



wwPDB X-ray Structure Validation Summary Report ⓘ

Sep 2, 2023 – 04:42 PM EDT

PDB ID : 3RHO
Title : Crystal structure of the E673Q MUTANT OF C-Terminal domain of 10'FOR MYLTETRAHYDROFOLATE DEHYDROGENASE in complex with NADP
Authors : Tsybovsky, Y.
Deposited on : 2011-04-11
Resolution : 2.26 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

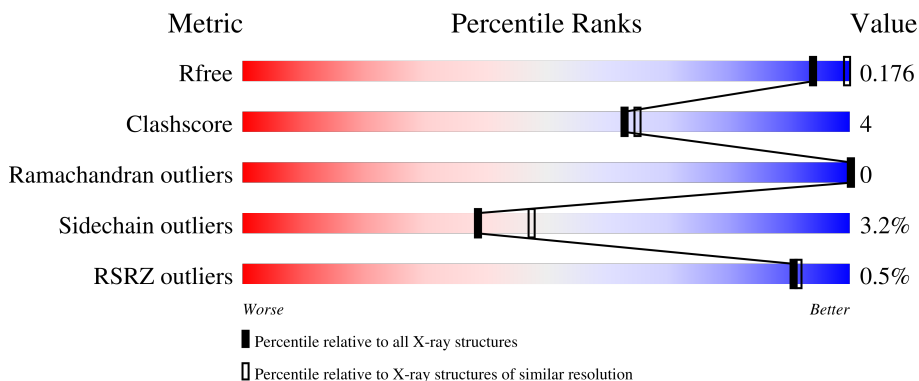
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



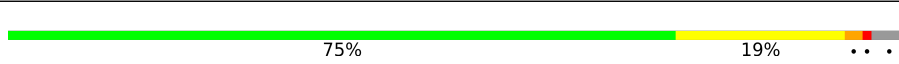
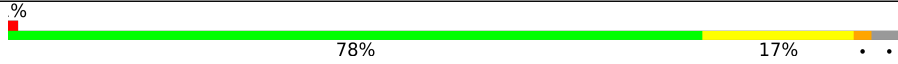
The reported resolution of this entry is 2.26 Å.

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	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	517	 78% 15% . .
1	B	517	 79% 16% . .
1	C	517	 75% 19% . .
1	D	517	 78% 17% . .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 17496 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 Aldehyde dehydrogenase 1 family, member L1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	498	3906	2478	672	735	21	4	19	0
1	B	498	3906	2478	672	735	21	0	19	0
1	C	498	3906	2478	672	735	21	0	19	0
1	D	498	3906	2478	672	735	21	0	19	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	386	MET	-	expression tag	UNP Q5HQB2
A	387	ARG	-	expression tag	UNP Q5HQB2
A	388	GLY	-	expression tag	UNP Q5HQB2
A	389	SER	-	expression tag	UNP Q5HQB2
A	390	HIS	-	expression tag	UNP Q5HQB2
A	391	HIS	-	expression tag	UNP Q5HQB2
A	392	HIS	-	expression tag	UNP Q5HQB2
A	393	HIS	-	expression tag	UNP Q5HQB2
A	394	HIS	-	expression tag	UNP Q5HQB2
A	395	THR	-	expression tag	UNP Q5HQB2
A	396	THR	-	expression tag	UNP Q5HQB2
A	673	GLN	GLU	engineered mutation	UNP Q5HQB2
B	386	MET	-	expression tag	UNP Q5HQB2
B	387	ARG	-	expression tag	UNP Q5HQB2
B	388	GLY	-	expression tag	UNP Q5HQB2
B	389	SER	-	expression tag	UNP Q5HQB2
B	390	HIS	-	expression tag	UNP Q5HQB2
B	391	HIS	-	expression tag	UNP Q5HQB2
B	392	HIS	-	expression tag	UNP Q5HQB2
B	393	HIS	-	expression tag	UNP Q5HQB2
B	394	HIS	-	expression tag	UNP Q5HQB2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	395	THR	-	expression tag	UNP Q5HZB2
B	396	THR	-	expression tag	UNP Q5HZB2
B	673	GLN	GLU	engineered mutation	UNP Q5HZB2
C	386	MET	-	expression tag	UNP Q5HZB2
C	387	ARG	-	expression tag	UNP Q5HZB2
C	388	GLY	-	expression tag	UNP Q5HZB2
C	389	SER	-	expression tag	UNP Q5HZB2
C	390	HIS	-	expression tag	UNP Q5HZB2
C	391	HIS	-	expression tag	UNP Q5HZB2
C	392	HIS	-	expression tag	UNP Q5HZB2
C	393	HIS	-	expression tag	UNP Q5HZB2
C	394	HIS	-	expression tag	UNP Q5HZB2
C	395	THR	-	expression tag	UNP Q5HZB2
C	396	THR	-	expression tag	UNP Q5HZB2
C	673	GLN	GLU	engineered mutation	UNP Q5HZB2
D	386	MET	-	expression tag	UNP Q5HZB2
D	387	ARG	-	expression tag	UNP Q5HZB2
D	388	GLY	-	expression tag	UNP Q5HZB2
D	389	SER	-	expression tag	UNP Q5HZB2
D	390	HIS	-	expression tag	UNP Q5HZB2
D	391	HIS	-	expression tag	UNP Q5HZB2
D	392	HIS	-	expression tag	UNP Q5HZB2
D	393	HIS	-	expression tag	UNP Q5HZB2
D	394	HIS	-	expression tag	UNP Q5HZB2
D	395	THR	-	expression tag	UNP Q5HZB2
D	396	THR	-	expression tag	UNP Q5HZB2
D	673	GLN	GLU	engineered mutation	UNP Q5HZB2

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total 56	C 21	N 7	O 23	P 5	3	1
2	B	1	Total 56	C 21	N 7	O 23	P 5	3	1
2	C	1	Total 56	C 21	N 7	O 23	P 5	3	1
2	D	1	Total 56	C 21	N 7	O 23	P 5	3	1

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



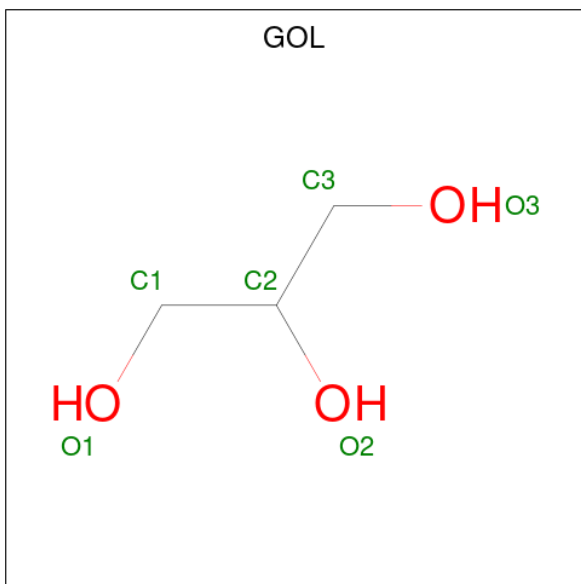
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

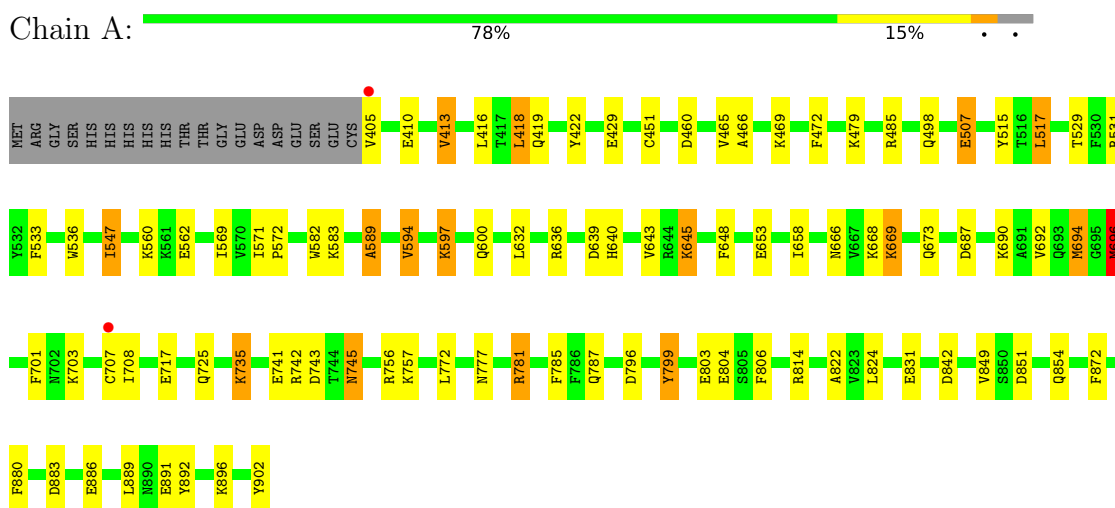
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	377	Total 377	O 377	0	4
5	B	336	Total 336	O 336	0	3
5	C	394	Total 394	O 394	0	3
5	D	377	Total 377	O 377	0	3

