



# Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 05:32 pm BST

PDB ID : 6VFZ  
Title : Crystal Structure of Human Mitochondrial Isocitrate Dehydrogenase (IDH2) R140Q Mutant Homodimer in Complex with NADPH and AG-881 (Vorasidenib) Inhibitor.  
Authors : Padyana, A.; Jin, L.  
Deposited on : 2020-01-07  
Resolution : 1.99 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](mailto: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.

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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

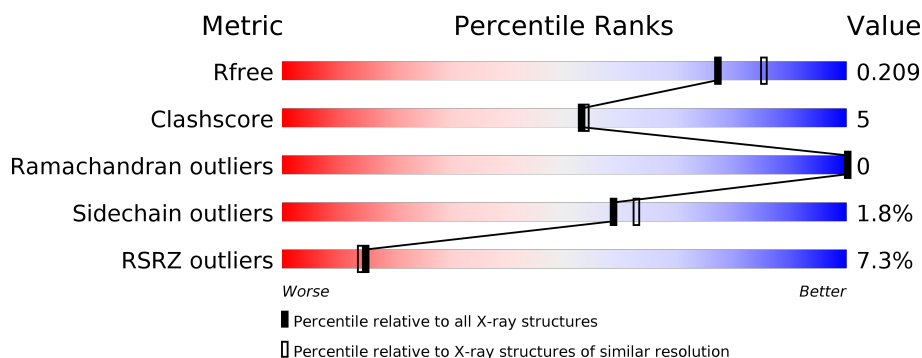
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 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	458	<div> <div>3%</div> <div> <div></div> <div>81%</div> <div>10%</div> <div>9%</div> </div> </div>
1	B	458	<div> <div>10%</div> <div> <div></div> <div>77%</div> <div>11%</div> <div>11%</div> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7356 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.

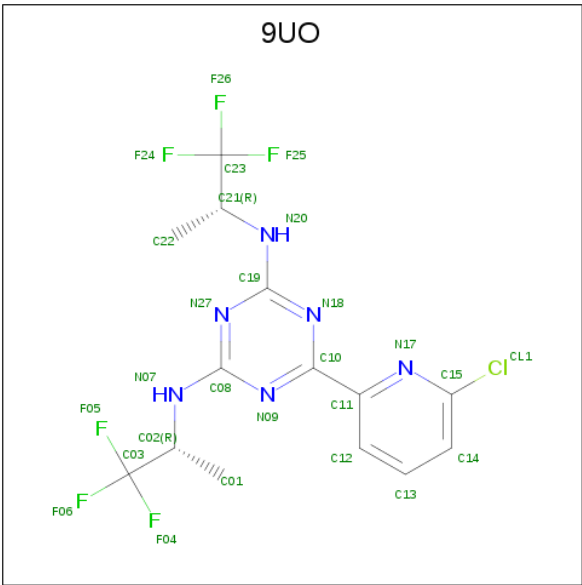
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP], mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	6	0
			3353	2132	583	619	19			
1	B	406	Total	C	N	O	S	0	1	0
			3233	2058	556	601	18			

There are 14 discrepancies between the modelled and reference sequences:

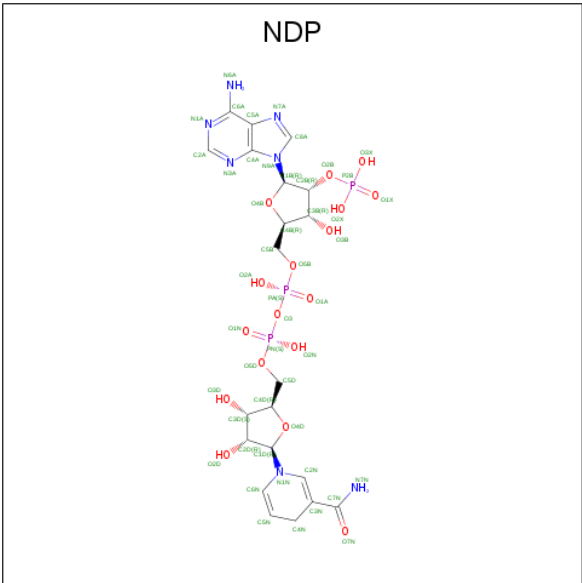
Chain	Residue	Modelled	Actual	Comment	Reference
A	140	GLN	ARG	engineered mutation	UNP P48735
A	453	HIS	-	expression tag	UNP P48735
A	454	HIS	-	expression tag	UNP P48735
A	455	HIS	-	expression tag	UNP P48735
A	456	HIS	-	expression tag	UNP P48735
A	457	HIS	-	expression tag	UNP P48735
A	458	HIS	-	expression tag	UNP P48735
B	140	GLN	ARG	engineered mutation	UNP P48735
B	453	HIS	-	expression tag	UNP P48735
B	454	HIS	-	expression tag	UNP P48735
B	455	HIS	-	expression tag	UNP P48735
B	456	HIS	-	expression tag	UNP P48735
B	457	HIS	-	expression tag	UNP P48735
B	458	HIS	-	expression tag	UNP P48735

- Molecule 2 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<sub>14</sub>H<sub>13</sub>ClF<sub>6</sub>N<sub>6</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Cl	F	N		
2	A	1	54	28	2	12	12	0	1

- Molecule 3 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C<sub>21</sub>H<sub>30</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).



