Full wwPDB X-ray Structure Validation Report (i)

Jan 2, 2024 – 09:03 pm GMT

PDB ID	:	4WAS
Title	:	STRUCTURE OF THE ETR1P/NADP/CROTONYL-COA COMPLEX
Authors	:	Quade, N.; Voegeli, B.; Rosenthal, R.; Capitani, G.; Erb, T.J.
Deposited on	:	2014-08-31
Resolution	:	1.70 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:

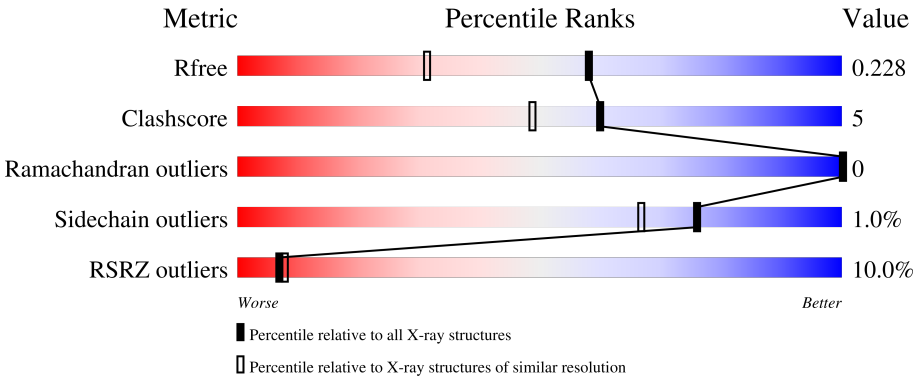
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
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 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	364	7% 90% 10%
1	B	364	9% 89% 11%
1	C	364	13% 88% 11%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	COO	A	402	X	-	-	-
3	COO	B	402[A]	X	-	-	-
3	COO	B	402[B]	X	-	-	-

2 Entry composition (i)

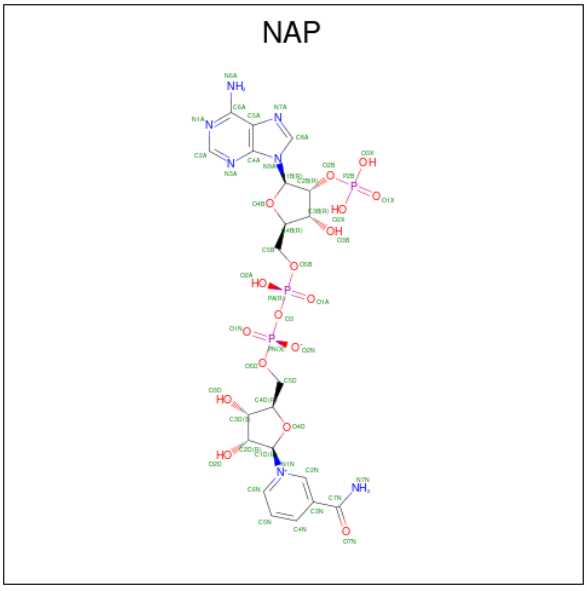
There are 4 unique types of molecules in this entry. The entry contains 10132 atoms, of which 1035 are hydrogens and 0 are deuteriums.

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 Enoyl-[acyl-carrier-protein] reductase [NADPH, B-specific] 1, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	364	Total	C	H	N	O	S	0	0	0
			3133	1768	345	466	548	6			
1	B	364	Total	C	H	N	O	S	0	2	0
			3153	1779	345	471	552	6			
1	C	364	Total	C	H	N	O	S	0	0	0
			3133	1768	345	466	548	6			

- 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
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

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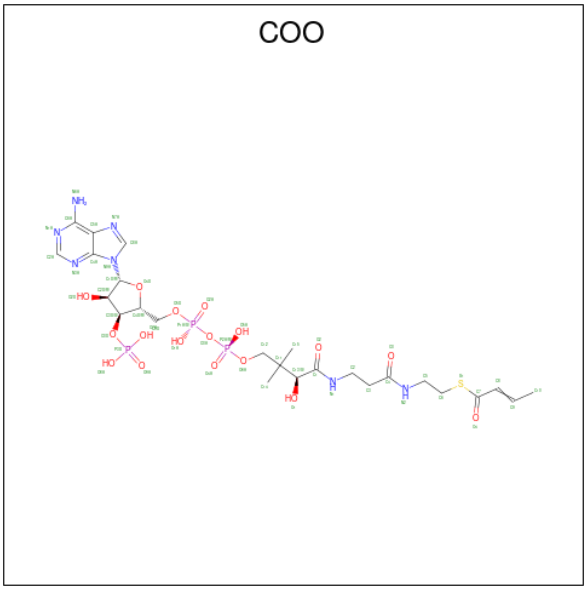
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is CROTONYL COENZYME A (three-letter code: COO) (formula: $C_{25}H_{40}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			31	15	2	11	2	1		
3	B	1	Total	C	N	O	P	S	0	1
			43	23	3	13	2	2		
3	C	1	Total	C	N	O	S		0	0
			12	8	1	2	1			

- Molecule 4 is water.

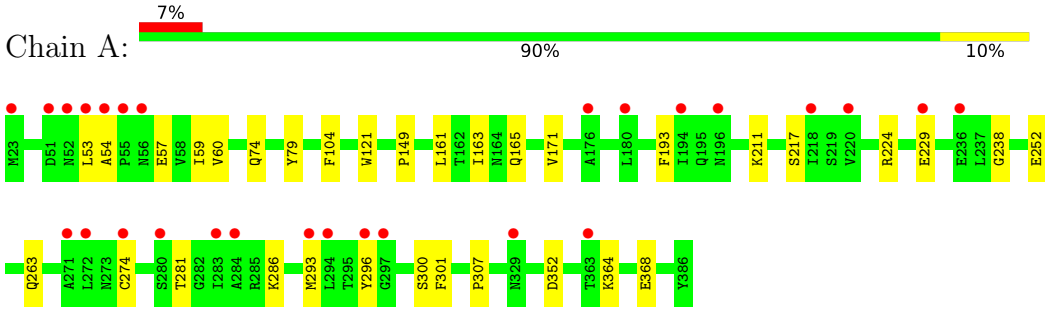
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	175	Total	O	0	0
			175	175		
4	B	147	Total	O	0	0
			147	147		
4	C	161	Total	O	0	0
			161	161		

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.

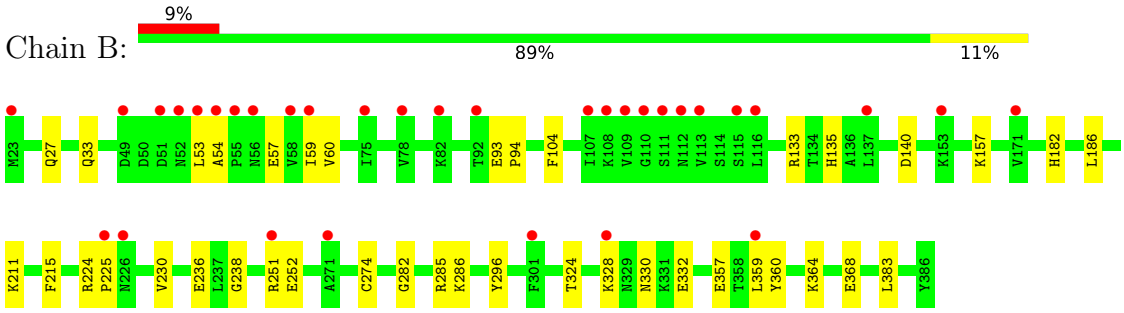
• Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADPH, B-specific] 1, mitochondrial

Chain A:



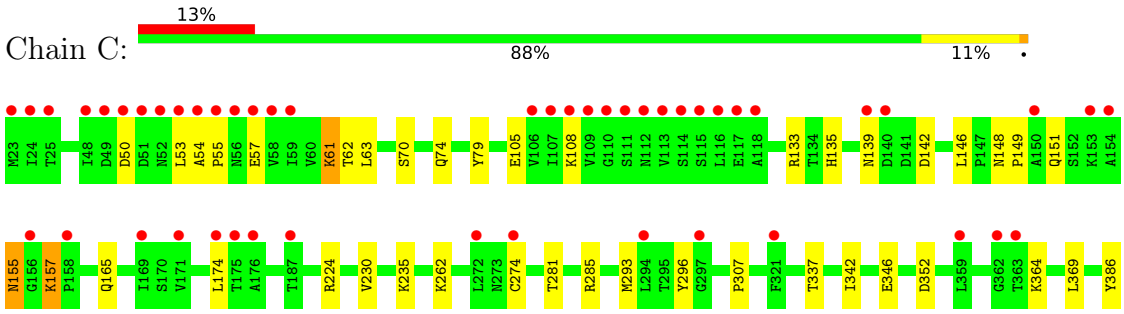
• Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADPH, B-specific] 1, mitochondrial

Chain B:



• Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADPH, B-specific] 1, mitochondrial

Chain C:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	147.92Å 106.14Å 94.41Å 90.00° 98.41° 90.00°	Depositor
Resolution (Å)	37.81 – 1.70 37.81 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (37.81-1.70) 99.1 (37.81-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	1.01 (at 1.70Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.188 , 0.224 0.194 , 0.228	Depositor DCC
R_{free} test set	1097 reflections (0.70%)	wwPDB-VP
Wilson B-factor (Å^2)	36.1	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.39 , 58.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	10132	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COO, 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	0.42	0/2847	0.56	0/3870
1	B	0.39	0/2867	0.53	0/3896
1	C	0.40	0/2847	0.54	1/3870 (0.0%)
All	All	0.41	0/8561	0.54	1/11636 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	293	MET	CG-SD-CE	-5.68	91.11	100.20

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2788	345	2767	22	0
1	B	2808	345	2784	25	0
1	C	2788	345	2767	31	0
2	A	48	0	25	2	0
2	B	48	0	25	1	0
2	C	48	0	25	1	0
3	A	31	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	43	0	28	1	0
3	C	12	0	10	1	0
4	A	175	0	0	4	0
4	B	147	0	0	4	0
4	C	161	0	0	3	0
All	All	9097	1035	8456	79	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 5.

All (79) 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:54:ALA:HB1	1:C:55:PRO:HD2	1.59	0.84
1:B:211:LYS:HE3	1:B:238:GLY:HA2	1.61	0.80
1:B:224:ARG:HD2	1:B:230:VAL:HG21	1.68	0.76
1:B:54:ALA:HB3	1:B:57:GLU:HG3	1.66	0.75
1:C:224:ARG:HD2	1:C:230:VAL:HG21	1.67	0.75
1:B:285[A]:ARG:NH2	4:B:641:HOH:O	2.19	0.74
1:A:296:TYR:O	2:A:401:NAP:H2N	1.93	0.68
1:C:364:LYS:HG3	1:C:369:LEU:HD11	1.80	0.63
1:B:211:LYS:HE3	1:B:238:GLY:CA	2.29	0.63
1:C:296:TYR:O	2:C:401:NAP:H2N	2.04	0.58
1:B:182:HIS:ND1	4:B:585:HOH:O	2.32	0.58
1:B:296:TYR:O	2:B:401:NAP:H2N	2.04	0.57
1:B:211:LYS:NZ	1:B:236:GLU:O	2.37	0.57
1:C:174:LEU:HD22	1:C:337:THR:HG22	1.87	0.57
1:C:54:ALA:HB3	1:C:57:GLU:HG3	1.89	0.55
1:A:257:ILE:HD12	1:A:286:LYS:HE3	1.88	0.55
3:B:402[B]:COO:H103	4:B:583:HOH:O	2.06	0.55
1:B:224:ARG:HB2	1:B:225:PRO:HD2	1.88	0.54
1:C:50:ASP:HB2	1:C:108:LYS:HD2	1.91	0.53
1:C:165:GLN:HG2	1:C:352:ASP:HB3	1.91	0.52
1:C:364:LYS:HG3	1:C:369:LEU:CD1	2.39	0.51
1:A:257:ILE:CD1	1:A:286:LYS:HE3	2.41	0.50
1:C:133:ARG:HD2	1:C:135:HIS:O	2.11	0.50
1:B:364:LYS:HE3	1:B:368:GLU:HB3	1.93	0.50
1:C:54:ALA:HB3	1:C:57:GLU:CG	2.42	0.50
1:C:155:ASN:O	1:C:155:ASN:ND2	2.34	0.50
1:A:53:LEU:HD21	1:A:59:ILE:HG13	1.94	0.49
1:A:364:LYS:HE2	1:A:368:GLU:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:SER:HB3	3:C:402:COO:C9	2.43	0.48
1:B:357[A]:GLU:HA	1:B:383:LEU:O	2.14	0.48
1:A:229:GLU:HG3	4:A:649:HOH:O	2.14	0.48
1:B:54:ALA:HB3	1:B:57:GLU:CG	2.39	0.48
1:A:193:PHE:CE1	1:A:217:SER:HB3	2.49	0.47
1:C:274:CYS:HB3	1:C:296:TYR:CZ	2.50	0.46
1:B:330:ASN:OD1	1:B:332:GLU:HG2	2.15	0.46
1:A:281:THR:HG23	1:A:307:PRO:HD3	1.97	0.46
1:A:54:ALA:HB3	1:A:57:GLU:HG3	1.98	0.46
1:A:259:GLU:O	1:A:263:GLN:HG2	2.16	0.46
1:A:364:LYS:HE3	1:A:368:GLU:OE1	2.16	0.46
1:B:133:ARG:HD2	1:B:135:HIS:O	2.16	0.46
1:A:274:CYS:HB3	1:A:296:TYR:CZ	2.50	0.46
1:C:61:LYS:HB3	1:C:61:LYS:HE3	1.76	0.45
1:C:63:LEU:HD11	1:C:105:GLU:HB2	1.99	0.45
1:B:252:GLU:OE1	4:B:600:HOH:O	2.21	0.45
1:C:61:LYS:HG2	4:C:660:HOH:O	2.15	0.45
1:B:60:VAL:HB	1:B:104:PHE:HB3	1.97	0.45
1:A:74:GLN:HA	1:A:79:TYR:HB3	1.99	0.44
1:C:148:ASN:HB2	1:C:149:PRO:CD	2.46	0.44
1:B:324:THR:O	1:B:328:LYS:HB2	2.18	0.44
1:C:139:ASN:O	1:C:142:ASP:HB2	2.17	0.44
1:B:274:CYS:HB3	1:B:296:TYR:CZ	2.53	0.43
1:C:54:ALA:HB1	1:C:55:PRO:CD	2.40	0.43
1:B:140:ASP:OD1	1:B:140:ASP:N	2.51	0.43
1:B:282:GLY:O	1:B:286:LYS:HG2	2.19	0.43
1:C:157:LYS:HZ2	1:C:157:LYS:HB3	1.84	0.43
1:A:171:VAL:HG12	2:A:401:NAP:H5N	2.01	0.42
1:C:146:LEU:CD2	1:C:342:ILE:HD11	2.49	0.42
1:A:300:SER:O	1:A:301:PHE:HB2	2.19	0.42
1:C:157:LYS:HD3	1:C:346:GLU:OE1	2.20	0.42
1:A:60:VAL:HB	1:A:104:PHE:HB3	2.02	0.42
1:A:165:GLN:HG2	1:A:352:ASP:HB3	2.01	0.42
1:C:62:THR:O	1:C:386:TYR:HB2	2.20	0.42
1:C:157:LYS:H	1:C:157:LYS:HG2	1.67	0.42
1:B:27:GLN:HG3	1:B:360:TYR:OH	2.20	0.41
1:B:332:GLU:H	1:B:332:GLU:CD	2.24	0.41
1:C:148:ASN:OD1	1:C:151:GLN:HG3	2.20	0.41
1:A:121:TRP:CH2	1:A:163:ILE:HD13	2.55	0.41
1:B:53:LEU:HD21	1:B:59:ILE:HG13	2.03	0.41
1:A:252:GLU:HG2	4:A:566:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:293:MET:HG2	4:A:579:HOH:O	2.21	0.41
1:C:262:LYS:HD3	4:C:587:HOH:O	2.21	0.41
1:C:281:THR:HG23	1:C:307:PRO:HD3	2.01	0.41
1:B:93:GLU:HG3	1:B:94:PRO:HD2	2.02	0.41
1:C:74:GLN:HA	1:C:79:TYR:HB3	2.01	0.41
4:A:523:HOH:O	1:C:235:LYS:HE3	2.21	0.40
1:C:285:ARG:NH1	4:C:503:HOH:O	2.54	0.40
1:A:149:PRO:HG3	1:A:161:LEU:O	2.22	0.40
1:A:211:LYS:HB2	1:A:238:GLY:HA3	2.02	0.40
1:B:186:LEU:HD12	1:B:215:PHE:CD2	2.57	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	362/364 (100%)	354 (98%)	8 (2%)	0	100	100
1	B	364/364 (100%)	353 (97%)	11 (3%)	0	100	100
1	C	362/364 (100%)	352 (97%)	10 (3%)	0	100	100
All	All	1088/1092 (100%)	1059 (97%)	29 (3%)	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	307/307 (100%)	306 (100%)	1 (0%)	92 89
1	B	309/307 (101%)	305 (99%)	4 (1%)	69 56
1	C	307/307 (100%)	303 (99%)	4 (1%)	69 56
All	All	923/921 (100%)	914 (99%)	9 (1%)	76 67

