



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 03:27 am BST

PDB ID : 6VEI
Title : Crystal Structure of Human Cytosolic Isocitrate Dehydrogenase (IDH1) R132H Mutant in Complex with NADPH and AG-881 (Vorasidenib) Inhibitor
Authors : Padyana, A.; Jin, L.
Deposited on : 2020-01-02
Resolution : 2.10 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

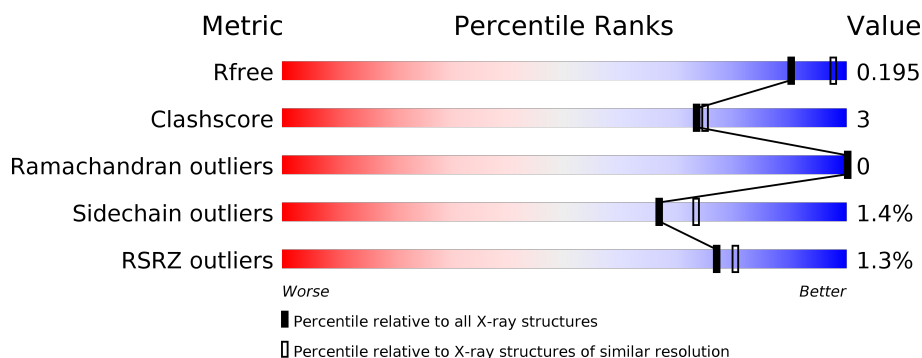
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 88%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 88% 9% </div> </div>
1	B	425	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 91%, grey 7%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 91% 7% </div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 7591 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.

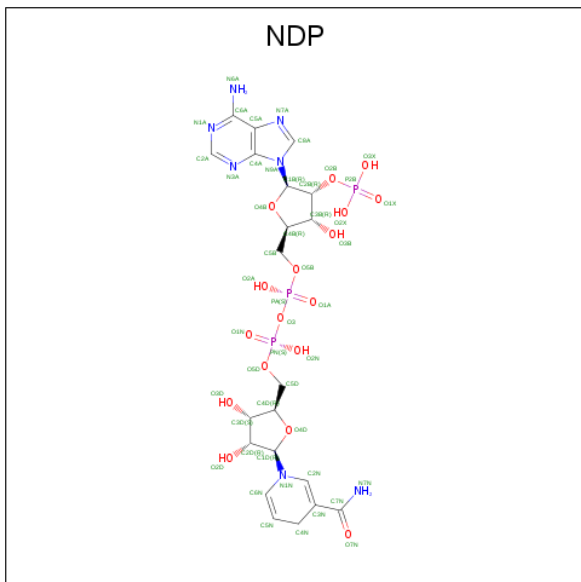
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	413	Total	C	N	O	S	0	11	0
			3320	2119	555	628	18			
1	B	417	Total	C	N	O	S	0	12	0
			3362	2144	566	633	19			

There are 24 discrepancies between the modelled and reference sequences:

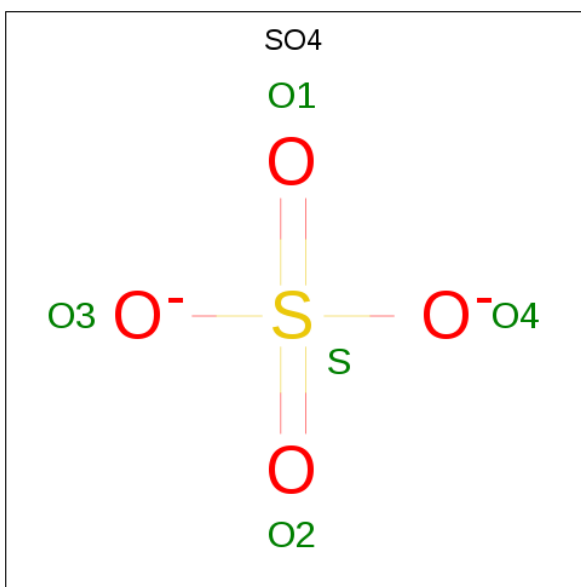
Chain	Residue	Modelled	Actual	Comment	Reference
A	132	HIS	ARG	engineered mutation	UNP O75874
A	415	SER	-	expression tag	UNP O75874
A	416	LEU	-	expression tag	UNP O75874
A	417	GLU	-	expression tag	UNP O75874
A	418	HIS	-	expression tag	UNP O75874
A	419	HIS	-	expression tag	UNP O75874
A	420	HIS	-	expression tag	UNP O75874
A	421	HIS	-	expression tag	UNP O75874
A	422	HIS	-	expression tag	UNP O75874
A	423	HIS	-	expression tag	UNP O75874
A	424	HIS	-	expression tag	UNP O75874
A	425	HIS	-	expression tag	UNP O75874
B	132	HIS	ARG	engineered mutation	UNP O75874
B	415	SER	-	expression tag	UNP O75874
B	416	LEU	-	expression tag	UNP O75874
B	417	GLU	-	expression tag	UNP O75874
B	418	HIS	-	expression tag	UNP O75874
B	419	HIS	-	expression tag	UNP O75874
B	420	HIS	-	expression tag	UNP O75874
B	421	HIS	-	expression tag	UNP O75874
B	422	HIS	-	expression tag	UNP O75874
B	423	HIS	-	expression tag	UNP O75874
B	424	HIS	-	expression tag	UNP O75874
B	425	HIS	-	expression tag	UNP O75874

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$).



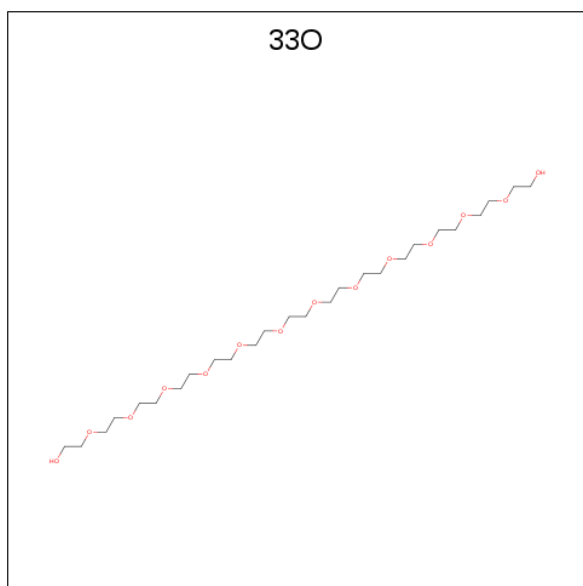
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	1
			96	42	14	34	6		
2	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



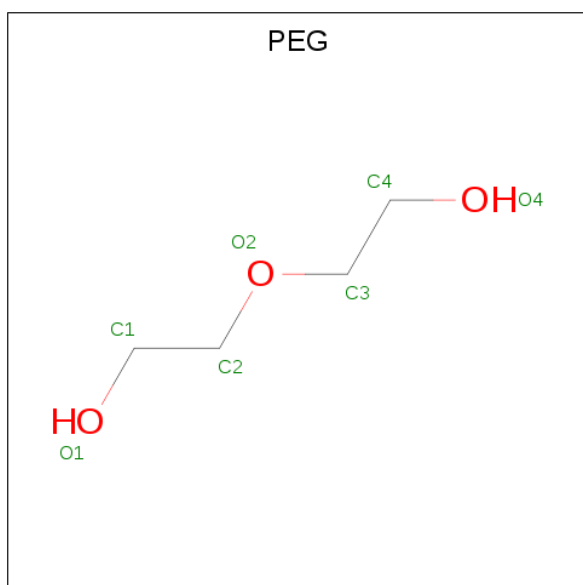
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is 3,6,9,12,15,18,21,24,27,30,33,36-dodecaoxaoctatriacontane-1,38-diol (three-letter code: 33O) (formula: $C_{26}H_{54}O_{14}$).



