Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 11:44 PM JST

PDB ID	:	6IHI
Title	:	Crystal structure of RasADH 3B3/I91V from Ralstonia.sp in complex with NADPH and A6O
Authors	:	Zhang, H.L.; Chen, X.; Liu, W.D.; Wu, Q.Q.; Zhu, D.M.
Deposited on	:	2018-09-30
Resolution	:	1.78 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 (i) 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 (i)) were used in the production of this report:

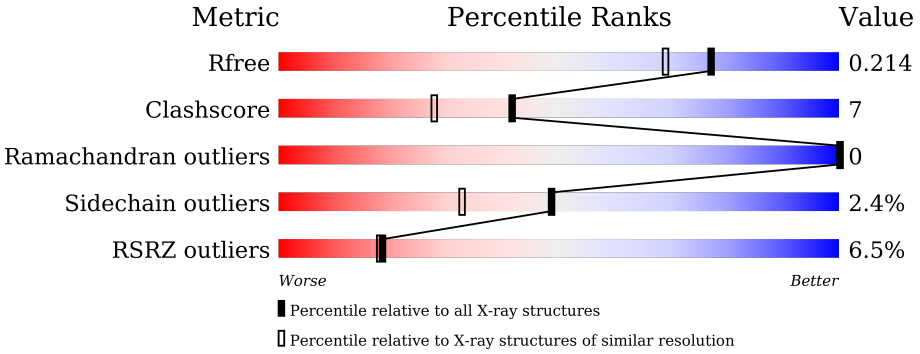
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (1.80-1.76)

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 <=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	249	8% 90% 9%
1	B	249	3% 86% 8% • 6%
1	C	249	6% 87% 8% • •
1	D	249	8% 85% 9% • 5%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 8124 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.

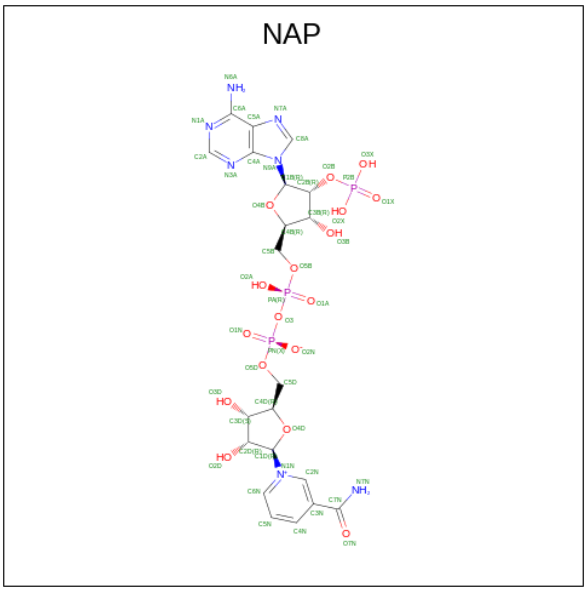
- Molecule 1 is a protein called Alclohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O				
1	A	248	1868	1168	335	365		0	0	0
			Total	C	N	O	S			
1	B	235	1765	1108	319	337	1	0	0	0
			Total	C	N	O	S			
1	C	239	1798	1129	323	345	1	0	0	0
			Total	C	N	O	S			
1	D	236	1773	1114	320	338	1	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

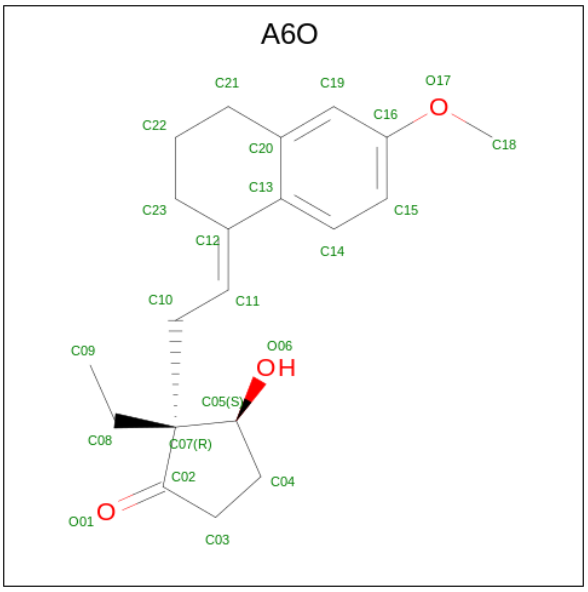
Chain	Residue	Modelled	Actual	Comment	Reference
A	91	VAL	ILE	engineered mutation	UNP C0IR58
A	187	SER	ILE	engineered mutation	UNP C0IR58
A	188	LEU	ILE	engineered mutation	UNP C0IR58
A	205	ALA	PHE	engineered mutation	UNP C0IR58
B	91	VAL	ILE	engineered mutation	UNP C0IR58
B	187	SER	ILE	engineered mutation	UNP C0IR58
B	188	LEU	ILE	engineered mutation	UNP C0IR58
B	205	ALA	PHE	engineered mutation	UNP C0IR58
C	91	VAL	ILE	engineered mutation	UNP C0IR58
C	187	SER	ILE	engineered mutation	UNP C0IR58
C	188	LEU	ILE	engineered mutation	UNP C0IR58
C	205	ALA	PHE	engineered mutation	UNP C0IR58
D	91	VAL	ILE	engineered mutation	UNP C0IR58
D	187	SER	ILE	engineered mutation	UNP C0IR58
D	188	LEU	ILE	engineered mutation	UNP C0IR58
D	205	ALA	PHE	engineered mutation	UNP C0IR58

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	48	21	7	17	3	0	0
			Total	C	N	O	P		
2	B	1	48	21	7	17	3	0	0
			Total	C	N	O	P		
2	C	1	48	21	7	17	3	0	0

- Molecule 3 is (2R,3S)-2-ethyl-2-[(2E)-2-(6-methoxy-3,4-dihydro-2H-naphthalen-1-ylidene)ethyl]-3-oxidanyl-cyclopentan-1-one (three-letter code: A6O) (formula: $C_{20}H_{26}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	23	20	3	0	0
			Total	C	O		
3	C	1	23	20	3	0	0

- Molecule 4 is water.

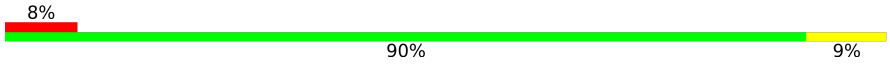
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	189	189	189	0	0
			Total	O		
4	B	196	196	196	0	0
			Total	O		
4	C	167	167	167	0	0
			Total	O		
4	D	178	178	178	0	0

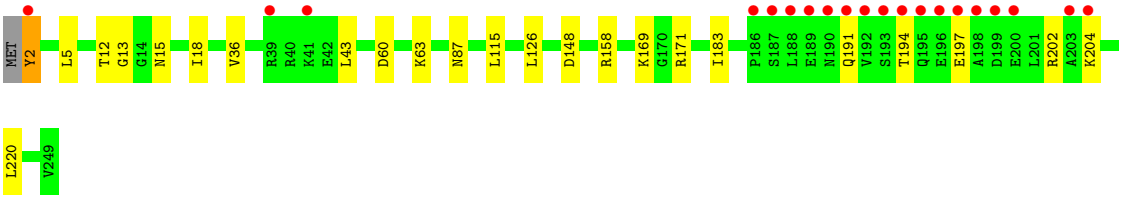
3 Residue-property plots (i)

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: Alclohol dehydrogenase

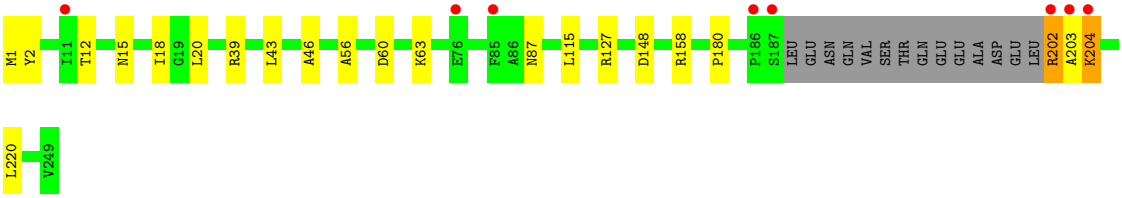
Chain A:





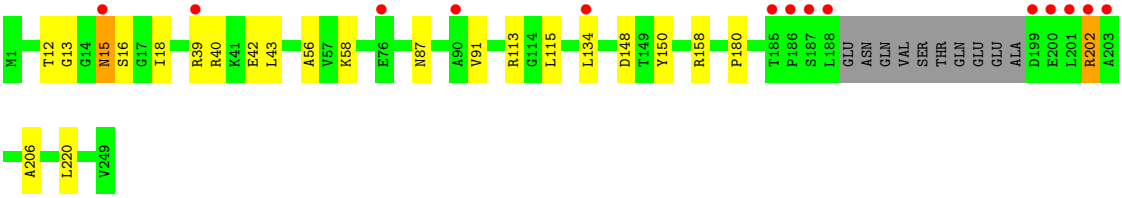
• Molecule 1: Alclohol dehydrogenase

Chain B:



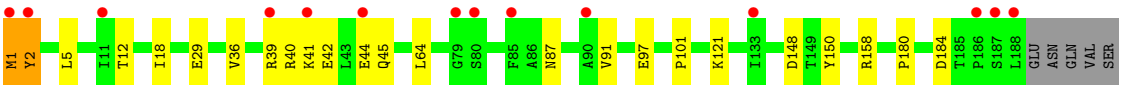
• Molecule 1: Alclohol dehydrogenase

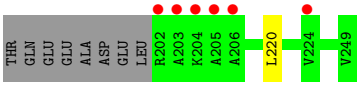
Chain C:



• Molecule 1: Alclohol dehydrogenase

Chain D:




THR
GLN
GLU
GLU
ALA
ASP
GLU
LEU
R202
A203
K204
A205
A206
L220
V224
V249

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	69.08Å 69.08Å 388.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.71 – 1.78 24.71 – 1.78	Depositor EDS
% Data completeness (in resolution range)	99.9 (24.71-1.78) 99.9 (24.71-1.78)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.08 (at 1.78Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.185 , 0.214 0.186 , 0.214	Depositor DCC
R_{free} test set	2007 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å^2)	26.7	Xtriage
Anisotropy	0.292	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 50.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.082 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8124	wwPDB-VP
Average B, all atoms (Å^2)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: A6O, NAP

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	0.29	0/1890	0.50	0/2561
1	B	0.29	0/1786	0.51	0/2417
1	C	0.30	0/1819	0.50	0/2462
1	D	0.33	0/1794	0.52	0/2428
All	All	0.30	0/7289	0.51	0/9868

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	1868	0	1902	18	0
1	B	1765	0	1815	20	0
1	C	1798	0	1847	28	0
1	D	1773	0	1826	34	0
2	A	48	0	23	4	0
2	B	48	0	23	4	0
2	C	48	0	25	3	0
3	A	23	0	0	6	0
3	C	23	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	189	0	0	3	0
4	B	196	0	0	2	0
4	C	167	0	0	3	0
4	D	178	0	0	2	0
All	All	8124	0	7461	104	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 7.

