



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:43 am BST

PDB ID : 6NJ7
Title : 11-BETA DEHYDROGENASE ISOZYME 1 IN COMPLEX WITH COL-
LETOIC ACID
Authors : Miller, D.J.; Rivas, F.
Deposited on : 2019-01-02
Resolution : 2.60 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

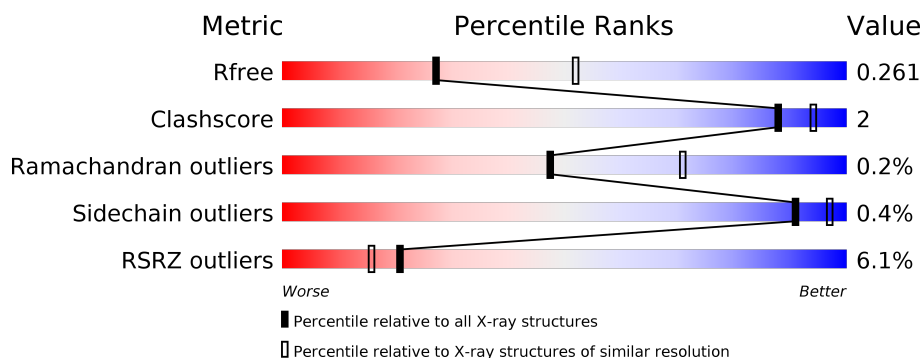
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 ≥ 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	289	<div> <div>6%</div> <div> <div></div> <div>77%</div> <div>7%</div> <div>16%</div> </div> </div>
1	B	289	<div> <div>7%</div> <div> <div></div> <div>85%</div> <div>•</div> <div>11%</div> </div> </div>
1	C	289	<div> <div>2%</div> <div> <div></div> <div>83%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	289	<div> <div>6%</div> <div> <div></div> <div>80%</div> <div>6%</div> <div>14%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7396 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 Corticosteroid 11-beta-dehydrogenase isozyme 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1734	1095	293	333	13			
1	B	257	Total	C	N	O	S	0	0	0
			1844	1164	314	353	13			
1	C	250	Total	C	N	O	S	0	0	0
			1805	1145	310	337	13			
1	D	248	Total	C	N	O	S	0	0	0
			1745	1103	300	332	10			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	MET	-	initiating methionine	UNP P28845
A	5	LYS	-	expression tag	UNP P28845
A	6	HIS	-	expression tag	UNP P28845
A	7	GLN	-	expression tag	UNP P28845
A	8	HIS	-	expression tag	UNP P28845
A	9	GLN	-	expression tag	UNP P28845
A	10	HIS	-	expression tag	UNP P28845
A	11	GLN	-	expression tag	UNP P28845
A	12	HIS	-	expression tag	UNP P28845
A	13	GLN	-	expression tag	UNP P28845
A	14	HIS	-	expression tag	UNP P28845
A	15	GLN	-	expression tag	UNP P28845
A	16	HIS	-	expression tag	UNP P28845
A	17	GLN	-	expression tag	UNP P28845
A	18	GLN	-	expression tag	UNP P28845
A	19	PRO	-	expression tag	UNP P28845
A	20	LEU	-	expression tag	UNP P28845
A	21	GLN	-	expression tag	UNP P28845
A	22	PRO	-	expression tag	UNP P28845
A	23	LEU	-	expression tag	UNP P28845
A	272	SER	CYS	engineered mutation	UNP P28845

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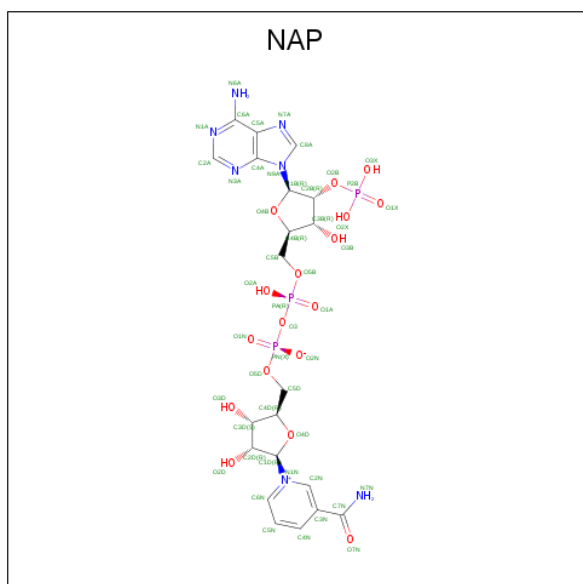
Chain	Residue	Modelled	Actual	Comment	Reference
B	4	MET	-	initiating methionine	UNP P28845
B	5	LYS	-	expression tag	UNP P28845
B	6	HIS	-	expression tag	UNP P28845
B	7	GLN	-	expression tag	UNP P28845
B	8	HIS	-	expression tag	UNP P28845
B	9	GLN	-	expression tag	UNP P28845
B	10	HIS	-	expression tag	UNP P28845
B	11	GLN	-	expression tag	UNP P28845
B	12	HIS	-	expression tag	UNP P28845
B	13	GLN	-	expression tag	UNP P28845
B	14	HIS	-	expression tag	UNP P28845
B	15	GLN	-	expression tag	UNP P28845
B	16	HIS	-	expression tag	UNP P28845
B	17	GLN	-	expression tag	UNP P28845
B	18	GLN	-	expression tag	UNP P28845
B	19	PRO	-	expression tag	UNP P28845
B	20	LEU	-	expression tag	UNP P28845
B	21	GLN	-	expression tag	UNP P28845
B	22	PRO	-	expression tag	UNP P28845
B	23	LEU	-	expression tag	UNP P28845
B	272	SER	CYS	engineered mutation	UNP P28845
C	4	MET	-	initiating methionine	UNP P28845
C	5	LYS	-	expression tag	UNP P28845
C	6	HIS	-	expression tag	UNP P28845
C	7	GLN	-	expression tag	UNP P28845
C	8	HIS	-	expression tag	UNP P28845
C	9	GLN	-	expression tag	UNP P28845
C	10	HIS	-	expression tag	UNP P28845
C	11	GLN	-	expression tag	UNP P28845
C	12	HIS	-	expression tag	UNP P28845
C	13	GLN	-	expression tag	UNP P28845
C	14	HIS	-	expression tag	UNP P28845
C	15	GLN	-	expression tag	UNP P28845
C	16	HIS	-	expression tag	UNP P28845
C	17	GLN	-	expression tag	UNP P28845
C	18	GLN	-	expression tag	UNP P28845
C	19	PRO	-	expression tag	UNP P28845
C	20	LEU	-	expression tag	UNP P28845
C	21	GLN	-	expression tag	UNP P28845
C	22	PRO	-	expression tag	UNP P28845
C	23	LEU	-	expression tag	UNP P28845
C	272	SER	CYS	engineered mutation	UNP P28845

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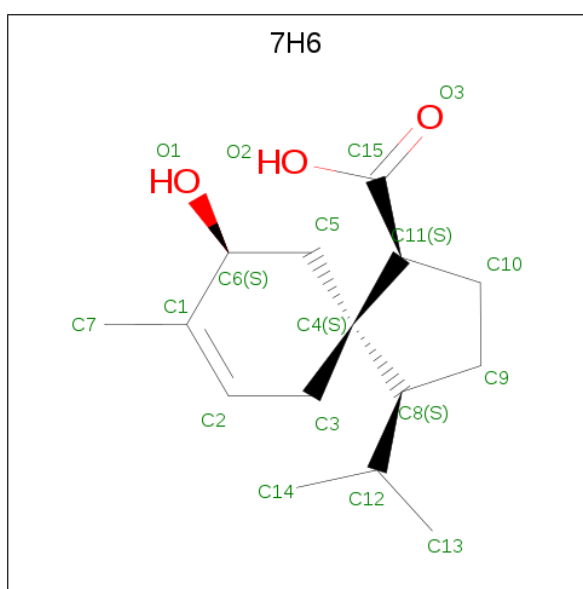
Chain	Residue	Modelled	Actual	Comment	Reference
D	4	MET	-	initiating methionine	UNP P28845
D	5	LYS	-	expression tag	UNP P28845
D	6	HIS	-	expression tag	UNP P28845
D	7	GLN	-	expression tag	UNP P28845
D	8	HIS	-	expression tag	UNP P28845
D	9	GLN	-	expression tag	UNP P28845
D	10	HIS	-	expression tag	UNP P28845
D	11	GLN	-	expression tag	UNP P28845
D	12	HIS	-	expression tag	UNP P28845
D	13	GLN	-	expression tag	UNP P28845
D	14	HIS	-	expression tag	UNP P28845
D	15	GLN	-	expression tag	UNP P28845
D	16	HIS	-	expression tag	UNP P28845
D	17	GLN	-	expression tag	UNP P28845
D	18	GLN	-	expression tag	UNP P28845
D	19	PRO	-	expression tag	UNP P28845
D	20	LEU	-	expression tag	UNP P28845
D	21	GLN	-	expression tag	UNP P28845
D	22	PRO	-	expression tag	UNP P28845
D	23	LEU	-	expression tag	UNP P28845
D	272	SER	CYS	engineered mutation	UNP P28845

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $\text{C}_{21}\text{H}_{28}\text{N}_7\text{O}_{17}\text{P}_3$).