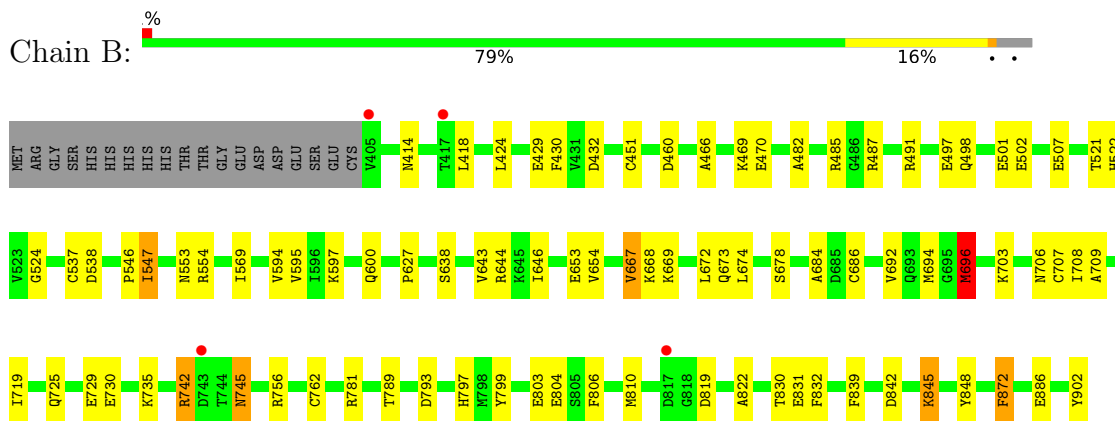
3 Residue-property plots [i](#)

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.

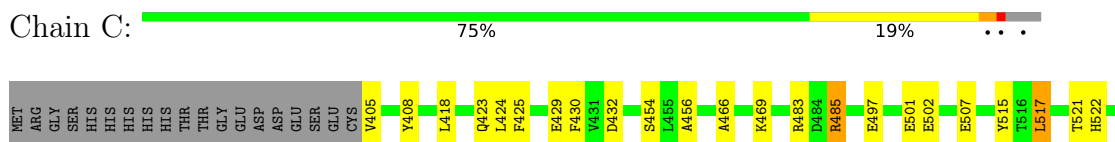
- Molecule 1: Aldehyde dehydrogenase 1 family, member L1

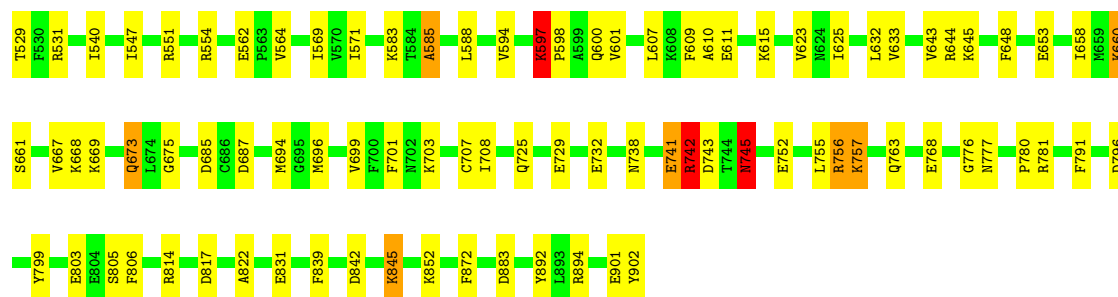


- Molecule 1: Aldehyde dehydrogenase 1 family, member L1

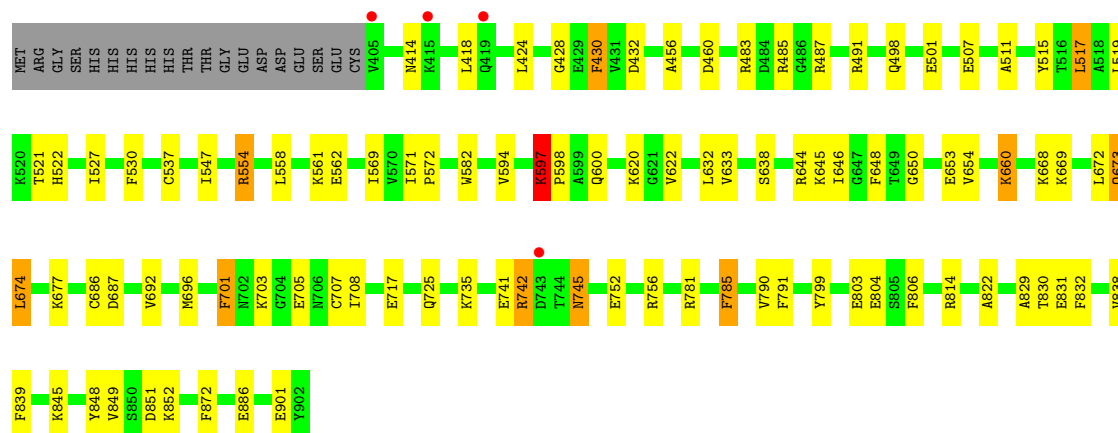
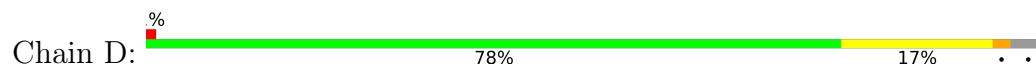


- Molecule 1: Aldehyde dehydrogenase 1 family, member L1





- Molecule 1: Aldehyde dehydrogenase 1 family, member L1



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	259.90Å 194.40Å 97.22Å 90.00° 109.09° 90.00°	Depositor
Resolution (Å)	49.15 – 2.26 49.12 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.15-2.26) 99.5 (49.12-2.26)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.74 (at 2.24Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.149 , 0.178 0.149 , 0.176	Depositor DCC
R_{free} test set	10652 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtrriage
Anisotropy	0.015	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.015 for -h-2*1,-k,l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17496	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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, GOL, SO4

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	1.63	39/4066 (1.0%)	1.09	15/5496 (0.3%)
1	B	1.63	45/4066 (1.1%)	1.08	17/5496 (0.3%)
1	C	1.74	55/4066 (1.4%)	1.19	26/5496 (0.5%)
1	D	1.67	53/4066 (1.3%)	1.07	12/5496 (0.2%)
All	All	1.66	192/16264 (1.2%)	1.11	70/21984 (0.3%)

The worst 5 of 192 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	501	GLU	CD-OE2	12.97	1.40	1.25
1	B	831	GLU	CG-CD	9.77	1.66	1.51
1	C	653	GLU	CG-CD	9.66	1.66	1.51
1	D	804	GLU	CG-CD	9.26	1.65	1.51
1	B	501	GLU	CD-OE2	8.98	1.35	1.25

The worst 5 of 70 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	742	ARG	NE-CZ-NH2	-18.10	111.25	120.30
1	C	742	ARG	NE-CZ-NH1	12.23	126.42	120.30
1	B	694	MET	CG-SD-CE	9.28	115.05	100.20
1	D	487	ARG	NE-CZ-NH1	9.03	124.82	120.30
1	D	485	ARG	NE-CZ-NH2	-8.65	115.97	120.30

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	3906	0	3923	35	0
1	B	3906	0	3922	29	0
1	C	3906	0	3923	42	0
1	D	3906	0	3923	31	0
2	A	56	0	14	1	0
2	B	56	0	14	1	0
2	C	56	0	14	1	0
2	D	56	0	14	1	0
3	A	35	0	0	0	0
3	B	40	0	0	1	0
3	C	40	0	0	2	0
3	D	25	0	0	1	0
4	A	6	0	8	0	0
4	B	6	0	8	2	0
4	C	6	0	8	0	0
4	D	6	0	8	0	0
5	A	377	0	0	6	0
5	B	336	0	0	5	0
5	C	394	0	0	6	0
5	D	377	0	0	3	0
All	All	17496	0	15779	138	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

The worst 5 of 138 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:669:LYS:CE	1:A:669:LYS:NZ	1.69	1.50
1:B:799:TYR:CE2	1:B:803[A]:GLU:HG3	2.14	0.83
1:D:742:ARG:HD2	5:D:1465:HOH:O	1.83	0.78
1:B:725[A]:GLN:NE2	1:B:729:GLU:OE2	2.19	0.76
1:A:569:ILE:HD12	1:A:594:VAL:HG21	1.69	0.74

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	515/517 (100%)	501 (97%)	14 (3%)	0	100	100
1	B	515/517 (100%)	495 (96%)	20 (4%)	0	100	100
1	C	515/517 (100%)	492 (96%)	23 (4%)	0	100	100
1	D	515/517 (100%)	497 (96%)	18 (4%)	0	100	100
All	All	2060/2068 (100%)	1985 (96%)	75 (4%)	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	428/426 (100%)	410 (96%)	18 (4%)	30	34
1	B	428/426 (100%)	415 (97%)	13 (3%)	41	50
1	C	428/426 (100%)	412 (96%)	16 (4%)	34	40
1	D	428/426 (100%)	414 (97%)	14 (3%)	38	46
All	All	1712/1704 (100%)	1651 (96%)	61 (4%)	39	42

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	806	PHE
1	D	725[A]	GLN
1	C	673	GLN

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Mol	Chain	Res	Type
1	D	707[B]	CYS
1	D	830	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 20 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	797	HIS
1	D	600	GLN
1	D	745	ASN
1	D	706	ASN
1	B	498	GLN

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](#)