All (104) 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)
2:B:301:NAP:C1D	2:B:301:NAP:O4D	1.63	1.10
1:D:1:MET:H3	1:D:5:LEU:HD13	1.14	1.08
1:D:1:MET:N	1:D:5:LEU:HD13	1.72	1.05
1:C:15:ASN:ND2	1:C:42:GLU:HG2	1.82	0.94
1:B:202:ARG:HG2	1:B:203:ALA:H	1.31	0.94
1:D:1:MET:N	1:D:5:LEU:HD22	1.91	0.85
1:C:15:ASN:HD21	1:C:42:GLU:HG2	1.38	0.85
1:C:15:ASN:HD21	1:C:42:GLU:CG	1.89	0.84
1:C:15:ASN:ND2	1:C:42:GLU:CG	2.42	0.81
1:A:191:GLN:HE22	3:A:302:A6O:C09	1.92	0.81
1:D:1:MET:H3	1:D:5:LEU:CD1	1.93	0.79
1:B:202:ARG:CG	1:B:203:ALA:H	1.95	0.79
1:D:1:MET:N	1:D:5:LEU:CD1	2.48	0.76
1:B:39:ARG:NH2	4:B:402:HOH:O	2.22	0.71
1:D:18:ILE:HG12	1:D:220:LEU:HD23	1.72	0.70
1:B:202:ARG:HG2	1:B:203:ALA:N	2.06	0.69
1:C:16:SER:HB2	4:C:429:HOH:O	1.93	0.67
1:D:1:MET:H2	1:D:5:LEU:HD22	1.57	0.67
1:A:194:THR:HG23	1:A:197:GLU:H	1.59	0.67
1:B:127:ARG:HD2	4:B:401:HOH:O	1.95	0.67
1:C:15:ASN:HD21	1:C:42:GLU:CB	2.07	0.66
1:A:191:GLN:HE22	3:A:302:A6O:C08	2.08	0.65
1:D:1:MET:N	1:D:5:LEU:CD2	2.62	0.63
1:A:191:GLN:NE2	3:A:302:A6O:C09	2.62	0.62
1:B:60:ASP:HB3	1:B:63:LYS:HD2	1.82	0.62
1:B:18:ILE:HG21	1:B:220:LEU:HD23	1.84	0.59
1:A:169:LYS:NZ	4:A:401:HOH:O	2.32	0.58
1:A:18:ILE:HG21	1:A:220:LEU:HD23	1.85	0.58
2:A:301:NAP:O1A	2:A:301:NAP:H4B	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:ARG:CG	1:B:203:ALA:N	2.62	0.58
1:B:87:ASN:ND2	2:B:301:NAP:H4D	2.18	0.58
4:C:535:HOH:O	1:D:97:GLU:HG3	2.03	0.57
1:B:180:PRO:HG3	1:B:220:LEU:HD21	1.85	0.57
1:A:183:ILE:H	2:A:301:NAP:H72N	1.54	0.56
1:D:41:LYS:O	1:D:45:GLN:HG3	2.06	0.56
1:C:115:LEU:HD11	1:C:134:LEU:HD22	1.88	0.56
1:C:91:VAL:HG22	1:C:150:TYR:CE1	2.41	0.56
1:D:180:PRO:HG3	1:D:220:LEU:HD21	1.88	0.55
1:A:204:LYS:O	1:A:204:LYS:HD2	2.07	0.55
3:A:302:A6O:C09	3:A:302:A6O:C04	2.86	0.54
1:C:113:ARG:NH2	1:D:101:PRO:HB3	2.23	0.53
1:C:180:PRO:HG3	1:C:220:LEU:HD21	1.90	0.53
1:D:1:MET:O	1:D:1:MET:HG3	2.05	0.53
3:A:302:A6O:C18	4:A:468:HOH:O	2.56	0.53
1:B:1:MET:N	1:B:1:MET:SD	2.80	0.53
1:D:91:VAL:HG22	1:D:150:TYR:CE1	2.44	0.53
1:A:87:ASN:OD1	2:A:301:NAP:H4D	2.09	0.53
1:A:191:GLN:NE2	4:A:408:HOH:O	2.42	0.53
1:D:2:TYR:HB3	1:D:29:GLU:O	2.10	0.52
1:C:202:ARG:NH2	4:C:404:HOH:O	2.43	0.52
1:D:40:ARG:O	1:D:44:GLU:HB2	2.09	0.52
1:B:15:ASN:HB3	1:B:43:LEU:HD23	1.89	0.52
1:B:15:ASN:OD1	2:B:301:NAP:O3B	2.28	0.51
1:C:43:LEU:HB3	1:C:56:ALA:HB1	1.93	0.51
1:C:202:ARG:HH12	1:C:206:ALA:HB2	1.76	0.50
1:C:15:ASN:HD21	1:C:42:GLU:C	2.15	0.49
2:C:301:NAP:O1A	2:C:301:NAP:H4B	2.12	0.49
1:A:126:LEU:O	1:A:171:ARG:NH2	2.46	0.49
1:D:18:ILE:CG1	1:D:220:LEU:HD23	2.41	0.49
1:A:12:THR:O	1:A:87:ASN:HB3	2.12	0.49
1:A:15:ASN:HB3	1:A:43:LEU:HD23	1.94	0.48
1:B:43:LEU:HB3	1:B:56:ALA:HB1	1.95	0.48
1:B:20:LEU:HD13	1:B:46:ALA:HB1	1.96	0.48
1:A:18:ILE:CG2	1:A:220:LEU:HD23	2.44	0.47
1:C:87:ASN:OD1	2:C:301:NAP:H4D	2.14	0.47
1:B:18:ILE:CG2	1:B:220:LEU:HD23	2.44	0.47
1:A:60:ASP:HB3	1:A:63:LYS:HG3	1.97	0.47
1:D:40:ARG:NE	1:D:44:GLU:OE2	2.47	0.47
1:A:12:THR:HA	1:A:36:VAL:HB	1.97	0.47
1:B:12:THR:O	1:B:87:ASN:HB3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:ILE:HG21	1:C:220:LEU:HD23	1.96	0.46
1:C:18:ILE:HG12	1:C:220:LEU:HD23	1.97	0.46
1:D:12:THR:O	1:D:87:ASN:HB3	2.16	0.46
1:B:204:LYS:O	1:B:204:LYS:HG3	2.15	0.45
1:C:13:GLY:HA2	2:C:301:NAP:H1B	1.96	0.45
1:C:39:ARG:HG2	1:C:39:ARG:HH11	1.80	0.45
1:C:18:ILE:CG2	1:C:220:LEU:HD23	2.47	0.45
1:D:1:MET:H1	1:D:5:LEU:HD13	1.72	0.45
1:D:1:MET:N	1:D:5:LEU:CG	2.80	0.45
1:C:18:ILE:CG1	1:C:220:LEU:HD23	2.47	0.44
1:D:184:ASP:OD2	4:D:301:HOH:O	2.21	0.44
1:C:18:ILE:HD13	1:C:220:LEU:CD2	2.48	0.44
1:D:220:LEU:HD12	1:D:220:LEU:HA	1.81	0.44
1:D:39:ARG:NE	1:D:42:GLU:OE1	2.51	0.44
1:D:39:ARG:CZ	1:D:42:GLU:OE1	2.67	0.43
1:C:113:ARG:HH22	1:D:101:PRO:HB3	1.82	0.43
1:A:2:TYR:N	1:A:5:LEU:HB2	2.34	0.43
1:D:91:VAL:HG22	1:D:150:TYR:CZ	2.54	0.42
1:C:39:ARG:HG2	1:C:39:ARG:NH1	2.34	0.42
1:C:15:ASN:ND2	1:C:42:GLU:CD	2.73	0.42
1:B:1:MET:HG2	1:B:2:TYR:CD2	2.54	0.42
3:A:302:A6O:C10	3:A:302:A6O:C14	2.98	0.42
1:D:1:MET:H1	1:D:5:LEU:CD1	2.27	0.42
1:C:12:THR:O	1:C:87:ASN:HB3	2.20	0.42
1:D:12:THR:HA	1:D:36:VAL:HB	2.02	0.42
1:D:1:MET:H3	1:D:5:LEU:CG	2.33	0.41
1:C:18:ILE:HD13	1:C:220:LEU:HD23	2.02	0.41
1:C:40:ARG:NE	1:C:58:LYS:HZ1	2.18	0.41
1:D:2:TYR:H	1:D:5:LEU:HB2	1.86	0.41
1:A:13:GLY:HA2	2:A:301:NAP:H1B	2.02	0.41
1:B:18:ILE:HD12	2:B:301:NAP:O1N	2.21	0.41
1:D:64:LEU:HG	1:D:121:LYS:HE3	2.03	0.41
1:D:1:MET:H1	1:D:5:LEU:CD2	2.32	0.40
1:D:45:GLN:NE2	4:D:303:HOH:O	2.34	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	246/249 (99%)	240 (98%)	6 (2%)	0	100	100
1	B	231/249 (93%)	225 (97%)	6 (3%)	0	100	100
1	C	235/249 (94%)	227 (97%)	8 (3%)	0	100	100
1	D	232/249 (93%)	228 (98%)	4 (2%)	0	100	100
All	All	944/996 (95%)	920 (98%)	24 (2%)	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	193/194 (100%)	188 (97%)	5 (3%)	46	29
1	B	181/194 (93%)	176 (97%)	5 (3%)	43	27
1	C	185/194 (95%)	181 (98%)	4 (2%)	52	36
1	D	182/194 (94%)	178 (98%)	4 (2%)	52	36
All	All	741/776 (96%)	723 (98%)	18 (2%)	49	33

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	TYR
1	A	115	LEU

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Mol	Chain	Res	Type
1	A	148	ASP
1	A	158	ARG
1	A	202	ARG
1	B	115	LEU
1	B	148	ASP
1	B	158	ARG
1	B	202	ARG
1	B	204	LYS
1	C	15	ASN
1	C	148	ASP
1	C	158	ARG
1	C	202	ARG
1	D	1	MET
1	D	2	TYR
1	D	148	ASP
1	D	158	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	ASN
1	C	15	ASN

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)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NAP	A	301	-	45,52,52	4.50	14 (31%)	56,80,80	2.07	9 (16%)
3	A6O	C	302	-	25,25,25	4.17	13 (52%)	28,36,36	1.78	5 (17%)
2	NAP	B	301	-	45,52,52	4.53	15 (33%)	56,80,80	2.04	9 (16%)
3	A6O	A	302	-	25,25,25	4.16	15 (60%)	28,36,36	1.69	4 (14%)
2	NAP	C	301	-	45,52,52	0.83	1 (2%)	56,80,80	1.09	3 (5%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	A	301	-	-	7/31/67/67	0/5/5/5
3	A6O	C	302	-	-	6/11/38/38	0/3/3/3
2	NAP	B	301	-	-	11/31/67/67	0/5/5/5
3	A6O	A	302	-	-	8/11/38/38	0/3/3/3
2	NAP	C	301	-	-	7/31/67/67	0/5/5/5