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		
2	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
2	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is (1S,4S,5S,9S)-9-hydroxy-8-methyl-4-(propan-2-yl)spiro[4.5]dec-7-ene-1-carboxylic acid (three-letter code: 7H6) (formula: C₁₅H₂₄O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			18	15	3		
3	B	1	Total	C	O	0	0
			18	15	3		
3	C	1	Total	C	O	0	0
			18	15	3		
3	D	1	Total	C	O	0	0
			18	15	3		

- Molecule 4 is water.

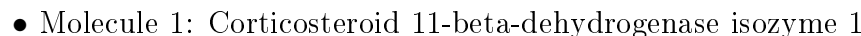
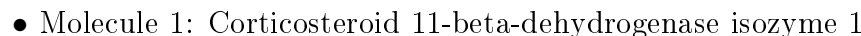
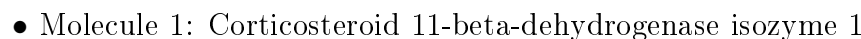
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	O	0	0
			2	2		

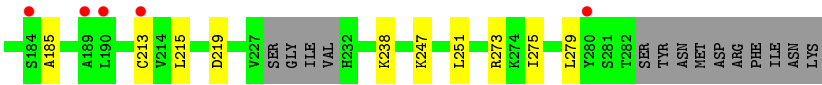
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	2	Total	O	0	0
			2	2		

- Molecule 1: Corticosteroid 11-beta-dehydrogenase isozyme 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.02Å 153.75Å 74.05Å 90.00° 93.24° 90.00°	Depositor
Resolution (Å)	76.87 – 2.60 29.98 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.3 (76.87-2.60) 98.4 (29.98-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0123	Depositor
R, R_{free}	0.213 , 0.264 0.213 , 0.261	Depositor DCC
R_{free} test set	1930 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	85.3	Xtriage
Anisotropy	0.434	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 82.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7396	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAP, 7H6

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	0/1758	0.61	0/2388
1	B	0.42	0/1870	0.62	0/2537
1	C	0.44	0/1830	0.63	0/2482
1	D	0.44	0/1770	0.62	0/2410
All	All	0.43	0/7228	0.62	0/9817

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	1734	0	1654	12	0
1	B	1844	0	1775	7	0
1	C	1805	0	1766	6	0
1	D	1745	0	1628	9	0
2	A	48	0	25	1	0
2	B	48	0	25	1	0
2	C	48	0	25	0	0
2	D	48	0	25	0	0
3	A	18	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	18	0	0	2	0
3	C	18	0	0	0	0
3	D	18	0	0	0	0
4	B	2	0	0	0	0
4	D	2	0	0	0	0
All	All	7396	0	6923	29	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 2.

All (29) 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:223:ALA:HB2	2:A:500:NAP:H72N	1.63	0.63
1:A:149:VAL:HG22	1:B:133:ILE:HD13	1.85	0.57
1:C:193:PHE:HB2	1:D:185:ALA:HB2	1.85	0.57
1:A:197:ILE:HA	1:A:200:GLU:HB2	1.87	0.57
1:B:180:VAL:HG11	3:B:501:7H6:C14	2.38	0.53
1:A:39:VAL:HG12	1:A:42:ALA:HB2	1.92	0.50
1:B:42:ALA:HB3	1:B:63:VAL:HB	1.93	0.50
1:D:219:ASP:OD1	1:D:238:LYS:N	2.43	0.50
1:A:185:ALA:HB2	1:B:193:PHE:HB2	1.93	0.49
1:A:194:PHE:HA	1:A:197:ILE:HG12	1.96	0.48
1:A:219:ASP:OD1	1:A:238:LYS:N	2.42	0.48
1:C:37:VAL:HG11	1:C:54:LEU:HD13	1.98	0.46
1:D:39:VAL:HG12	1:D:42:ALA:HB2	1.98	0.46
1:A:34:GLY:N	1:A:58:GLY:O	2.49	0.45
1:C:133:ILE:HD13	1:D:149:VAL:HG22	1.98	0.44
1:D:37:VAL:HG22	1:D:115:MET:HB3	2.01	0.43
1:C:175:VAL:HG12	1:D:273:ARG:HG3	2.01	0.43
1:A:155:LEU:HB3	1:A:156:PRO:HD3	2.02	0.42
1:D:213:CYS:HB3	1:D:215:LEU:HD21	2.01	0.42
1:C:37:VAL:HG11	1:C:54:LEU:CD1	2.50	0.42
1:D:247:LYS:O	1:D:251:LEU:HG	2.20	0.42
1:A:56:LYS:HA	1:A:81:LEU:O	2.20	0.41
1:B:36:LYS:HA	1:B:60:HIS:HB2	2.03	0.41
1:D:275:ILE:HG22	1:D:279:LEU:HD11	2.02	0.41
1:A:193:PHE:HB2	1:B:185:ALA:HB2	2.02	0.41
1:C:64:THR:HB	1:C:102:PHE:CE1	2.55	0.41
2:B:500:NAP:C5N	3:B:501:7H6:C15	2.99	0.40
1:B:96:MET:CE	1:B:141:GLU:HG3	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:ASP:C	1:A:261:SER:H	2.25	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	239/289 (83%)	225 (94%)	13 (5%)	1 (0%)	34	57
1	B	255/289 (88%)	244 (96%)	11 (4%)	0	100	100
1	C	246/289 (85%)	228 (93%)	18 (7%)	0	100	100
1	D	242/289 (84%)	223 (92%)	18 (7%)	1 (0%)	34	57
All	All	982/1156 (85%)	920 (94%)	60 (6%)	2 (0%)	47	71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	65	ALA
1	A	260	SER

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	170/246 (69%)	170 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	182/246 (74%)	181 (100%)	1 (0%)	88	96
1	C	180/246 (73%)	178 (99%)	2 (1%)	73	88
1	D	164/246 (67%)	164 (100%)	0	100	100
All	All	696/984 (71%)	693 (100%)	3 (0%)	91	97

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	145	LEU
1	C	137	ARG
1	C	260	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	162	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	D	500	-	45,52,52	0.65	0	56,80,80	0.89	3 (5%)
3	7H6	B	501	-	16,19,19	0.80	1 (6%)	17,29,29	2.98	9 (52%)
3	7H6	D	501	-	16,19,19	0.82	1 (6%)	17,29,29	3.40	7 (41%)
3	7H6	C	501	-	16,19,19	0.89	0	17,29,29	3.25	9 (52%)
3	7H6	A	501	-	16,19,19	0.81	1 (6%)	17,29,29	3.41	6 (35%)
2	NAP	B	500	-	45,52,52	0.71	1 (2%)	56,80,80	1.05	3 (5%)
2	NAP	A	500	-	45,52,52	0.63	0	56,80,80	0.88	2 (3%)
2	NAP	C	500	-	45,52,52	0.68	0	56,80,80	0.94	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	D	500	-	-	7/31/67/67	0/5/5/5
3	7H6	B	501	-	-	0/4/40/40	0/2/2/2
3	7H6	D	501	-	-	0/4/40/40	0/2/2/2
3	7H6	C	501	-	-	0/4/40/40	0/2/2/2
3	7H6	A	501	-	-	0/4/40/40	0/2/2/2
2	NAP	B	500	-	-	2/31/67/67	0/5/5/5
2	NAP	A	500	-	-	7/31/67/67	0/5/5/5
2	NAP	C	500	-	-	4/31/67/67	0/5/5/5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	501	7H6	C3-C4	2.42	1.58	1.54
3	B	501	7H6	C3-C4	2.26	1.58	1.54
2	B	500	NAP	C8A-N7A	-2.06	1.31	1.34
3	A	501	7H6	C3-C4	2.03	1.57	1.54

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	7H6	C7-C1-C6	9.07	122.24	115.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	501	7H6	C7-C1-C6	9.06	122.22	115.93
3	A	501	7H6	C6-C1-C2	-7.88	115.39	121.27
3	B	501	7H6	C7-C1-C6	7.58	121.20	115.93
3	C	501	7H6	C6-C1-C2	-7.17	115.92	121.27
3	D	501	7H6	C6-C1-C2	-7.16	115.93	121.27
3	C	501	7H6	C7-C1-C6	7.03	120.82	115.93
3	D	501	7H6	C9-C8-C4	5.81	108.98	104.22
3	C	501	7H6	C9-C8-C4	5.60	108.81	104.22
3	B	501	7H6	C6-C1-C2	-5.43	117.22	121.27
3	B	501	7H6	C9-C8-C4	5.35	108.60	104.22
3	A	501	7H6	C9-C8-C4	4.67	108.04	104.22
2	B	500	NAP	C3N-C7N-N7N	3.95	122.49	117.75
3	C	501	7H6	O1-C6-C5	3.45	116.89	109.88
2	B	500	NAP	O7N-C7N-N7N	-2.84	118.54	122.58
2	C	500	NAP	C3N-C7N-N7N	2.79	121.10	117.75
3	C	501	7H6	C3-C4-C8	2.66	118.04	111.42
3	A	501	7H6	O1-C6-C5	2.56	115.09	109.88
2	A	500	NAP	O7N-C7N-N7N	-2.54	118.97	122.58
2	D	500	NAP	O7N-C7N-N7N	-2.52	119.00	122.58
3	C	501	7H6	C9-C8-C12	-2.52	108.40	113.66
2	D	500	NAP	C3N-C7N-N7N	2.46	120.70	117.75
2	B	500	NAP	O2N-PN-O1N	2.44	124.29	112.24
3	C	501	7H6	C4-C8-C12	-2.40	116.45	119.55
3	D	501	7H6	C3-C4-C8	2.38	117.34	111.42
2	C	500	NAP	O7N-C7N-N7N	-2.35	119.23	122.58
3	C	501	7H6	C3-C2-C1	-2.34	119.48	124.30
3	B	501	7H6	C3-C2-C1	-2.31	119.54	124.30
3	B	501	7H6	C4-C8-C12	-2.29	116.59	119.55
3	D	501	7H6	C10-C11-C4	2.27	107.93	103.67
3	D	501	7H6	C3-C2-C1	-2.25	119.66	124.30
2	C	500	NAP	PN-O3-PA	-2.24	125.15	132.83
3	B	501	7H6	C14-C12-C8	-2.22	106.47	112.42
3	C	501	7H6	C14-C12-C8	-2.21	106.50	112.42
3	A	501	7H6	C3-C4-C8	2.19	116.87	111.42
2	A	500	NAP	O2N-PN-O1N	2.17	122.97	112.24
3	B	501	7H6	C3-C4-C8	2.17	116.81	111.42
3	A	501	7H6	C9-C8-C12	-2.15	109.15	113.66
2	D	500	NAP	O2N-PN-O1N	2.09	122.57	112.24
3	B	501	7H6	O1-C6-C5	2.07	114.08	109.88
3	B	501	7H6	C10-C11-C4	2.04	107.50	103.67
3	D	501	7H6	C4-C8-C12	-2.02	116.95	119.55

