



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 06:21 am BST

PDB ID : 6AJ6
Title : Crystal structure of Trypanosoma brucei glycosomal isocitrate dehydrogenase in complex with NADP+
Authors : Wang, X.; Inaoka, D.K.; Shiba, T.; Balogun, E.O.; Ziebart, N.; Allman, S.; Watanabe, Y.; Nozaki, T.; Boshart, M.; Bringaud, F.; Harada, S.; Kita, K.
Deposited on : 2018-08-27
Resolution : 3.20 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 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

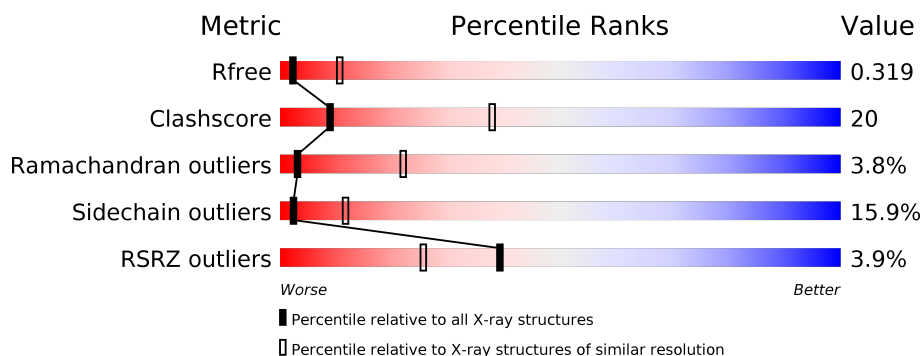
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	413	<div> <div>4%</div> <div> <div></div> <div>54%</div> <div>36%</div> <div>9%</div> </div> <div></div> </div>
1	C	413	<div> <div>4%</div> <div> <div></div> <div>68%</div> <div>26%</div> <div>6%</div> </div> <div></div> </div>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 6660 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.

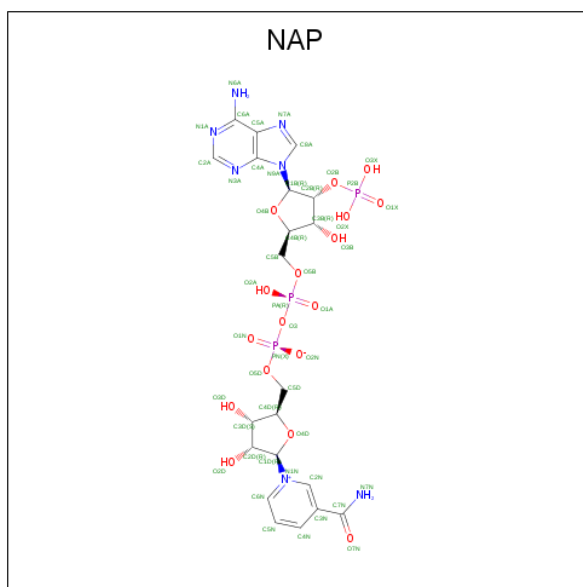
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	0	0
			3282	2084	563	613	22			
1	C	413	Total	C	N	O	S	0	0	0
			3282	2084	563	613	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	MET	engineered mutation	UNP Q387G0
C	1	SER	MET	engineered mutation	UNP Q387G0

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by author).



Continued from previous page...

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

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). 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.

Chain A:

4% 54% 36% 9%

Index	Value	Category
V345	D143	Green
D346	K72	Green
N347	C73	Green
N348	A74	Green
L351	T75	Green
L352	T76	Red
E353	T77	Green
R357	D79	Green
I362	R82	Green
I363	W83	Green
A364	E85	Green
T365	L88	Green
I366	K89	Green
A370	K90	Green
T371	W91	Green
E372	M92	Green
D374	R93	Green
I377	S94	Green
G378	P95	Green
I379	N96	Green
K380	G97	Green
G381	T98	Green
E382	T101	Green
D889	L103	Green
Y390	G104	Green
L391	G105	Green
N392	R109	Green
T393	I112	Green
D394	C114	Green
I397	A199	Green
V400	R202	Red
K405	K203	Red
V413	W204	Green
	P205	Green
	L206	Green
	E233	Green
	T324	Green
	S325	Green
	T326	Green
	N327	Green
	K216	Green
	H217	Green
	Y218	Green
	P226	Green
	E235	Green
	H247	Red
	R342	Green
	A344	Green
	T345	Green
	D143	Green
	K72	Green
	C73	Green
	A74	Green
	T75	Green
	T76	Red
	T77	Green
	D79	Green
	R82	Green
	W83	Green
	E85	Green
	L88	Green
	K89	Green
	K90	Green
	W91	Green
	M92	Green
	R93	Green
	S94	Green
	P95	Green
	N96	Green
	G97	Green
	T98	Green
	T101	Green
	L103	Green
	G104	Green
	G105	Green
	R109	Green
	I112	Green
	C114	Green
	A199	Green
	R202	Red
	K203	Red
	W204	Green
	P205	Green
	L206	Green
	E233	Green
	T324	Green
	S325	Green
	T326	Green
	N327	Green
	K216	Green
	H217	Green
	Y218	Green
	P226	Green
	E235	Green
	H247	Red
	R342	Green
	A344	Green
	T345	Green
	D143	Green
	K72	Green
	C73	Green
	A74	Green
	T75	Green
	T76	Red
	T77	Green
	D79	Green
	R82	Green
	W83	Green
	E85	Green
	L88	Green
	K89	Green
	K90	Green
	W91	Green
	M92	Green
	R93	Green
	S94	Green
	P95	Green
	N96	Green
	G97	Green
	T98	Green
	T101	Green
	L103	Green
	G104	Green
	G105	Green
	R109	Green
	I112	Green
	C114	Green
	A199	Green
	R202	Red
	K203	Red
	W204	Green
	P205	Green
	L206	Green
	E233	Green
	T324	Green
	S325	Green
	T326	Green
	N327	Green
	K216	Green
	H217	Green
	Y218	Green
	P226	Green
	E235	Green
	H247	Red
	R342	Green
	A344	Green
	T345	Green
	D143	Green
	K72	Green
	C73	Green
	A74	Green
	T75	Green
	T76	Red
	T77	Green
	D79	Green
	R82	Green
	W83	Green
	E85	Green
	L88	Green
	K89	Green
	K90	Green
	W91	Green
	M92	Green
	R93	Green
	S94	Green
	P95	Green
	N96	Green
	G97	Green
	T98	Green

Chain C:

68% 26% 6%

4%

Chain C: A horizontal bar chart showing a distribution of values. The bar is primarily green, with a small yellow segment at the end. The values are 68%, 26%, and 6%.

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.44Å 109.75Å 64.96Å 90.00° 89.98° 90.00°	Depositor
Resolution (Å)	45.02 – 3.20 45.03 – 3.20	Depositor EDS
% Data completeness (in resolution range)	83.9 (45.02-3.20) 87.0 (45.03-3.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.212 , 0.309 0.228 , 0.319	Depositor DCC
R_{free} test set	582 reflections (4.60%)	wwPDB-VP
Wilson B-factor (Å ²)	68.6	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -l,k,h 0.226 for h,-k,-l 0.008 for -l,-k,-h	Xtriage
Reported twinning fraction	0.609 for H, K, L 0.391 for h,-k,-l	Depositor
Outliers	0 of 12657 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6660	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.78	0/3350	1.05	10/4522 (0.2%)
1	C	0.68	1/3350 (0.0%)	0.88	5/4522 (0.1%)
All	All	0.73	1/6700 (0.0%)	0.97	15/9044 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	C	0	4
All	All	0	9

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	186	ASP	CB-CG	5.50	1.63	1.51

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	137	ASP	N-CA-C	-8.48	88.11	111.00
1	C	55	ARG	N-CA-C	7.20	130.43	111.00
1	A	90	LYS	N-CA-C	6.43	128.37	111.00
1	A	53	GLU	N-CA-C	-6.29	94.02	111.00
1	A	55	ARG	N-CA-C	6.26	127.90	111.00

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	GLY	Peptide
1	A	136	GLY	Peptide
1	A	139	TYR	Peptide
1	A	320	LYS	Peptide
1	A	71	VAL	Peptide

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	3282	0	3274	167	0
1	C	3282	0	3271	107	0
2	A	48	0	25	4	0
2	C	48	0	25	3	0
All	All	6660	0	6595	269	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 20.