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	NAP	O4D-C1D	16.35	1.63	1.41
2	B	301	NAP	O4B-C1B	15.45	1.62	1.41
2	A	301	NAP	C2D-C1D	-15.26	1.30	1.53
2	A	301	NAP	O4B-C1B	15.23	1.62	1.41
2	A	301	NAP	O4D-C1D	15.14	1.62	1.41
2	B	301	NAP	C2D-C1D	-14.05	1.32	1.53
3	A	302	A6O	C04-C05	-9.96	1.34	1.53
3	C	302	A6O	C04-C05	-9.96	1.34	1.53
3	C	302	A6O	C10-C07	-8.63	1.42	1.55
3	A	302	A6O	C10-C07	-8.09	1.43	1.55
3	A	302	A6O	C21-C20	7.34	1.63	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	302	A6O	C21-C20	7.28	1.63	1.51
3	C	302	A6O	C03-C02	-7.09	1.40	1.51
3	A	302	A6O	C03-C02	-7.07	1.40	1.51
2	B	301	NAP	C7N-N7N	6.73	1.45	1.33
2	A	301	NAP	C7N-N7N	6.53	1.45	1.33
2	B	301	NAP	O4B-C4B	-6.30	1.30	1.45
2	A	301	NAP	O4D-C4D	-6.09	1.31	1.45
2	A	301	NAP	O4B-C4B	-6.05	1.31	1.45
3	C	302	A6O	C04-C03	5.82	1.66	1.53
3	A	302	A6O	C04-C03	5.81	1.66	1.53
3	A	302	A6O	C08-C07	5.75	1.63	1.54
3	C	302	A6O	C08-C07	5.65	1.63	1.54
2	B	301	NAP	O4D-C4D	-5.32	1.33	1.45
3	A	302	A6O	C13-C12	4.51	1.56	1.46
3	C	302	A6O	C13-C12	4.30	1.55	1.46
3	C	302	A6O	O01-C02	-4.14	1.14	1.21
3	A	302	A6O	O01-C02	-4.06	1.14	1.21
3	C	302	A6O	C07-C05	3.68	1.62	1.55
3	A	302	A6O	C07-C05	3.64	1.62	1.55
3	C	302	A6O	C10-C11	3.32	1.58	1.50
3	A	302	A6O	C10-C11	3.26	1.58	1.50
2	A	301	NAP	P2B-O2B	3.24	1.65	1.59
2	B	301	NAP	C6A-N6A	3.23	1.45	1.34
2	A	301	NAP	C6A-N6A	3.20	1.45	1.34
2	B	301	NAP	O3D-C3D	-3.15	1.35	1.43
2	B	301	NAP	P2B-O2B	3.11	1.65	1.59
2	B	301	NAP	O3B-C3B	-3.00	1.35	1.43
3	A	302	A6O	C07-C02	2.97	1.57	1.52
2	A	301	NAP	O3B-C3B	-2.95	1.36	1.43
2	A	301	NAP	O2D-C2D	2.86	1.49	1.43
2	A	301	NAP	O3D-C3D	-2.85	1.36	1.43
2	A	301	NAP	O7N-C7N	-2.69	1.19	1.24
2	A	301	NAP	C5A-C4A	-2.67	1.33	1.40
3	C	302	A6O	C07-C02	2.64	1.56	1.52
2	B	301	NAP	C5A-C4A	-2.63	1.34	1.40
2	B	301	NAP	O7N-C7N	-2.62	1.19	1.24
3	A	302	A6O	C14-C13	2.54	1.43	1.39
3	C	302	A6O	C14-C13	2.46	1.43	1.39
3	A	302	A6O	C11-C12	2.41	1.35	1.34
2	C	301	NAP	C5A-C4A	2.41	1.47	1.40
3	A	302	A6O	C19-C16	2.39	1.43	1.38
2	B	301	NAP	C3N-C7N	2.38	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	NAP	C2A-N3A	2.35	1.35	1.32
2	A	301	NAP	C2A-N3A	2.35	1.35	1.32
3	C	302	A6O	C19-C16	2.32	1.43	1.38
2	B	301	NAP	O2D-C2D	2.29	1.48	1.43
3	A	302	A6O	C19-C20	2.00	1.42	1.39

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	301	NAP	C5A-C6A-N6A	10.08	135.67	120.35
2	B	301	NAP	C5A-C6A-N6A	9.65	135.02	120.35
2	A	301	NAP	N6A-C6A-N1A	-7.08	103.88	118.57
2	B	301	NAP	N6A-C6A-N1A	-6.57	104.95	118.57
3	A	302	A6O	C03-C02-C07	-6.24	103.86	109.25
3	C	302	A6O	C03-C02-C07	-6.00	104.07	109.25
2	A	301	NAP	N3A-C2A-N1A	-5.54	120.02	128.68
2	B	301	NAP	N3A-C2A-N1A	-5.39	120.25	128.68
3	C	302	A6O	C23-C12-C13	3.84	119.40	114.92
2	B	301	NAP	O4D-C1D-C2D	-3.57	101.72	106.93
2	C	301	NAP	PN-O3-PA	-3.53	120.71	132.83
3	C	302	A6O	C03-C04-C05	3.51	108.23	104.06
2	C	301	NAP	N3A-C2A-N1A	-3.32	123.48	128.68
2	A	301	NAP	PN-O3-PA	-3.28	121.56	132.83
2	A	301	NAP	C1B-N9A-C4A	-3.28	120.88	126.64
3	A	302	A6O	C03-C04-C05	3.11	107.76	104.06
2	B	301	NAP	C1B-N9A-C4A	-3.04	121.29	126.64
2	B	301	NAP	PN-O3-PA	-2.93	122.76	132.83
3	A	302	A6O	O01-C02-C07	2.93	128.88	125.32
3	A	302	A6O	C23-C12-C11	-2.68	117.14	122.86
2	C	301	NAP	C4A-C5A-N7A	-2.61	106.68	109.40
2	A	301	NAP	C3N-C7N-N7N	2.48	120.72	117.75
2	A	301	NAP	O4D-C1D-C2D	-2.46	103.33	106.93
2	B	301	NAP	C3N-C7N-N7N	2.43	120.66	117.75
2	B	301	NAP	O7N-C7N-N7N	-2.41	119.15	122.58
2	A	301	NAP	C6N-N1N-C2N	-2.38	119.80	121.97
3	C	302	A6O	O01-C02-C07	2.23	128.03	125.32
2	A	301	NAP	O7N-C7N-N7N	-2.16	119.51	122.58
2	B	301	NAP	C6N-N1N-C2N	-2.15	120.02	121.97
3	C	302	A6O	C14-C13-C12	-2.06	119.93	123.78

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	NAP	C4B-C5B-O5B-PA
2	A	301	NAP	C5D-O5D-PN-O3
2	B	301	NAP	C5B-O5B-PA-O1A
2	B	301	NAP	C5B-O5B-PA-O2A
2	B	301	NAP	C5D-O5D-PN-O3
2	B	301	NAP	C5D-O5D-PN-O2N
2	C	301	NAP	C4B-C5B-O5B-PA
2	C	301	NAP	C5D-O5D-PN-O3
2	C	301	NAP	C5D-O5D-PN-O2N
3	A	302	A6O	C10-C11-C12-C13
3	A	302	A6O	C05-C07-C08-C09
3	A	302	A6O	C02-C07-C10-C11
3	A	302	A6O	C08-C07-C10-C11
3	C	302	A6O	C10-C11-C12-C13
3	C	302	A6O	C10-C11-C12-C23
3	C	302	A6O	C02-C07-C08-C09
3	C	302	A6O	C19-C16-O17-C18
3	C	302	A6O	C15-C16-O17-C18
2	B	301	NAP	O4B-C4B-C5B-O5B
2	B	301	NAP	C3B-C4B-C5B-O5B
3	A	302	A6O	C19-C16-O17-C18
3	A	302	A6O	C15-C16-O17-C18
3	A	302	A6O	C10-C07-C08-C09
2	B	301	NAP	C5B-O5B-PA-O3
2	A	301	NAP	C5D-O5D-PN-O1N
2	B	301	NAP	C5D-O5D-PN-O1N
3	A	302	A6O	C05-C07-C10-C11
2	A	301	NAP	PN-O3-PA-O1A
2	B	301	NAP	PN-O3-PA-O5B
2	A	301	NAP	C2B-O2B-P2B-O1X
3	C	302	A6O	C10-C07-C08-C09
2	C	301	NAP	O4D-C4D-C5D-O5D
2	A	301	NAP	C2B-O2B-P2B-O2X
2	B	301	NAP	C2B-O2B-P2B-O2X
2	B	301	NAP	C2B-O2B-P2B-O3X
2	C	301	NAP	C2B-O2B-P2B-O2X
2	C	301	NAP	C2B-O2B-P2B-O3X
2	A	301	NAP	O4B-C4B-C5B-O5B
2	C	301	NAP	O4B-C4B-C5B-O5B