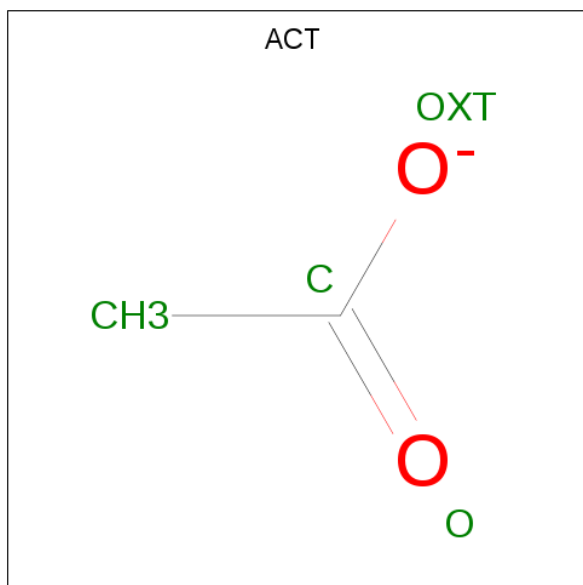
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			40	26	14		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



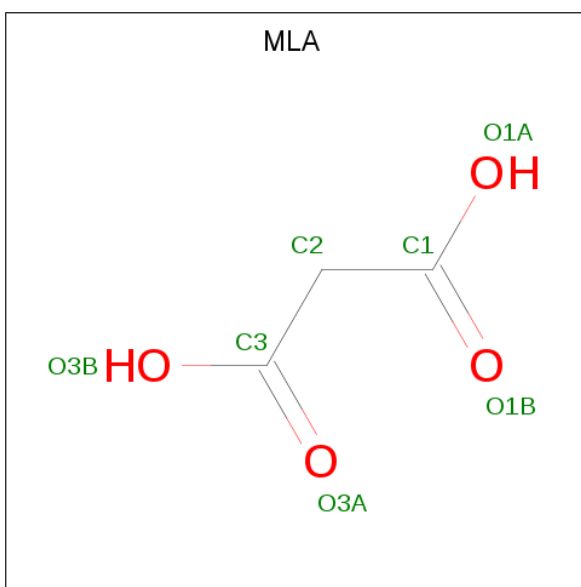
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	1
			14	8	6		
5	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



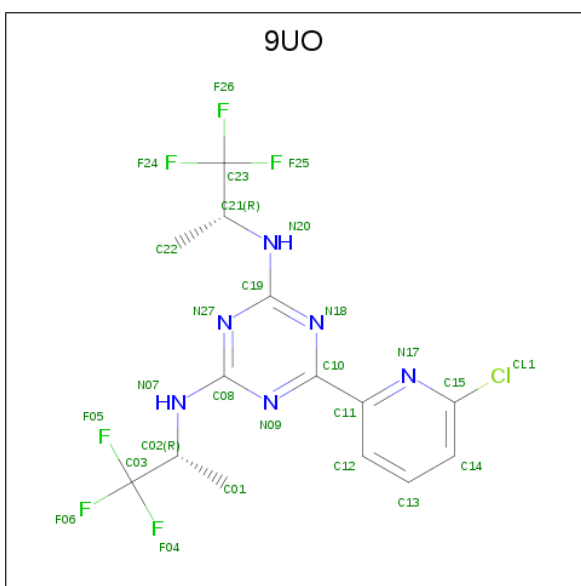
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is MALONIC ACID (three-letter code: MLA) (formula: $C_3H_4O_4$).



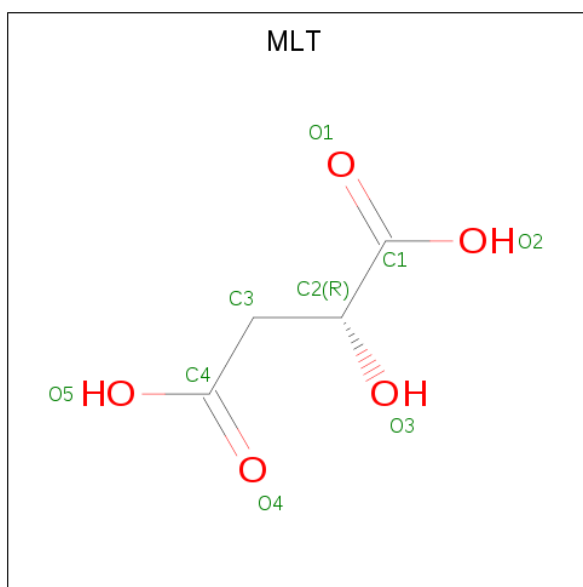
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	O		0	0
			7	3	4			

- Molecule 8 is 6-(6-chloropyridin-2-yl)-N2,N4-bis[(2R)-1,1,1-trifluoropropan-2-yl]-1,3,5-triazine-2,4-diamine (three-letter code: 9UO) (formula: $C_{14}H_{13}ClF_6N_6$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	B	1	Total	C	Cl	F	N	0	0
			27	14	1	6	6		

- Molecule 9 is D-MALATE (three-letter code: MLT) (formula: $C_4H_6O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			9	4	5		

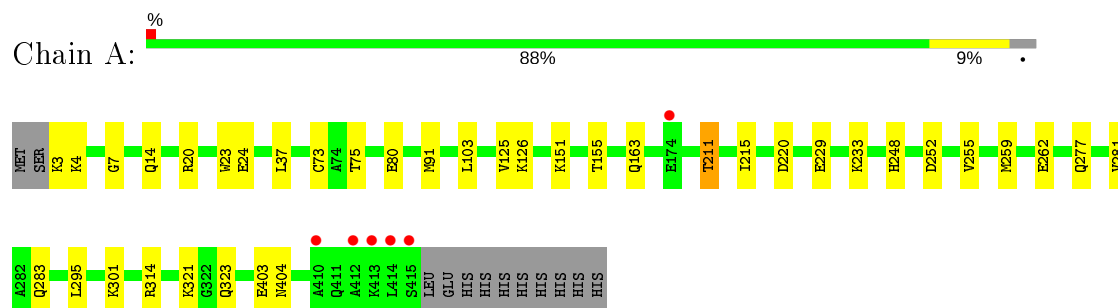
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	292	Total	O	0	0
			292	292		
10	B	346	Total	O	0	0
			346	346		