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

Mol	Chain	Res	Type
1	A	224	ARG
1	B	33	GLN
1	B	157	LYS
1	B	251	ARG
1	B	359	LEU
1	C	53	LEU
1	C	61	LYS
1	C	155	ASN
1	C	157	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

7 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	C	401	-	45,52,52	2.11	10 (22%)	56,80,80	1.33	7 (12%)
3	COO	B	402[B]	-	23,30,55	1.74	3 (13%)	33,42,81	1.72	3 (9%)
2	NAP	A	401	-	45,52,52	2.08	9 (20%)	56,80,80	1.31	7 (12%)
3	COO	B	402[A]	-	23,30,55	1.73	3 (13%)	33,42,81	1.63	4 (12%)
3	COO	A	402	-	23,30,55	1.76	3 (13%)	33,42,81	1.65	4 (12%)
3	COO	C	402	-	9,11,55	1.65	2 (22%)	8,12,81	3.26	3 (37%)
2	NAP	B	401	-	45,52,52	2.08	11 (24%)	56,80,80	1.50	9 (16%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	C	401	-	-	4/31/67/67	0/5/5/5
3	COO	B	402[B]	-	1/1/7/16	13/39/39/70	-
2	NAP	A	401	-	-	2/31/67/67	0/5/5/5
3	COO	B	402[A]	-	1/1/7/16	11/39/39/70	-
3	COO	A	402	-	1/1/7/16	5/39/39/70	-
3	COO	C	402	-	-	4/10/10/70	-
2	NAP	B	401	-	-	5/31/67/67	0/5/5/5

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAP	O4B-C1B	6.38	1.50	1.41
2	C	401	NAP	O4B-C1B	6.30	1.49	1.41
2	B	401	NAP	O4B-C1B	5.64	1.49	1.41
3	A	402	COO	C4-N2	5.24	1.45	1.33
2	A	401	NAP	C2N-N1N	5.21	1.41	1.35
2	C	401	NAP	C7N-N7N	5.16	1.42	1.33
3	B	402[A]	COO	C4-N2	5.16	1.45	1.33
3	B	402[B]	COO	C4-N2	5.15	1.45	1.33
2	A	401	NAP	C7N-N7N	5.02	1.42	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	NAP	C7N-N7N	4.89	1.42	1.33
3	A	402	COO	C1-N1	4.77	1.44	1.33
3	B	402[A]	COO	C1-N1	4.70	1.43	1.33
3	B	402[B]	COO	C1-N1	4.70	1.43	1.33
2	C	401	NAP	C2N-N1N	4.63	1.40	1.35
2	C	401	NAP	O4D-C1D	4.63	1.47	1.41
2	B	401	NAP	O4D-C1D	4.54	1.47	1.41
2	B	401	NAP	C2N-N1N	4.45	1.40	1.35
2	C	401	NAP	C2D-C1D	-4.08	1.47	1.53
2	A	401	NAP	C2D-C1D	-4.06	1.47	1.53
2	A	401	NAP	O4D-C1D	3.90	1.46	1.41
2	B	401	NAP	C6N-N1N	3.87	1.44	1.35
3	C	402	COO	C4-N2	3.83	1.45	1.34
2	C	401	NAP	C6N-N1N	3.61	1.44	1.35
2	B	401	NAP	C2D-C1D	-3.60	1.48	1.53
2	A	401	NAP	C6N-N1N	3.35	1.43	1.35
2	C	401	NAP	C3B-C4B	-2.97	1.45	1.53
2	A	401	NAP	C3B-C4B	-2.85	1.45	1.53
2	B	401	NAP	C3B-C4B	-2.84	1.45	1.53
2	A	401	NAP	C2D-C3D	-2.40	1.46	1.53
2	B	401	NAP	C2D-C3D	-2.35	1.46	1.53
2	C	401	NAP	C2D-C3D	-2.34	1.46	1.53
2	B	401	NAP	C6A-N6A	2.33	1.42	1.34
2	C	401	NAP	C6A-N6A	2.29	1.42	1.34
2	B	401	NAP	C8A-N7A	2.27	1.38	1.34
3	B	402[A]	COO	P1A-O5X	2.20	1.63	1.54
3	B	402[B]	COO	P1A-O5X	2.20	1.63	1.54
3	A	402	COO	P1A-O5X	2.16	1.63	1.54
2	C	401	NAP	C8A-N7A	2.14	1.38	1.34
2	B	401	NAP	C2N-C3N	2.14	1.42	1.39
3	C	402	COO	C3-C4	2.08	1.54	1.50
2	A	401	NAP	C8A-N7A	2.07	1.38	1.34

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	402	COO	C6-S1-C7	7.91	109.74	99.80
3	B	402[B]	COO	C6-S1-C7	7.38	109.07	99.80
3	A	402	COO	C6-S1-C7	6.74	108.26	99.80
3	B	402[A]	COO	C6-S1-C7	6.20	107.59	99.80
2	B	401	NAP	C1B-N9A-C4A	-4.09	119.46	126.64
2	A	401	NAP	N3A-C2A-N1A	-4.04	122.37	128.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	401	NAP	N3A-C2A-N1A	-3.94	122.52	128.68
2	B	401	NAP	N3A-C2A-N1A	-3.69	122.91	128.68
2	B	401	NAP	C4A-C5A-N7A	-3.30	105.96	109.40
2	B	401	NAP	O7N-C7N-N7N	-3.29	117.90	122.58
2	B	401	NAP	C6N-N1N-C2N	-3.23	119.03	121.97
2	A	401	NAP	C1B-N9A-C4A	-3.21	121.00	126.64
2	A	401	NAP	C6N-N1N-C2N	-3.17	119.08	121.97
3	B	402[A]	COO	C10-C9-C8	-3.14	119.15	125.34
3	A	402	COO	C6-C5-N2	-3.07	105.96	112.42
3	A	402	COO	C10-C9-C8	-2.96	119.50	125.34
3	C	402	COO	C3-C4-N2	2.94	121.29	116.09
2	C	401	NAP	C6N-N1N-C2N	-2.91	119.32	121.97
2	B	401	NAP	O7N-C7N-C3N	2.87	123.07	119.63
2	C	401	NAP	C4A-C5A-N7A	-2.80	106.48	109.40
2	B	401	NAP	C2A-N1A-C6A	2.57	123.16	118.75
2	C	401	NAP	C2B-C3B-C4B	2.55	107.54	101.99
3	B	402[B]	COO	C10-C9-C8	-2.52	120.36	125.34
2	A	401	NAP	C4A-C5A-N7A	-2.52	106.77	109.40
3	B	402[A]	COO	C3-C2-N1	-2.45	106.95	111.90
2	C	401	NAP	O2X-P2B-O2B	2.32	116.40	105.99
2	A	401	NAP	C2N-C3N-C4N	2.23	120.78	118.26
3	C	402	COO	C5-N2-C4	-2.22	119.14	122.56
2	A	401	NAP	O4D-C1D-C2D	-2.20	103.72	106.93
2	B	401	NAP	C6N-C5N-C4N	2.18	122.61	119.44
2	C	401	NAP	C1B-N9A-C4A	-2.18	122.81	126.64
3	A	402	COO	O1A-P1A-O3A	2.16	111.88	104.64
2	B	401	NAP	C3N-C2N-N1N	2.14	122.52	120.43
2	A	401	NAP	C2B-C3B-C4B	2.12	106.61	101.99
3	B	402[A]	COO	P2A-O3A-P1A	-2.09	125.67	132.83
3	B	402[B]	COO	P2A-O3A-P1A	-2.09	125.67	132.83
2	C	401	NAP	C3B-C2B-C1B	2.05	106.75	102.89