40 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	903[A]	-	45,33,52	3.24	6 (13%)	56,52,80	1.93	8 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	3015	-	4,4,4	0.51	0	6,6,6	1.53	1 (16%)
3	SO4	A	3001	-	4,4,4	0.38	0	6,6,6	0.52	0
3	SO4	B	3019	-	4,4,4	0.48	0	6,6,6	0.59	0
3	SO4	C	3002	-	4,4,4	0.27	0	6,6,6	0.37	0
4	GOL	B	3026	-	5,5,5	0.55	0	5,5,5	0.57	0
3	SO4	A	3018	-	4,4,4	0.55	0	6,6,6	0.67	0
2	NAP	A	903[B]	-	45,52,52	3.65	5 (11%)	56,80,80	1.21	5 (8%)
3	SO4	B	3003	-	4,4,4	0.43	0	6,6,6	0.47	0
3	SO4	B	3009	-	4,4,4	0.26	0	6,6,6	0.51	0
3	SO4	C	3007	-	4,4,4	0.33	0	6,6,6	1.53	1 (16%)
3	SO4	A	3024	-	4,4,4	0.19	0	6,6,6	0.45	0
3	SO4	D	3016	-	4,4,4	0.72	0	6,6,6	1.16	1 (16%)
3	SO4	B	3022	-	4,4,4	0.33	0	6,6,6	0.92	0
2	NAP	D	903[B]	-	45,52,52	2.23	5 (11%)	56,80,80	1.27	4 (7%)
2	NAP	B	903[A]	-	45,33,52	5.37	6 (13%)	56,52,80	2.57	9 (16%)
3	SO4	B	3005	-	4,4,4	0.33	0	6,6,6	1.14	1 (16%)
2	NAP	C	903[B]	-	45,52,52	2.40	5 (11%)	56,80,80	1.75	7 (12%)
3	SO4	B	3025	-	4,4,4	0.36	0	6,6,6	0.72	0
3	SO4	C	3013	-	4,4,4	0.40	0	6,6,6	1.47	1 (16%)
3	SO4	C	3023	-	4,4,4	0.24	0	6,6,6	0.70	0
3	SO4	C	3026	-	4,4,4	0.28	0	6,6,6	1.68	2 (33%)
3	SO4	C	3028	-	4,4,4	0.48	0	6,6,6	1.28	1 (16%)
3	SO4	D	3020	-	4,4,4	0.71	0	6,6,6	0.87	0
4	GOL	A	3025	-	5,5,5	0.69	0	5,5,5	1.16	1 (20%)
3	SO4	C	3010	-	4,4,4	0.15	0	6,6,6	0.48	0
3	SO4	A	3021	-	4,4,4	0.40	0	6,6,6	1.00	0
3	SO4	D	3027	-	4,4,4	0.24	0	6,6,6	0.81	0
3	SO4	D	3012	-	4,4,4	0.17	0	6,6,6	0.45	0
2	NAP	A	903[A]	-	45,33,52	4.75	6 (13%)	56,52,80	1.66	6 (10%)
2	NAP	B	903[B]	-	45,52,52	3.48	5 (11%)	56,80,80	2.28	8 (14%)
3	SO4	A	3006	-	4,4,4	0.24	0	6,6,6	1.21	0
4	GOL	C	3029	-	5,5,5	0.66	0	5,5,5	1.28	1 (20%)
3	SO4	B	3014	-	4,4,4	0.32	0	6,6,6	1.53	2 (33%)
4	GOL	D	3028	-	5,5,5	0.48	0	5,5,5	1.39	0
3	SO4	B	3004	-	4,4,4	0.39	0	6,6,6	0.38	0
3	SO4	C	3017	-	4,4,4	0.32	0	6,6,6	0.82	0
2	NAP	D	903[A]	-	45,33,52	3.20	6 (13%)	56,52,80	1.67	6 (10%)
3	SO4	D	3008	-	4,4,4	0.58	0	6,6,6	1.52	1 (16%)
3	SO4	A	3011	-	4,4,4	0.18	0	6,6,6	0.49	0

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
4	GOL	C	3029	-	-	3/4/4/4	-
2	NAP	A	903[A]	-	-	5/31/37/67	0/5/3/5
2	NAP	C	903[A]	-	-	4/31/37/67	0/5/3/5
2	NAP	B	903[B]	-	-	9/31/67/67	0/5/5/5
4	GOL	D	3028	-	-	2/4/4/4	-
2	NAP	D	903[B]	-	-	9/31/67/67	0/5/5/5
2	NAP	B	903[A]	-	-	5/31/37/67	0/5/3/5
4	GOL	A	3025	-	-	4/4/4/4	-
2	NAP	D	903[A]	-	-	7/31/37/67	0/5/3/5
2	NAP	A	903[B]	-	-	9/31/67/67	0/5/5/5
2	NAP	C	903[B]	-	-	9/31/67/67	0/5/5/5
4	GOL	B	3026	-	-	1/4/4/4	-

The worst 5 of 44 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	903[A]	NAP	O5D-C5D	27.51	2.50	1.44
2	A	903[A]	NAP	C3N-C7N	-21.66	1.18	1.50
2	A	903[B]	NAP	C3N-C7N	-21.66	1.18	1.50
2	A	903[A]	NAP	O5D-C5D	20.41	2.23	1.44
2	B	903[A]	NAP	C3N-C7N	-20.39	1.19	1.50

The worst 5 of 66 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	903[A]	NAP	PN-O5D-C5D	-8.87	69.68	121.68
2	B	903[A]	NAP	O7N-C7N-C3N	8.73	130.08	119.63
2	B	903[B]	NAP	O7N-C7N-C3N	8.73	130.08	119.63
2	A	903[A]	NAP	PN-O5D-C5D	-8.56	71.49	121.68
2	B	903[A]	NAP	C3N-C7N-N7N	-8.15	107.97	117.75

There are no chirality outliers.

5 of 67 torsion outliers are listed below:

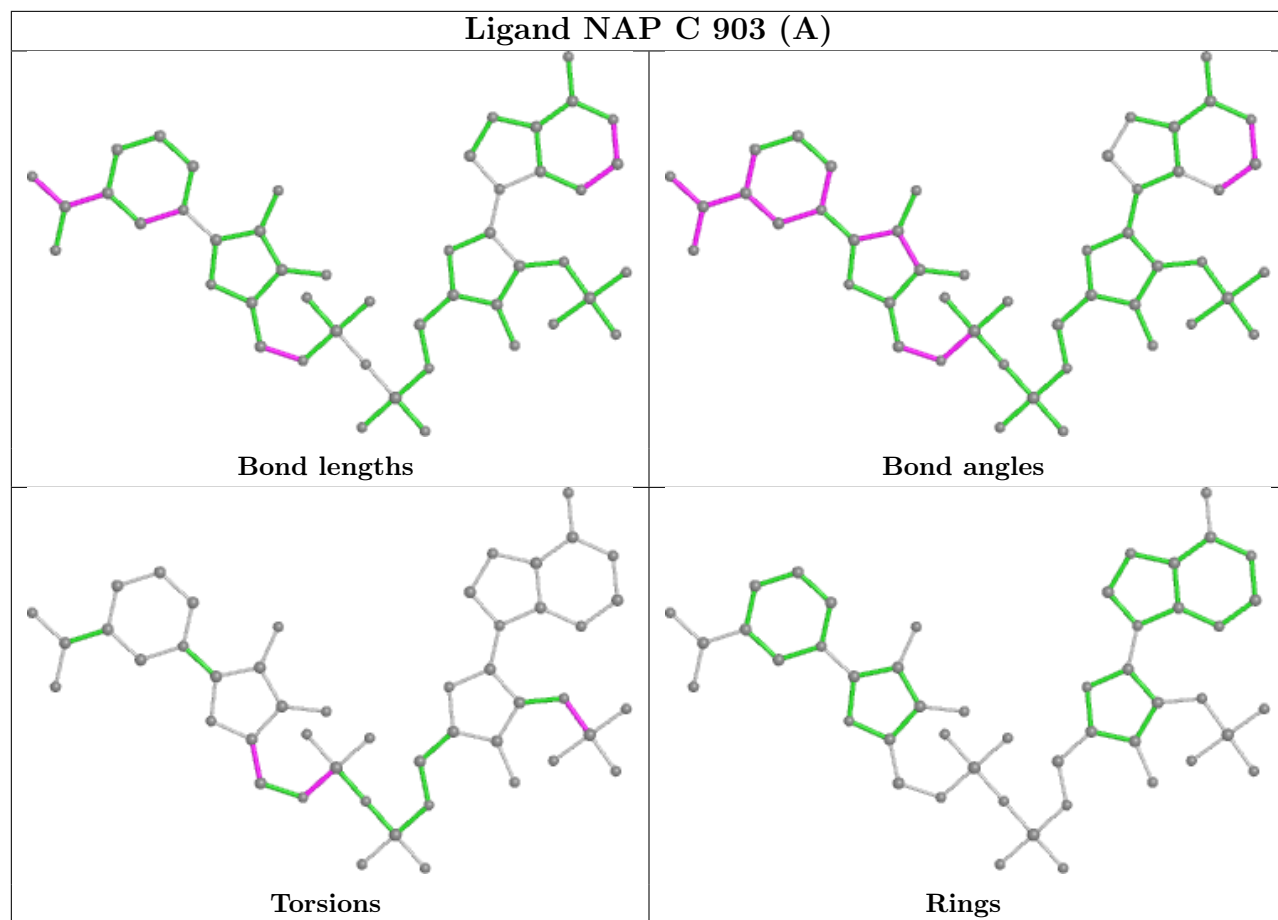
Mol	Chain	Res	Type	Atoms
2	A	903[A]	NAP	C5D-O5D-PN-O1N
2	A	903[B]	NAP	C5B-O5B-PA-O1A
2	A	903[B]	NAP	C3D-C4D-C5D-O5D
2	B	903[A]	NAP	C5D-O5D-PN-O1N
2	B	903[B]	NAP	C5B-O5B-PA-O1A

