.2
1	C	164	GLU	2.2
1	D	52	GLY	2.2
1	D	4	GLY	2.2
1	B	150	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	28	LEU	2.2
1	A	18	GLN	2.1
1	D	12	ILE	2.1
1	A	41	GLN	2.1
1	B	190	SER	2.1
1	D	51	GLY	2.1
1	A	183	LEU	2.1
1	B	184	ASP	2.1
1	B	165	PRO	2.1
1	A	62	ASN	2.1
1	B	134	ILE	2.0
1	D	261	GLY	2.0
1	D	8	GLY	2.0
1	D	23	THR	2.0
1	C	56	VAL	2.0
1	B	159	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 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(Å <sup>2</sup> )	Q<0.9
4	MG	B	303	1/1	0.88	0.09	40,40,40,40	0
3	N7E	C	302	26/26	0.91	0.16	24,26,35,36	0
4	MG	A	303	1/1	0.92	0.13	34,34,34,34	0
4	MG	D	303	1/1	0.92	0.06	39,39,39,39	0
3	N7E	D	302	26/26	0.92	0.18	34,40,50,53	0
2	ADP	D	301	27/27	0.94	0.13	24,26,33,33	0
4	MG	C	303	1/1	0.94	0.09	31,31,31,31	0

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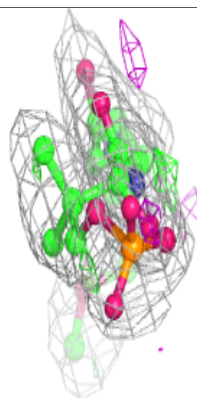
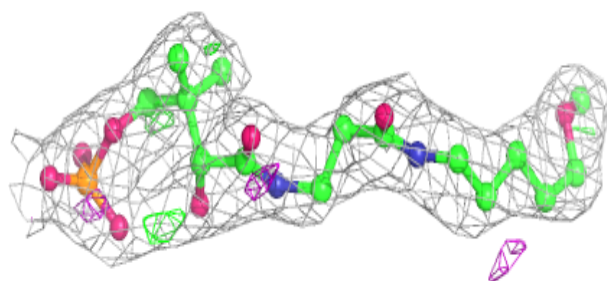
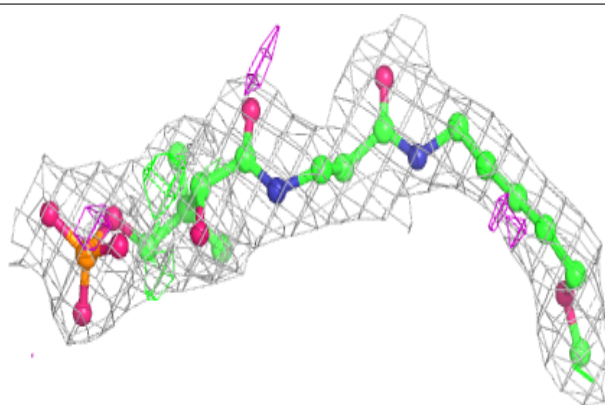
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ADP	C	301	27/27	0.94	0.14	23,25,28,28	0
3	N7E	A	302	26/26	0.95	0.15	16,18,20,22	0
3	N7E	B	302	26/26	0.96	0.17	21,23,27,27	0
2	ADP	B	301	27/27	0.96	0.13	25,26,28,30	0
2	ADP	A	301	27/27	0.97	0.13	19,22,26,27	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.

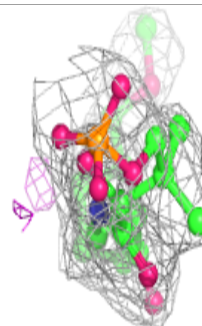
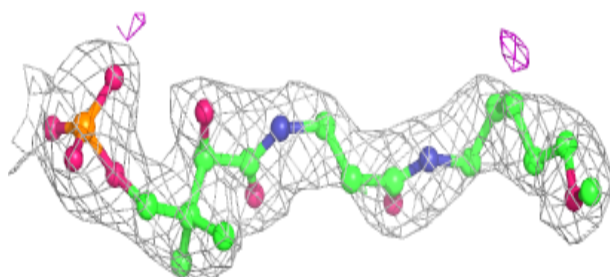
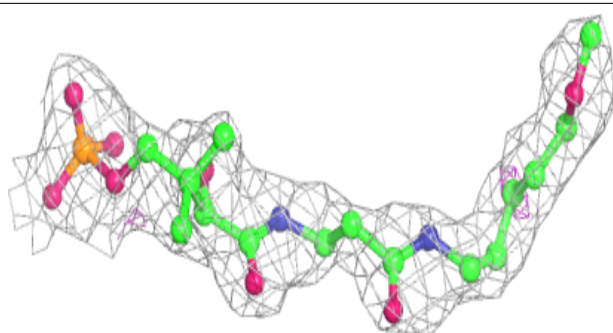
**Electron density around N7E C 302:**