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	402	COO	C13
3	B	402[A]	COO	C13
3	B	402[B]	COO	C13

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	NAP	C2B-O2B-P2B-O1X
2	B	401	NAP	C2B-O2B-P2B-O1X
2	B	401	NAP	O4D-C1D-N1N-C6N
3	A	402	COO	O4-C7-S1-C6
3	A	402	COO	C8-C7-S1-C6
3	B	402[A]	COO	C14-C11-C13-O1
3	B	402[A]	COO	C12-C11-C13-O1
3	B	402[A]	COO	C15-C11-C13-O1
3	B	402[A]	COO	C14-C11-C13-C1
3	B	402[A]	COO	C12-C11-C13-C1
3	B	402[A]	COO	C15-C11-C13-C1
3	B	402[A]	COO	S1-C7-C8-C9
3	B	402[B]	COO	C14-C11-C13-O1
3	B	402[B]	COO	C12-C11-C13-O1
3	B	402[B]	COO	C15-C11-C13-O1
3	B	402[B]	COO	C14-C11-C13-C1
3	B	402[B]	COO	C12-C11-C13-C1
3	B	402[B]	COO	C15-C11-C13-C1
3	B	402[B]	COO	N2-C5-C6-S1
3	B	402[B]	COO	O4-C7-S1-C6
3	C	402	COO	N2-C5-C6-S1
3	C	402	COO	O4-C7-S1-C6
3	C	402	COO	C8-C7-S1-C6
3	B	402[A]	COO	P1A-O3A-P2A-O4A
3	B	402[B]	COO	P1A-O3A-P2A-O4A
3	B	402[A]	COO	N2-C5-C6-S1
3	B	402[A]	COO	P1A-O3A-P2A-O6A
3	B	402[B]	COO	P1A-O3A-P2A-O6A
3	B	402[B]	COO	C8-C7-S1-C6
2	C	401	NAP	C2B-O2B-P2B-O1X
3	A	402	COO	P1A-O3A-P2A-O4A
3	A	402	COO	N2-C5-C6-S1
3	B	402[A]	COO	O4-C7-C8-C9
2	B	401	NAP	O4B-C4B-C5B-O5B
2	C	401	NAP	O4B-C4B-C5B-O5B
3	B	402[B]	COO	C6-C5-N2-C4
3	A	402	COO	C5-C6-S1-C7
3	C	402	COO	C5-C6-S1-C7
3	B	402[B]	COO	C3-C2-N1-C1
2	B	401	NAP	PA-O3-PN-O1N
2	B	401	NAP	PA-O3-PN-O2N
2	C	401	NAP	PN-O3-PA-O2A
2	A	401	NAP	O4B-C4B-C5B-O5B

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Mol	Chain	Res	Type	Atoms
2	C	401	NAP	C3B-C4B-C5B-O5B

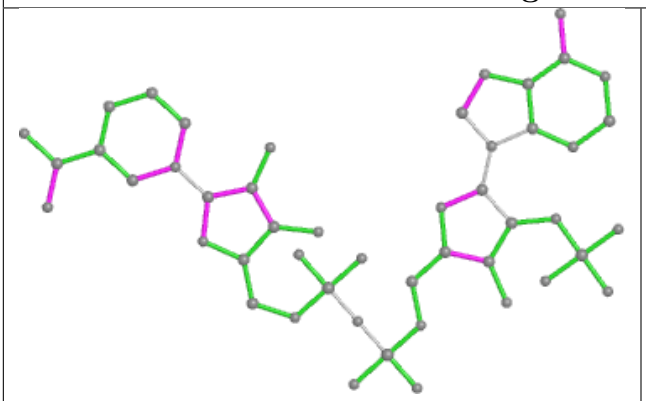
There are no ring outliers.

5 monomers are involved in 6 short contacts:

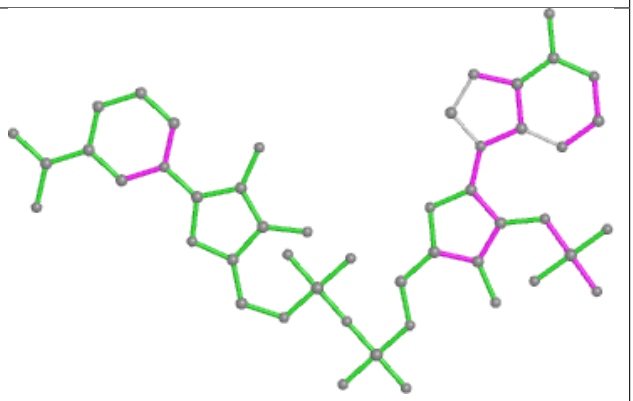
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	NAP	1	0
3	B	402[B]	COO	1	0
2	A	401	NAP	2	0
3	C	402	COO	1	0
2	B	401	NAP	1	0

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 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.

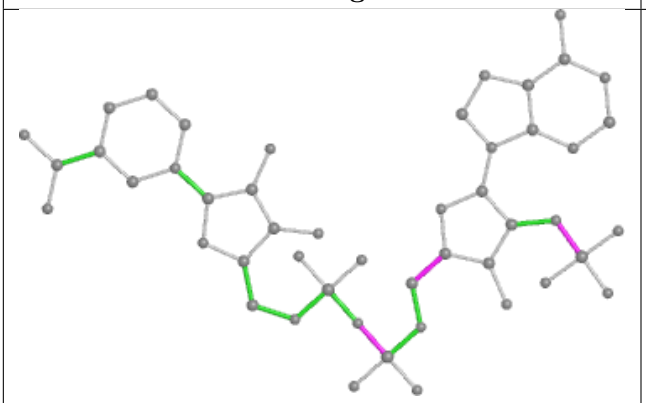
Ligand NAP C 401



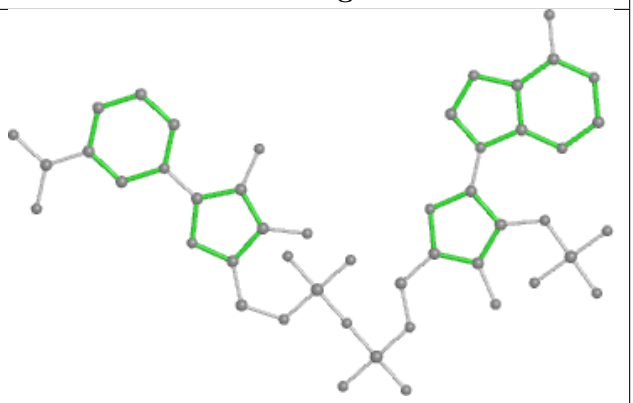
Bond lengths



Bond angles

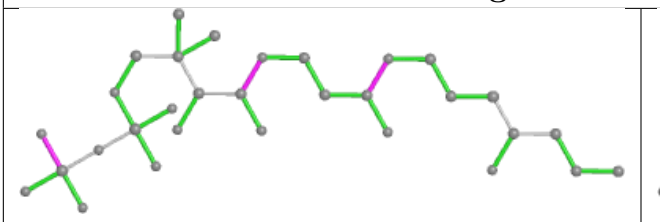


Torsions



Rings

Ligand COO B 402 (B)



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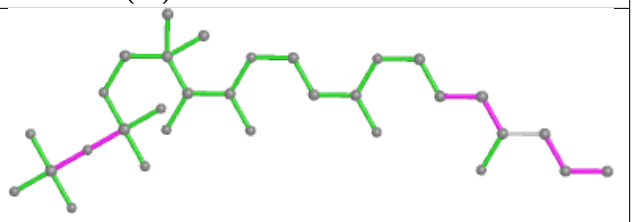