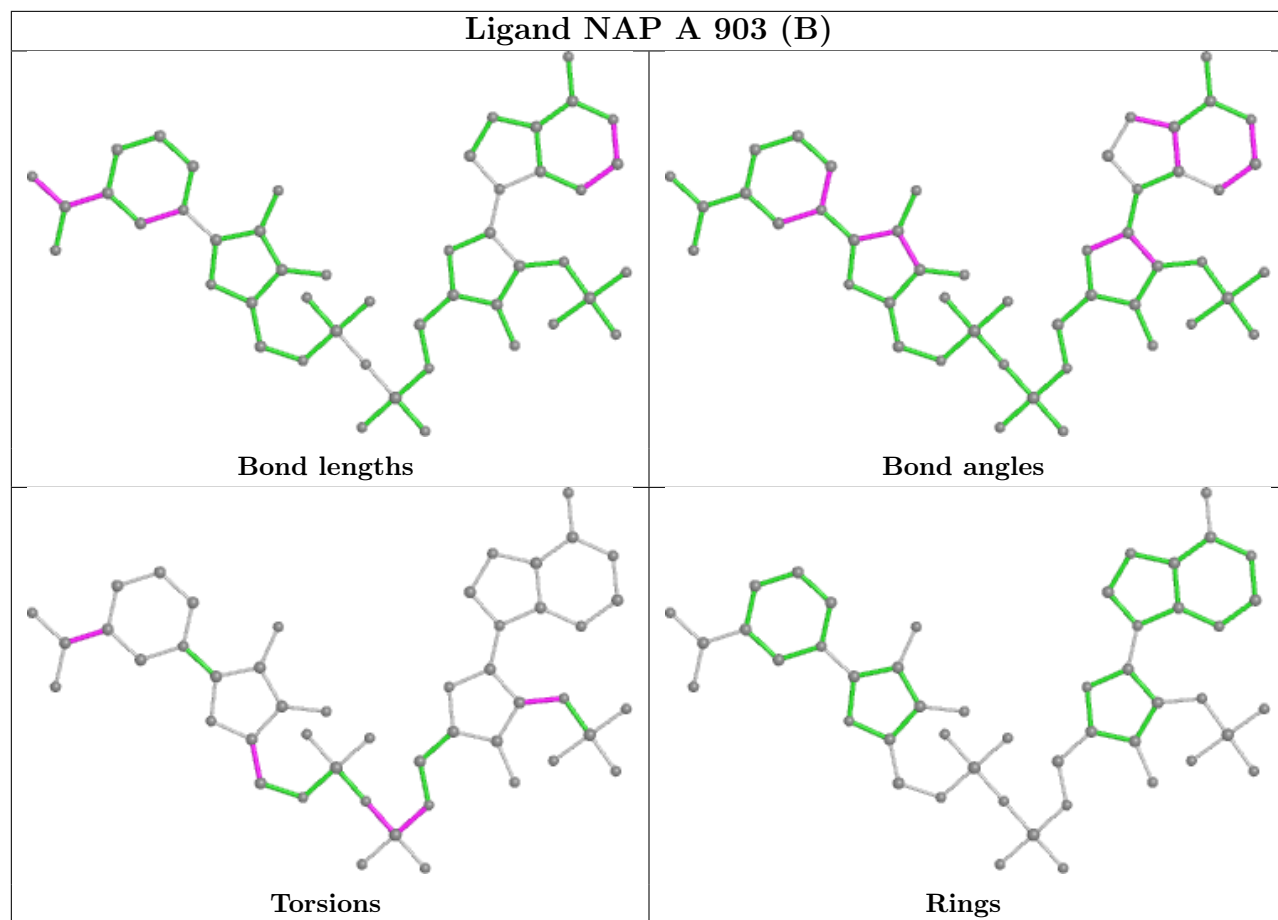
There are no ring outliers.

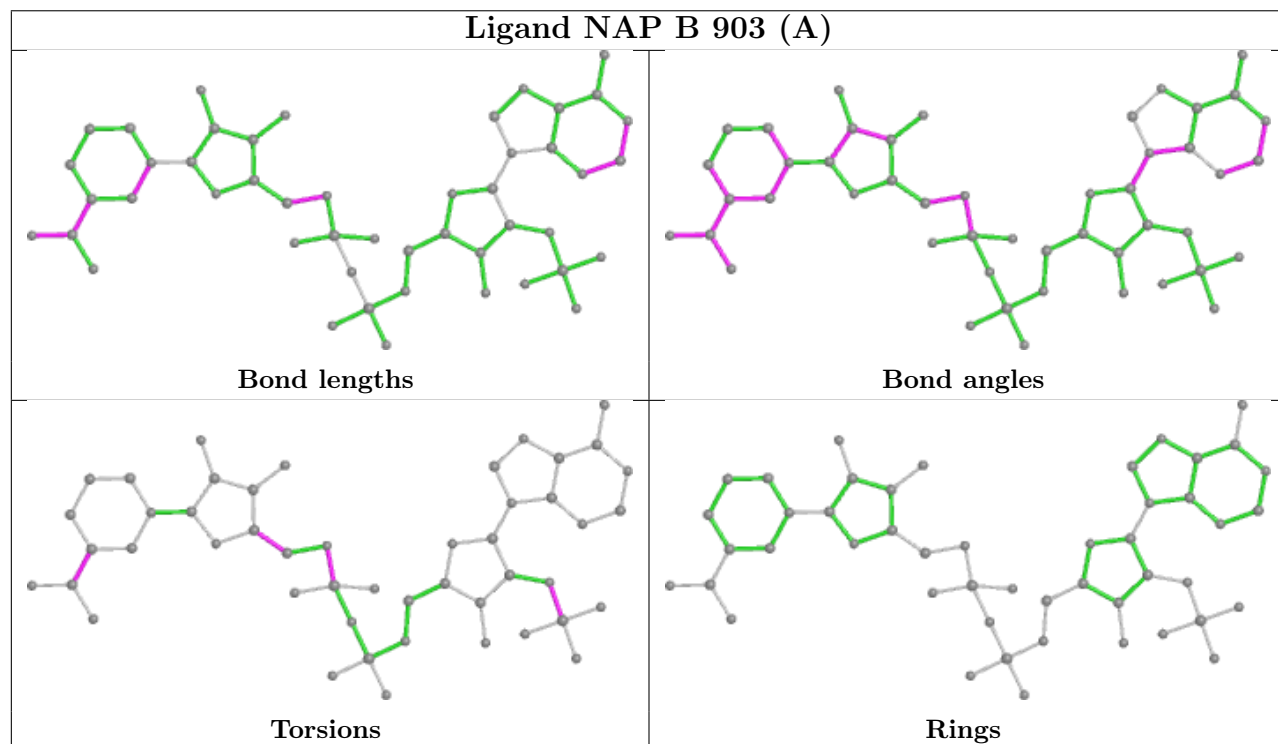
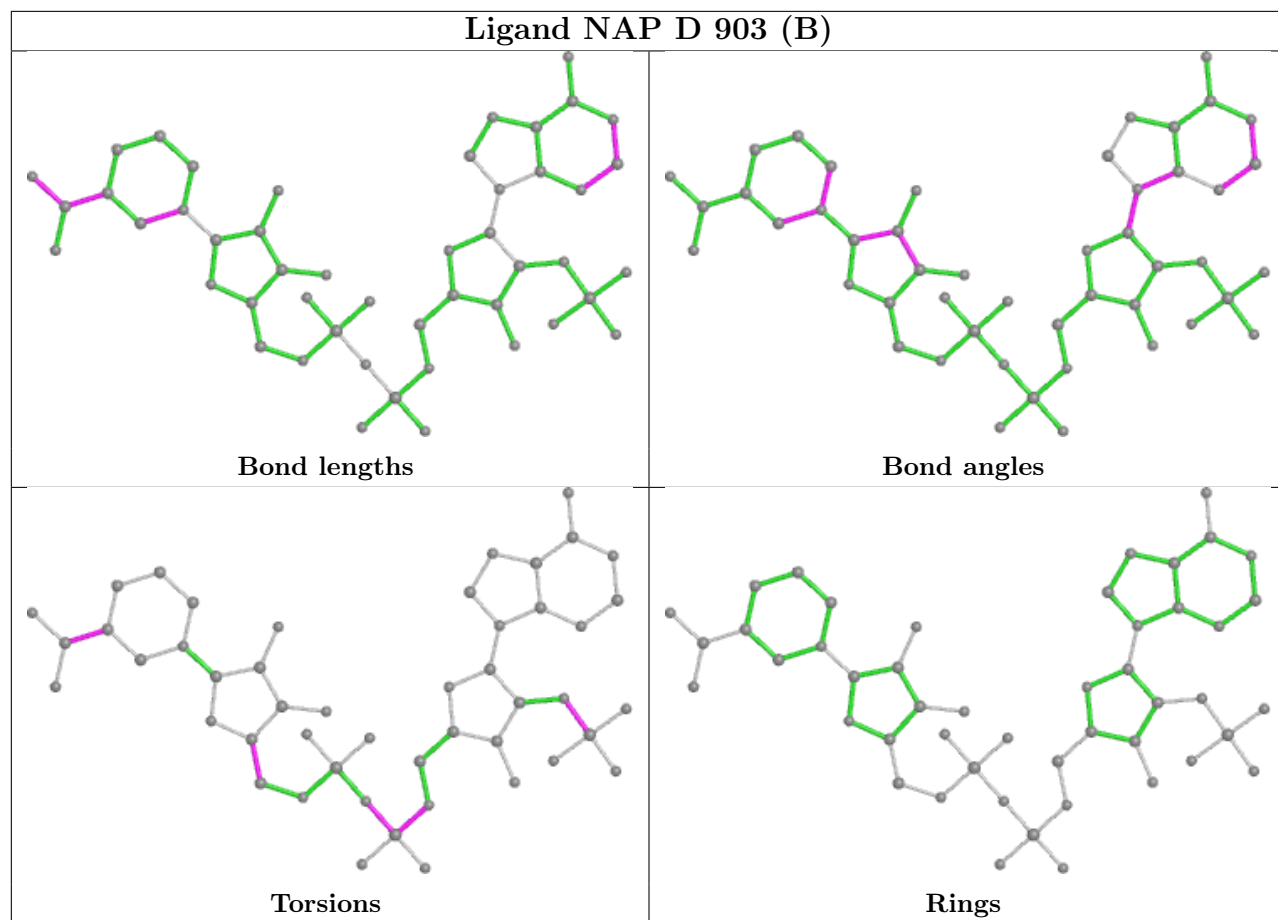
9 monomers are involved in 10 short contacts:

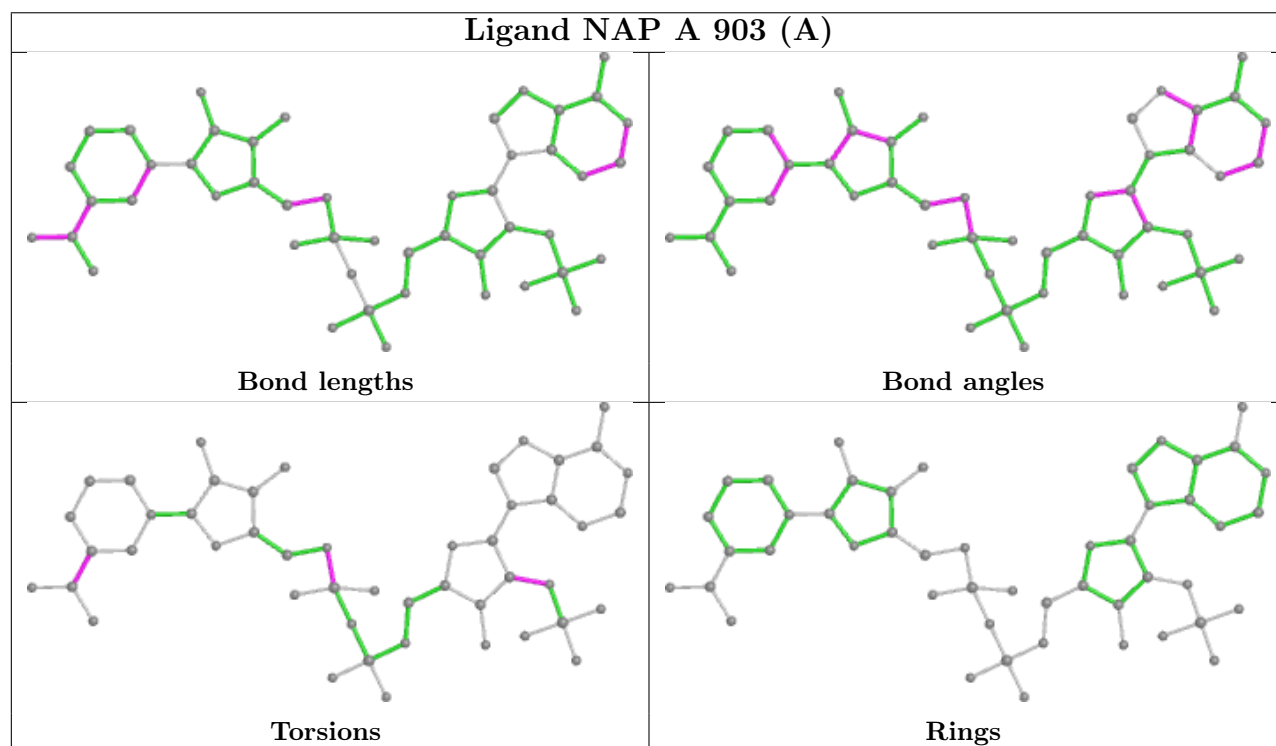
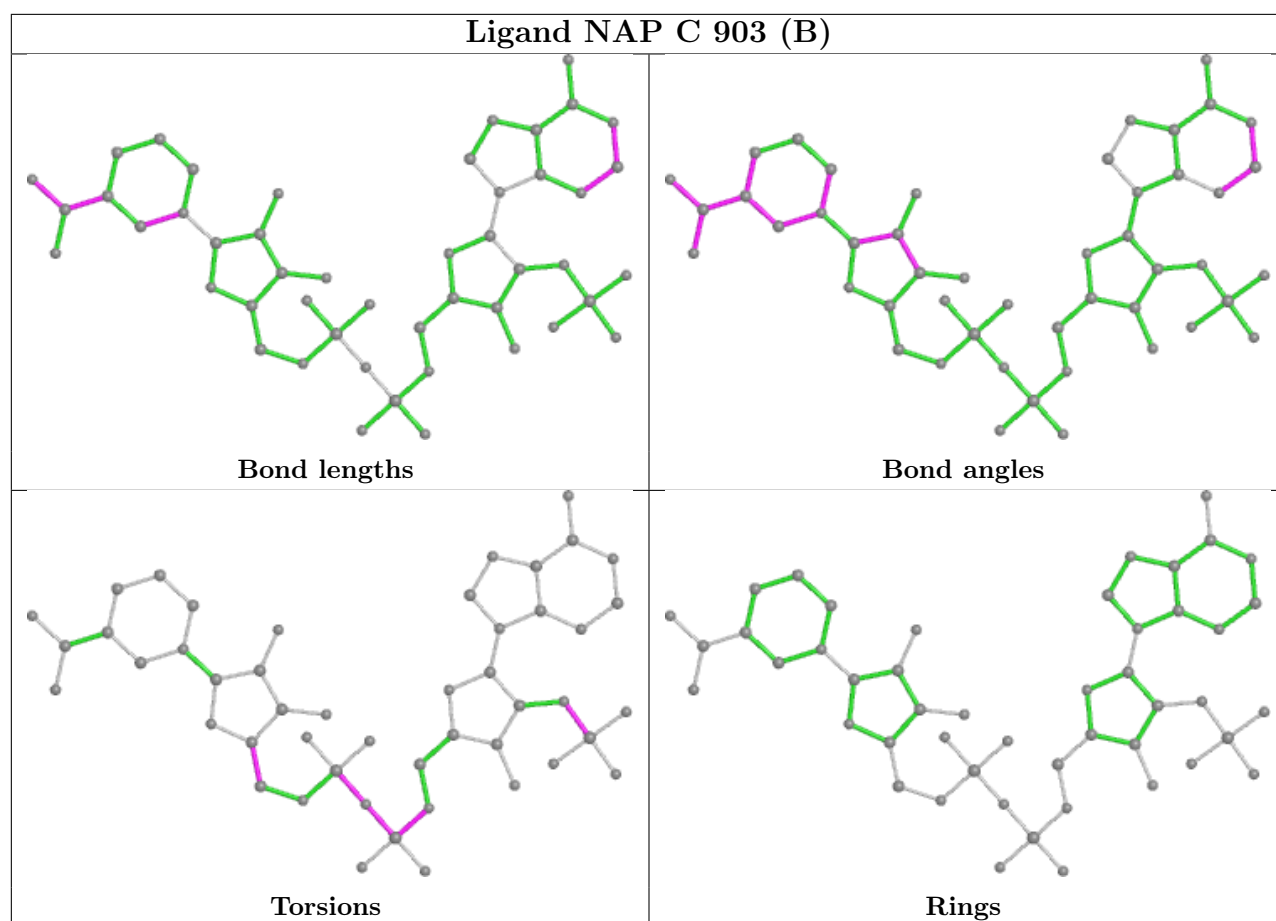
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	3002	SO4	1	0
4	B	3026	GOL	2	0
2	A	903[B]	NAP	1	0
3	B	3003	SO4	1	0
2	D	903[B]	NAP	1	0
2	C	903[B]	NAP	1	0
3	C	3028	SO4	1	0
3	D	3020	SO4	1	0
2	B	903[B]	NAP	1	0