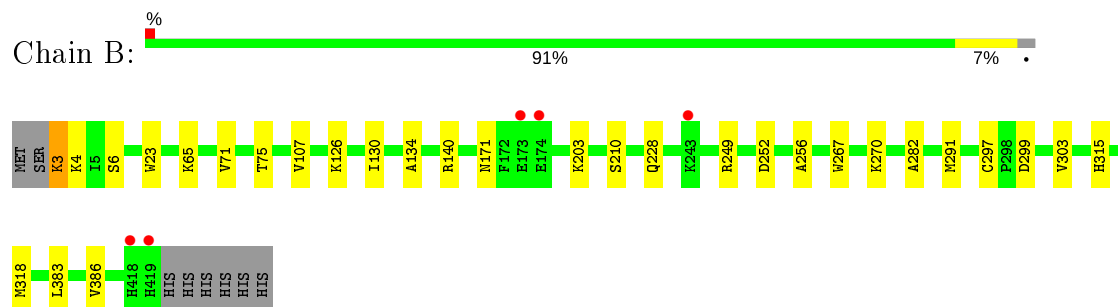
3 Residue-property plots

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.

- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	117.21Å 85.10Å 96.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.51 – 2.10 35.51 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.51-2.10) 99.8 (35.51-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	25.84 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.158 , 0.195 0.157 , 0.195	Depositor DCC
R_{free} test set	2886 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.1	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7591	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.26% 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: MLA, 9UO, 33O, SO4, MLT, ACT, NDP, PEG

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	0.25	0/3419	0.42	0/4614
1	B	0.24	0/3463	0.42	0/4670
All	All	0.24	0/6882	0.42	0/9284

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

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	3320	0	3321	25	0
1	B	3362	0	3358	19	0
2	A	96	0	52	2	0
2	B	48	0	26	4	0
3	A	5	0	0	0	0
4	A	40	0	54	0	0
5	A	14	0	20	0	0
5	B	21	0	30	4	0
6	A	4	0	3	0	0
7	A	7	0	2	0	0
8	B	27	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	9	0	4	0	0
10	A	292	0	0	5	0
10	B	346	0	0	3	0
All	All	7591	0	6870	46	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 3.

All (46) 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:3:LYS:HD2	1:A:4:LYS:HG3	1.52	0.90
1:A:73:CYS:SG	10:A:601:HOH:O	2.41	0.78
2:B:501:NDP:H42N	5:B:503[B]:PEG:H12	1.67	0.75
1:A:301:LYS:NZ	10:A:602:HOH:O	2.23	0.71
1:A:20:ARG:HA	10:A:601:HOH:O	1.91	0.68
1:A:163:GLN:NE2	10:A:603:HOH:O	2.27	0.67
1:B:4[B]:LYS:NZ	10:B:604:HOH:O	2.36	0.59
1:B:270:LYS:NZ	10:B:603:HOH:O	2.32	0.58
1:A:163:GLN:NE2	10:A:609:HOH:O	2.37	0.58
1:A:103:LEU:HD12	1:A:295:LEU:HD23	1.88	0.56
1:B:75:THR:O	2:B:501:NDP:H2N	2.07	0.55
1:A:283:GLN:OE1	1:B:256:ALA:HB1	2.06	0.55
1:A:211[A]:THR:HG23	1:A:220:ASP:HB3	1.89	0.53
1:B:23:TRP:HH2	1:B:71[B]:VAL:HG22	1.74	0.52
1:A:14:GLN:NE2	1:A:24:GLU:OE2	2.37	0.52
1:B:228:GLN:NE2	10:B:602:HOH:O	2.32	0.50
1:A:80:GLU:HG2	1:A:91:MET:HE1	1.93	0.50
1:A:314:ARG:NE	2:A:501[B]:NDP:O1X	2.34	0.50
1:A:403:GLU:HG3	1:A:404:ASN:N	2.27	0.49
1:B:23:TRP:HH2	1:B:71[A]:VAL:HG12	1.76	0.49
1:A:321:LYS:NZ	1:A:323:GLN:NE2	2.60	0.49
1:A:215:ILE:HD11	5:B:503[B]:PEG:H42	1.94	0.49
1:A:23:TRP:CD2	1:A:73:CYS:HB2	2.48	0.49
1:B:23:TRP:CH2	1:B:71[B]:VAL:HG22	2.48	0.48
2:B:501:NDP:C4N	5:B:503[B]:PEG:H12	2.41	0.48
1:A:75:THR:O	2:A:501[B]:NDP:H2N	2.14	0.47
1:A:229:GLU:OE2	1:A:233:LYS:HE2	2.15	0.46
1:A:125:VAL:HG13	1:A:126:LYS:HG3	1.97	0.46
1:B:130[B]:ILE:HD13	1:B:267:TRP:HB3	1.99	0.45
1:A:211[A]:THR:HG22	1:A:248:HIS:NE2	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LYS:HE2	1:B:65:LYS:HB3	1.64	0.44
1:A:125:VAL:HG12	1:A:262:GLU:HB3	2.00	0.44
2:B:501:NDP:C5N	5:B:503[A]:PEG:H11	2.48	0.44
1:B:107:VAL:HG23	1:B:134:ALA:HB2	1.98	0.44
1:B:383:LEU:O	1:B:386[A]:VAL:HG12	2.17	0.43
1:B:23:TRP:CH2	1:B:71[A]:VAL:HG12	2.54	0.43
1:B:297:CYS:HB2	1:B:299:ASP:OD2	2.18	0.43
1:B:3:LYS:C	1:B:3:LYS:HD2	2.39	0.43
1:A:7:GLY:HA3	1:A:37:LEU:HD23	2.00	0.43
1:A:277:GLN:O	1:A:281[A]:VAL:HG13	2.19	0.42
1:B:282:ALA:HB2	1:B:291:MET:SD	2.59	0.42
1:B:315:HIS:HA	1:B:318:MET:HE2	2.00	0.42
1:A:252:ASP:N	1:A:252:ASP:OD1	2.53	0.42
1:B:210:SER:HA	1:B:249:ARG:O	2.20	0.42
1:A:255:VAL:O	1:A:259:MET:HG3	2.21	0.41
1:B:203:LYS:HB3	1:B:203:LYS:HE3	1.75	0.41

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	422/425 (99%)	411 (97%)	11 (3%)	0	100	100
1	B	427/425 (100%)	413 (97%)	14 (3%)	0	100	100
All	All	849/850 (100%)	824 (97%)	25 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	360/361 (100%)	355 (99%)	5 (1%)	67	73
1	B	365/361 (101%)	357 (98%)	8 (2%)	52	57
All	All	725/722 (100%)	712 (98%)	13 (2%)	67	65

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	151	LYS
1	A	155[A]	THR
1	A	155[B]	THR
1	A	211[A]	THR
1	A	211[B]	THR
1	B	3	LYS
1	B	6	SER
1	B	126	LYS
1	B	140	ARG
1	B	171	ASN
1	B	252	ASP
1	B	303[A]	VAL
1	B	303[B]	VAL