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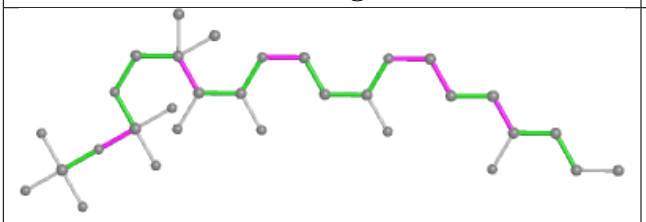

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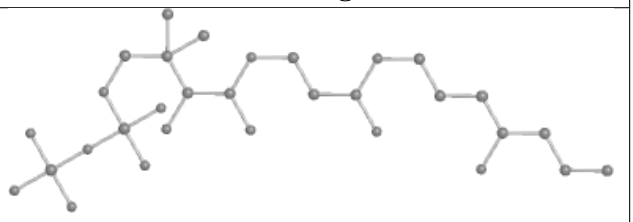
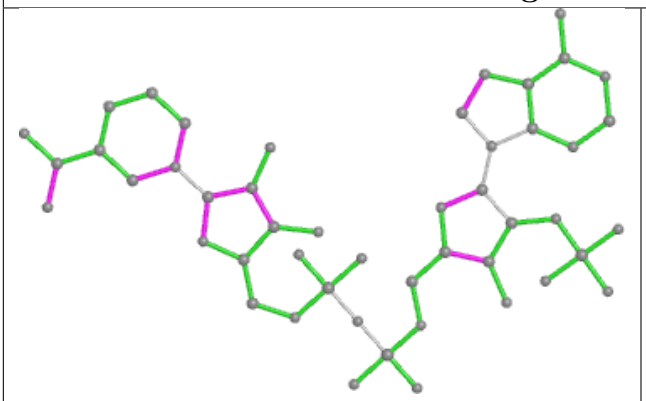
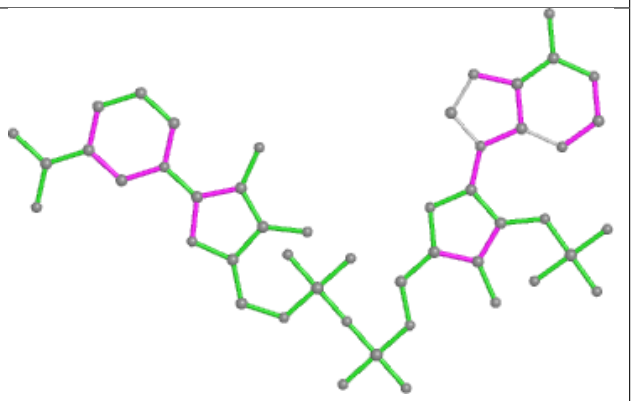


Rings

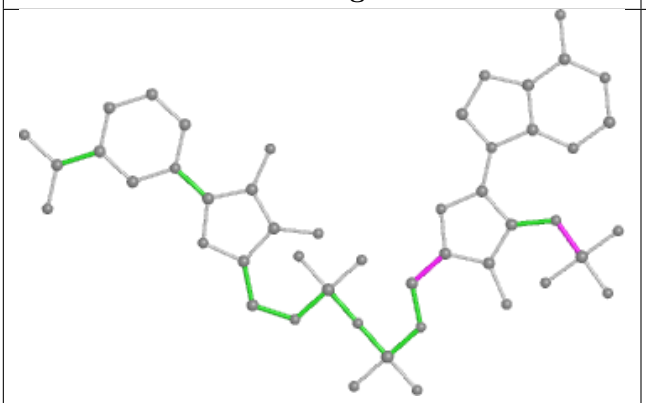
Ligand NAP A 401



Bond lengths



Bond angles



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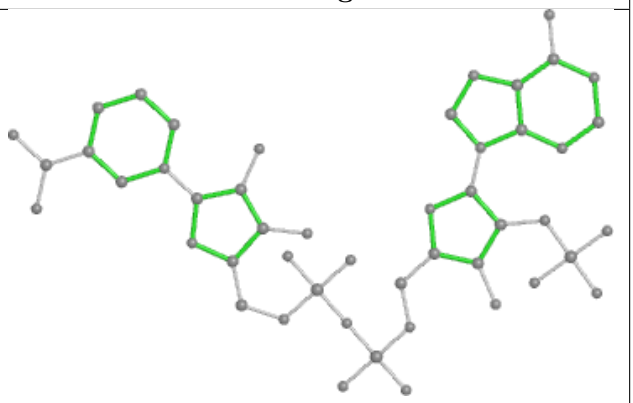
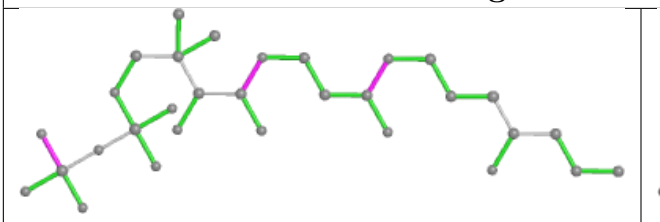
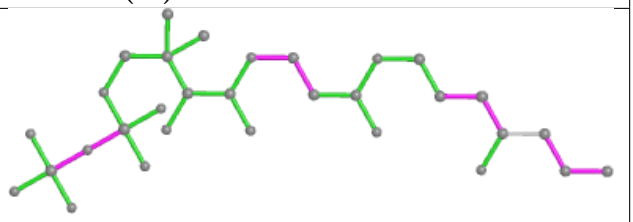


Rings

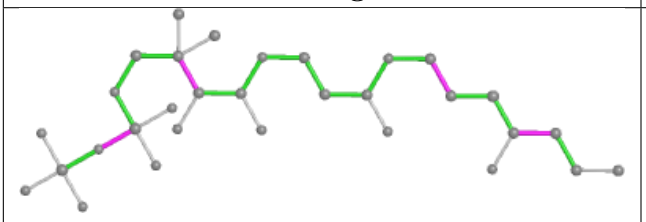
Ligand COO B 402 (A)



Bond lengths



Bond angles



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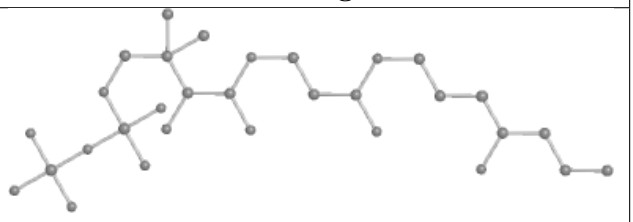
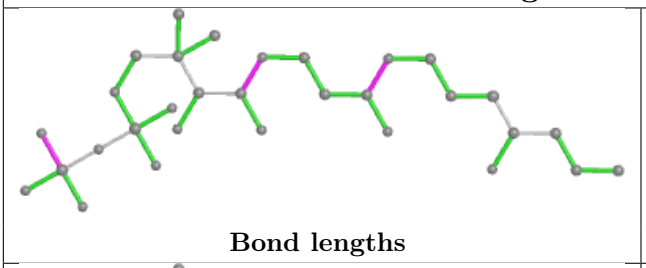


Rings

Ligand COO A 402



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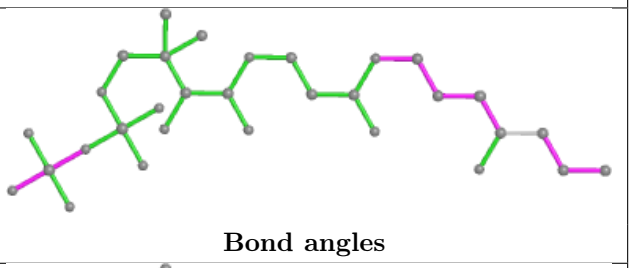
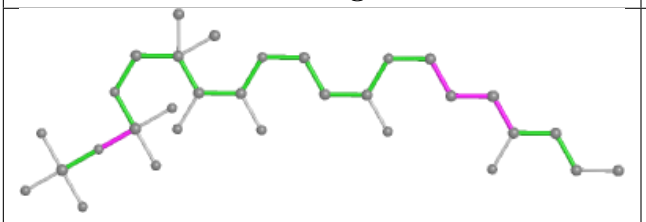