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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0
4	A	1	Total 1	Ca 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total 1	Na 1	0	0
5	A	1	Total 1	Na 1	0	0

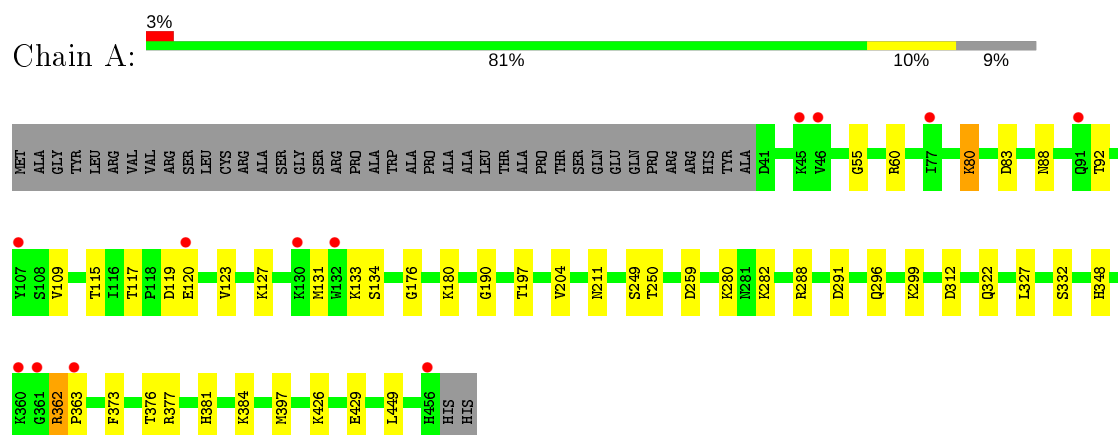
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	360	Total 360	O 360	0	0
6	B	256	Total 256	O 256	0	0

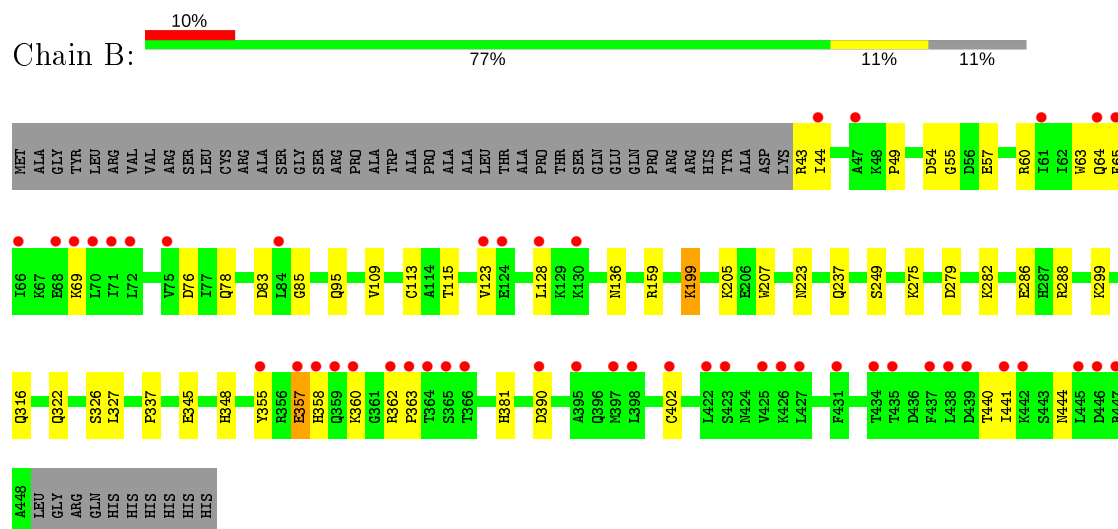
### 3 Residue-property plots [i](#)

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 ( $RSRZ > 2$ ). 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], mitochondrial



- Molecule 1: Isocitrate dehydrogenase [NADP], mitochondrial



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.07Å 119.80Å 130.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.06 – 1.99 49.11 – 1.99	Depositor EDS
% Data completeness (in resolution range)	96.3 (42.06-1.99) 96.3 (49.11-1.99)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.20 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, $R_{free}$	0.177 , 0.209 0.177 , 0.209	Depositor DCC
$R_{free}$ test set	3046 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.3	Xtriage
Anisotropy	0.094	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 57.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7356	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: NDP, CA, 9UO, NA

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	0.29	0/3440	0.48	0/4646
1	B	0.29	0/3307	0.49	0/4469
All	All	0.29	0/6747	0.49	0/9115

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	3353	0	3321	33	0
1	B	3233	0	3202	43	0
2	A	54	0	0	2	0
3	A	48	0	26	3	0
3	B	48	0	26	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	360	0	0	1	0
6	B	256	0	0	6	0
All	All	7356	0	6575	72	0