The worst 5 of 269 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:341:HIS:O	1:C:345:VAL:HG23	1.31	1.25
1:A:79:ASP:O	1:A:83:VAL:HG13	1.61	1.01
1:C:136:GLY:HA3	1:C:140:SER:OG	1.61	1.00
1:C:341:HIS:O	1:C:345:VAL:CG2	2.10	1.00
1:A:79:ASP:CB	1:A:82:ARG:HB2	1.96	0.96

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	411/413 (100%)	318 (77%)	74 (18%)	19 (5%)	2	18
1	C	411/413 (100%)	355 (86%)	44 (11%)	12 (3%)	4	28
All	All	822/826 (100%)	673 (82%)	118 (14%)	31 (4%)	3	22

5 of 31 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	PHE
1	A	235	GLU
1	A	308	HIS
1	A	323	GLU
1	C	72	LYS

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	353/353 (100%)	288 (82%)	65 (18%)	1	8
1	C	353/353 (100%)	306 (87%)	47 (13%)	4	18
All	All	706/706 (100%)	594 (84%)	112 (16%)	2	12

5 of 112 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	346	ASP
1	A	394	ASP
1	C	357	ARG
1	A	353	GLU
1	A	374	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	HIS
1	C	392	ASN
1	A	392	ASN
1	A	67	HIS
1	C	270	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	501	-	45,52,52	0.93	2 (4%)	56,80,80	1.60	7 (12%)
2	NAP	A	501	-	45,52,52	0.97	3 (6%)	56,80,80	1.42	7 (12%)

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	501	-	-	10/31/67/67	0/5/5/5
2	NAP	A	501	-	-	6/31/67/67	0/5/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	NAP	O4D-C1D	2.90	1.45	1.41
2	C	501	NAP	O4D-C1D	2.79	1.45	1.41
2	A	501	NAP	C5A-C4A	2.58	1.47	1.40
2	C	501	NAP	C5A-C4A	2.19	1.46	1.40
2	A	501	NAP	C2A-N3A	2.06	1.35	1.32

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	NAP	C6N-N1N-C2N	-3.88	118.44	121.97
2	C	501	NAP	PN-O3-PA	-3.76	119.91	132.83
2	C	501	NAP	C3D-C2D-C1D	3.72	106.59	100.98
2	A	501	NAP	C3N-C7N-N7N	3.68	122.17	117.75
2	A	501	NAP	PN-O3-PA	-3.61	120.45	132.83

There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	NAP	C5D-O5D-PN-O3
2	C	501	NAP	C5D-O5D-PN-O1N
2	C	501	NAP	C5D-O5D-PN-O2N
2	C	501	NAP	C2D-C1D-N1N-C2N
2	C	501	NAP	C2D-C1D-N1N-C6N

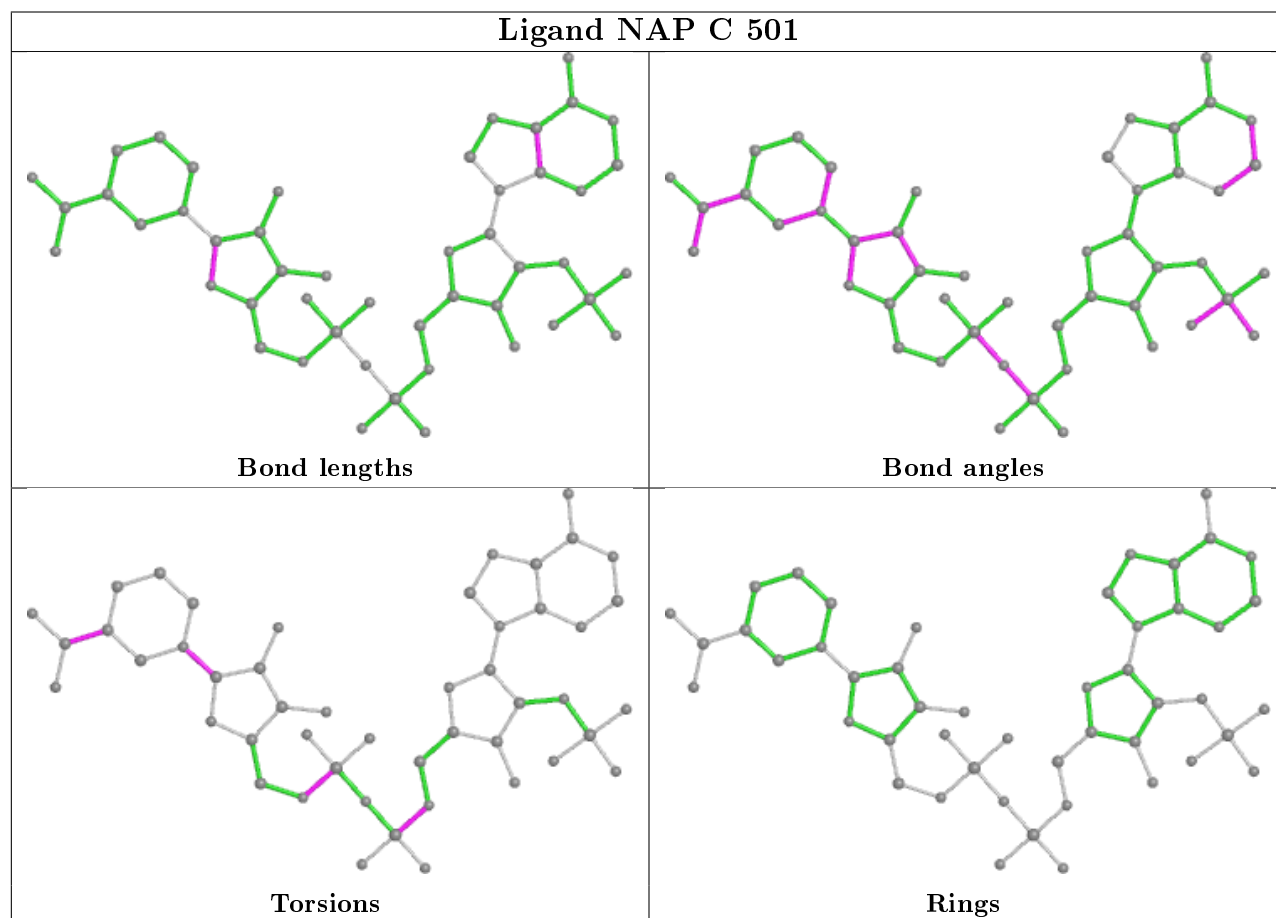
There are no ring outliers.