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

Mol	Chain	Res	Type
1	A	323	GLN
1	A	411	GLN

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 ⓘ

14 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	NDP	A	501[B]	-	45,52,52	0.51	0	53,80,80	0.51	1 (1%)
5	PEG	B	503[B]	-	6,6,6	0.48	0	5,5,5	0.28	0
2	NDP	A	501[A]	-	45,52,52	0.51	0	53,80,80	0.52	1 (1%)
8	9UO	B	502	-	28,28,28	2.45	3 (10%)	32,42,42	3.13	10 (31%)
4	33O	A	503	-	39,39,39	0.55	0	38,38,38	0.37	0
6	ACT	A	506	-	1,3,3	6.13	1 (100%)	0,3,3	0.00	-
2	NDP	B	501	-	45,52,52	0.49	0	53,80,80	0.55	1 (1%)
5	PEG	A	505	-	6,6,6	0.49	0	5,5,5	0.28	0
7	MLA	A	507	-	0,6,6	0.00	-	0,7,7	0.00	-
5	PEG	B	503[A]	-	6,6,6	0.48	0	5,5,5	0.26	0
5	PEG	A	504	-	6,6,6	0.48	0	5,5,5	0.28	0
9	MLT	B	505	-	2,8,8	0.50	0	3,10,10	0.91	0
5	PEG	B	504	-	6,6,6	0.49	0	5,5,5	0.28	0
3	SO4	A	502	-	4,4,4	0.13	0	6,6,6	0.09	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
2	NDP	A	501[B]	-	-	11/30/77/77	0/5/5/5
5	PEG	B	503[B]	-	-	1/4/4/4	-
5	PEG	B	503[A]	-	-	1/4/4/4	-
2	NDP	A	501[A]	-	-	6/30/77/77	0/5/5/5
8	9UO	B	502	-	-	0/24/24/24	0/2/2/2
4	33O	A	503	-	-	18/37/37/37	-
2	NDP	B	501	-	-	9/30/77/77	0/5/5/5
5	PEG	A	505	-	-	1/4/4/4	-
5	PEG	B	504	-	-	2/4/4/4	-
5	PEG	A	504	-	-	0/4/4/4	-
9	MLT	B	505	-	-	0/2/8/8	-
7	MLA	A	507	-	-	0/0/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	502	9UO	C08-N07	8.53	1.45	1.34
8	B	502	9UO	C19-N20	8.24	1.45	1.34
6	A	506	ACT	CH3-C	6.13	1.56	1.48
8	B	502	9UO	C15-CL1	2.03	1.78	1.74

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	502	9UO	C08-N09-C10	8.23	119.85	114.60
8	B	502	9UO	C19-N18-C10	7.90	119.64	114.60
8	B	502	9UO	C19-N20-C21	-7.72	118.19	124.53
8	B	502	9UO	C08-N07-C02	-5.01	120.42	124.53
8	B	502	9UO	N27-C08-N09	-4.06	119.80	126.23
8	B	502	9UO	N27-C19-N18	-3.77	120.26	126.23
8	B	502	9UO	C19-N27-C08	3.64	120.06	113.89
8	B	502	9UO	N18-C10-N09	-3.00	120.38	125.23
2	A	501[B]	NDP	C5A-C6A-N6A	2.34	123.92	120.35
2	A	501[A]	NDP	C5A-C6A-N6A	2.32	123.88	120.35
2	B	501	NDP	C5A-C6A-N6A	2.30	123.85	120.35
8	B	502	9UO	C14-C15-N17	-2.21	121.43	125.50
8	B	502	9UO	N07-C08-N09	2.10	120.33	117.18

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501[B]	NDP	C5B-O5B-PA-O1A
2	A	501[B]	NDP	C5D-O5D-PN-O2N
2	A	501[A]	NDP	C5B-O5B-PA-O1A
2	A	501[A]	NDP	C5D-O5D-PN-O2N
2	B	501	NDP	C5B-O5B-PA-O1A
2	B	501	NDP	C5D-O5D-PN-O1N
2	B	501	NDP	C5D-O5D-PN-O2N
4	A	503	33O	O21-C43-C44-O22
4	A	503	33O	O22-C45-C46-O23
4	A	503	33O	O27-C55-C56-O28
4	A	503	33O	O25-C51-C52-O26
5	A	505	PEG	O1-C1-C2-O2
4	A	503	33O	O19-C39-C40-O20
4	A	503	33O	C61-C62-O31-C63
4	A	503	33O	O20-C41-C42-O21
2	A	501[B]	NDP	C2D-C1D-N1N-C2N
4	A	503	33O	O28-C57-C58-O29
2	A	501[B]	NDP	C2D-C1D-N1N-C6N
4	A	503	33O	O24-C49-C50-O25
5	B	504	PEG	C1-C2-O2-C3
4	A	503	33O	C45-C46-O23-C47
2	A	501[B]	NDP	O4D-C1D-N1N-C2N
2	A	501[B]	NDP	C5B-O5B-PA-O3
2	A	501[B]	NDP	C5D-O5D-PN-O3
2	A	501[A]	NDP	C5B-O5B-PA-O3
2	A	501[A]	NDP	C5D-O5D-PN-O3
2	B	501	NDP	C5B-O5B-PA-O3
2	B	501	NDP	C5D-O5D-PN-O3
4	A	503	33O	C52-C51-O25-C50
5	B	503[A]	PEG	C4-C3-O2-C2
2	A	501[A]	NDP	O4D-C1D-N1N-C6N
2	A	501[B]	NDP	C5D-O5D-PN-O1N
2	A	501[B]	NDP	C2N-C3N-C7N-N7N
2	A	501[A]	NDP	C5D-O5D-PN-O1N
4	A	503	33O	O23-C47-C48-O24
2	A	501[B]	NDP	C2N-C3N-C7N-O7N
4	A	503	33O	C46-C45-O22-C44
4	A	503	33O	C53-C54-O27-C55
2	A	501[B]	NDP	O4D-C1D-N1N-C6N
2	B	501	NDP	O4D-C1D-N1N-C2N
2	B	501	NDP	C2D-C1D-N1N-C2N
4	A	503	33O	C42-C41-O20-C40
4	A	503	33O	C50-C49-O24-C48