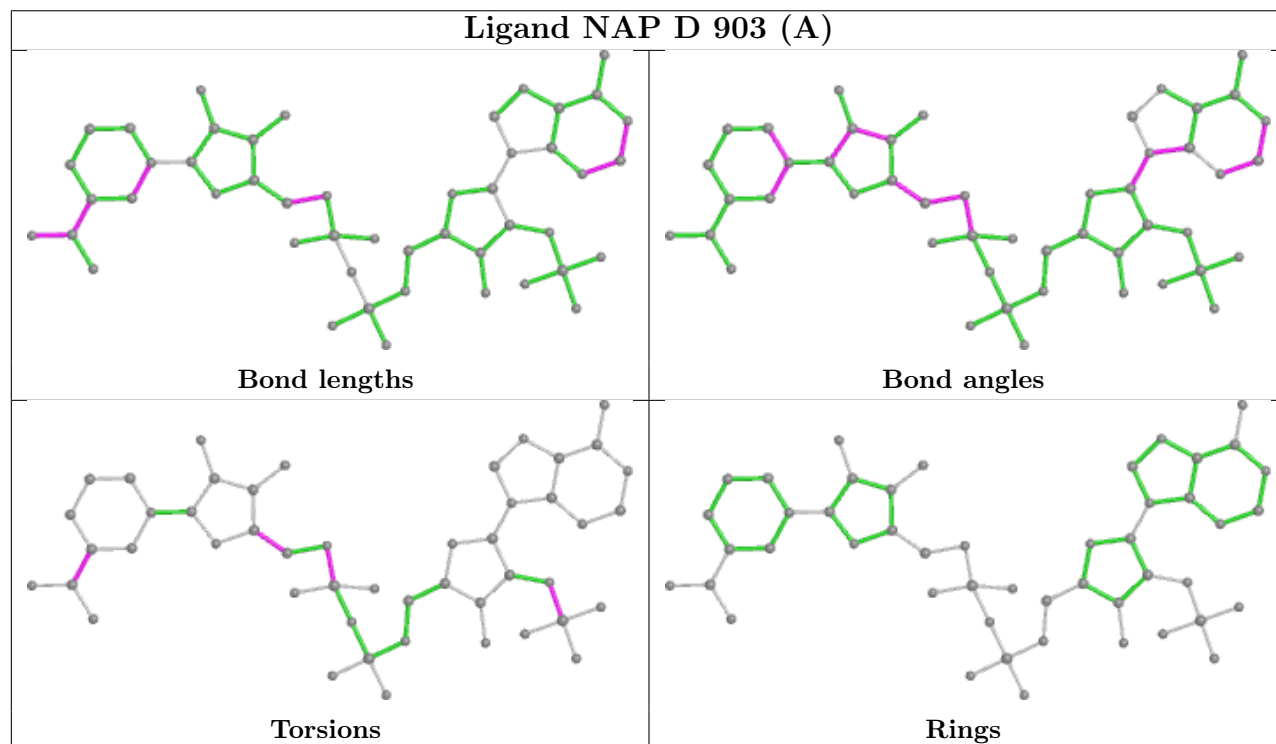
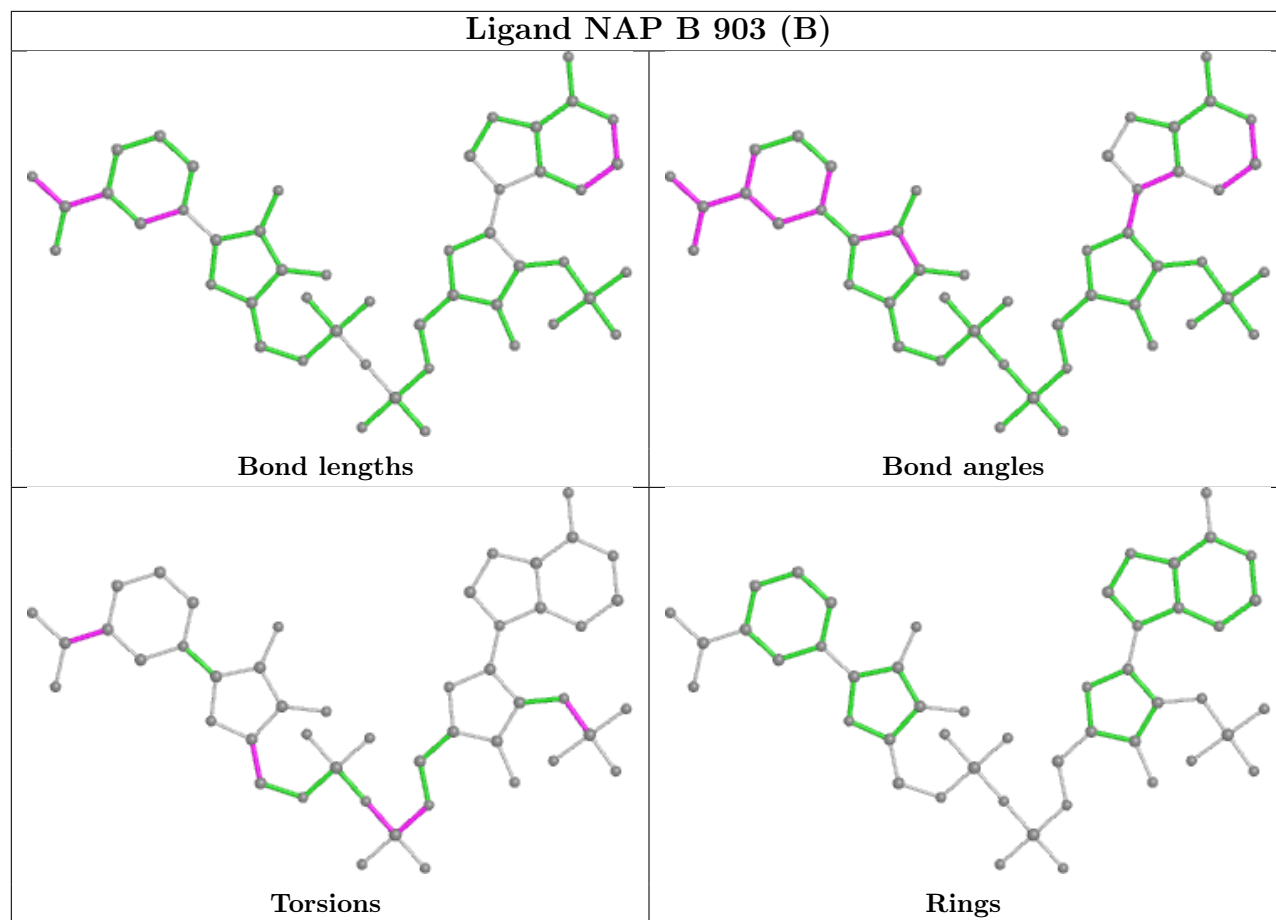
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











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	498/517 (96%)	-0.53	2 (0%) 92 93	13, 26, 38, 56	12 (2%)
1	B	498/517 (96%)	-0.50	4 (0%) 86 87	15, 28, 40, 63	14 (2%)
1	C	498/517 (96%)	-0.65	0 100 100	12, 24, 35, 61	17 (3%)
1	D	498/517 (96%)	-0.62	4 (0%) 86 87	13, 27, 41, 60	15 (3%)
All	All	1992/2068 (96%)	-0.57	10 (0%) 91 91	12, 26, 39, 63	58 (2%)

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	405	VAL	3.8
1	D	405	VAL	3.8
1	B	405	VAL	2.7
1	D	415	LYS	2.4
1	B	417	THR	2.2

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.