Bond angles



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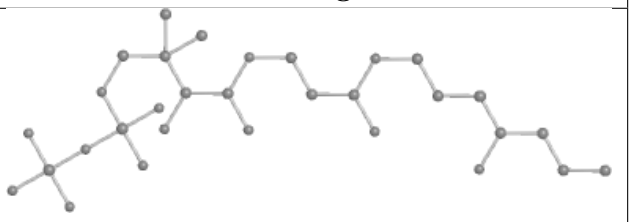
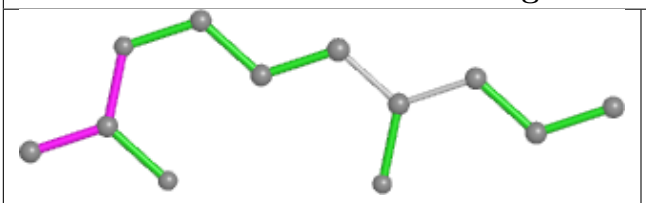
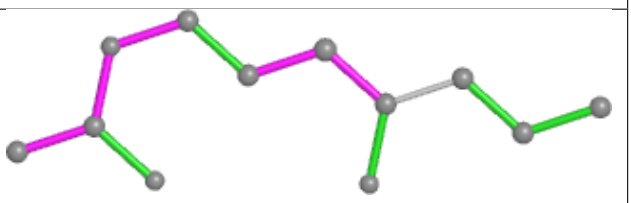


Rings

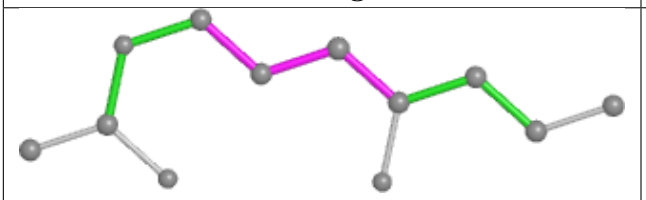
Ligand COO C 402



Bond lengths



Bond angles



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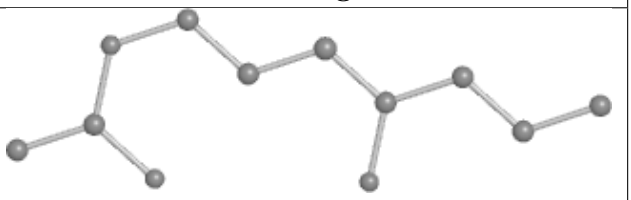
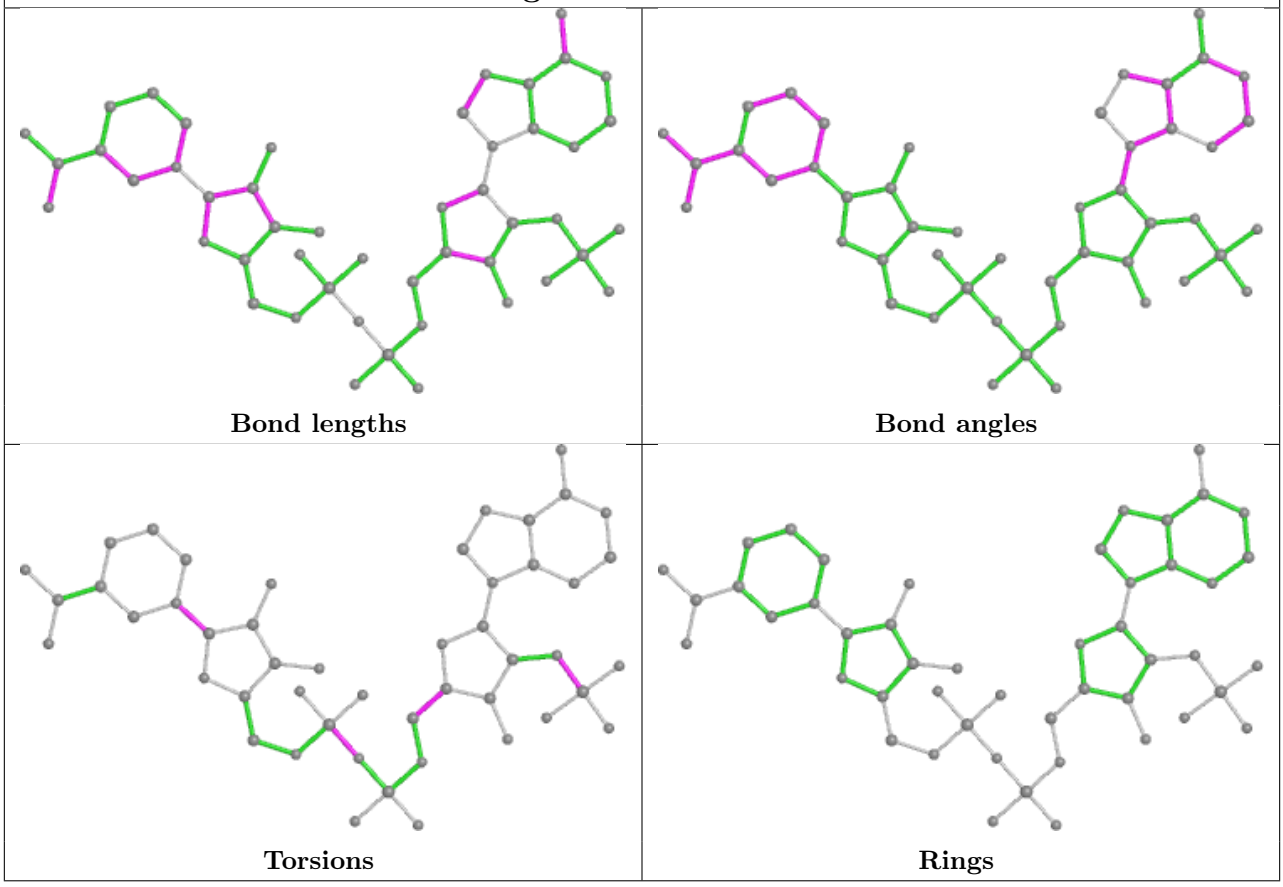


Rings

Ligand NAP B 401



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, 95^{th} 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	364/364 (100%)	0.37	27 (7%) 14 16	34, 43, 71, 101	0
1	B	364/364 (100%)	0.58	33 (9%) 9 10	33, 47, 91, 159	0
1	C	364/364 (100%)	0.58	49 (13%) 3 3	30, 47, 97, 129	0
All	All	1092/1092 (100%)	0.51	109 (9%) 7 8	30, 45, 89, 159	0

All (109) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	111	SER	7.6
1	B	53	LEU	7.6
1	C	109	VAL	7.0
1	C	53	LEU	6.9
1	C	112	ASN	6.8
1	B	54	ALA	6.3
1	C	113	VAL	6.1
1	B	52	ASN	6.1
1	B	109	VAL	5.9
1	C	55	PRO	5.4
1	A	272	LEU	5.2
1	C	56	ASN	5.1
1	B	23	MET	5.0
1	B	51	ASP	4.9
1	C	54	ALA	4.8
1	A	54	ALA	4.7
1	C	110	GLY	4.6
1	C	50	ASP	4.5
1	C	116	LEU	4.5
1	B	78	VAL	4.5
1	C	108	LYS	4.4
1	C	154	ALA	4.3
1	B	301	PHE	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	56	ASN	4.1
1	B	110	GLY	4.1
1	C	150	ALA	4.0
1	B	113	VAL	4.0
1	A	52	ASN	4.0
1	C	107	ILE	4.0
1	A	53	LEU	3.9
1	B	112	ASN	3.8
1	A	271	ALA	3.7
1	B	116	LEU	3.7
1	C	49	ASP	3.6
1	C	363	THR	3.6
1	C	359	LEU	3.6
1	C	114	SER	3.6
1	C	25	THR	3.6
1	A	51	ASP	3.4
1	B	55	PRO	3.4
1	C	58	VAL	3.4
1	B	111	SER	3.3
1	C	153	LYS	3.3
1	A	55	PRO	3.2
1	A	274	CYS	3.2
1	C	139	ASN	3.2
1	A	363	THR	3.2
1	C	51	ASP	3.1
1	A	220	VAL	3.1
1	C	117	GLU	3.0
1	C	156	GLY	3.0
1	B	171	VAL	3.0
1	C	171	VAL	3.0
1	C	118	ALA	2.9
1	A	283	ILE	2.9
1	C	115	SER	2.9
1	C	140	ASP	2.9
1	A	293	MET	2.9
1	A	236	GLU	2.9
1	B	153	LYS	2.9
1	A	23	MET	2.9
1	B	225	PRO	2.9
1	C	272	LEU	2.8
1	B	75	ILE	2.7
1	B	92	THR	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	59	ILE	2.7
1	A	294	LEU	2.7
1	C	24	ILE	2.6
1	A	296	TYR	2.6
1	B	251	ARG	2.6
1	A	176	ALA	2.6
1	C	48	ILE	2.6
1	A	329	ASN	2.5
1	B	49	ASP	2.5
1	B	328	LYS	2.5
1	B	137	LEU	2.5
1	C	297	GLY	2.5
1	C	106	VAL	2.4
1	A	56	ASN	2.4
1	C	158	PRO	2.4
1	A	284	ALA	2.4
1	C	175	THR	2.3
1	A	180	LEU	2.3
1	B	226	ASN	2.3
1	C	187	THR	2.3
1	C	294	LEU	2.3
1	C	169	ILE	2.2
1	B	108	LYS	2.2
1	A	218	ILE	2.2
1	B	82	LYS	2.2
1	C	52	ASN	2.2
1	B	107	ILE	2.2
1	A	229	GLU	2.1
1	C	57	GLU	2.1
1	B	115	SER	2.1
1	C	362	GLY	2.1
1	C	174	LEU	2.1
1	A	194	ILE	2.1
1	A	297	GLY	2.1
1	B	59	ILE	2.1
1	A	196	ASN	2.1
1	A	280	SER	2.1
1	B	271	ALA	2.1
1	C	176	ALA	2.1
1	B	58	VAL	2.0
1	C	321	PHE	2.0
1	B	359	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	274	CYS	2.0
1	C	23	MET	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)