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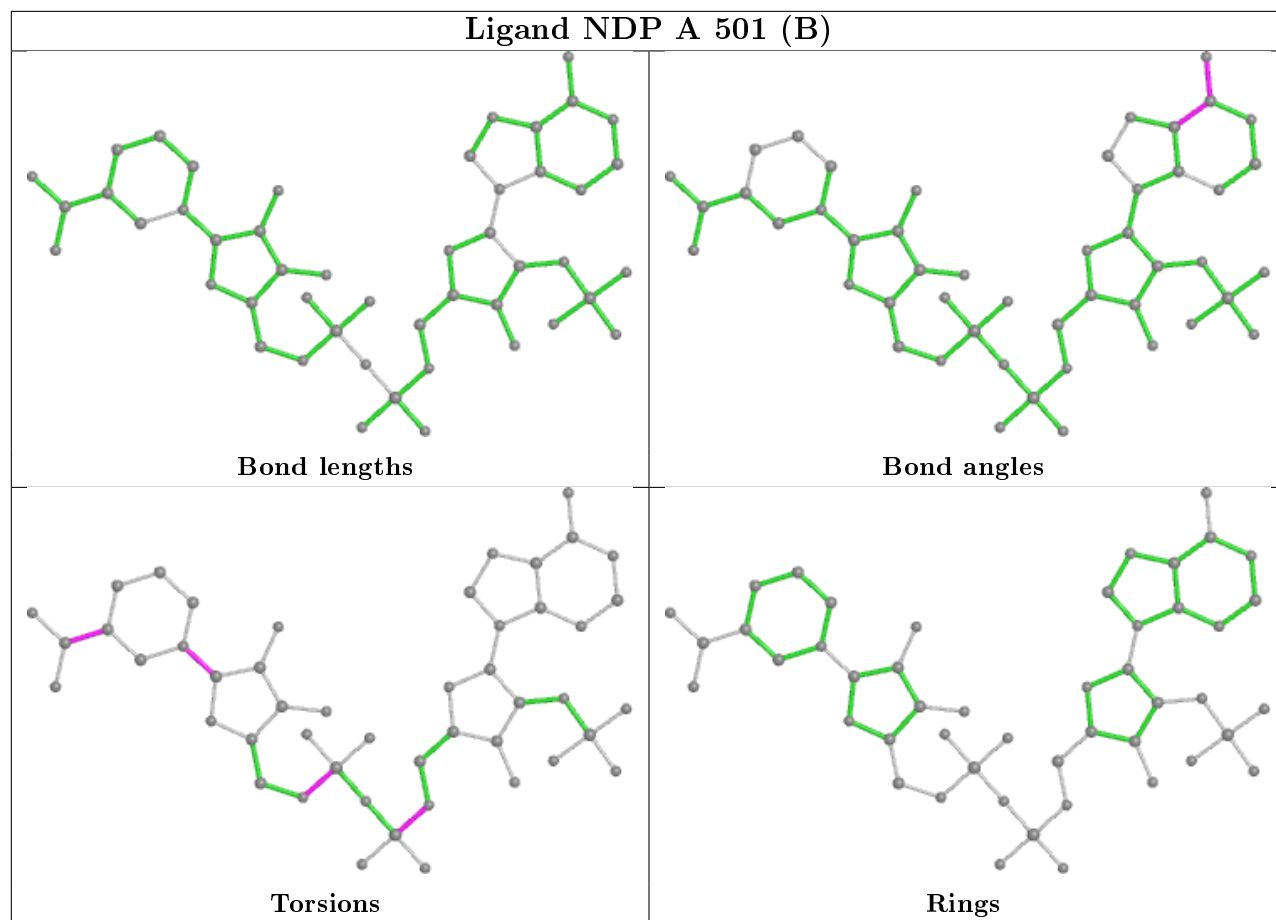
Mol	Chain	Res	Type	Atoms
5	B	504	PEG	C4-C3-O2-C2
2	B	501	NDP	O4D-C1D-N1N-C6N
2	B	501	NDP	C2D-C1D-N1N-C6N
5	B	503[B]	PEG	C4-C3-O2-C2
4	A	503	33O	O30-C61-C62-O31
4	A	503	33O	C51-C52-O26-C53

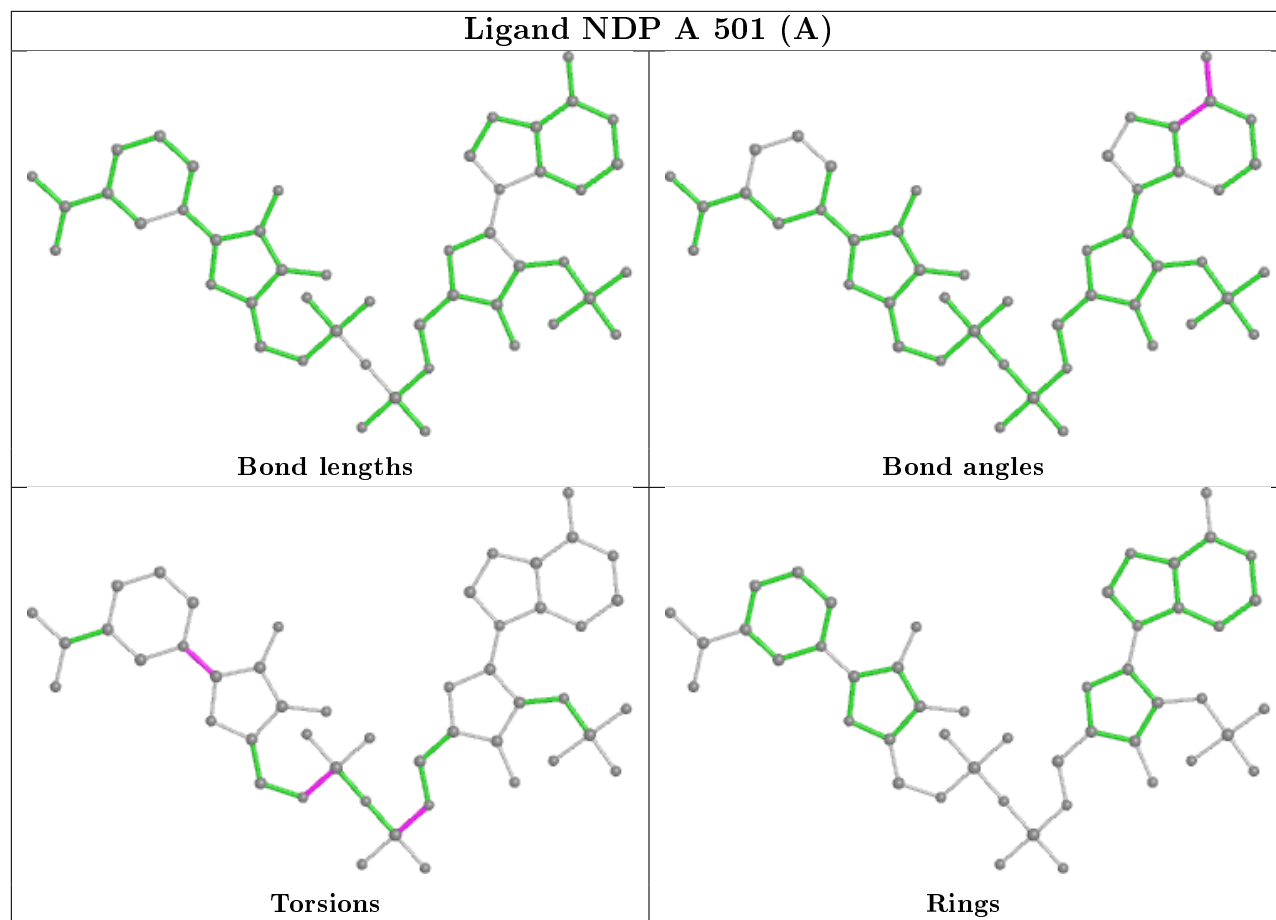
There are no ring outliers.