The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (72) 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:B:43:ARG:NH2	1:B:76:ASP:OD2	1.87	1.07
1:A:362:ARG:HG3	1:A:363:PRO:HD2	1.49	0.94
1:B:54:ASP:OD1	6:B:601:HOH:O	2.00	0.78
1:B:44:ILE:HD12	1:B:390:ASP:HB3	1.66	0.77
1:A:362:ARG:CG	1:A:363:PRO:HD2	2.16	0.75
1:A:120:GLU:HG2	1:A:131:MET:HE1	1.75	0.67
1:A:60:ARG:NH2	1:A:83:ASP:OD1	2.28	0.67
1:B:63:TRP:CD2	1:B:113:CYS:HB2	2.31	0.66
1:B:60:ARG:O	1:B:64:GLN:HG2	1.97	0.64
1:A:88:ASN:O	1:A:92:THR:HG22	1.98	0.64
1:B:109:VAL:HG22	1:B:381:HIS:HD2	1.64	0.62
1:B:205:LYS:NZ	6:B:605:HOH:O	2.32	0.62
1:B:64:GLN:NE2	6:B:601:HOH:O	2.32	0.62
1:B:286:GLU:OE1	1:B:288:ARG:NH1	2.29	0.60
1:B:136:ASN:ND2	1:B:345:GLU:OE2	2.36	0.59
1:B:440:THR:HG22	1:B:444:ASN:HD21	1.70	0.57
1:B:49:PRO:HB3	1:B:78:GLN:HG3	1.88	0.56
1:A:117:THR:OG1	3:A:502:NDP:O2D	2.25	0.55
1:B:440:THR:O	1:B:444:ASN:ND2	2.40	0.55
1:A:176:GLY:H	1:A:180:LYS:HD3	1.72	0.54
1:B:60:ARG:HH21	1:B:85:GLY:HA3	1.73	0.54
1:A:197:THR:HG23	1:A:204:VAL:HG22	1.89	0.53
1:B:63:TRP:CE3	1:B:113:CYS:HB2	2.44	0.53
1:B:237:GLN:NE2	1:B:337:PRO:O	2.42	0.52
1:A:362:ARG:HG3	1:A:363:PRO:CD	2.33	0.52
1:B:360:LYS:HD3	1:B:362:ARG:NH2	2.25	0.52
1:A:117:THR:HG22	1:A:134:SER:HA	1.91	0.51
1:B:440:THR:HG22	1:B:444:ASN:ND2	2.26	0.51
1:A:327:LEU:HB3	1:A:348:HIS:HB3	1.93	0.50
1:B:327:LEU:HB3	1:B:348:HIS:HB3	1.95	0.49
1:B:282:LYS:HE2	1:B:282:LYS:HA	1.95	0.49
1:A:211:ASN:OD1	1:B:199:LYS:NZ	2.46	0.49
1:A:322:GLN:OE1	1:B:299:LYS:HD2	2.12	0.48
1:A:299:LYS:HE2	1:B:322:GLN:HG3	1.94	0.48
1:B:60:ARG:NH2	1:B:83:ASP:OD1	2.45	0.48
1:B:249:SER:HA	1:B:288:ARG:O	2.13	0.48
1:A:190:GLY:HA2	1:B:199:LYS:HD3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:THR:H	3:A:502:NDP:H71N	1.61	0.48
1:A:119:ASP:O	1:A:123:VAL:HG23	2.13	0.48
1:B:358:HIS:ND1	6:B:607:HOH:O	2.35	0.48
1:A:249:SER:HA	1:A:288:ARG:O	2.15	0.46
1:A:55:GLY:O	1:A:60:ARG:NE	2.44	0.46
1:B:95:GLN:NE2	6:B:602:HOH:O	2.23	0.45
1:A:80:LYS:HA	1:A:80:LYS:HD2	1.78	0.45
1:B:65:PHE:HA	1:B:69:LYS:NZ	2.32	0.44
1:B:123:VAL:HA	1:B:128:LEU:HD12	1.99	0.44
1:A:426:LYS:HE2	1:A:429:GLU:OE1	2.18	0.44
1:B:402:CYS:HA	1:B:441:ILE:HD11	1.99	0.44
1:A:373:PHE:HA	1:A:376:THR:OG1	2.18	0.43
1:A:373:PHE:O	1:A:377:ARG:HG2	2.18	0.43
1:A:109:VAL:HG22	1:A:381:HIS:HD2	1.82	0.43
1:A:312:ASP:OD1	2:A:501[A]:9UO:CL1	2.73	0.43
1:B:357:GLU:CD	1:B:362:ARG:HD2	2.38	0.43
1:A:115:THR:O	3:A:502:NDP:H2N	2.18	0.43
1:A:299:LYS:HE2	1:B:322:GLN:CG	2.48	0.43
1:A:280:LYS:HA	1:A:280:LYS:HD3	1.84	0.43
1:B:205:LYS:HG2	1:B:207:TRP:CZ2	2.53	0.43
1:B:43:ARG:NE	1:B:76:ASP:HB2	2.34	0.43
1:A:176:GLY:N	1:A:180:LYS:HD3	2.34	0.42
1:B:360:LYS:HD3	1:B:362:ARG:CZ	2.48	0.42
1:A:250:THR:HB	1:A:259:ASP:HB3	2.02	0.42
2:A:501[A]:9UO:C12	1:B:316:GLN:NE2	2.83	0.42
1:B:55:GLY:O	1:B:60:ARG:NE	2.52	0.42
1:A:397[A]:MET:SD	1:A:449:LEU:HD23	2.59	0.42
3:B:501:NDP:O3X	3:B:501:NDP:O3B	2.33	0.41
1:B:362:ARG:HB2	1:B:363:PRO:HD2	2.02	0.41
1:B:275:LYS:HE3	1:B:279:ASP:OD2	2.20	0.41
1:B:60:ARG:HG3	6:B:601:HOH:O	2.20	0.41
1:A:176:GLY:HA2	1:A:180:LYS:HD3	2.04	0.40
1:B:115:THR:O	3:B:501:NDP:H2N	2.22	0.40
1:B:57:GLU:HG2	1:B:355:TYR:CD2	2.57	0.40
1:A:296[A]:GLN:NE2	6:A:626:HOH:O	2.54	0.40

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	420/458 (92%)	408 (97%)	12 (3%)	0	100	100
1	B	405/458 (88%)	392 (97%)	13 (3%)	0	100	100
All	All	825/916 (90%)	800 (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	363/389 (93%)	354 (98%)	9 (2%)	47	49
1	B	349/389 (90%)	344 (99%)	5 (1%)	67	72
All	All	712/778 (92%)	698 (98%)	14 (2%)	59	58

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LYS
1	A	127	LYS
1	A	133	LYS
1	A	282	LYS
1	A	291	ASP
1	A	332[A]	SER
1	A	332[B]	SER
1	A	362	ARG

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Mol	Chain	Res	Type
1	A	384	LYS
1	B	159	ARG
1	B	199	LYS
1	B	223	ASN
1	B	326	SER
1	B	357	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	424	ASN
1	B	88	ASN
1	B	141	ASN
1	B	316	GLN
1	B	444	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](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	9UO	A	501[A]	-	28,28,28	2.31	3 (10%)	32,42,42	2.88	10 (31%)
2	9UO	A	501[B]	-	28,28,28	2.35	3 (10%)	32,42,42	2.93	13 (40%)
3	NDP	B	501	-	45,52,52	0.51	0	53,80,80	0.54	1 (1%)
3	NDP	A	502	-	45,52,52	0.53	0	53,80,80	0.62	1 (1%)