There are no chirality outliers.

All (20) torsion outliers are listed below:

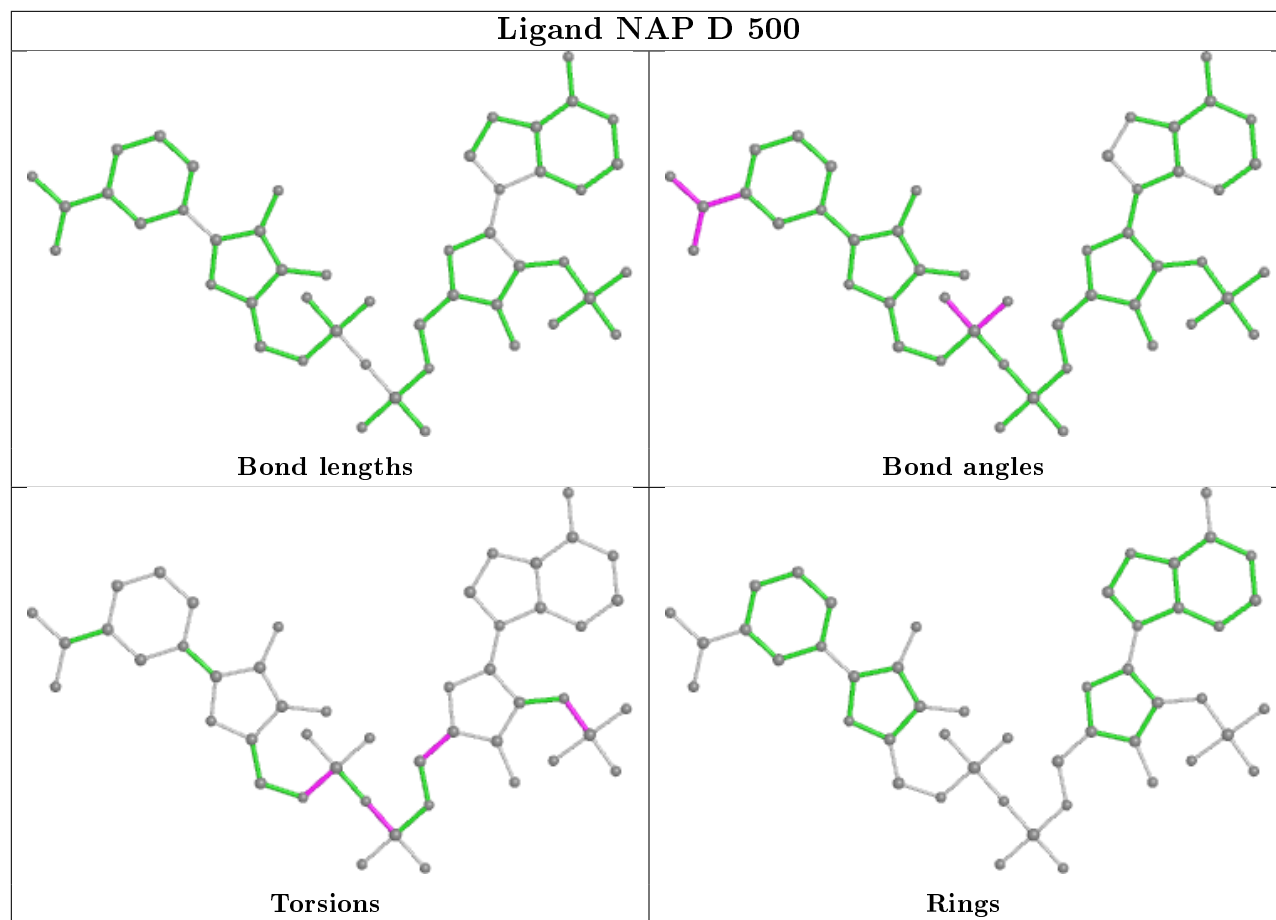
Mol	Chain	Res	Type	Atoms
2	D	500	NAP	C2B-O2B-P2B-O1X
2	D	500	NAP	C5D-O5D-PN-O1N
2	B	500	NAP	C2B-O2B-P2B-O1X
2	A	500	NAP	C2B-O2B-P2B-O1X
2	A	500	NAP	C5D-O5D-PN-O1N
2	A	500	NAP	O4D-C1D-N1N-C2N
2	C	500	NAP	C2B-O2B-P2B-O2X
2	C	500	NAP	O4D-C4D-C5D-O5D
2	C	500	NAP	C3D-C4D-C5D-O5D
2	D	500	NAP	C5D-O5D-PN-O3
2	A	500	NAP	C5D-O5D-PN-O3
2	D	500	NAP	C5D-O5D-PN-O2N
2	A	500	NAP	C5D-O5D-PN-O2N
2	D	500	NAP	PN-O3-PA-O1A
2	A	500	NAP	C2B-O2B-P2B-O3X
2	B	500	NAP	O4B-C4B-C5B-O5B
2	A	500	NAP	O4B-C4B-C5B-O5B
2	D	500	NAP	PN-O3-PA-O2A
2	D	500	NAP	O4B-C4B-C5B-O5B
2	C	500	NAP	O4B-C4B-C5B-O5B

There are no ring outliers.

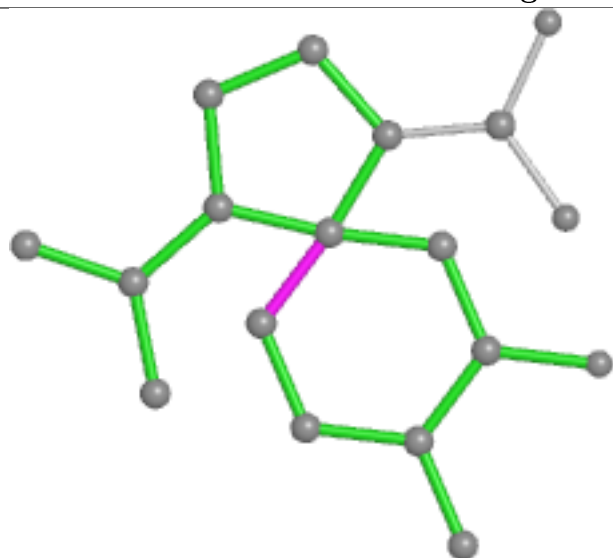
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	7H6	2	0
2	B	500	NAP	1	0
2	A	500	NAP	1	0

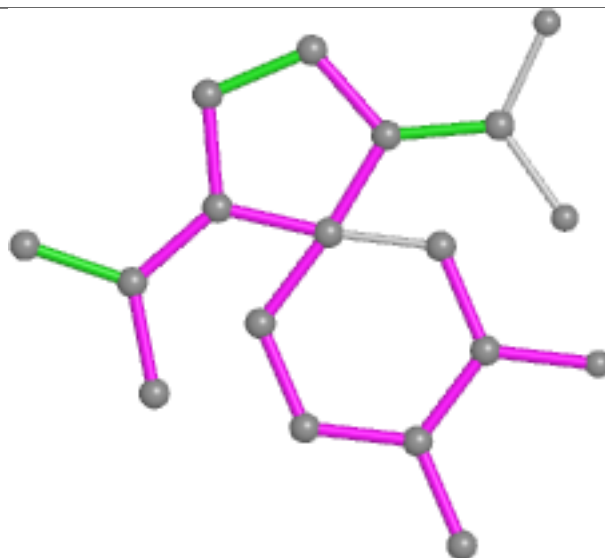
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