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)

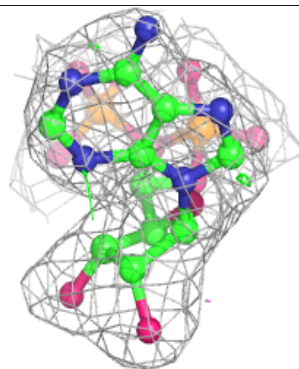
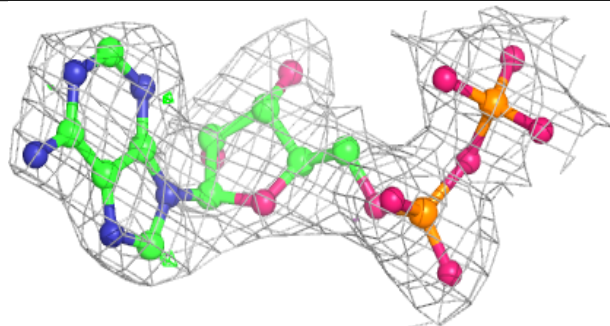
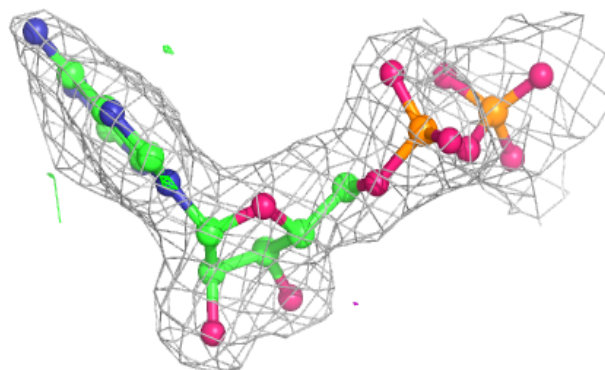


**Electron density around N7E D 302:**

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

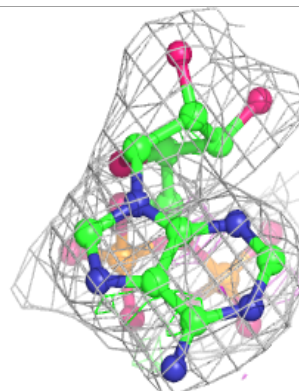
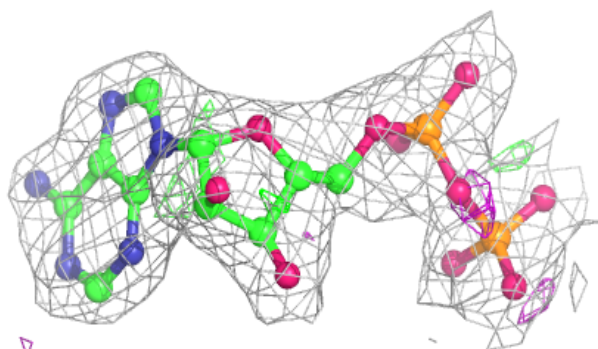
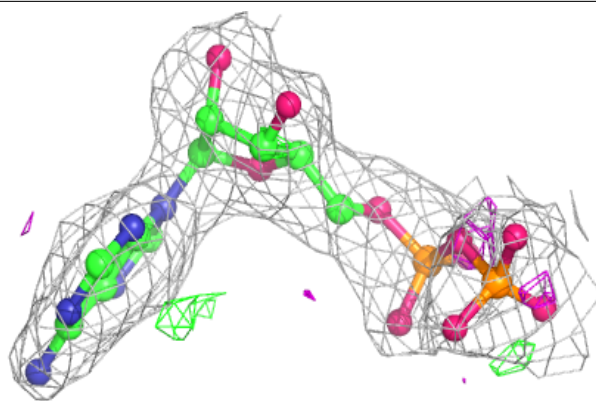
**Electron density around ADP D 301:**