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	9UO	A	501[A]	-	-	0/24/24/24	0/2/2/2
2	9UO	A	501[B]	-	-	3/24/24/24	0/2/2/2
3	NDP	B	501	-	-	8/30/77/77	0/5/5/5
3	NDP	A	502	-	-	8/30/77/77	0/5/5/5

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501[B]	9UO	C08-N07	8.14	1.45	1.34
2	A	501[A]	9UO	C08-N07	7.94	1.45	1.34
2	A	501[B]	9UO	C19-N20	7.88	1.44	1.34
2	A	501[A]	9UO	C19-N20	7.70	1.44	1.34
2	A	501[A]	9UO	C11-N17	-2.60	1.31	1.34
2	A	501[B]	9UO	C11-N17	-2.29	1.31	1.34

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501[A]	9UO	C08-N09-C10	9.21	120.47	114.60
2	A	501[B]	9UO	C08-N09-C10	8.29	119.89	114.60
2	A	501[B]	9UO	C19-N18-C10	8.25	119.86	114.60
2	A	501[A]	9UO	C19-N18-C10	7.61	119.45	114.60
2	A	501[B]	9UO	C19-N20-C21	-5.42	120.08	124.53
2	A	501[A]	9UO	C19-N20-C21	-4.86	120.54	124.53
2	A	501[B]	9UO	N27-C08-N09	-3.70	120.39	126.23
2	A	501[A]	9UO	N27-C08-N09	-3.67	120.43	126.23
2	A	501[B]	9UO	N27-C19-N18	-3.56	120.60	126.23
2	A	501[A]	9UO	N07-C08-N27	3.46	122.37	117.18
2	A	501[A]	9UO	N27-C19-N18	-3.37	120.90	126.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501[A]	9UO	N18-C10-N09	-3.22	120.02	125.23
2	A	501[B]	9UO	N18-C10-N09	-3.19	120.08	125.23
2	A	501[B]	9UO	C11-C10-N09	3.14	121.74	117.48
2	A	501[B]	9UO	C19-N27-C08	3.12	119.19	113.89
2	A	501[B]	9UO	C08-N07-C02	-2.93	122.13	124.53
2	A	501[A]	9UO	C19-N27-C08	2.84	118.71	113.89
2	A	501[A]	9UO	C11-C10-N18	2.74	121.19	117.48
2	A	501[B]	9UO	C14-C15-N17	-2.71	120.51	125.50
3	B	501	NDP	C5A-C6A-N6A	2.30	123.84	120.35
2	A	501[B]	9UO	C15-N17-C11	2.27	121.67	117.02
3	A	502	NDP	C5A-C6A-N6A	2.24	123.75	120.35
2	A	501[A]	9UO	C14-C15-N17	-2.19	121.47	125.50
2	A	501[B]	9UO	N20-C19-N27	2.14	120.39	117.18
2	A	501[B]	9UO	N07-C08-N27	2.01	120.19	117.18

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	501	NDP	C5B-O5B-PA-O1A
3	B	501	NDP	C5D-O5D-PN-O2N
3	B	501	NDP	C2D-C1D-N1N-C2N
3	B	501	NDP	C5D-O5D-PN-O3
3	B	501	NDP	C5D-O5D-PN-O1N
3	A	502	NDP	O4B-C4B-C5B-O5B
3	B	501	NDP	O4D-C1D-N1N-C2N
2	A	501[B]	9UO	N20-C21-C23-F25
3	A	502	NDP	O4D-C1D-N1N-C2N
3	A	502	NDP	C2D-C1D-N1N-C2N
3	B	501	NDP	O4D-C1D-N1N-C6N
3	B	501	NDP	C2D-C1D-N1N-C6N
3	A	502	NDP	C3B-C4B-C5B-O5B
2	A	501[B]	9UO	N20-C21-C23-F24
2	A	501[B]	9UO	N20-C21-C23-F26
3	A	502	NDP	C5D-O5D-PN-O3
3	A	502	NDP	C5B-O5B-PA-O1A
3	A	502	NDP	C5D-O5D-PN-O1N
3	A	502	NDP	C5D-O5D-PN-O2N

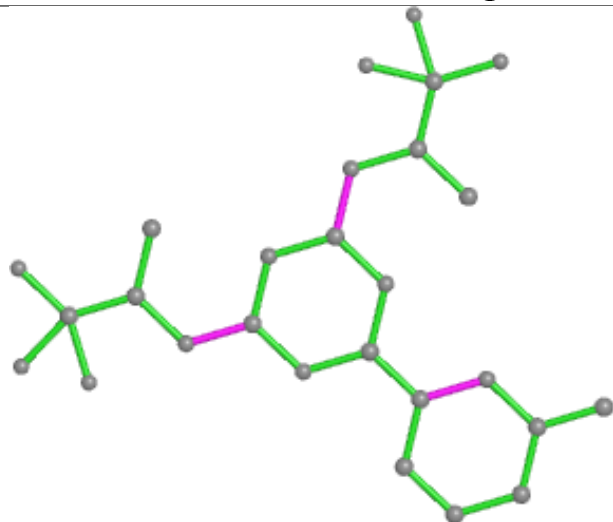
There are no ring outliers.