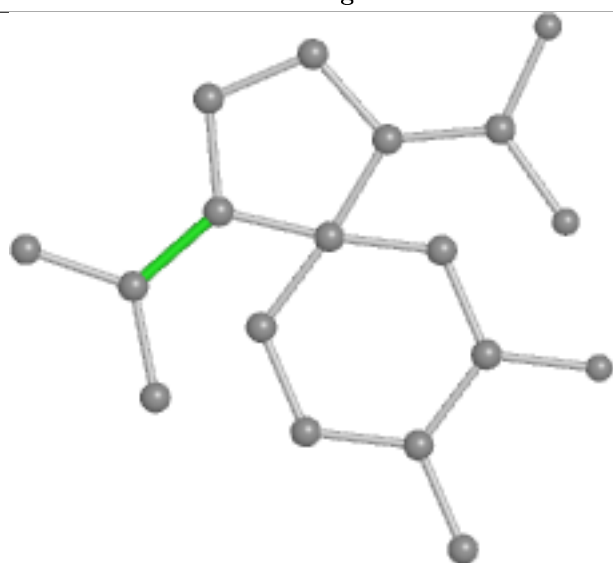
Ligand 7H6 B 501



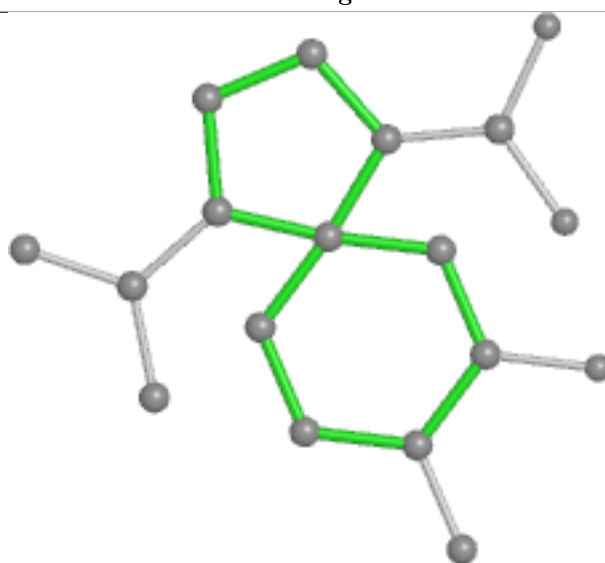
Bond lengths



Bond angles

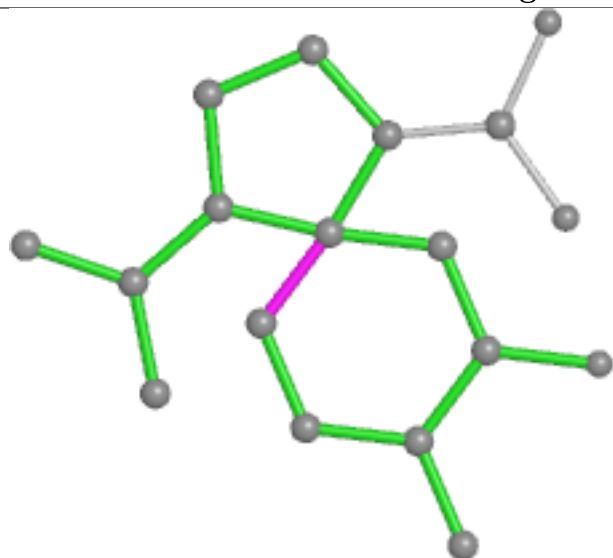


Torsions

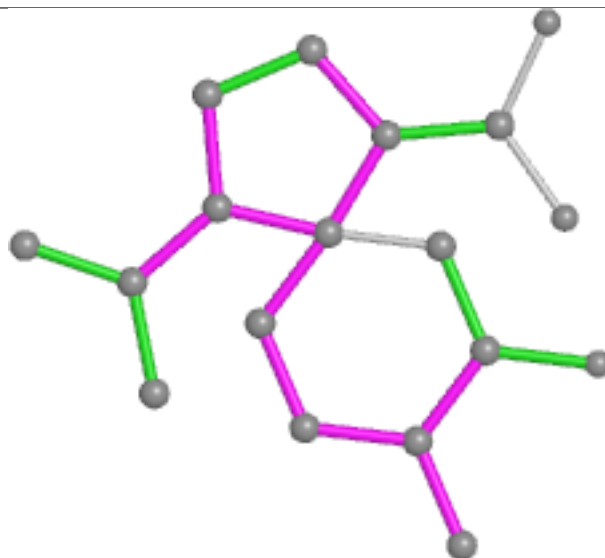


Rings

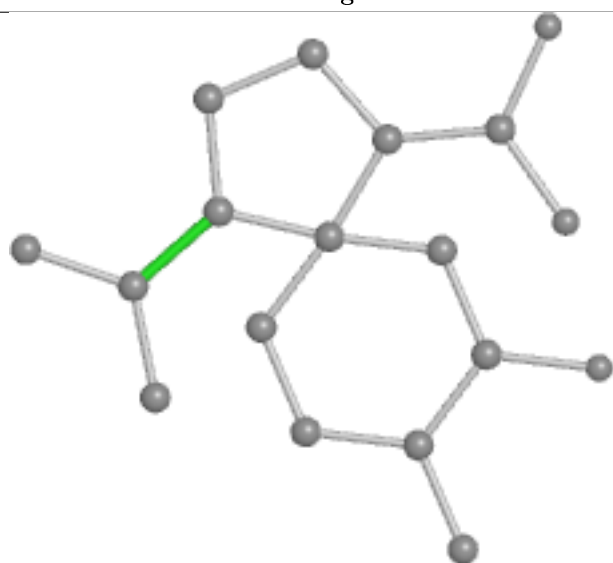
Ligand 7H6 D 501



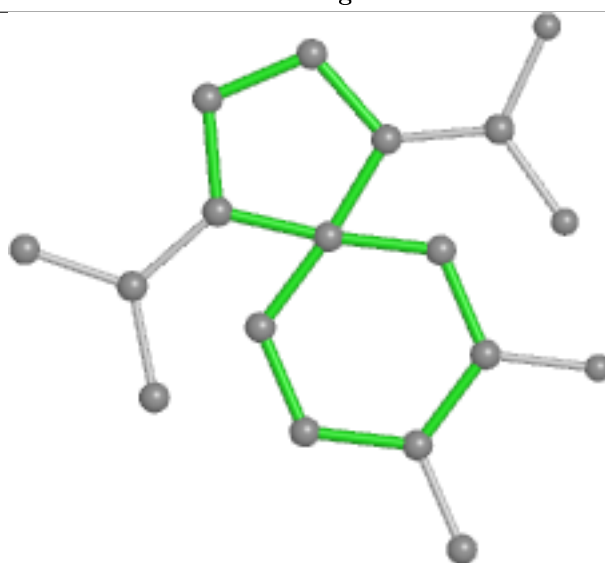
Bond lengths



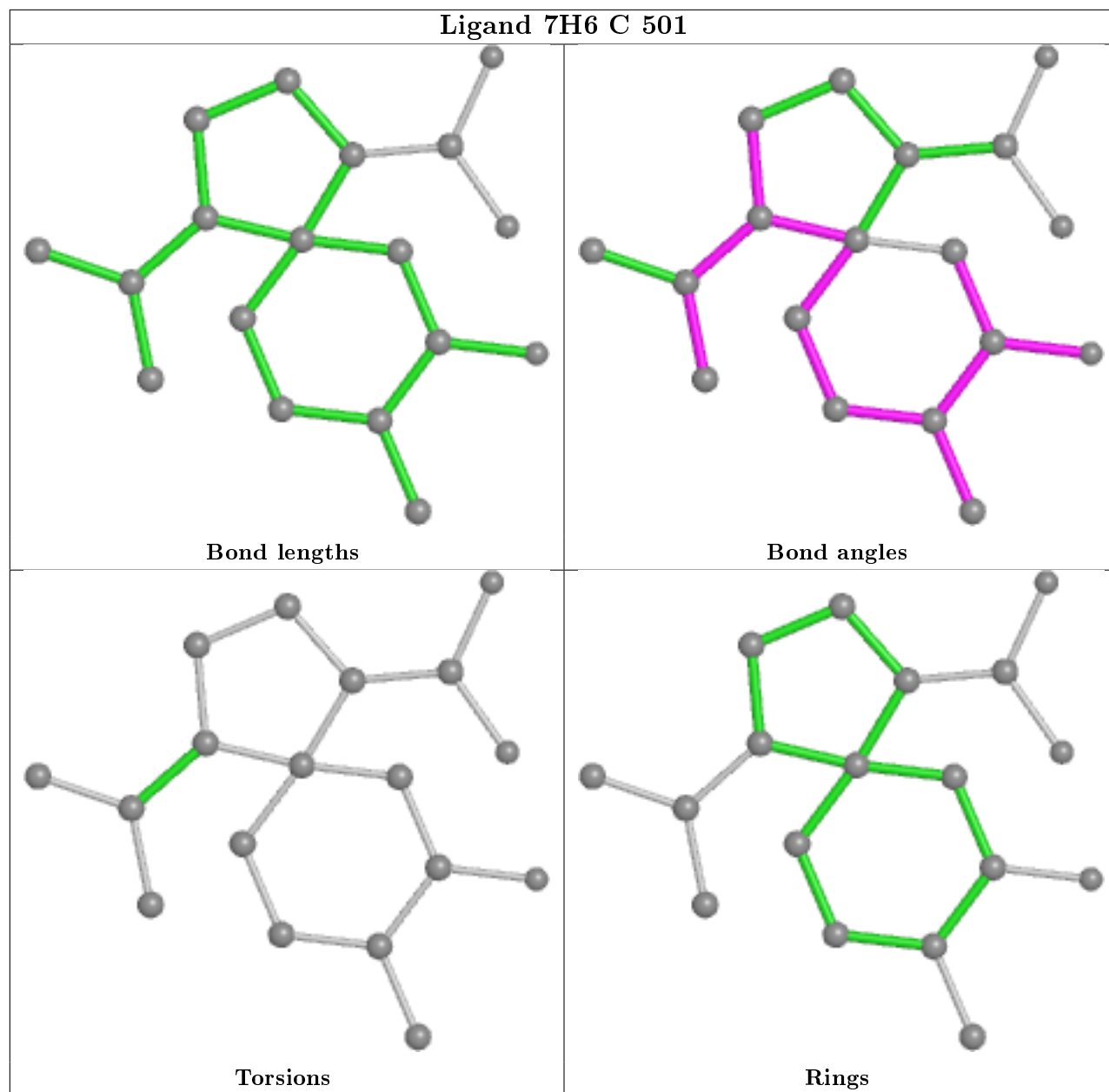
Bond angles

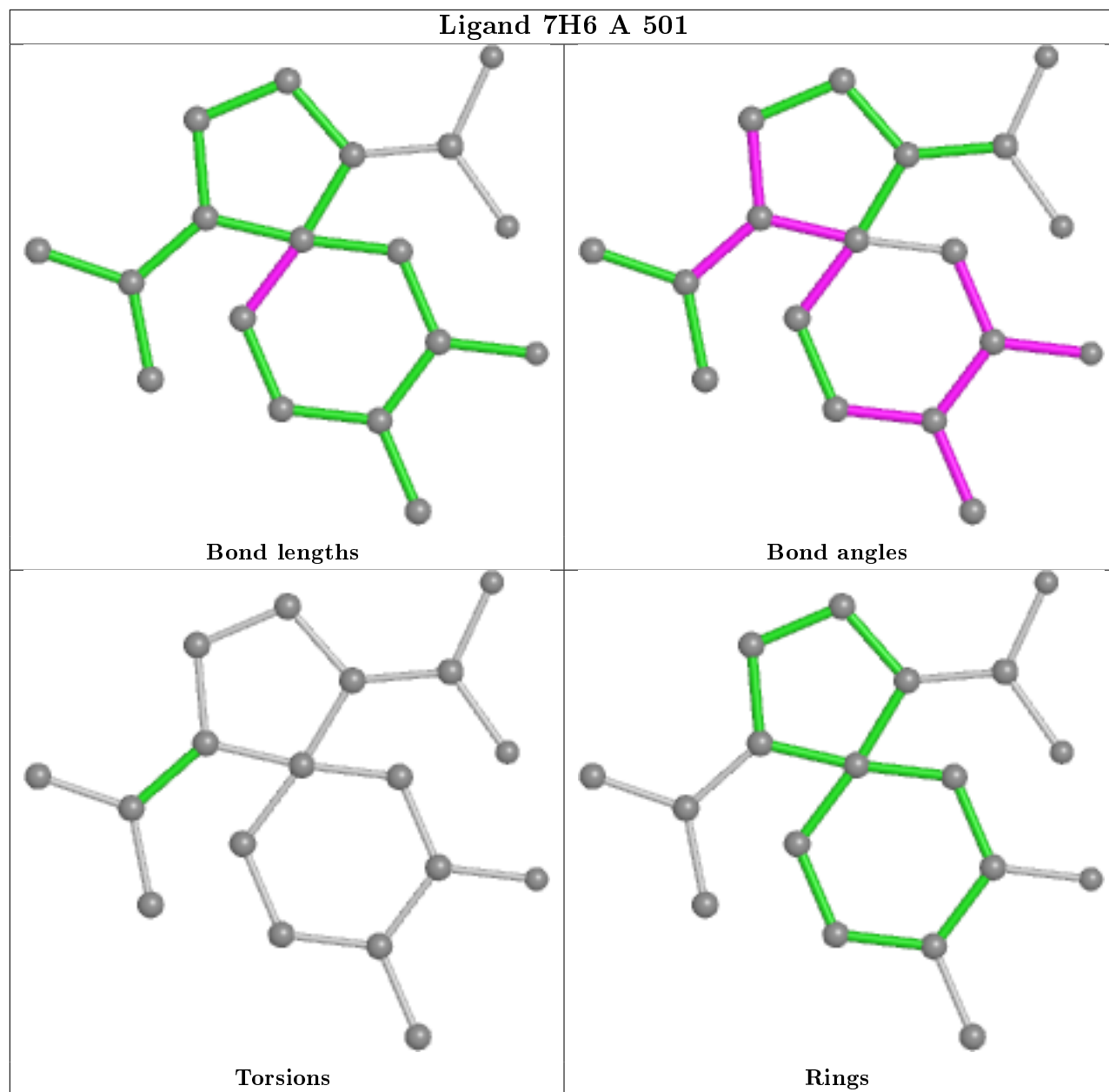


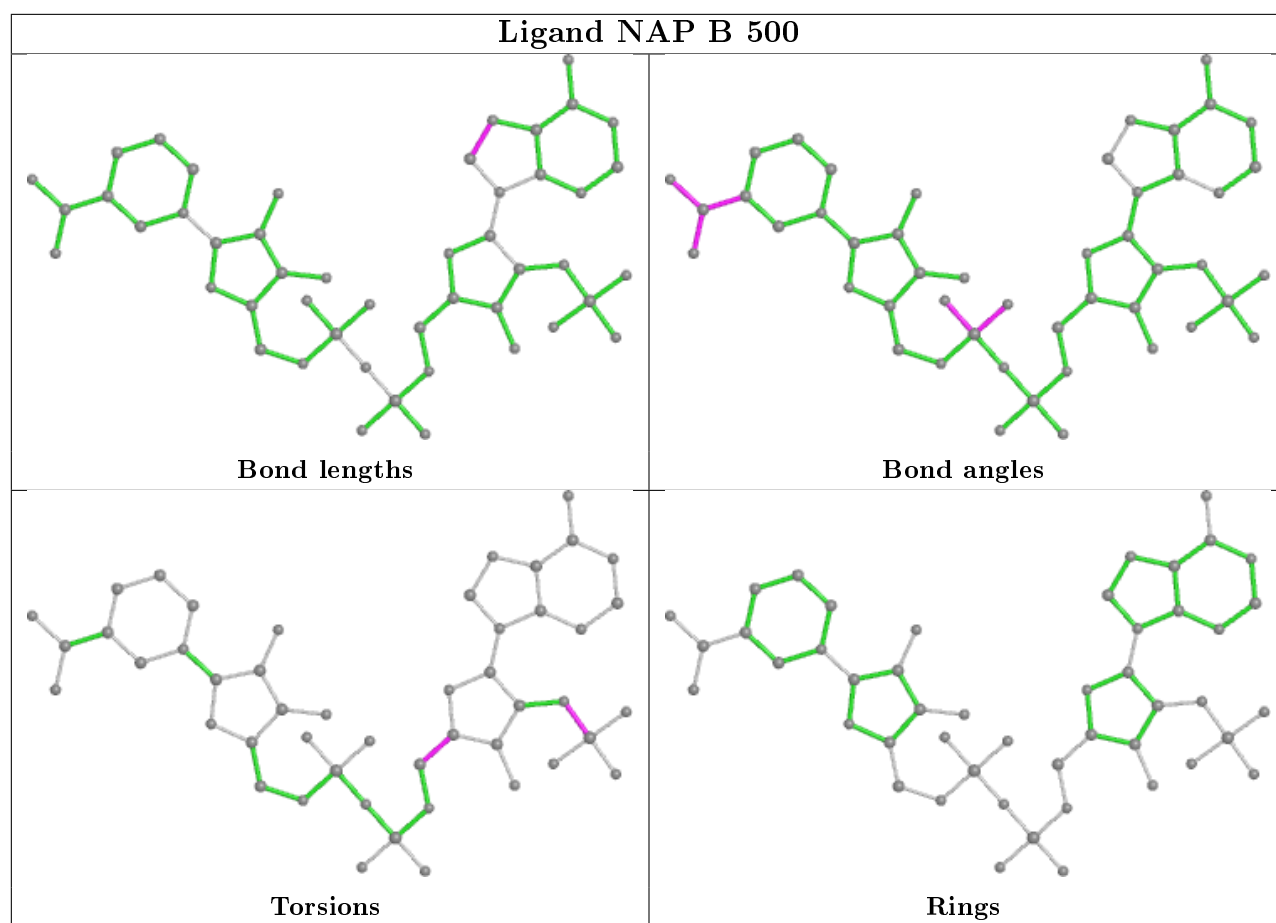
Torsions

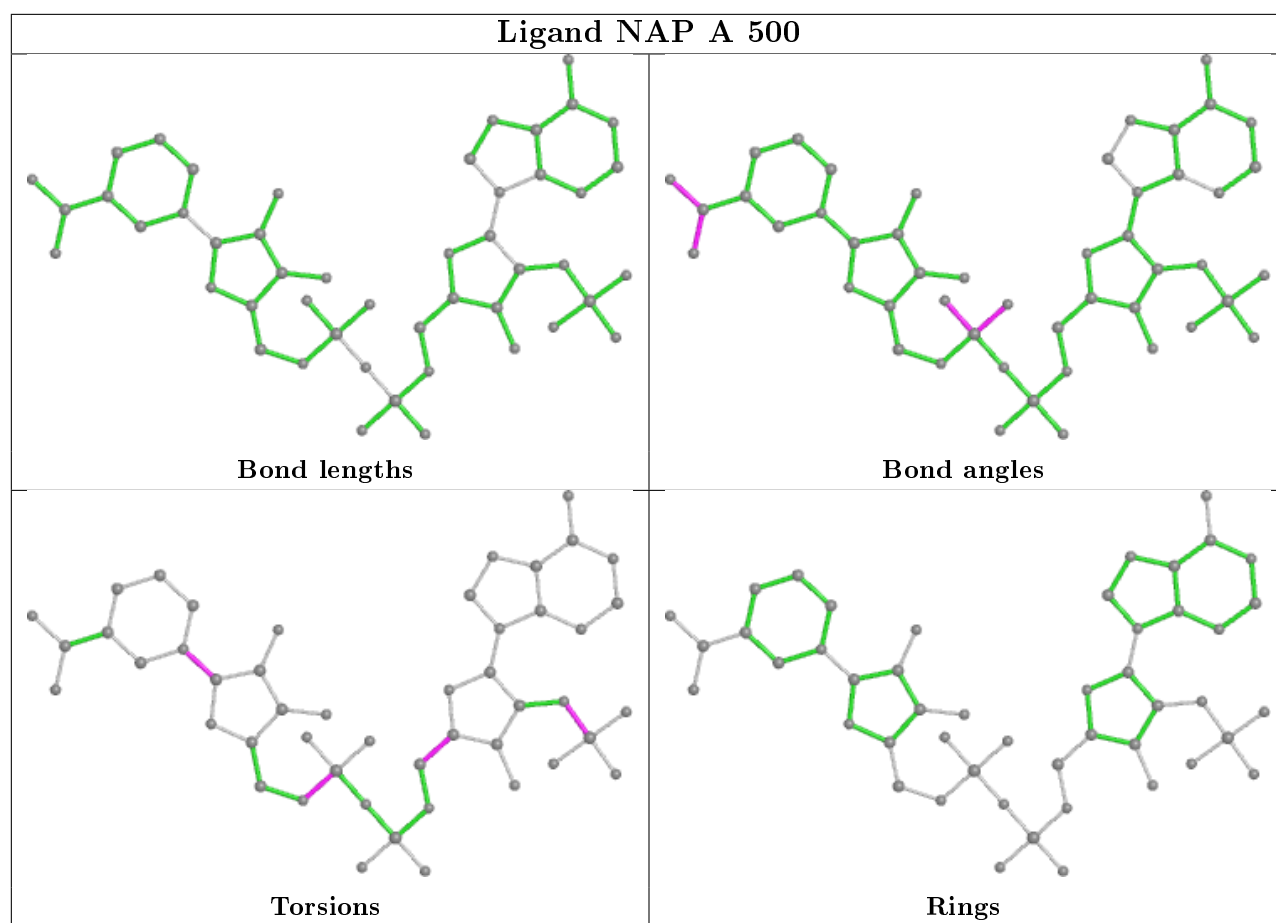


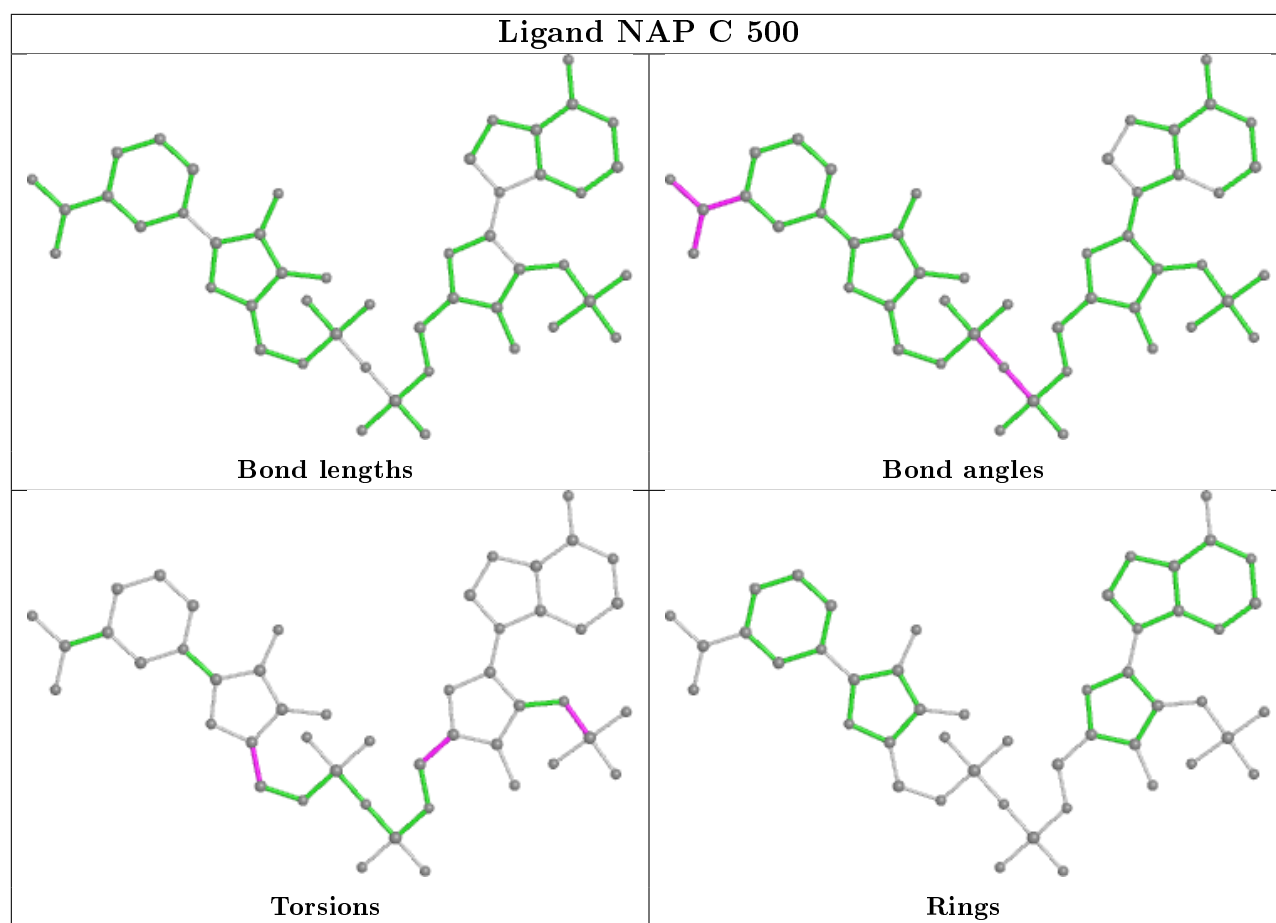
Rings











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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	243/289 (84%)	0.14	16 (6%) 18 13	70, 104, 129, 150	0