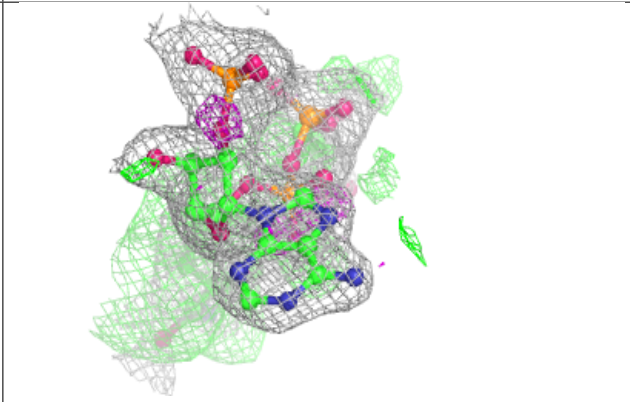
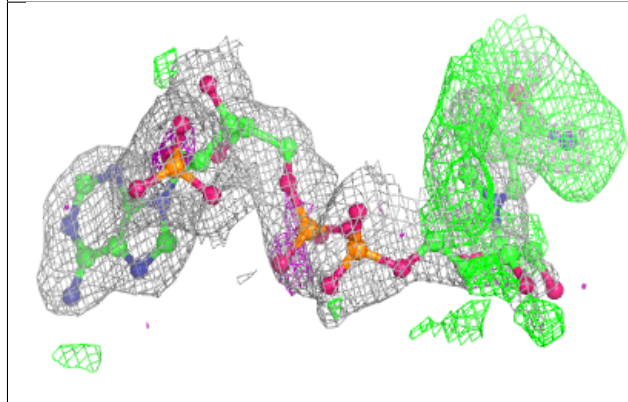
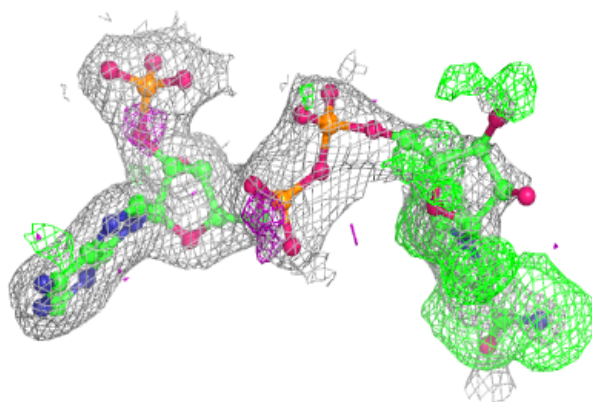
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SO4	D	3020	5/5	0.75	0.19	48,49,54,56	5
3	SO4	A	3021	5/5	0.79	0.27	42,50,55,56	5
3	SO4	A	3011	5/5	0.80	0.28	58,59,62,64	5
3	SO4	A	3018	5/5	0.83	0.17	50,51,54,55	5
3	SO4	B	3022	5/5	0.85	0.24	37,46,48,50	5
3	SO4	D	3012	5/5	0.86	0.16	49,54,56,58	5
4	GOL	C	3029	6/6	0.88	0.17	59,66,72,75	0
3	SO4	C	3010	5/5	0.89	0.16	47,48,53,55	5
3	SO4	C	3017	5/5	0.89	0.19	52,53,55,56	5
4	GOL	A	3025	6/6	0.89	0.21	59,65,68,70	0
3	SO4	C	3023	5/5	0.89	0.19	46,52,54,56	5
3	SO4	B	3019	5/5	0.90	0.18	50,50,54,56	5
4	GOL	D	3028	6/6	0.90	0.21	68,73,75,76	0
4	GOL	B	3026	6/6	0.91	0.17	50,58,60,63	0
3	SO4	C	3026	5/5	0.92	0.16	45,47,49,51	5
3	SO4	D	3027	5/5	0.92	0.20	42,45,48,49	5
3	SO4	A	3024	5/5	0.92	0.23	57,58,60,60	5
2	NAP	B	903[B]	48/48	0.93	0.15	31,37,39,46	25
3	SO4	B	3009	5/5	0.93	0.16	46,47,50,53	5
2	NAP	C	903[A]	31/48	0.93	0.15	26,29,34,35	8
2	NAP	C	903[B]	48/48	0.93	0.15	26,29,36,42	25
3	SO4	B	3025	5/5	0.93	0.14	47,52,53,53	5
2	NAP	A	903[A]	31/48	0.93	0.15	27,31,35,35	8
2	NAP	A	903[B]	48/48	0.93	0.15	27,35,38,45	25
2	NAP	B	903[A]	31/48	0.93	0.15	31,34,39,39	8
2	NAP	D	903[B]	48/48	0.94	0.13	26,29,36,40	25
2	NAP	D	903[A]	31/48	0.94	0.13	26,29,37,38	8
3	SO4	B	3005	5/5	0.95	0.17	37,42,49,50	5
3	SO4	C	3002	5/5	0.96	0.11	40,42,47,47	5
3	SO4	D	3008	5/5	0.96	0.11	34,35,39,41	5
3	SO4	A	3001	5/5	0.96	0.09	59,60,62,62	5
3	SO4	C	3007	5/5	0.97	0.11	38,40,46,49	5
3	SO4	B	3003	5/5	0.98	0.12	28,28,32,33	5
3	SO4	A	3006	5/5	0.98	0.11	31,34,41,45	5
3	SO4	A	3015	5/5	0.98	0.14	30,42,48,51	0
3	SO4	C	3028	5/5	0.98	0.10	28,36,38,39	5
3	SO4	B	3014	5/5	0.99	0.20	34,45,48,54	0
3	SO4	C	3013	5/5	0.99	0.12	32,41,45,47	0
3	SO4	B	3004	5/5	0.99	0.08	35,41,42,42	5
3	SO4	D	3016	5/5	0.99	0.17	30,43,45,50	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.

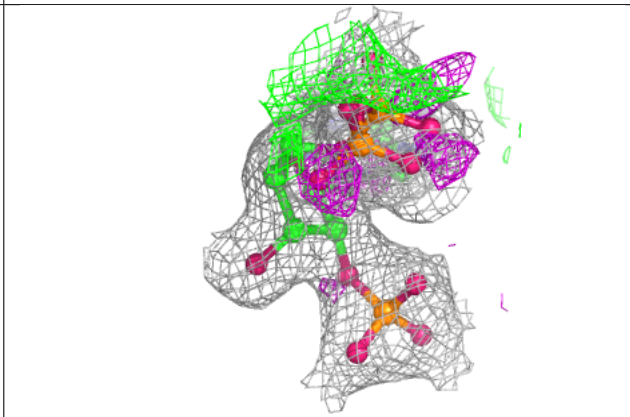
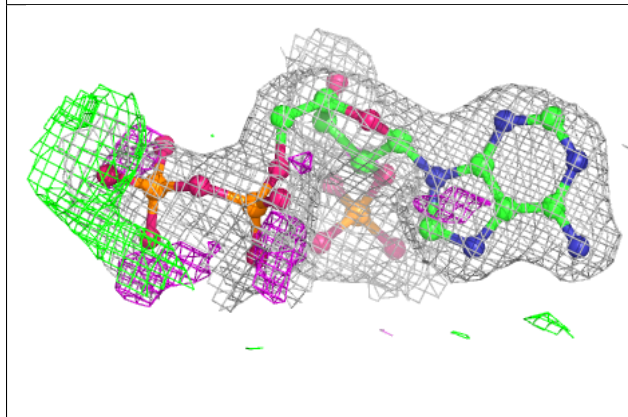
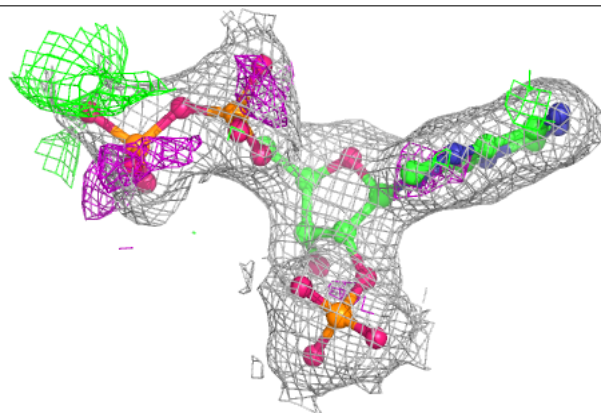
Electron density around NAP B 903 (B):

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



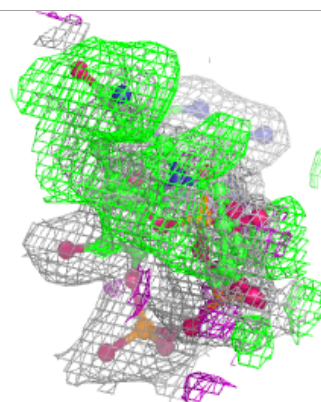
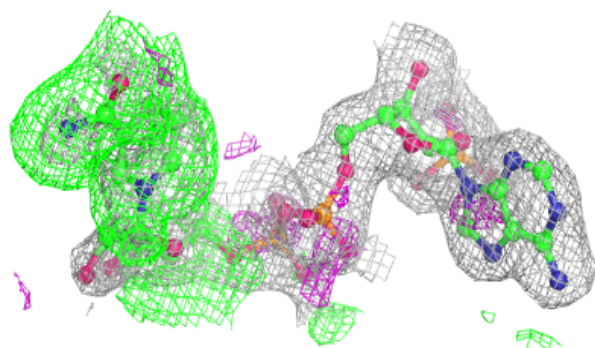
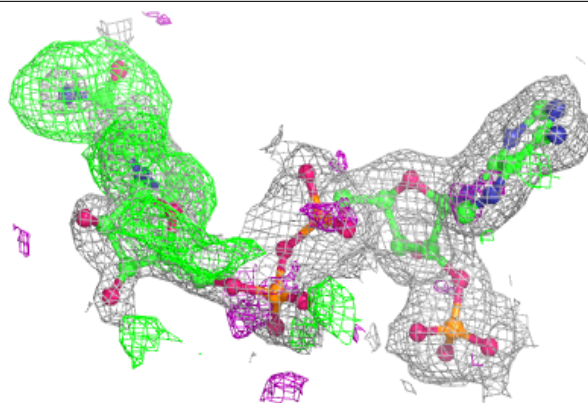
Electron density around NAP C 903 (A):