3 monomers are involved in 7 short contacts:

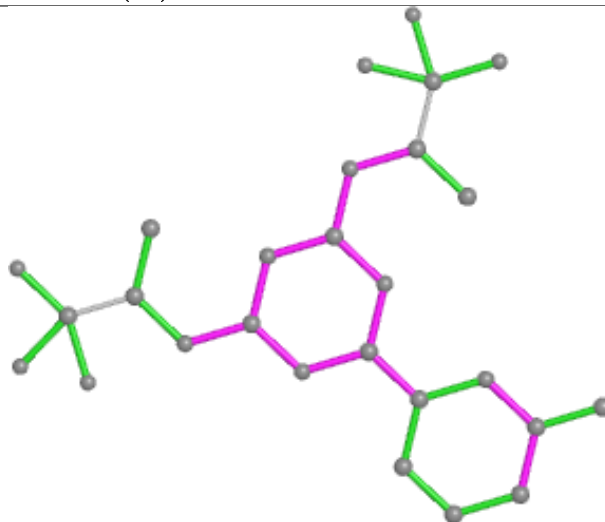
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501[A]	9UO	2	0
3	B	501	NDP	2	0
3	A	502	NDP	3	0

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.

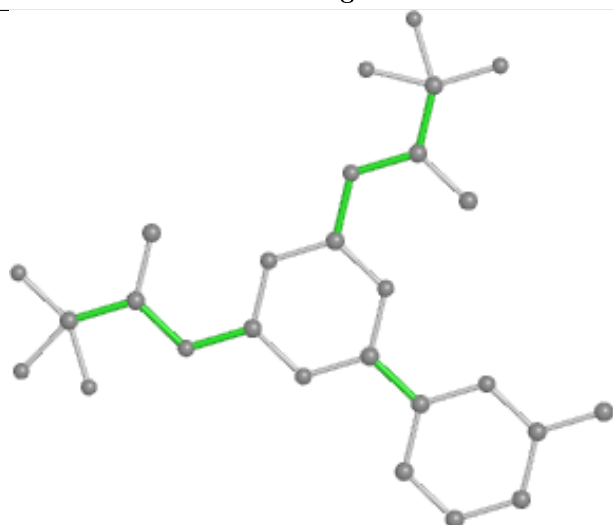
## Ligand 9UO A 501 (A)