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

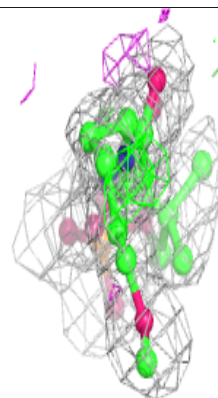
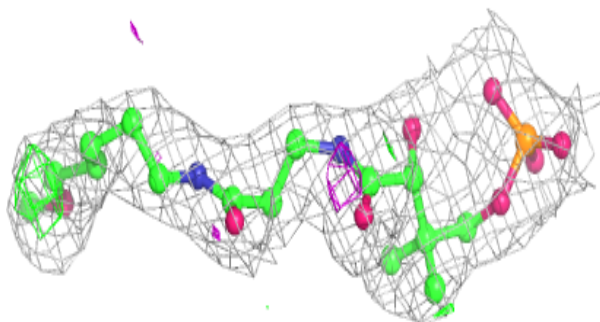
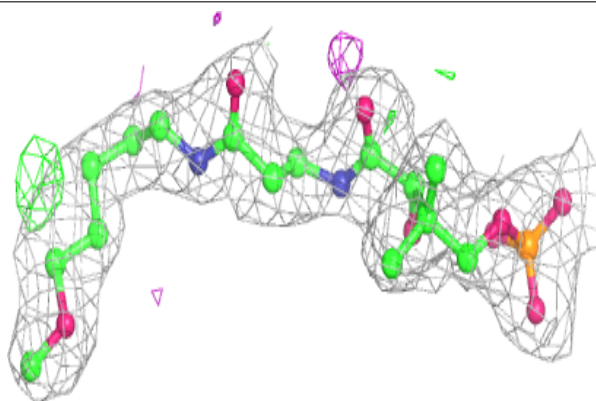


**Electron density around ADP C 301:**

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

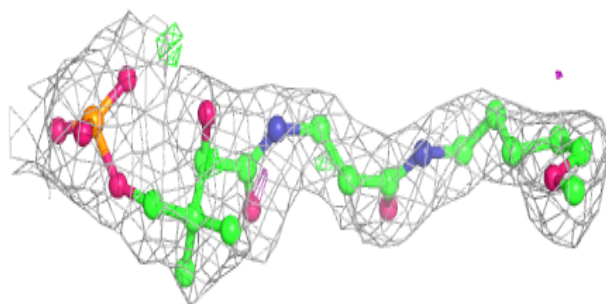
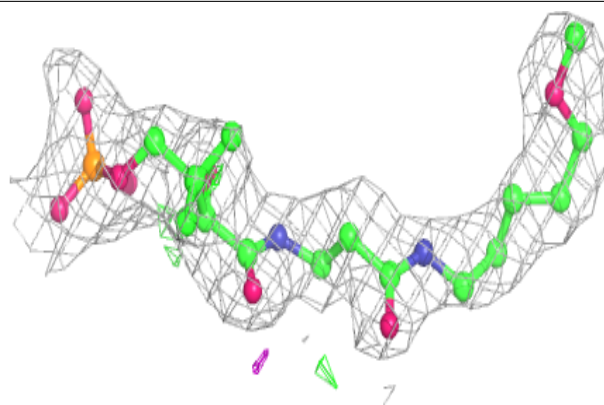
**Electron density around N7E A 302:**

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

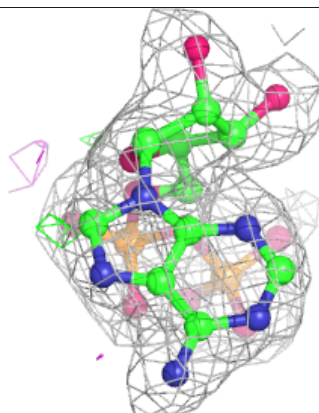
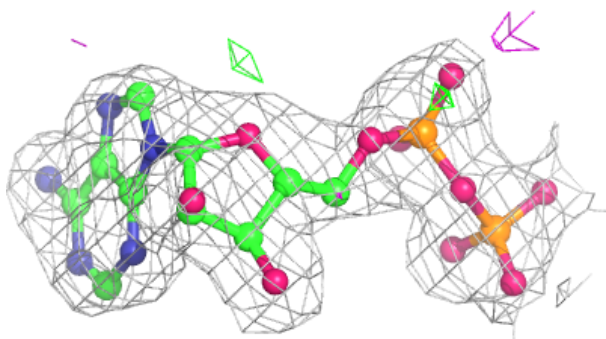
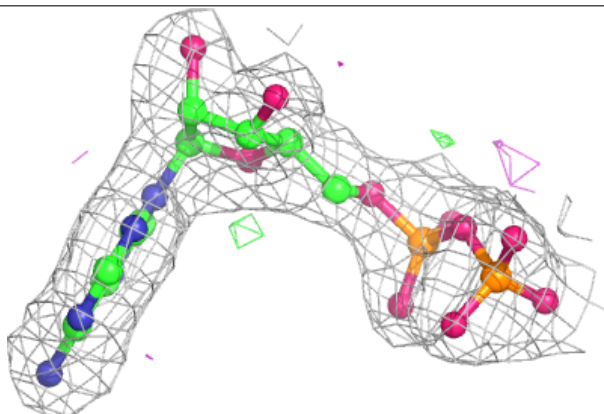