4 monomers are involved in 7 short contacts:

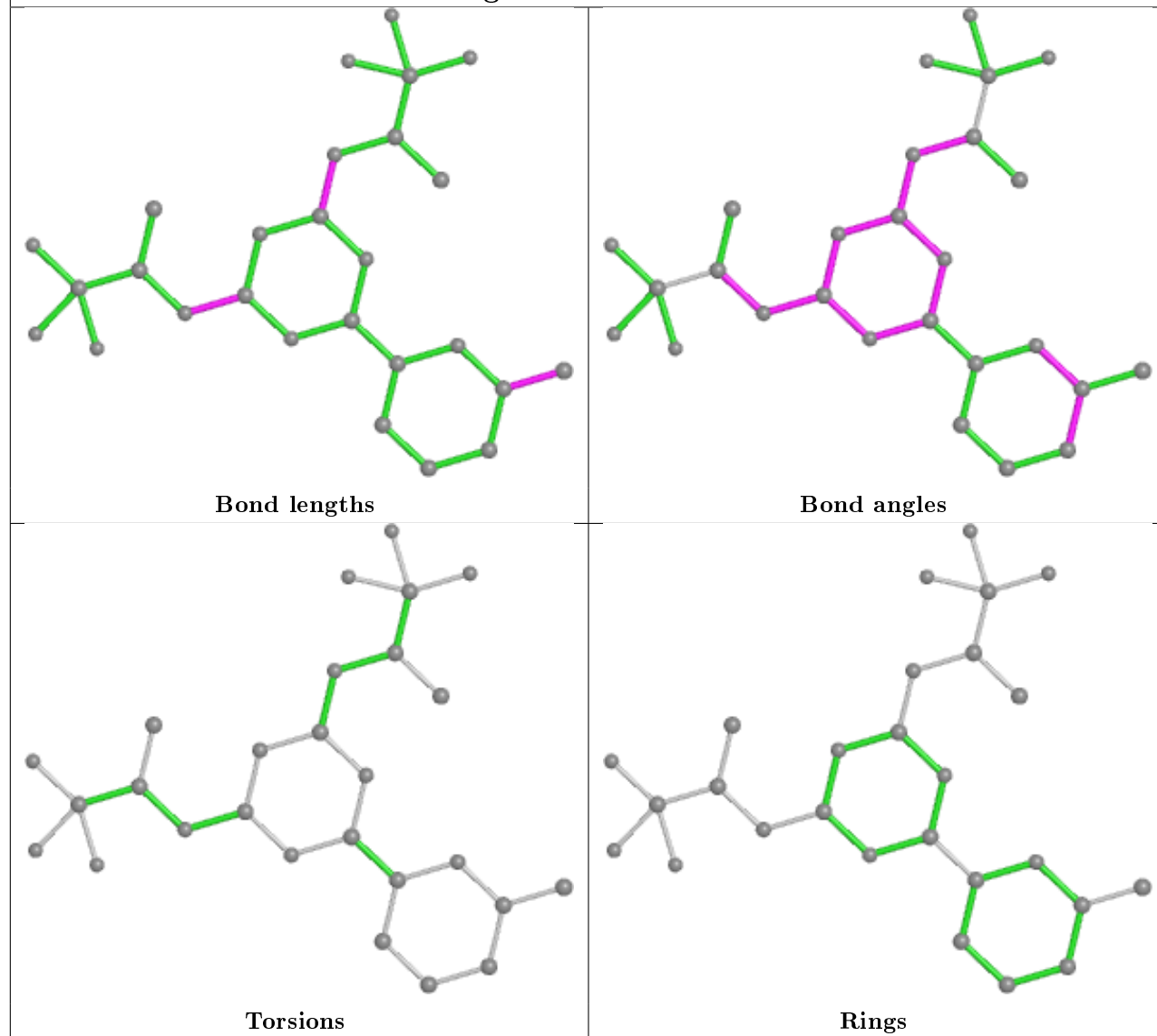
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501[B]	NDP	2	0
5	B	503[B]	PEG	3	0
2	B	501	NDP	4	0
5	B	503[A]	PEG	1	0

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.