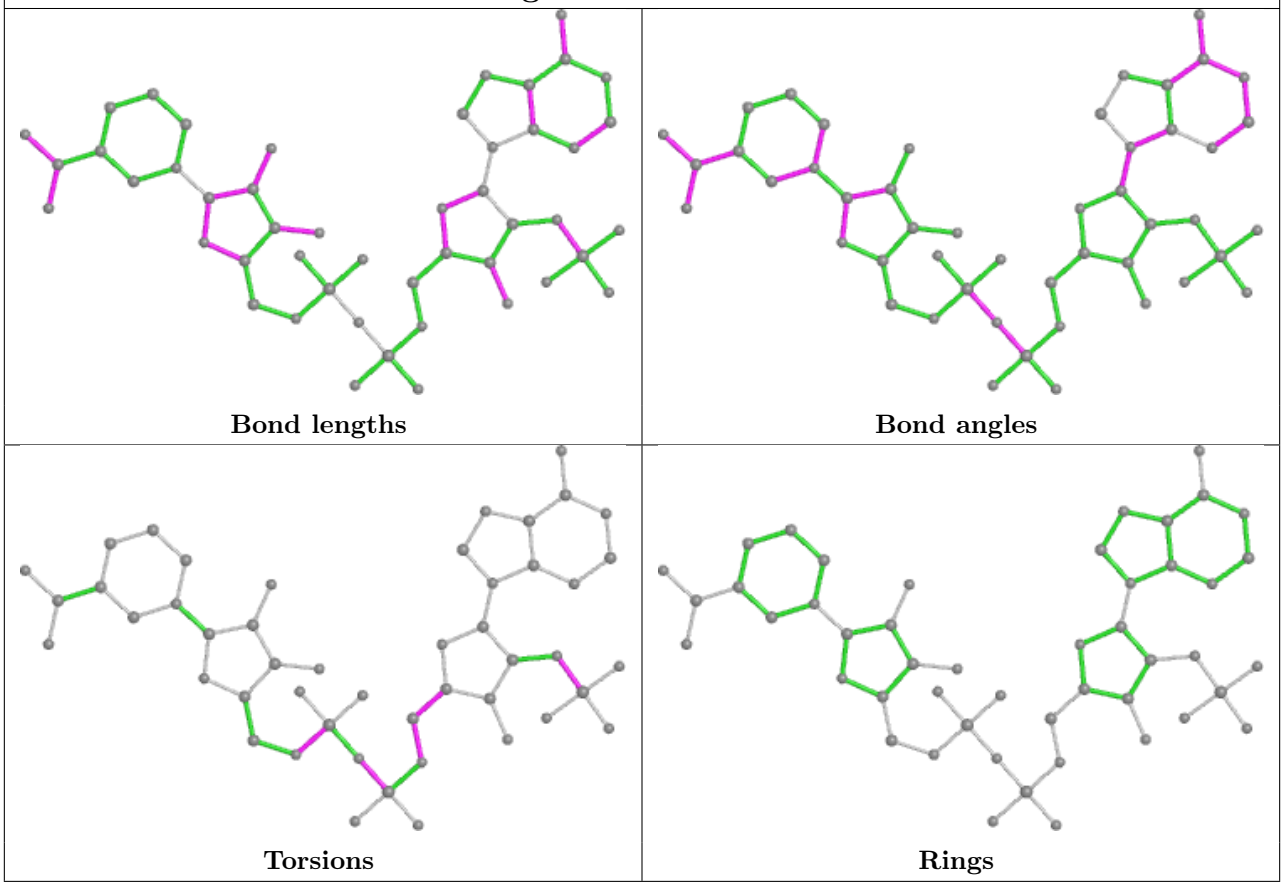
There are no ring outliers.

4 monomers are involved in 17 short contacts:

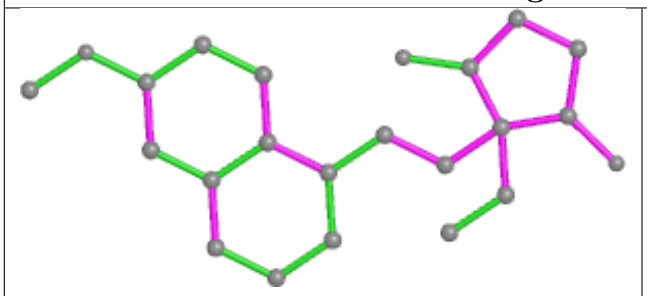
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NAP	4	0
2	B	301	NAP	4	0
3	A	302	A6O	6	0
2	C	301	NAP	3	0

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 then 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.