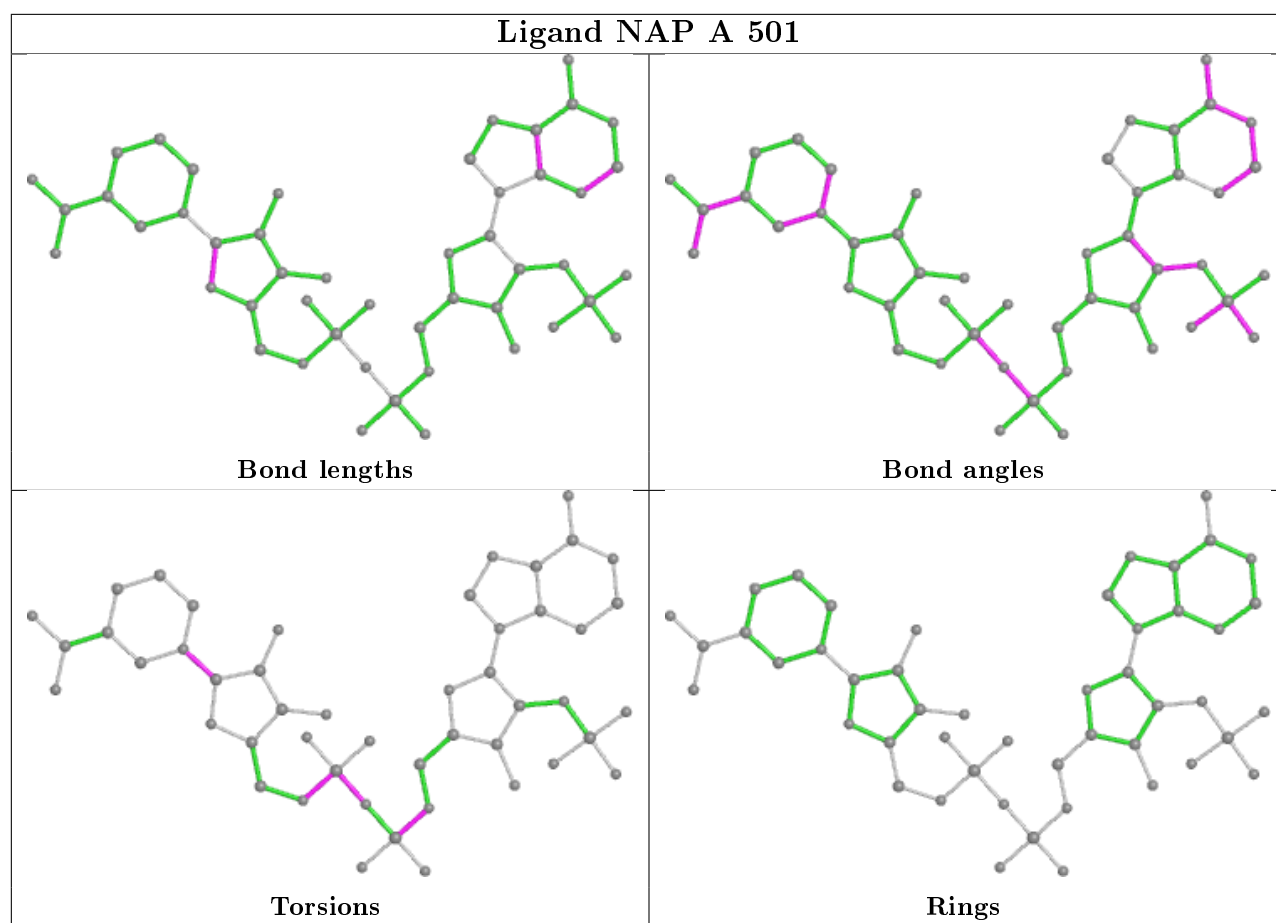
2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	NAP	3	0
2	A	501	NAP	4	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	413/413 (100%)	-0.05	17 (4%) 37 24	43, 84, 130, 163	0
1	C	413/413 (100%)	-0.08	15 (3%) 42 27	27, 67, 128, 197	0
All	All	826/826 (100%)	-0.06	32 (3%) 39 25	27, 76, 129, 197	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	319	GLN	8.5
1	C	284	PHE	8.4
1	C	283	GLY	7.1
1	C	139	TYR	6.4
1	A	257	CYS	4.8