**Electron density around N7E B 302:**

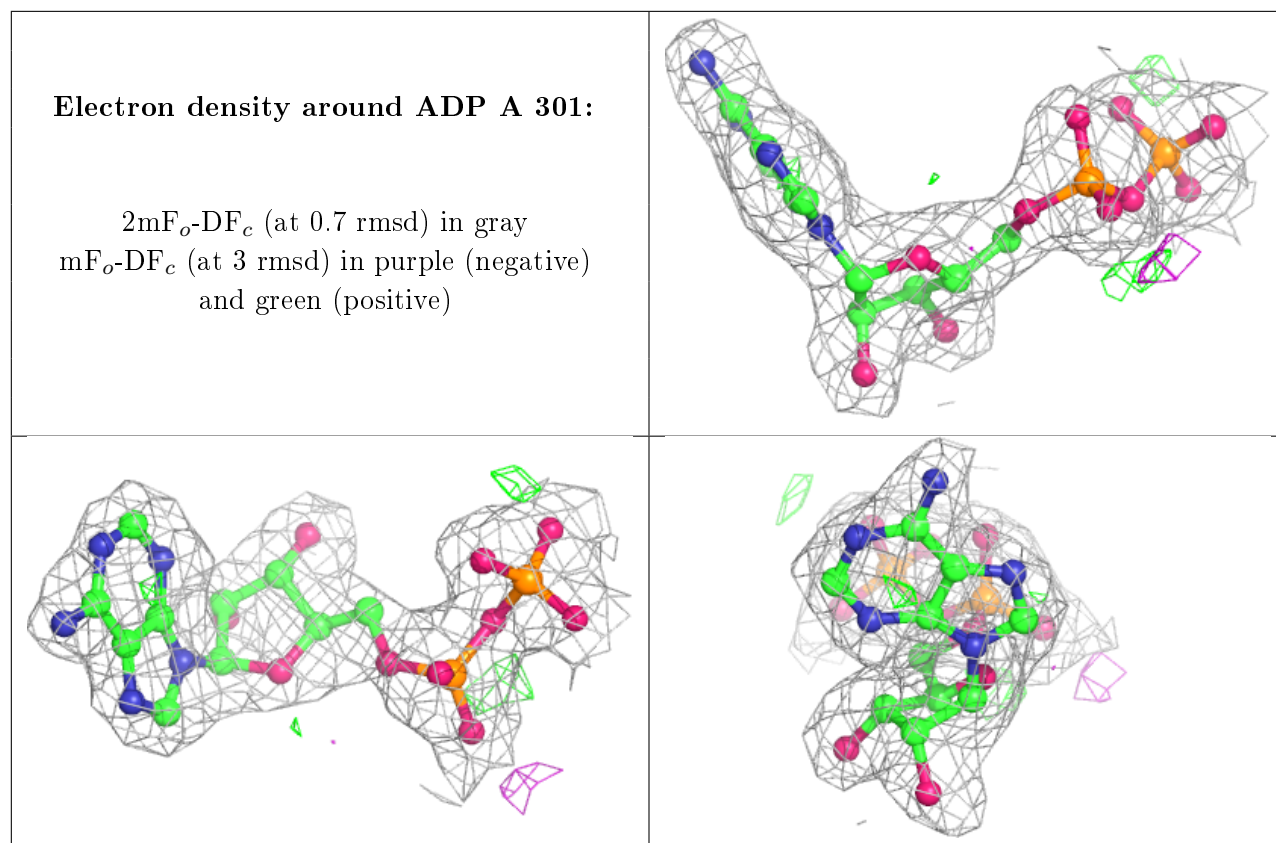
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around ADP B 301:**

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







## 6.5 Other polymers [i](#)

There are no such residues in this entry.