1	B	257/289 (88%)	0.16	21 (8%) 11 8	67, 95, 120, 152	0
1	C	250/289 (86%)	0.06	6 (2%) 59 53	60, 92, 126, 147	0
1	D	248/289 (85%)	0.25	18 (7%) 15 11	62, 105, 133, 155	0
All	All	998/1156 (86%)	0.15	61 (6%) 21 16	60, 99, 129, 155	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	118	LEU	4.4
1	C	168	VAL	4.3
1	D	118	LEU	4.2
1	D	78	CYS	4.2
1	B	167	VAL	4.0
1	B	190	LEU	3.8
1	B	119	ASN	3.8
1	A	118	LEU	3.6
1	B	117	ILE	3.6
1	B	168	VAL	3.5
1	B	189	ALA	3.5
1	A	189	ALA	3.2
1	A	168	VAL	3.1
1	B	177	TYR	3.1
1	D	213	CYS	3.1
1	D	117	ILE	3.0
1	D	169	SER	2.9
1	D	166	VAL	2.9
1	B	186	SER	2.8
1	D	90	ALA	2.8
1	C	192	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	119	ASN	2.8
1	A	117	ILE	2.7
1	C	187	LYS	2.7
1	B	147	TYR	2.7
1	B	40	THR	2.6
1	A	173	GLY	2.6
1	C	169	SER	2.6
1	D	190	LEU	2.6
1	A	190	LEU	2.5
1	B	27	PHE	2.5
1	A	169	SER	2.5
1	D	50	MET	2.5
1	D	40	THR	2.5
1	A	188	PHE	2.4
1	A	185	ALA	2.4
1	C	119	ASN	2.4
1	A	119	ASN	2.4
1	D	167	VAL	2.3
1	C	184	SER	2.3
1	D	168	VAL	2.3
1	B	187	LYS	2.3
1	B	229	GLY	2.2
1	D	184	SER	2.2
1	B	166	VAL	2.2
1	D	147	TYR	2.2
1	B	39	VAL	2.2
1	B	231	VAL	2.2
1	A	166	VAL	2.2
1	A	147	TYR	2.1
1	A	69	GLU	2.1
1	A	186	SER	2.1
1	D	39	VAL	2.1
1	B	213	CYS	2.1
1	D	280	TYR	2.1
1	D	189	ALA	2.1
1	A	203	VAL	2.0
1	B	192	GLY	2.0
1	B	185	ALA	2.0
1	B	235	ALA	2.0
1	A	187	LYS	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](#)