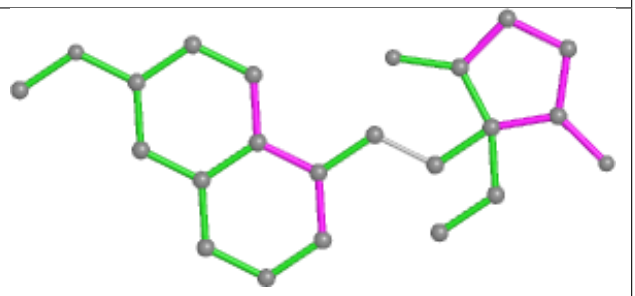
Ligand NAP A 301



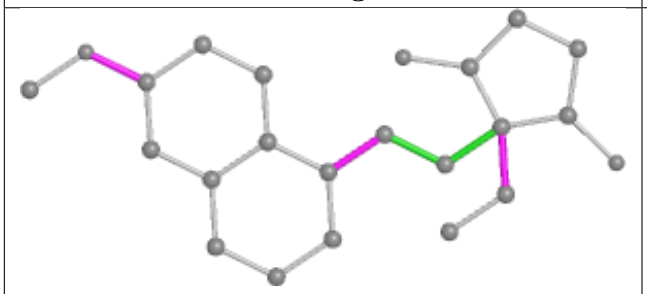
Ligand A6O C 302



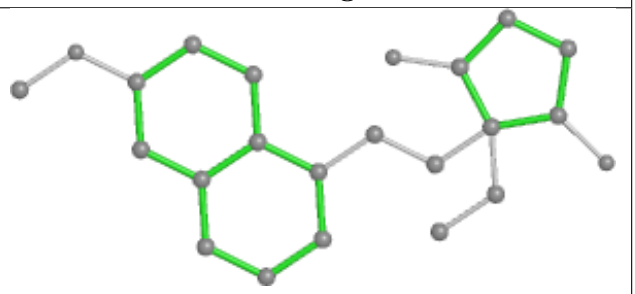
Bond lengths



Bond angles

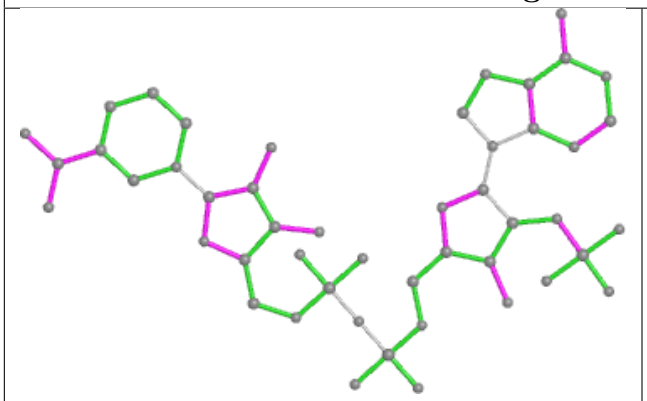


Torsions

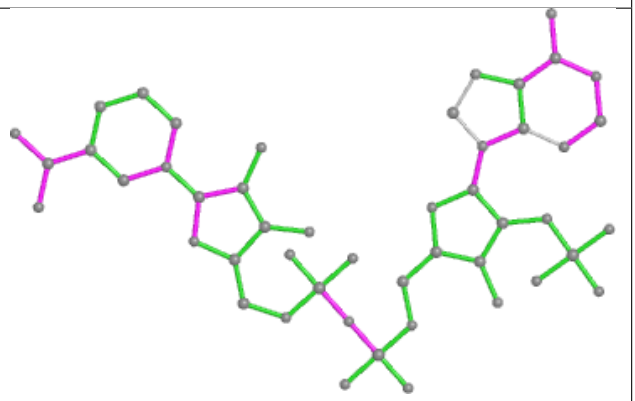


Rings

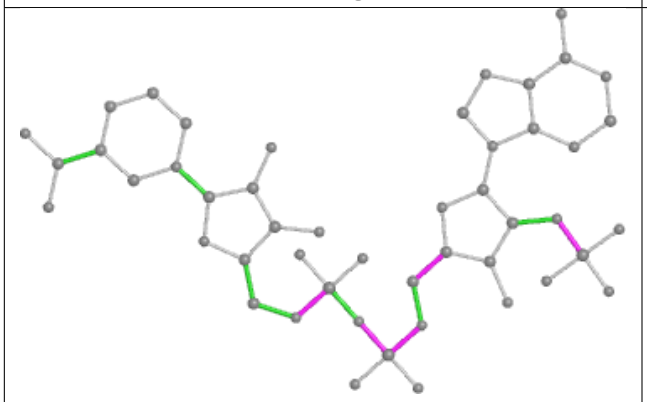
Ligand NAP B 301



Bond lengths



Bond angles



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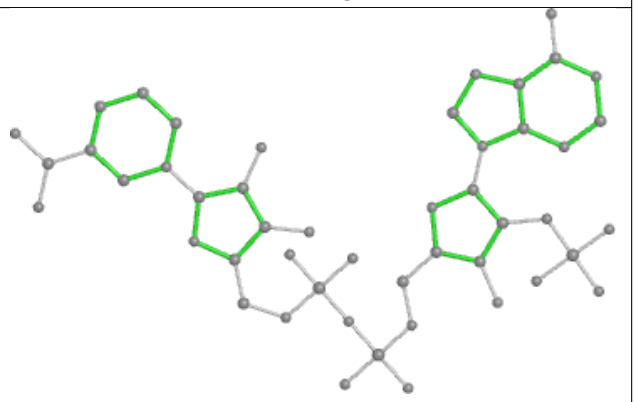
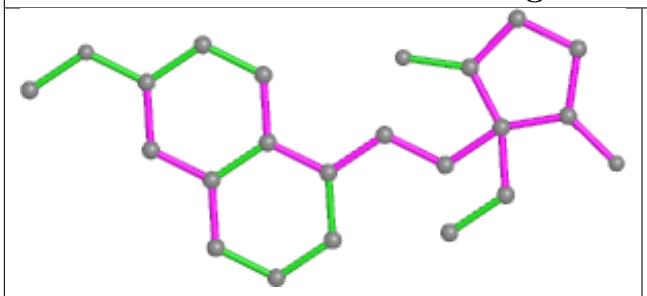
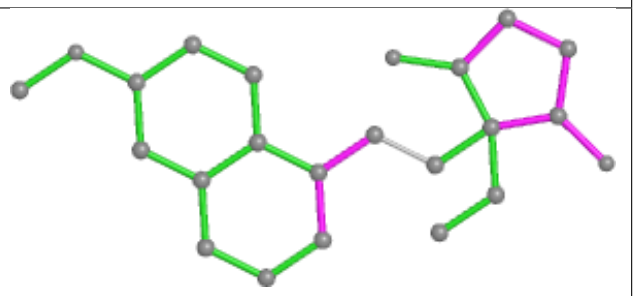