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](#)

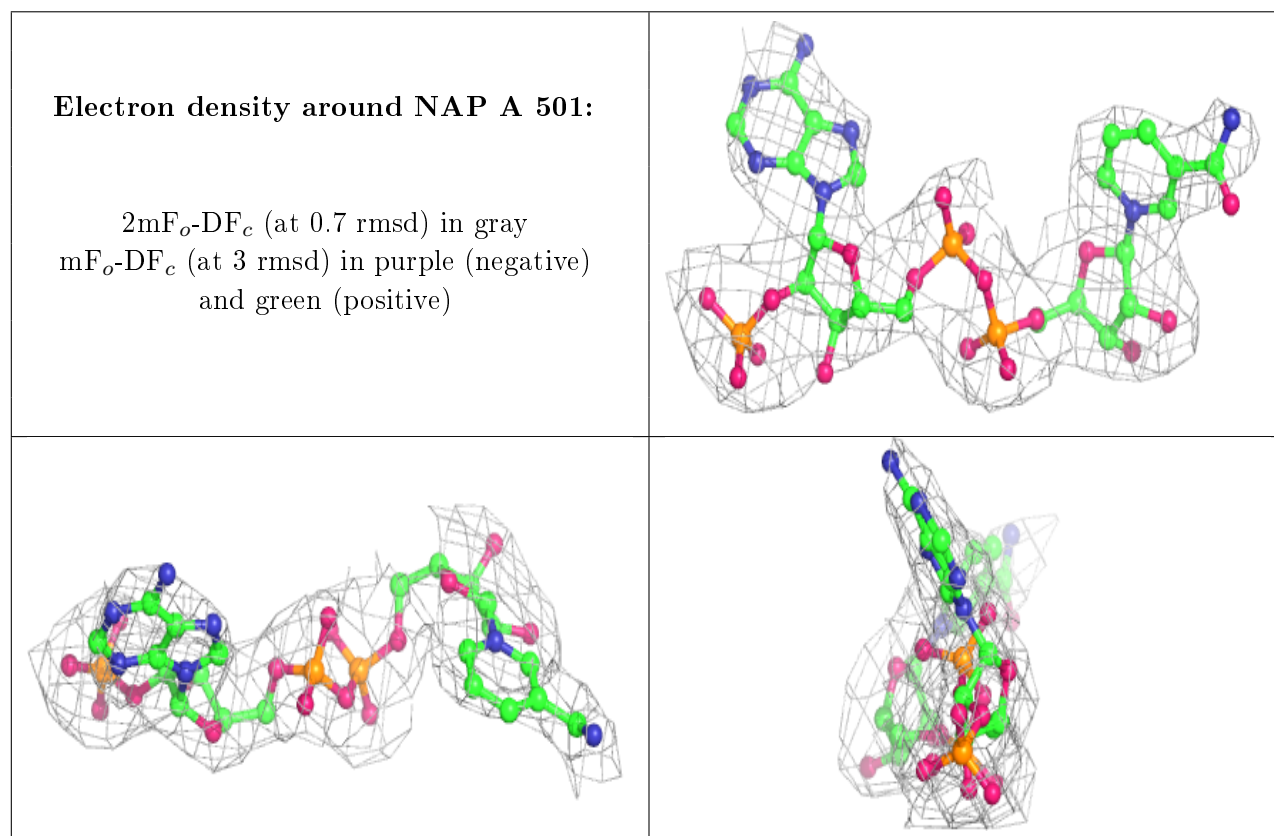
There are no carbohydrates in this entry.