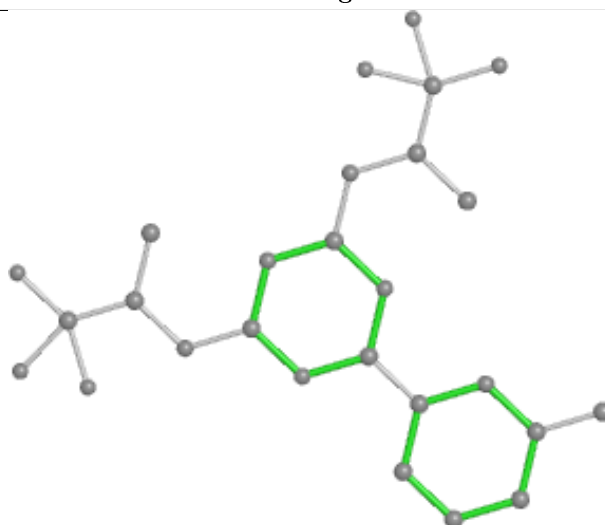
Bond lengths



Bond angles



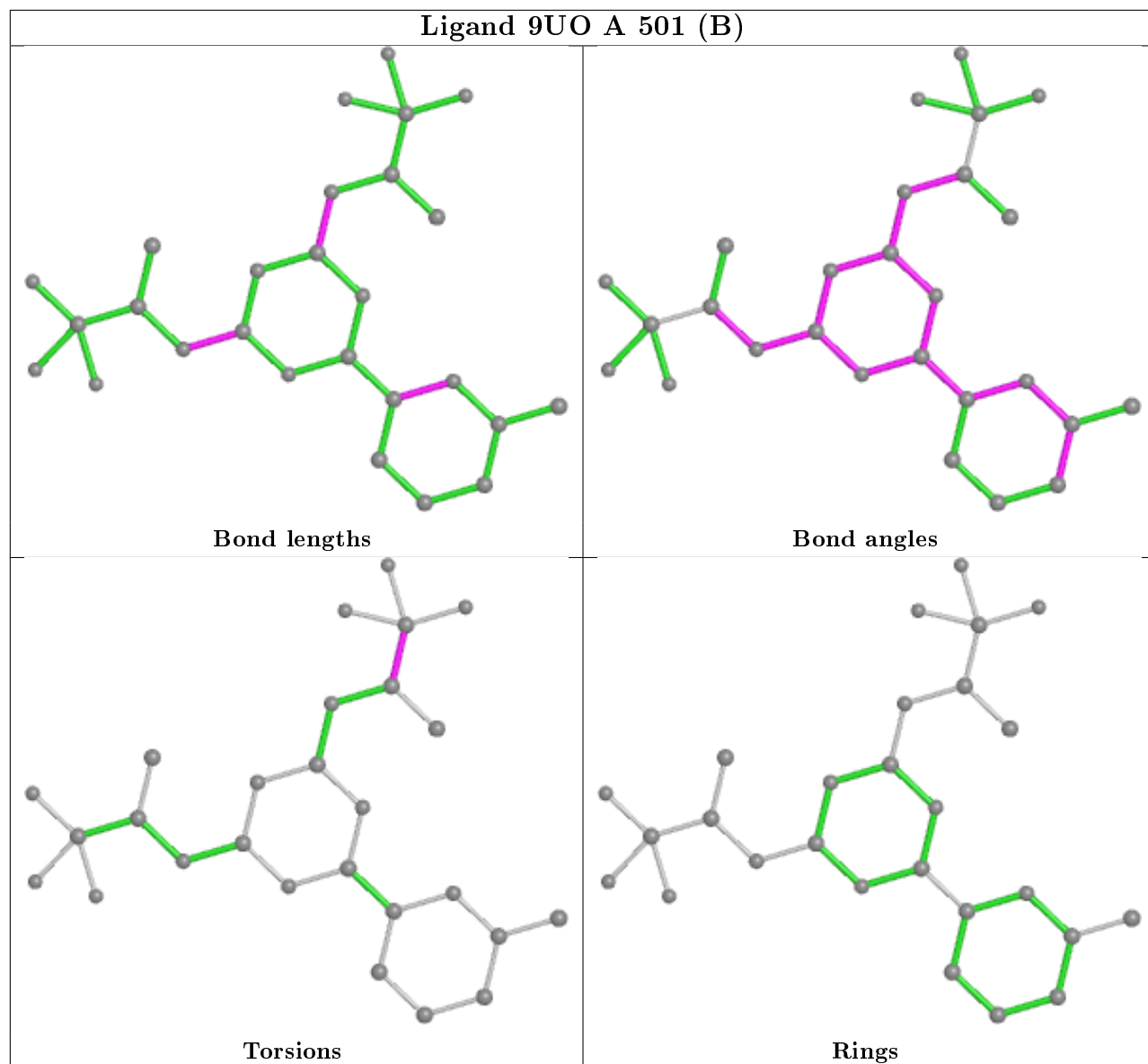
Torsions

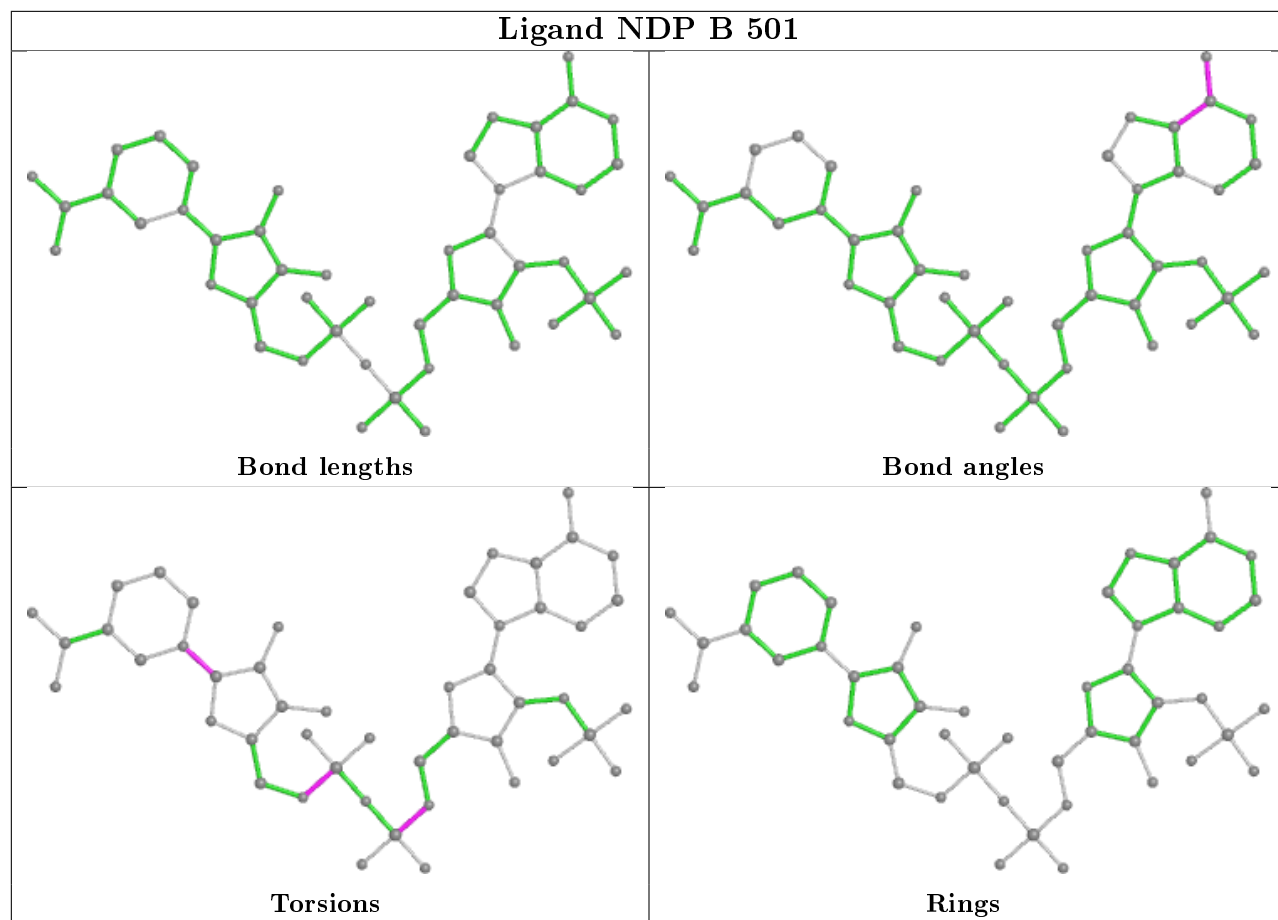


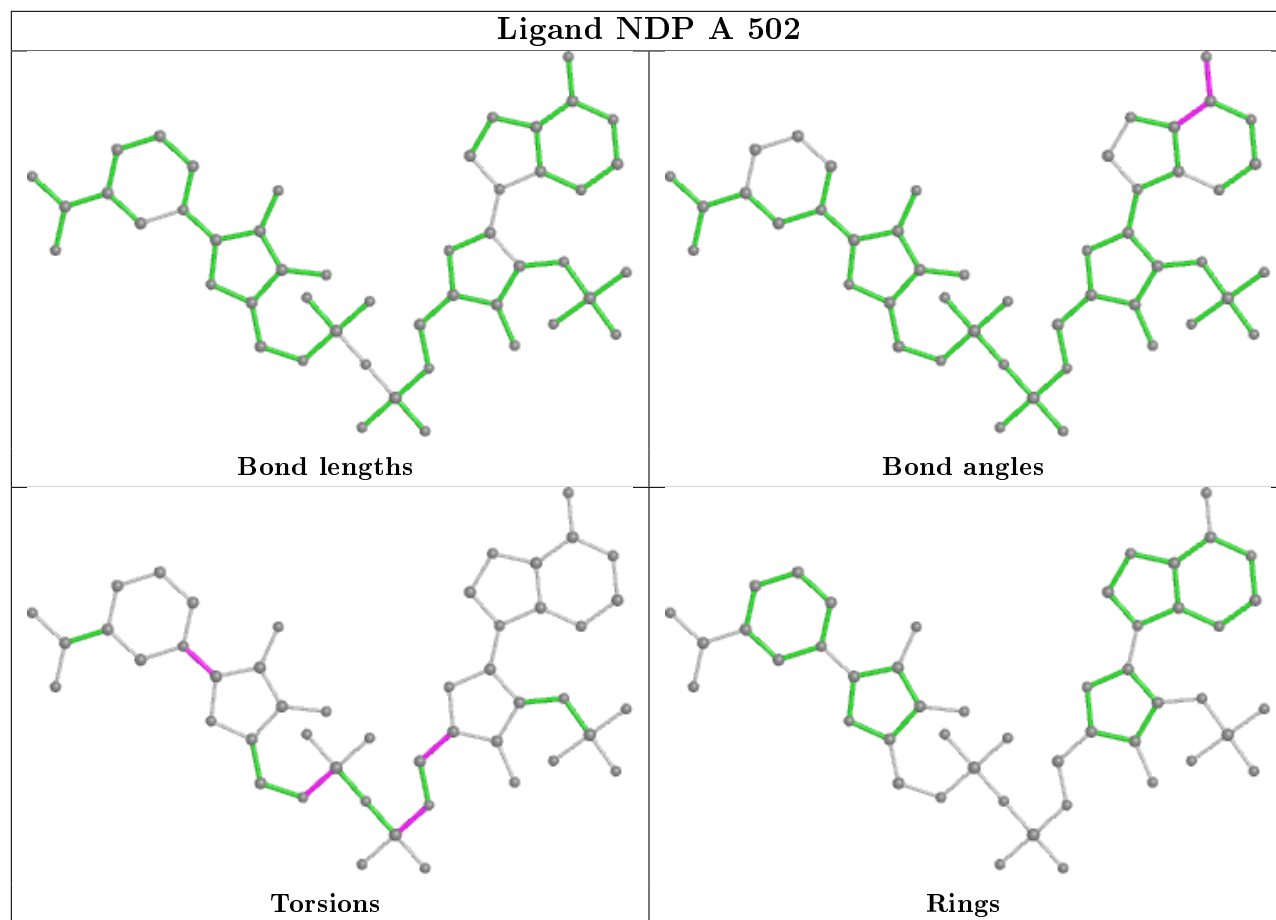
Rings



## Ligand 9UO A 501 (B)







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	416/458 (90%)	0.28	12 (2%) 51 50	9, 27, 51, 82	0
1	B	406/458 (88%)	0.63	48 (11%) 4 4	11, 36, 70, 94	0
All	All	822/916 (89%)	0.45	60 (7%) 15 14	9, 30, 65, 94	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	72	LEU	5.3
1	B	363	PRO	4.3
1	B	425	VAL	4.3
1	B	439	ASP	4.2
1	B	355	TYR	4.1
1	B	435	THR	4.1
1	B	426	LYS	3.9
1	B	362	ARG	3.8
1	B	437	PHE	3.7
1	B	446	ASP	3.7
1	B	441	ILE	3.6
1	B	438	LEU	3.5
1	B	357	GLU	3.5
1	A	91	GLN	3.5
1	B	442	LYS	3.4
1	A	107	TYR	3.3
1	B	431	PHE	3.2
1	B	75	VAL	3.1
1	B	447	ARG	2.9