Rings

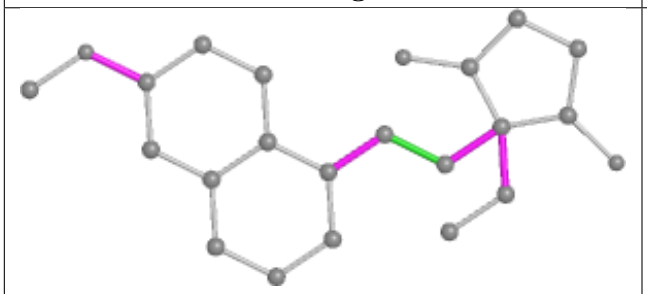
Ligand A6O A 302



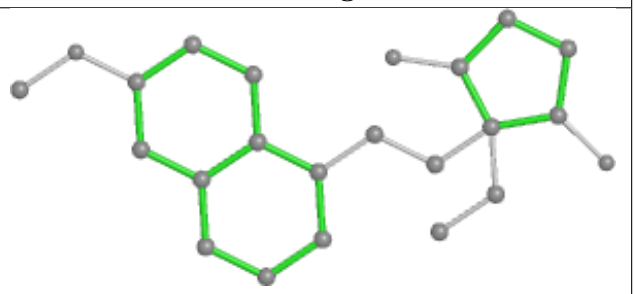
Bond lengths



Bond angles

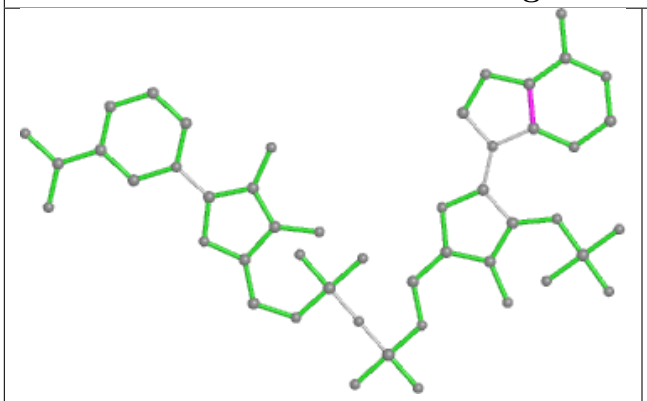


Torsions

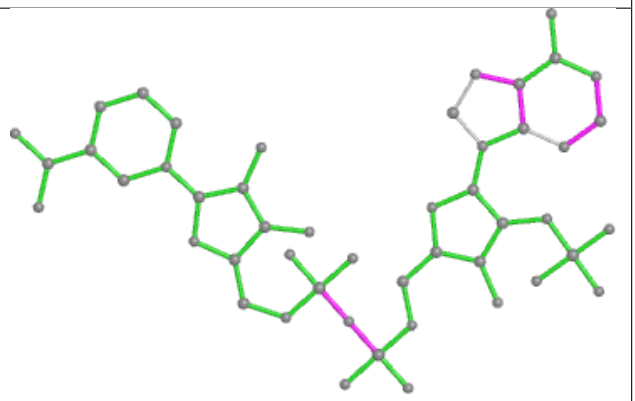


Rings

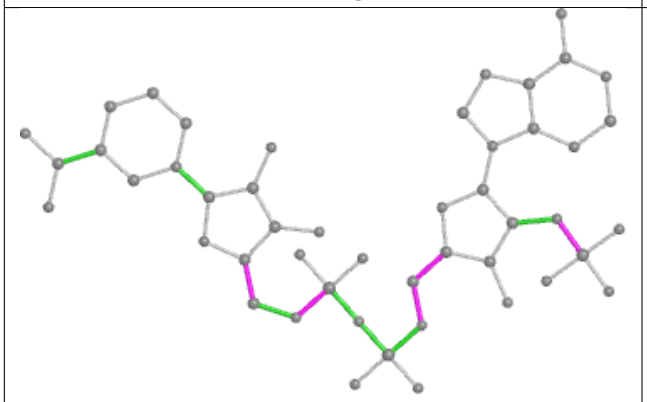
Ligand NAP C 301



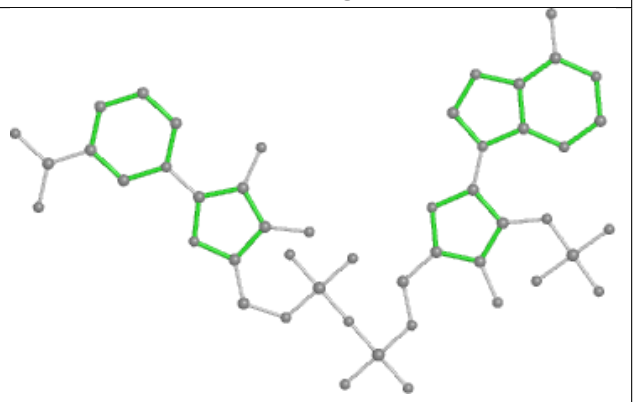
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	248/249 (99%)	0.14	20 (8%)	12	11	18, 27, 62, 80	0
1	B	235/249 (94%)	0.05	8 (3%)	45	44	18, 26, 44, 77	0
1	C	239/249 (95%)	0.29	14 (5%)	22	21	18, 30, 53, 98	0
1	D	236/249 (94%)	0.31	20 (8%)	10	10	20, 31, 54, 81	0
All	All	958/996 (96%)	0.20	62 (6%)	18	18	18, 28, 54, 98	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	186	PRO	9.8
1	D	1	MET	9.4
1	C	188	LEU	9.1
1	C	186	PRO	9.0
1	D	188	LEU	8.0
1	C	201	LEU	6.5
1	C	199	ASP	6.5
1	A	198	ALA	5.8
1	B	204	LYS	5.6
1	A	193	SER	5.5
1	B	186	PRO	5.4
1	D	187	SER	5.1
1	D	203	ALA	5.1
1	C	187	SER	4.9
1	B	187	SER	4.9
1	D	2	TYR	4.7
1	A	186	PRO	4.7
1	B	203	ALA	4.7
1	D	202	ARG	4.6
1	A	196	GLU	4.6
1	A	192	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
1	C	202	ARG	4.1
1	A	190	ASN	4.0
1	A	195	GLN	3.9
1	C	200	GLU	3.9
1	D	39	ARG	3.7
1	A	187	SER	3.7
1	A	200	GLU	3.7
1	A	2	TYR	3.6
1	D	90	ALA	3.5
1	A	189	GLU	3.4
1	D	205	ALA	3.4
1	D	41	LYS	3.4
1	C	185	THR	3.3
1	B	202	ARG	3.3
1	D	85	PHE	3.2
1	A	41	LYS	3.1
1	D	133	ILE	3.0
1	A	191	GLN	3.0
1	B	11	ILE	2.9
1	A	199	ASP	2.9
1	A	188	LEU	2.8
1	A	197	GLU	2.8
1	D	224	VAL	2.8
1	C	39	ARG	2.7
1	C	134	LEU	2.7
1	C	90	ALA	2.7
1	C	15	ASN	2.7
1	C	76	GLU	2.7
1	C	203	ALA	2.7
1	A	39	ARG	2.6
1	D	44	GLU	2.5
1	A	194	THR	2.4
1	D	11	ILE	2.4
1	D	204	LYS	2.3
1	B	76	GLU	2.2
1	B	85	PHE	2.2
1	D	80	SER	2.2
1	A	203	ALA	2.2
1	A	204	LYS	2.2
1	D	206	ALA	2.1
1	D	79	GLY	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)

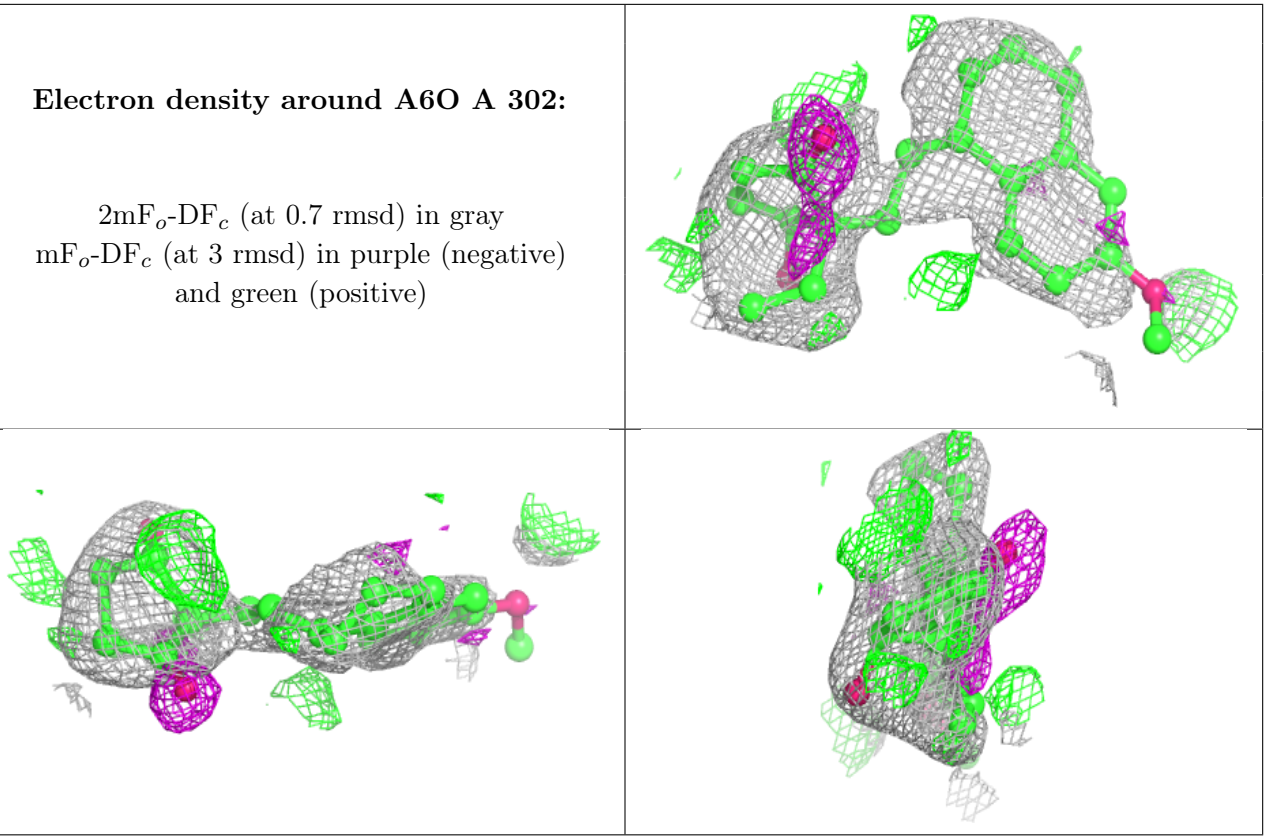
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	A6O	A	302	23/23	0.47	0.31	53,68,72,89	0
3	A6O	C	302	23/23	0.49	0.29	61,77,86,92	0
2	NAP	C	301	48/48	0.75	0.22	42,59,77,92	0
2	NAP	B	301	48/48	0.85	0.16	33,47,76,92	0
2	NAP	A	301	48/48	0.87	0.17	39,56,81,89	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.