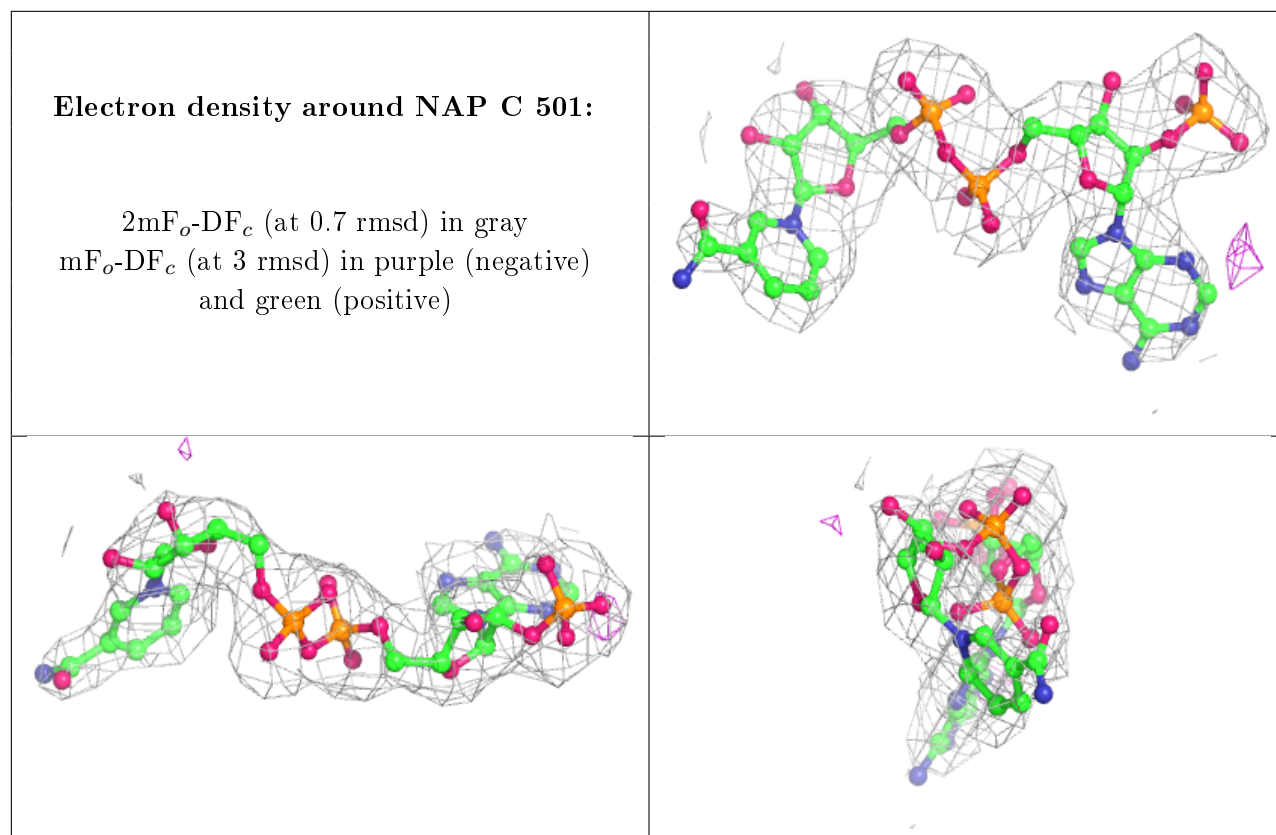
6.4 Ligands [i](#)

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
2	NAP	A	501	48/48	0.93	0.20	56,70,90,97	0
2	NAP	C	501	48/48	0.94	0.19	52,65,81,91	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.