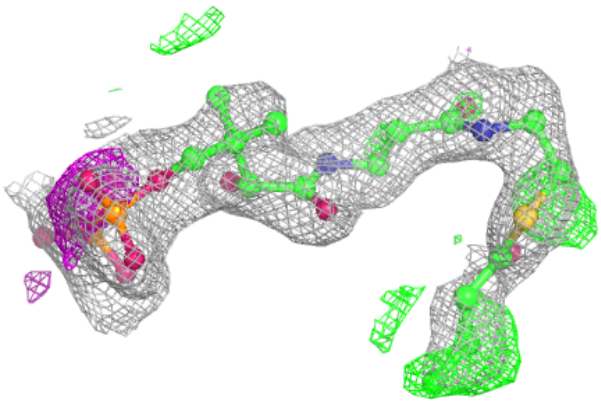
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

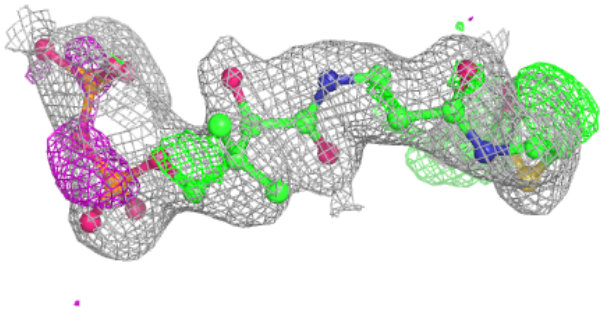
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	COO	B	402[A]	31/53	0.58	0.34	70,78,89,94	12
3	COO	B	402[B]	31/53	0.58	0.34	65,78,89,94	12
3	COO	A	402	31/53	0.80	0.18	68,72,77,86	0
3	COO	C	402	12/53	0.85	0.23	71,73,78,88	0
2	NAP	B	401	48/48	0.96	0.10	37,43,55,59	0
2	NAP	C	401	48/48	0.96	0.13	34,43,63,80	0
2	NAP	A	401	48/48	0.96	0.11	33,43,53,56	0

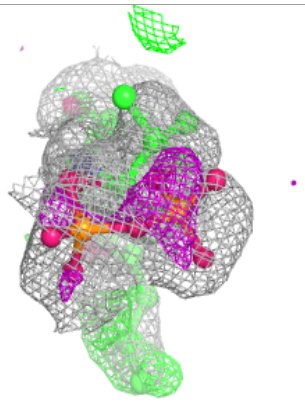
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around COO B 402 (A):

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

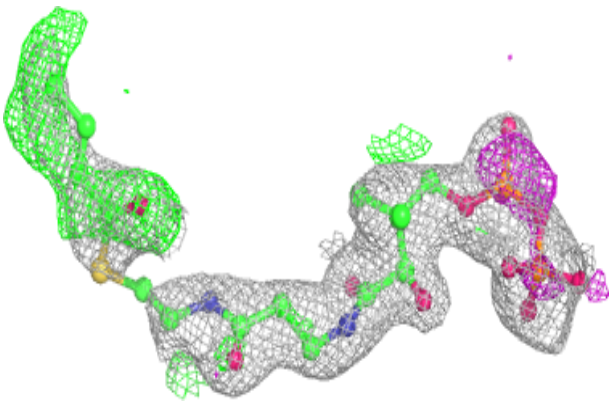


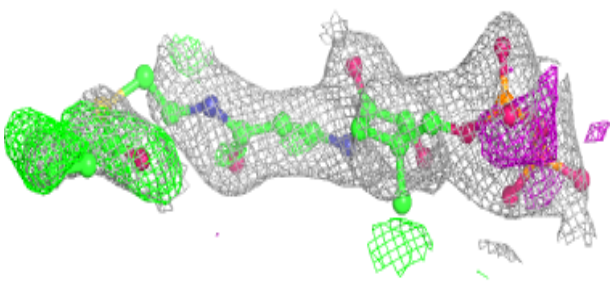


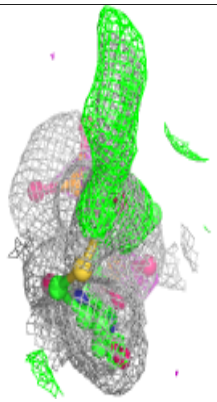


Electron density around COO B 402 (B):