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

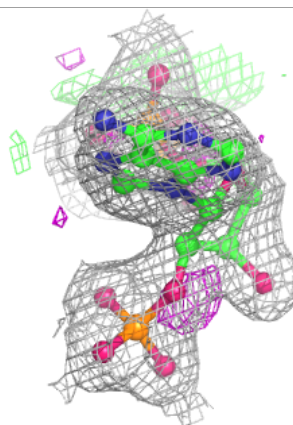
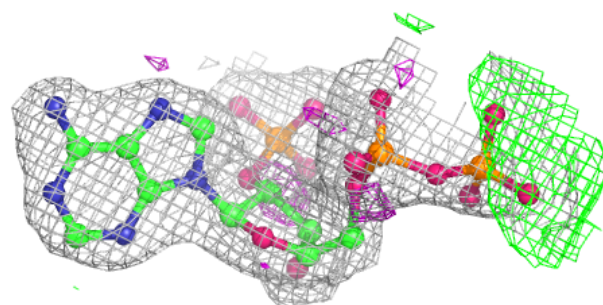
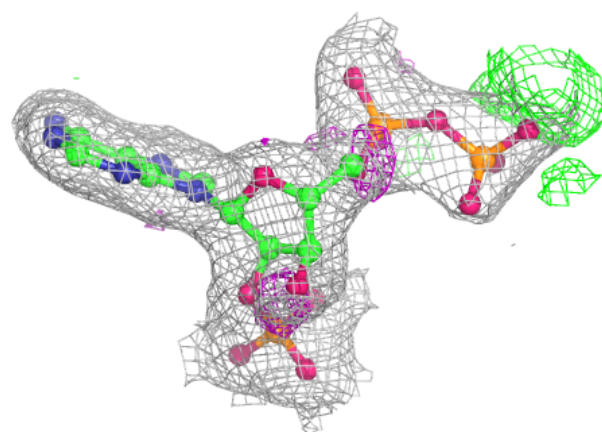


Electron density around NAP C 903 (B):

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

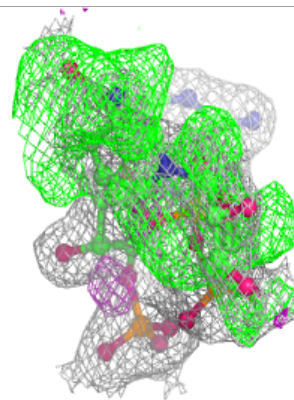
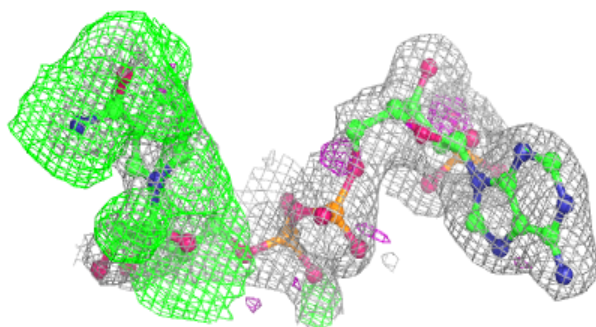
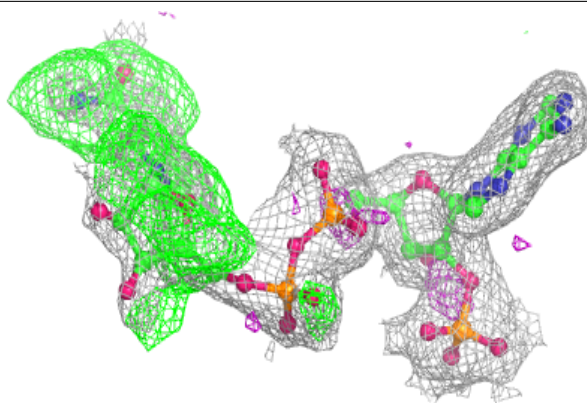
**Electron density around NAP A 903 (A):**

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

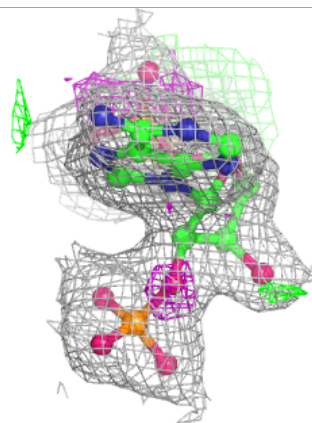
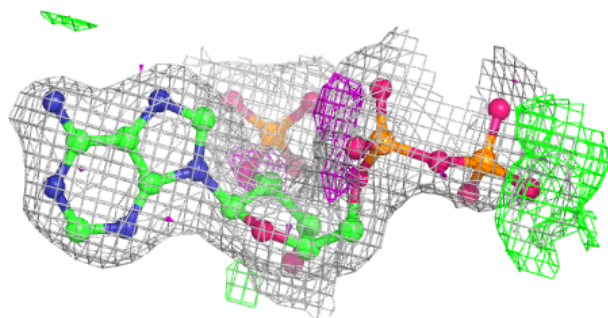
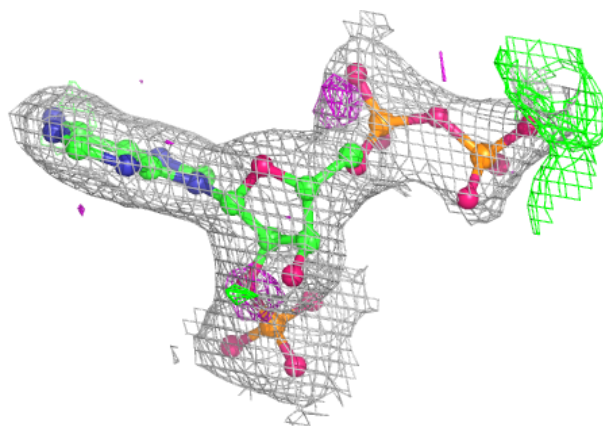


Electron density around NAP A 903 (B):

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

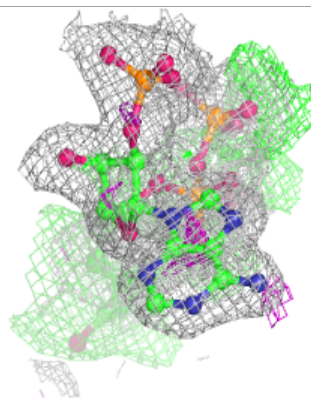
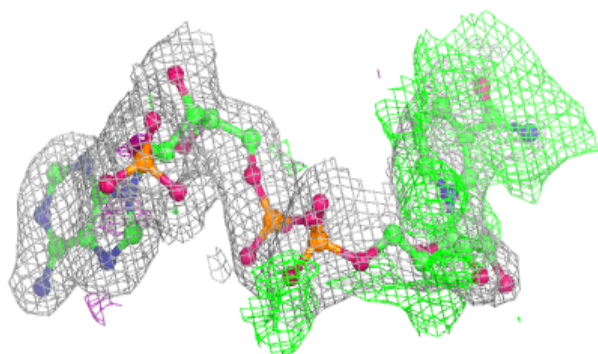
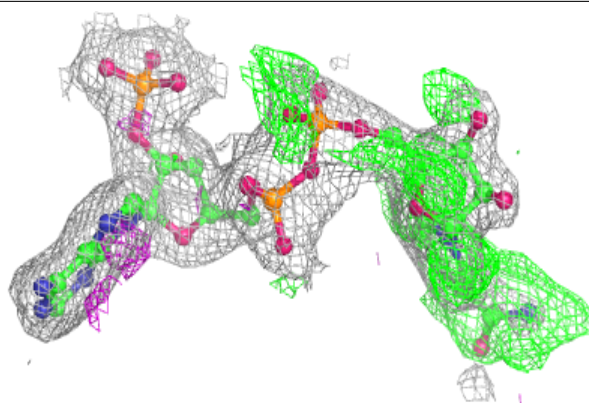
**Electron density around NAP B 903 (A):**

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

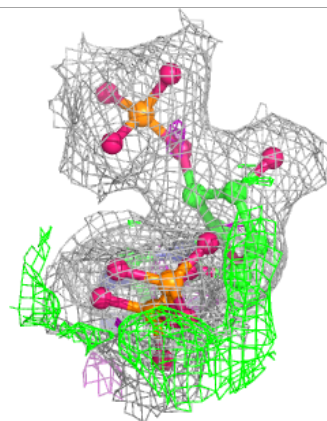
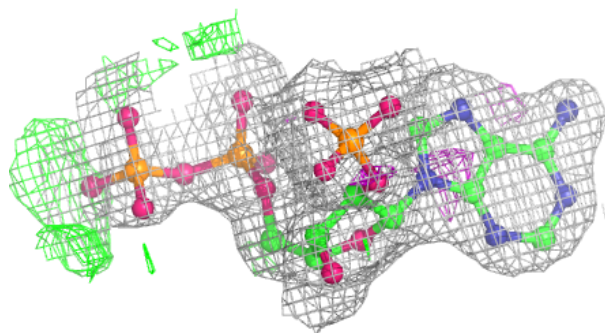
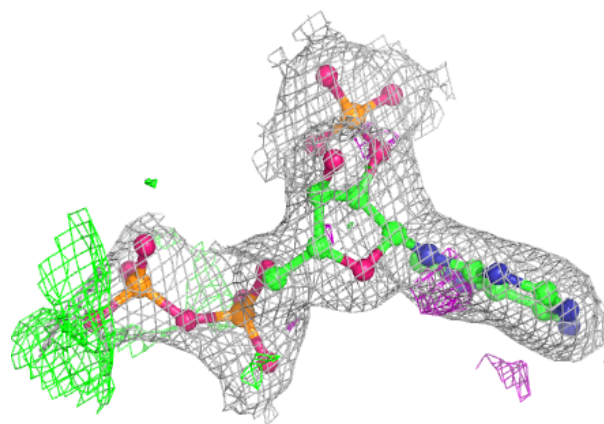


Electron density around NAP D 903 (B):

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

**Electron density around NAP D 903 (A):**

$2mF_o-DF_c$ (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.