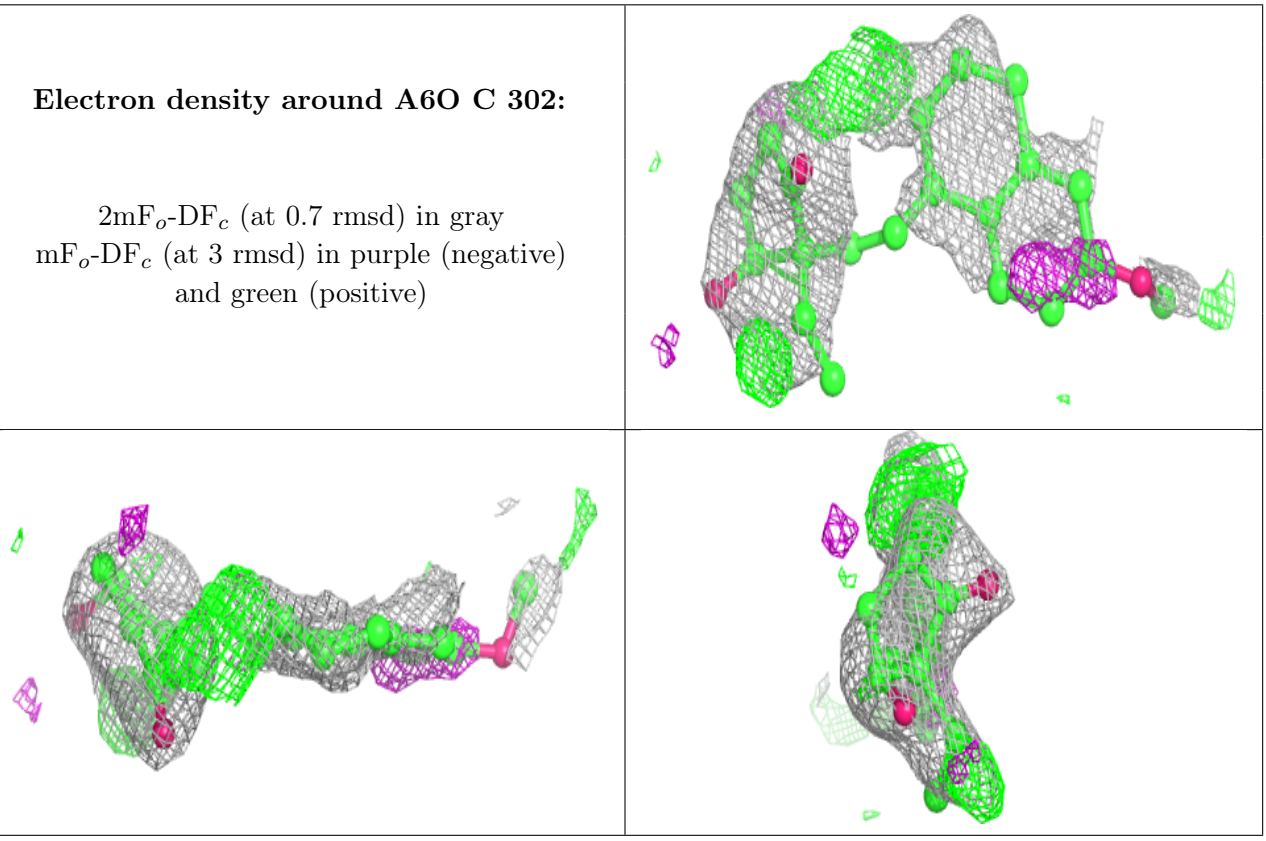
Electron density around A6O A 302:

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



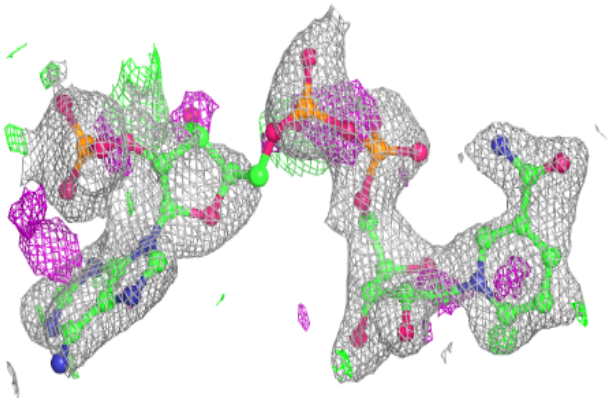
Electron density around A6O C 302:

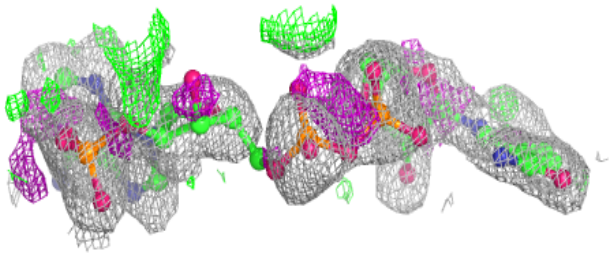
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

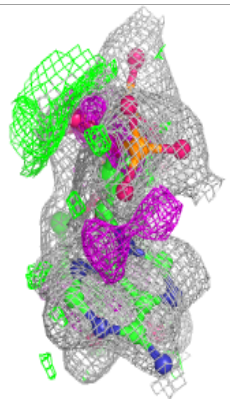


Electron density around NAP C 301:

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

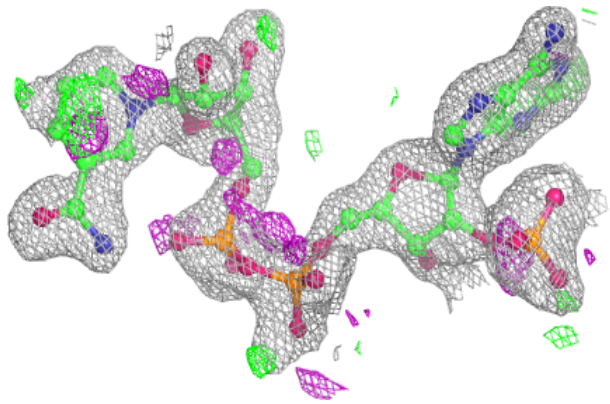


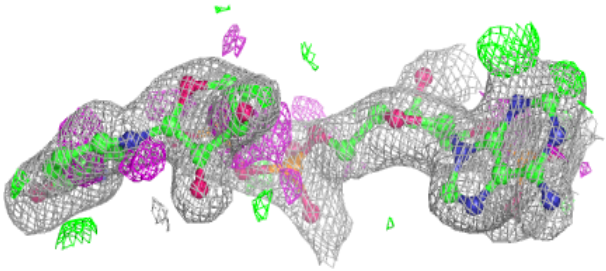


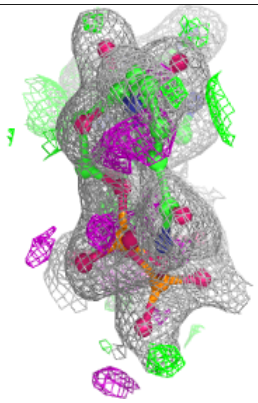


Electron density around NAP B 301:

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

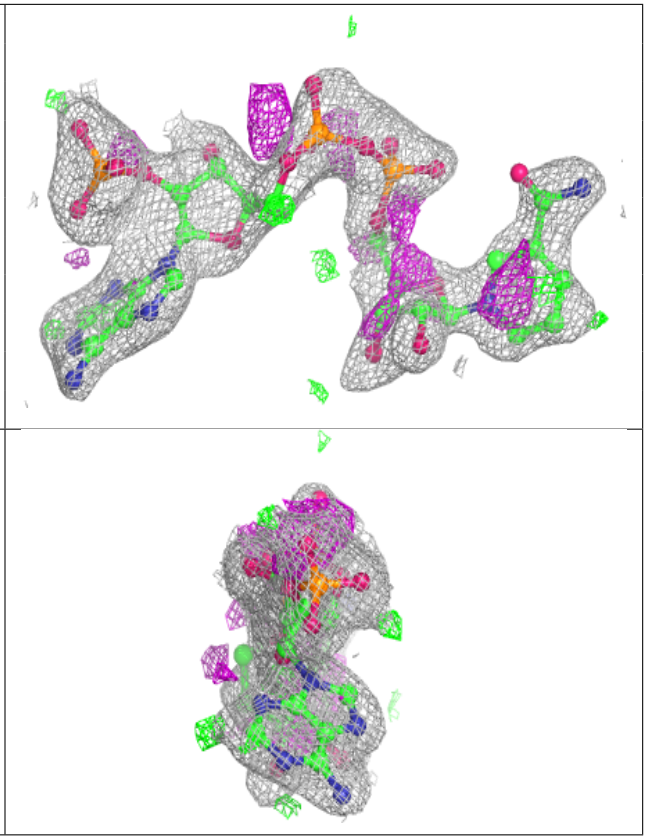






Electron density around NAP A 301:

$2mF_o$-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.