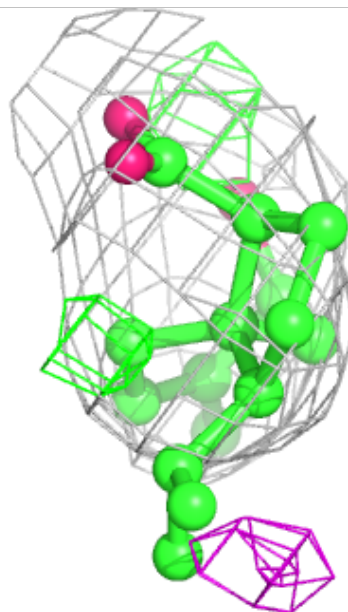
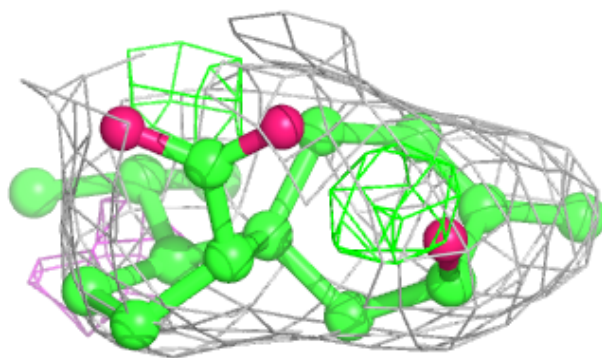
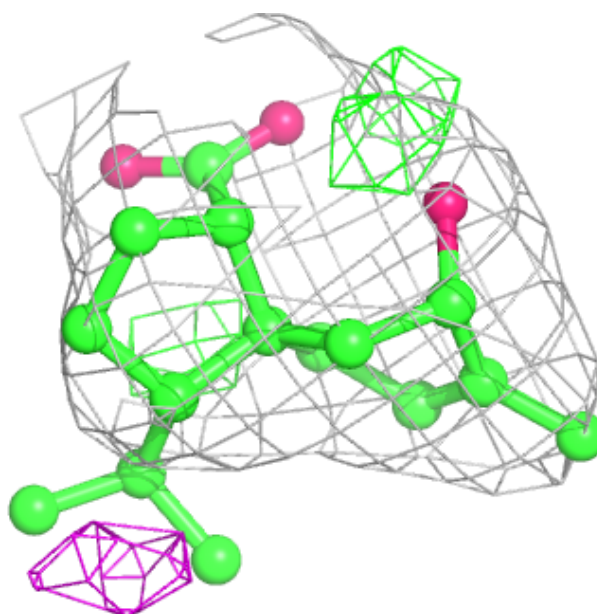
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	7H6	D	501	18/18	0.88	0.28	93,117,153,160	0
3	7H6	C	501	18/18	0.92	0.21	86,100,116,127	0
2	NAP	A	500	48/48	0.92	0.14	73,99,109,118	0
2	NAP	B	500	48/48	0.93	0.14	64,78,90,96	0
2	NAP	D	500	48/48	0.93	0.13	74,96,112,127	0
3	7H6	B	501	18/18	0.95	0.23	97,108,121,127	0
3	7H6	A	501	18/18	0.95	0.23	92,111,129,132	0
2	NAP	C	500	48/48	0.97	0.16	70,79,90,100	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.

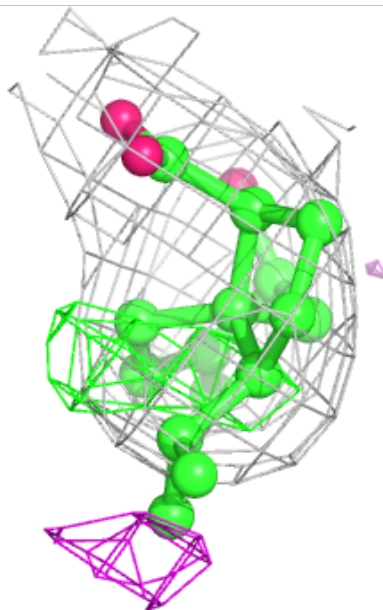
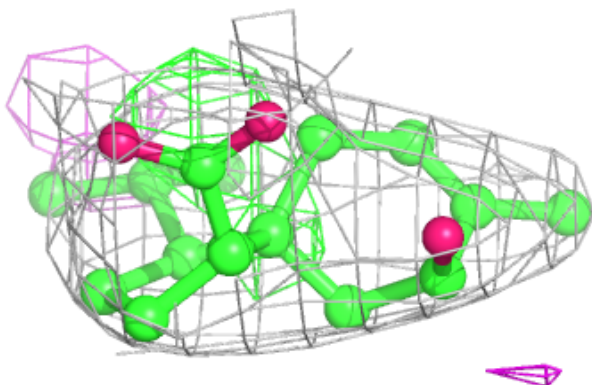
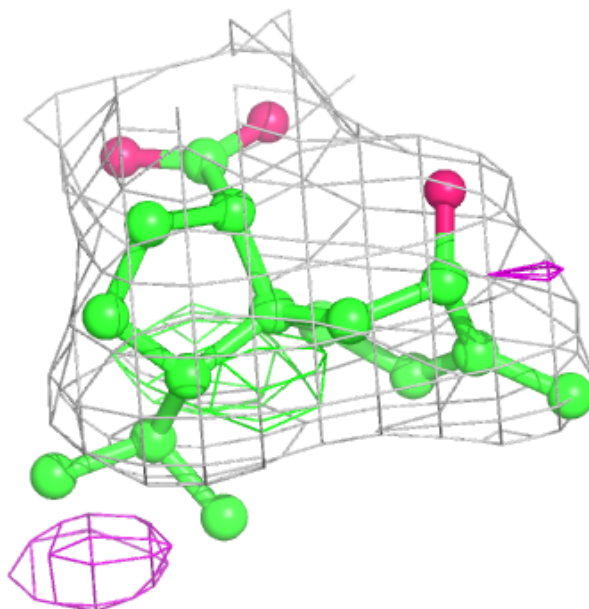
Electron density around 7H6 D 501:

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



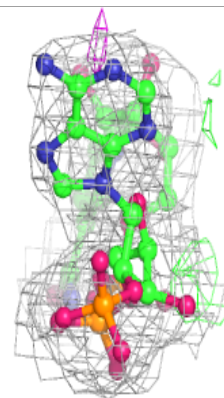
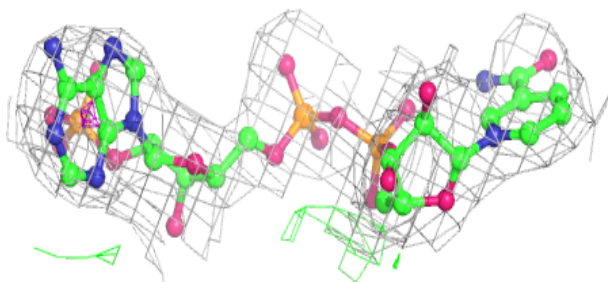
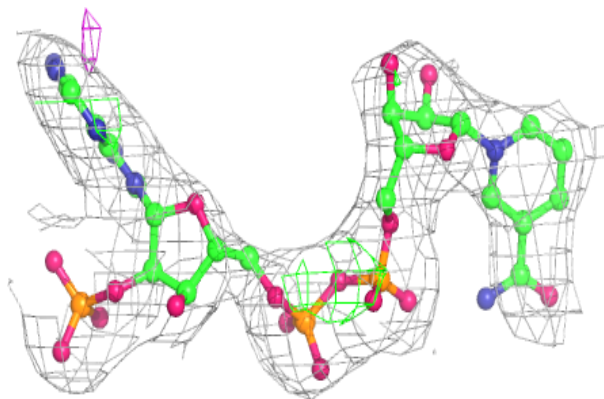
Electron density around 7H6 C 501:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

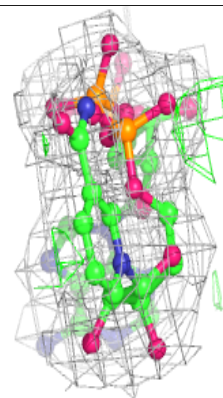
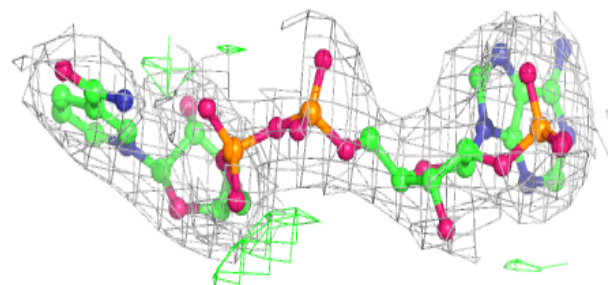


Electron density around NAP A 500:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

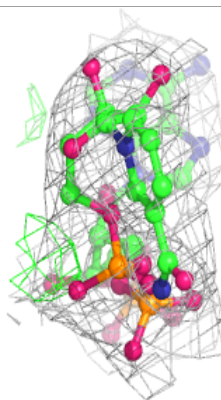
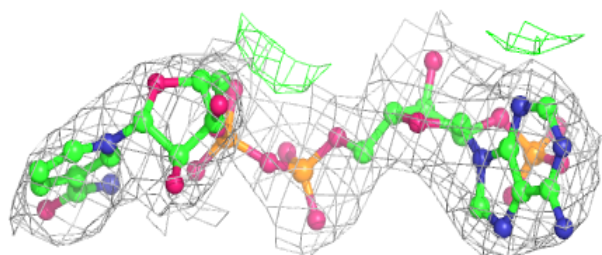
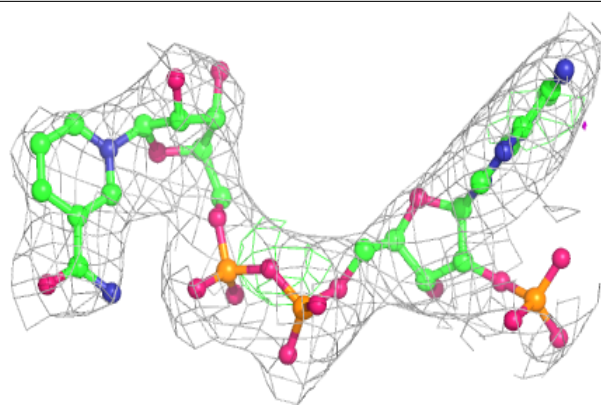
**Electron density around NAP B 500:**