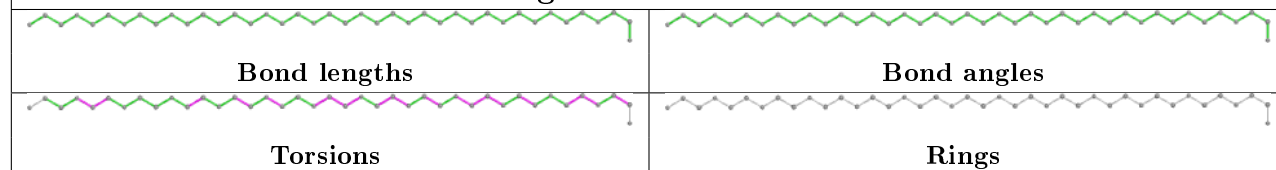


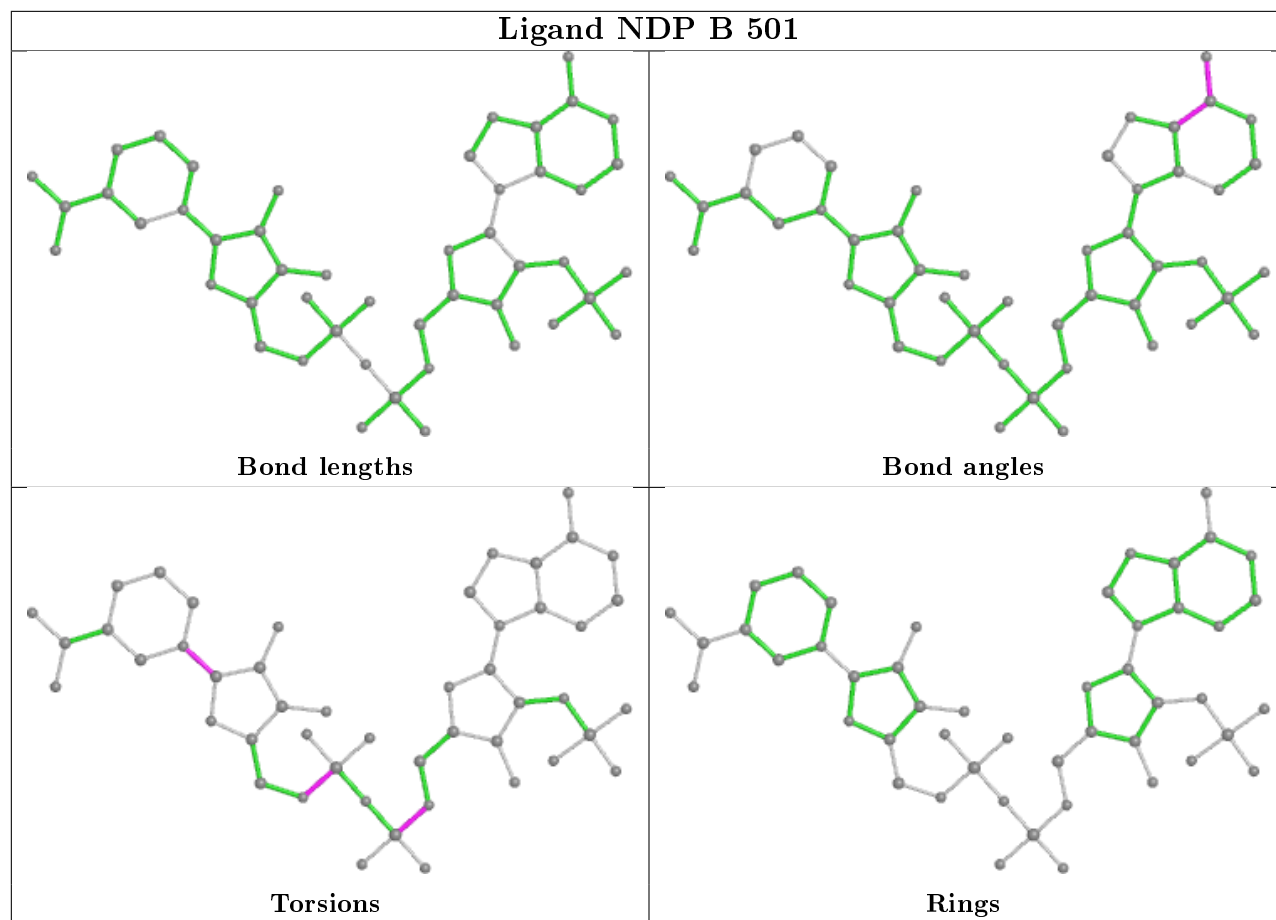


Ligand 9UO B 502



Ligand 33O A 503





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	413/425 (97%)	-0.32	6 (1%) 73 77	16, 25, 46, 73	0
1	B	417/425 (98%)	-0.37	5 (1%) 79 82	15, 24, 47, 74	0
All	All	830/850 (97%)	-0.35	11 (1%) 77 80	15, 25, 46, 74	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	414	LEU	7.4
1	B	174	GLU	5.3
1	B	419	HIS	5.1
1	A	413	LYS	5.0
1	A	412	ALA	4.1
1	A	415	SER	3.3
1	A	174	GLU	3.3
1	A	410	ALA	2.5
1	B	243	LYS	2.1
1	B	173	GLU	2.0
1	B	418	HIS	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

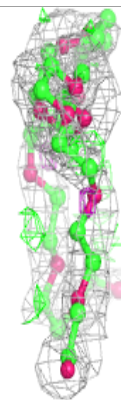
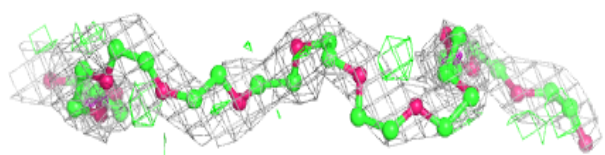
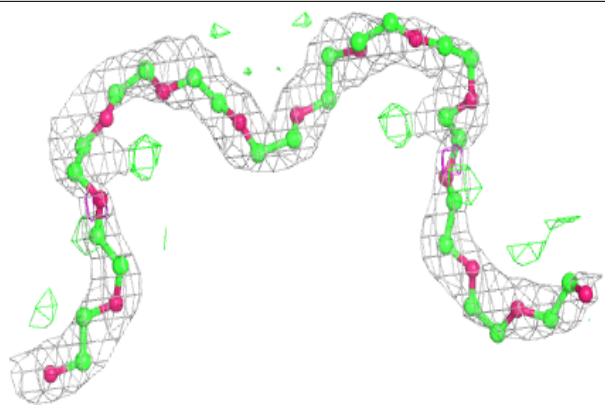
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	ACT	A	506	4/4	0.71	0.36	53,65,65,71	0
5	PEG	A	505	7/7	0.74	0.19	51,55,65,66	0
7	MLA	A	507	7/7	0.77	0.16	58,62,74,74	0
9	MLT	B	505	9/9	0.77	0.14	35,50,60,61	0
5	PEG	B	503[A]	7/7	0.81	0.27	21,32,43,44	7
5	PEG	B	503[B]	7/7	0.81	0.27	19,32,44,44	7
5	PEG	B	504	7/7	0.82	0.19	57,61,66,70	0
4	33O	A	503	40/40	0.84	0.20	43,52,56,57	40
5	PEG	A	504	7/7	0.85	0.15	52,54,57,61	0
8	9UO	B	502	27/27	0.96	0.13	14,20,37,41	0
2	NDP	A	501[B]	48/48	0.96	0.09	16,28,36,41	48
2	NDP	A	501[A]	48/48	0.96	0.09	20,29,34,41	48
2	NDP	B	501	48/48	0.97	0.09	15,21,27,31	0
3	SO4	A	502	5/5	0.97	0.10	29,32,38,53	5

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.

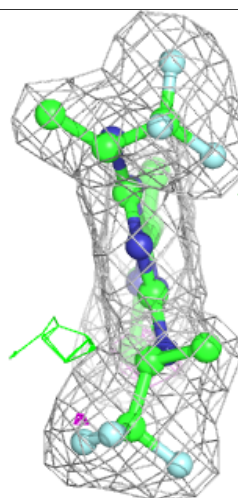
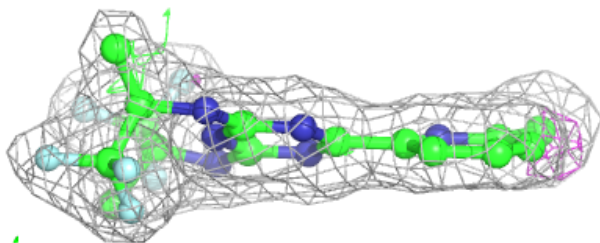
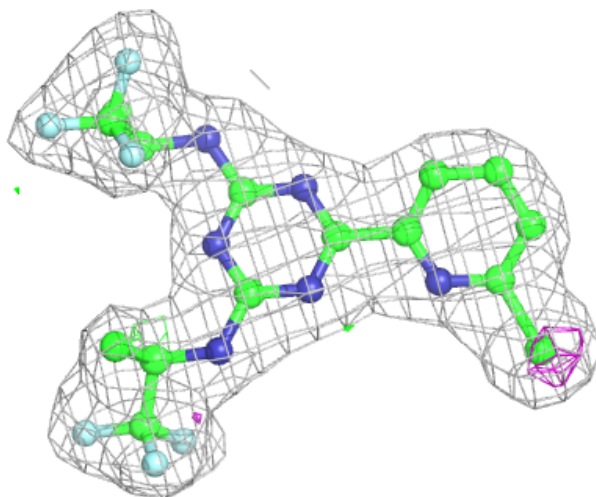
Electron density around 33O A 503:

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



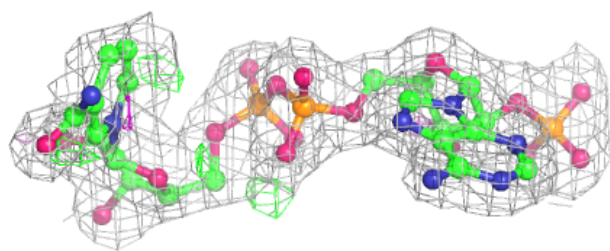
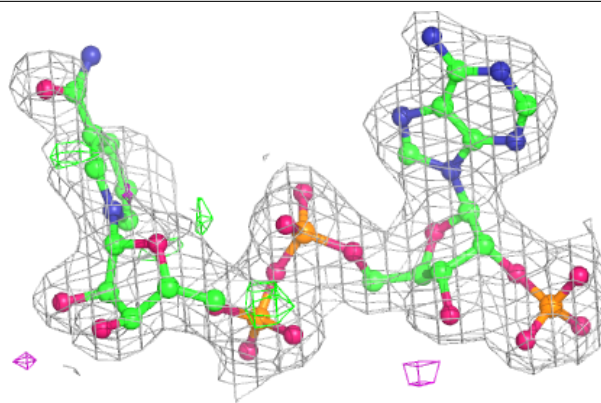
Electron density around 9UO B 502:

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

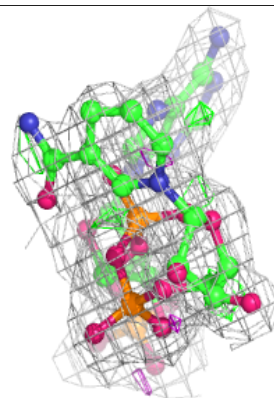
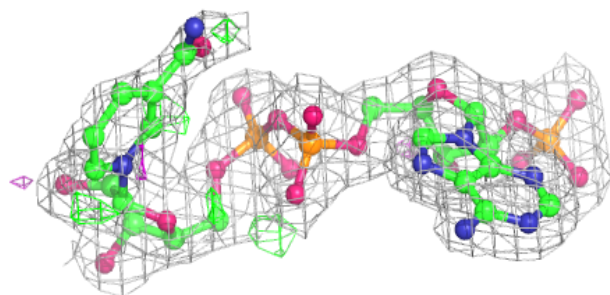
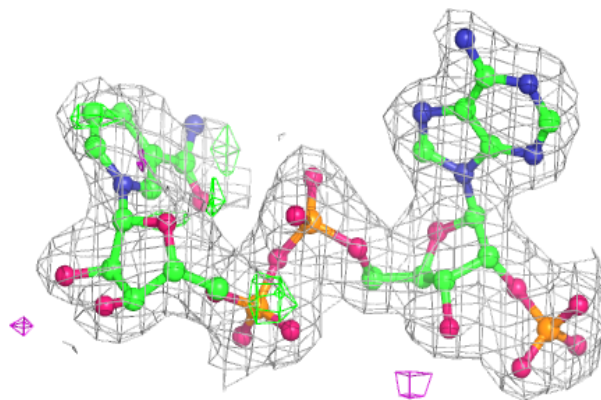


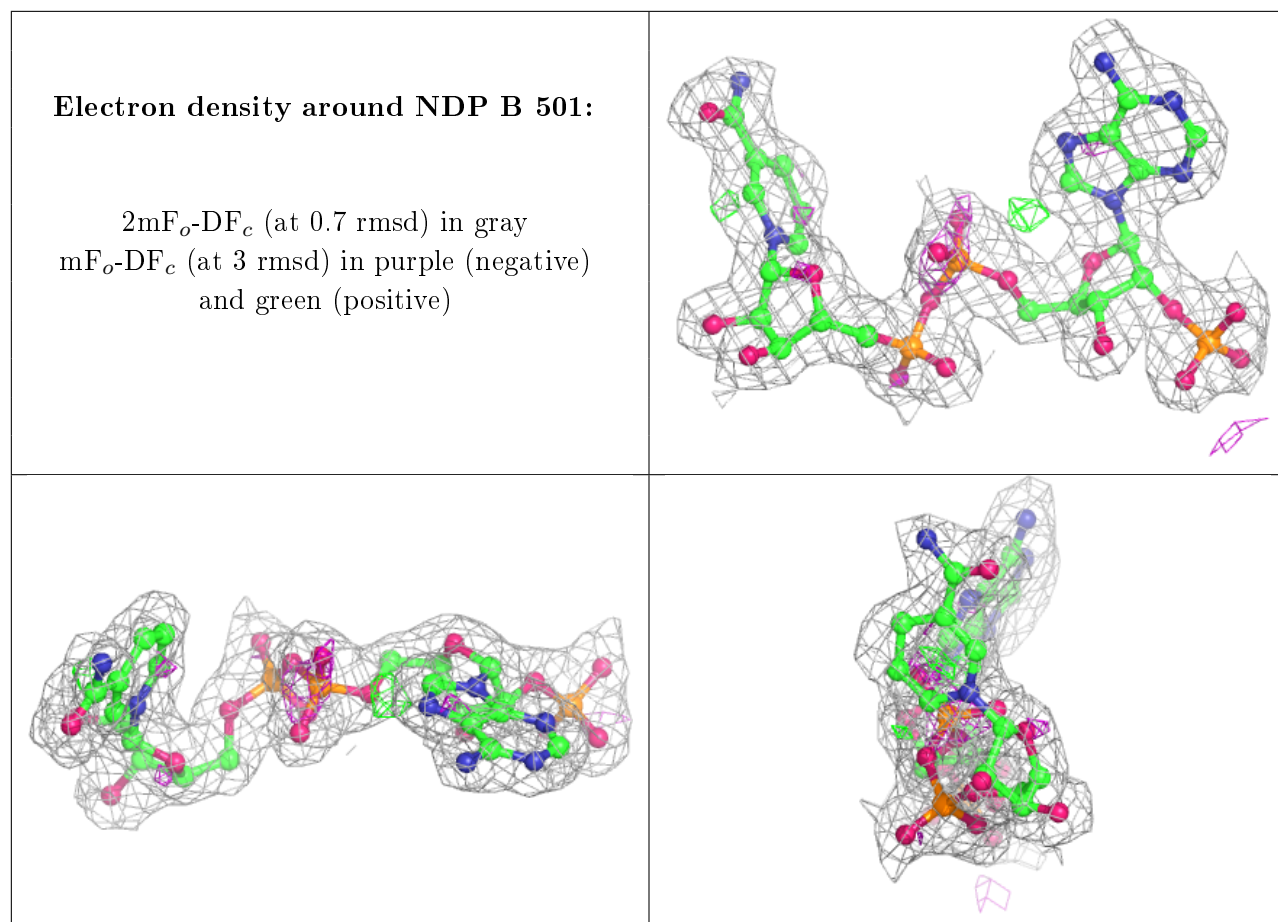
Electron density around NDP A 501 (B):

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

**Electron density around NDP A 501 (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 ⓘ

There are no such residues in this entry.