1	B	397	MET	2.9
1	A	130	LYS	2.9
1	B	358	HIS	2.9
1	B	422	LEU	2.9
1	A	360	LYS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	77	ILE	2.8
1	A	456	HIS	2.8
1	B	366	THR	2.8
1	B	365	SER	2.8
1	B	395	ALA	2.7
1	B	434	THR	2.7
1	A	363	PRO	2.7
1	B	364	THR	2.7
1	B	124	GLU	2.7
1	B	71	ILE	2.7
1	B	65	PHE	2.7
1	B	44	ILE	2.6
1	A	46	VAL	2.6
1	A	45	LYS	2.6
1	B	128	LEU	2.5
1	B	130	LYS	2.5
1	B	47	ALA	2.5
1	A	132	TRP	2.5
1	A	120	GLU	2.4
1	B	84	LEU	2.4
1	B	390	ASP	2.4
1	B	61	ILE	2.4
1	B	123	VAL	2.4
1	B	398	LEU	2.3
1	B	427	LEU	2.3
1	B	445	LEU	2.3
1	B	64	GLN	2.3
1	B	68	GLU	2.3
1	B	69	LYS	2.3
1	B	360	LYS	2.2
1	A	361	GLY	2.2
1	B	423	SER	2.1
1	B	359	GLN	2.1
1	B	66	ILE	2.1
1	B	402	CYS	2.0
1	B	70	LEU	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 6.3 Carbohydrates

There are no carbohydrates in this entry.