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

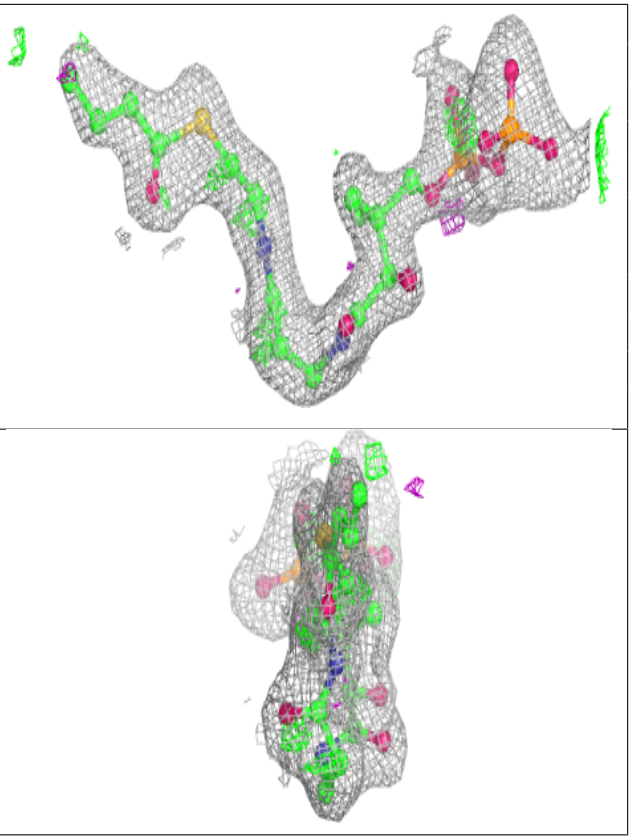






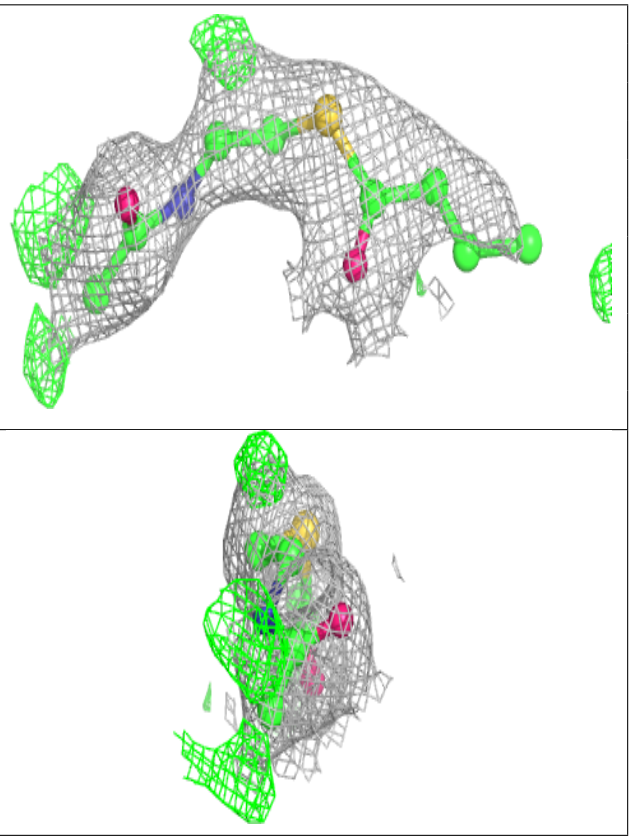
Electron density around COO A 402:

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



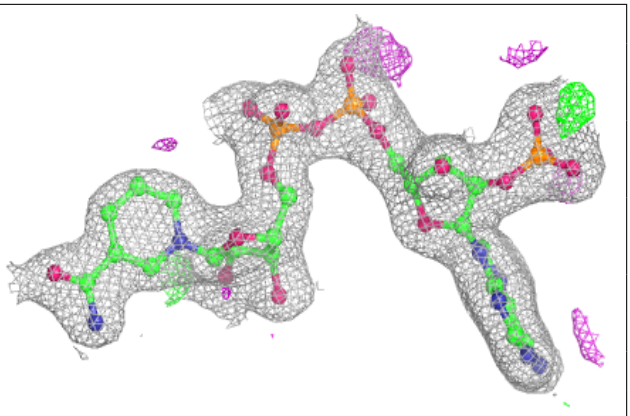
Electron density around COO C 402:

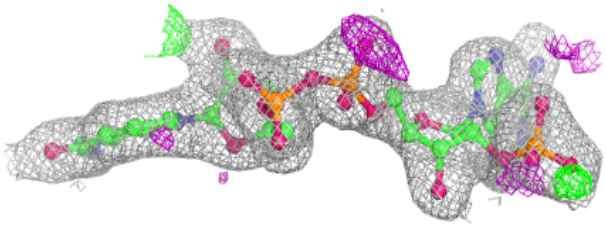
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

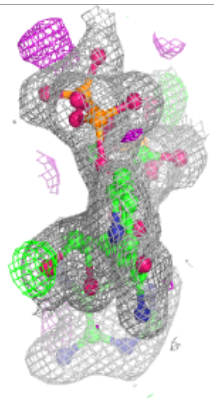


Electron density around NAP B 401:

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

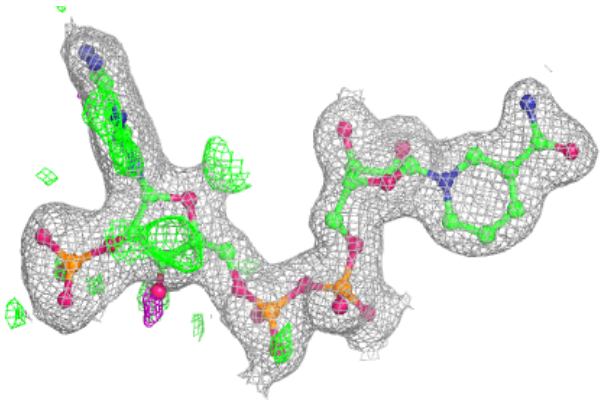


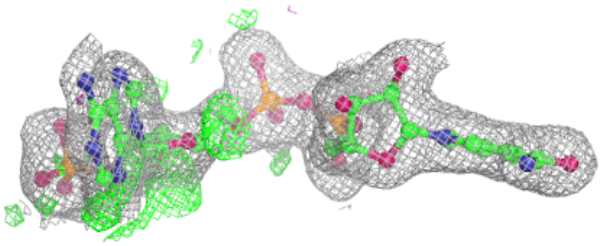


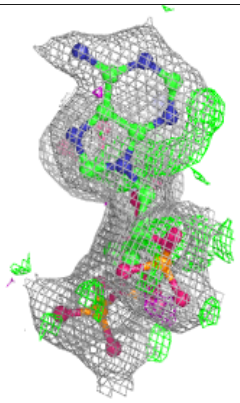


Electron density around NAP C 401:

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

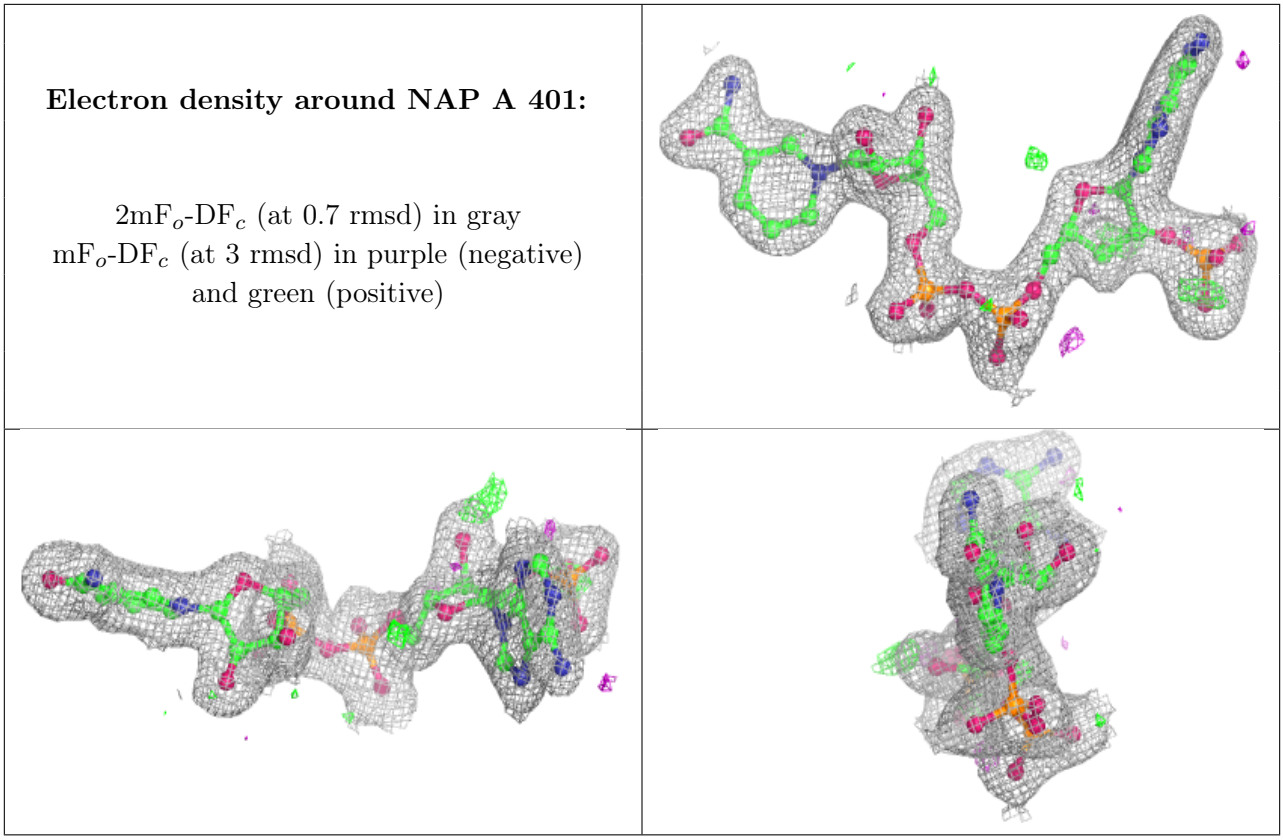






Electron density around NAP A 401:

$2mF_o$-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.