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



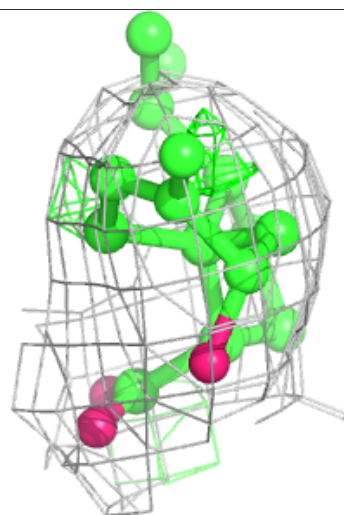
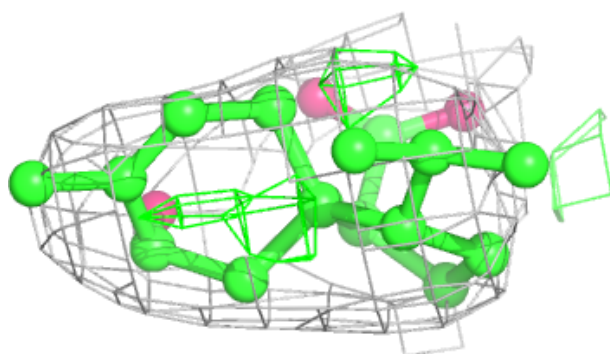
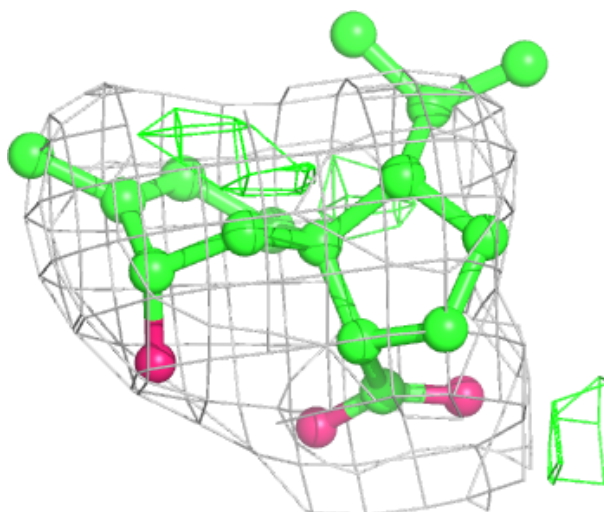
Electron density around NAP D 500:

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



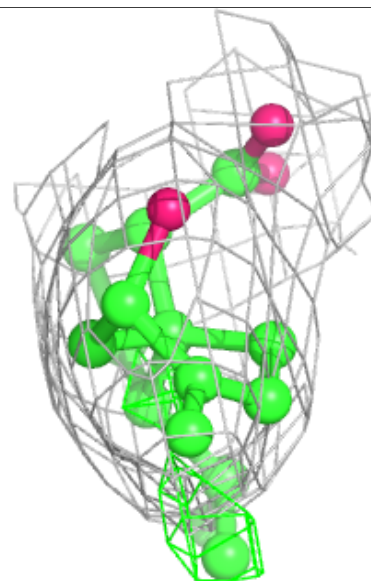
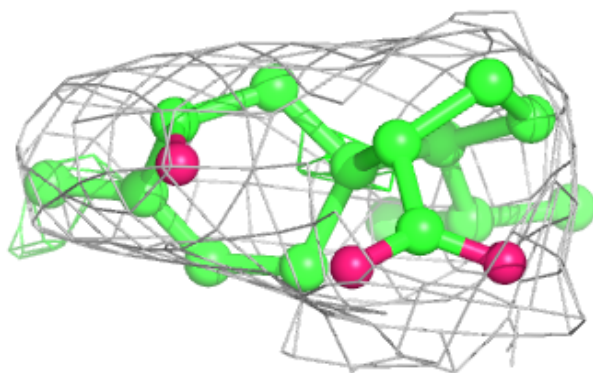
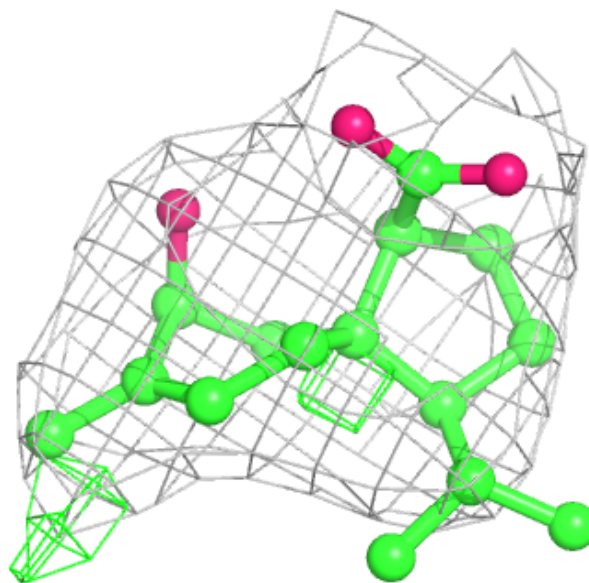
Electron density around 7H6 B 501:

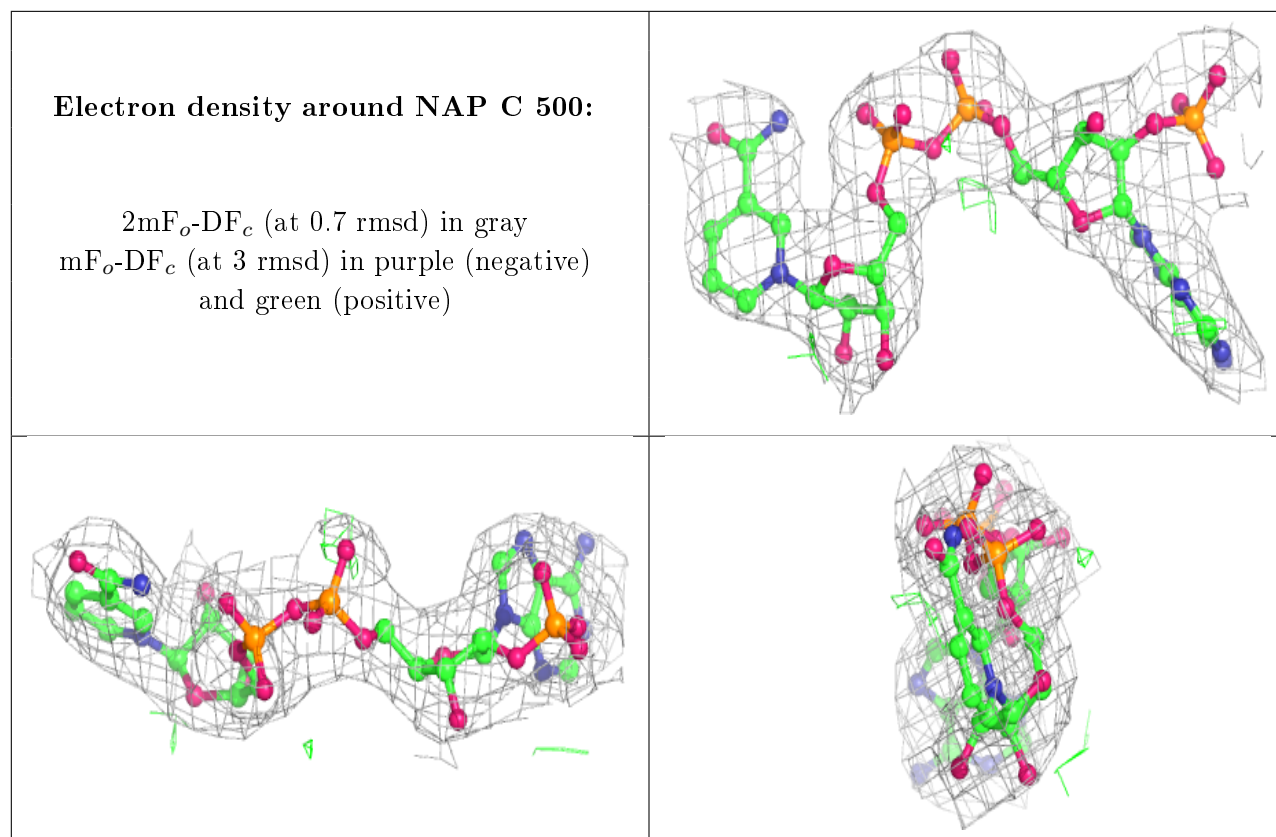
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 7H6 A 501:

$2mF_o-DF_c$ (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.