### 6.4 Ligands

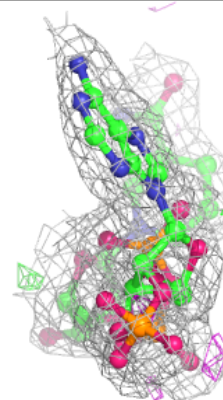
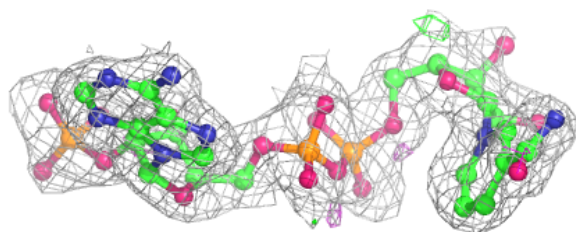
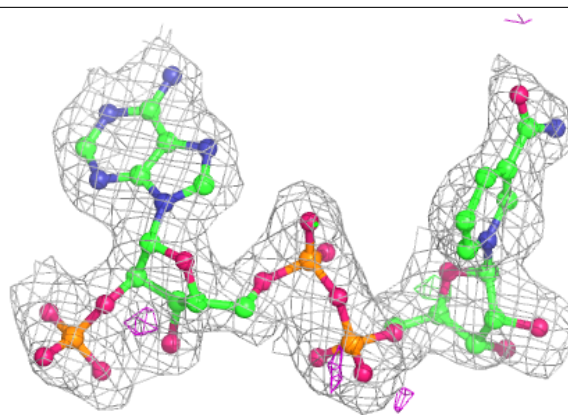
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	NA	A	504	1/1	0.87	0.29	52,52,52,52	0
5	NA	B	503	1/1	0.93	0.15	47,47,47,47	0
3	NDP	B	501	48/48	0.95	0.11	28,37,45,52	0
2	9UO	A	501[A]	27/27	0.97	0.15	8,12,16,19	27
2	9UO	A	501[B]	27/27	0.97	0.15	7,12,16,17	27
3	NDP	A	502	48/48	0.97	0.09	18,26,36,39	0
4	CA	A	503	1/1	0.98	0.08	20,20,20,20	0
4	CA	B	502	1/1	0.99	0.05	22,22,22,22	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.

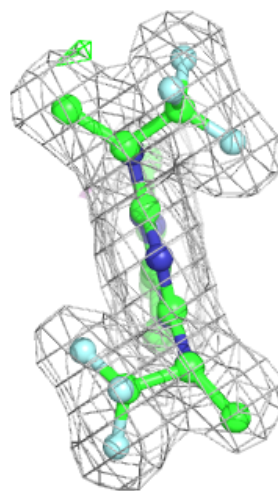
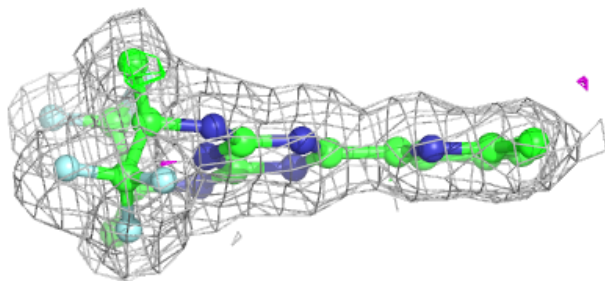
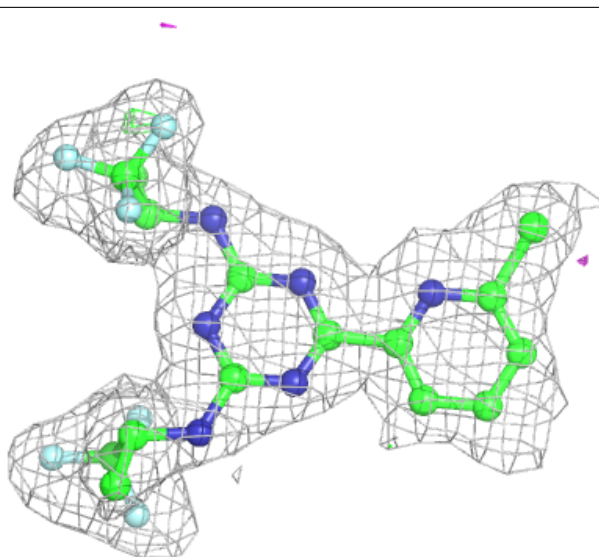
**Electron density around NDP B 501:**

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



**Electron density around 9UO A 501 (A):**

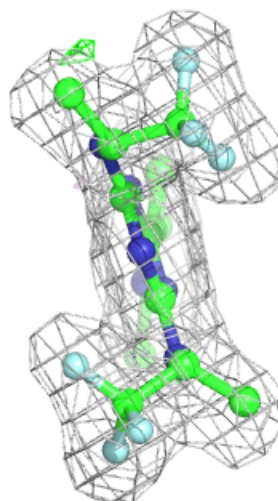
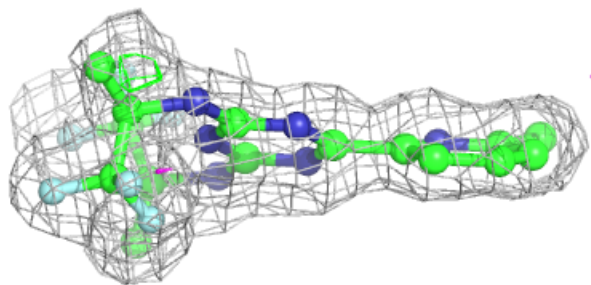
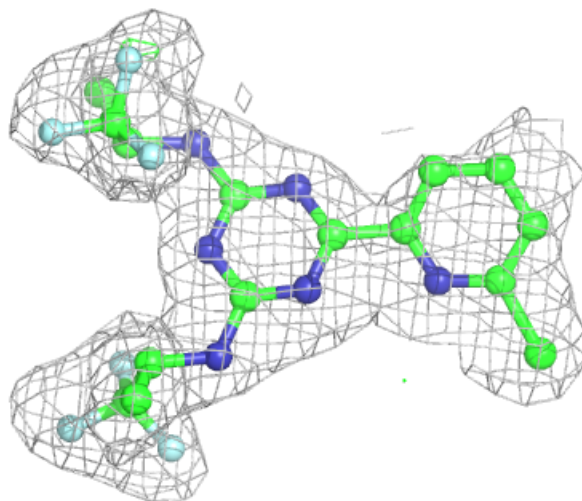
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

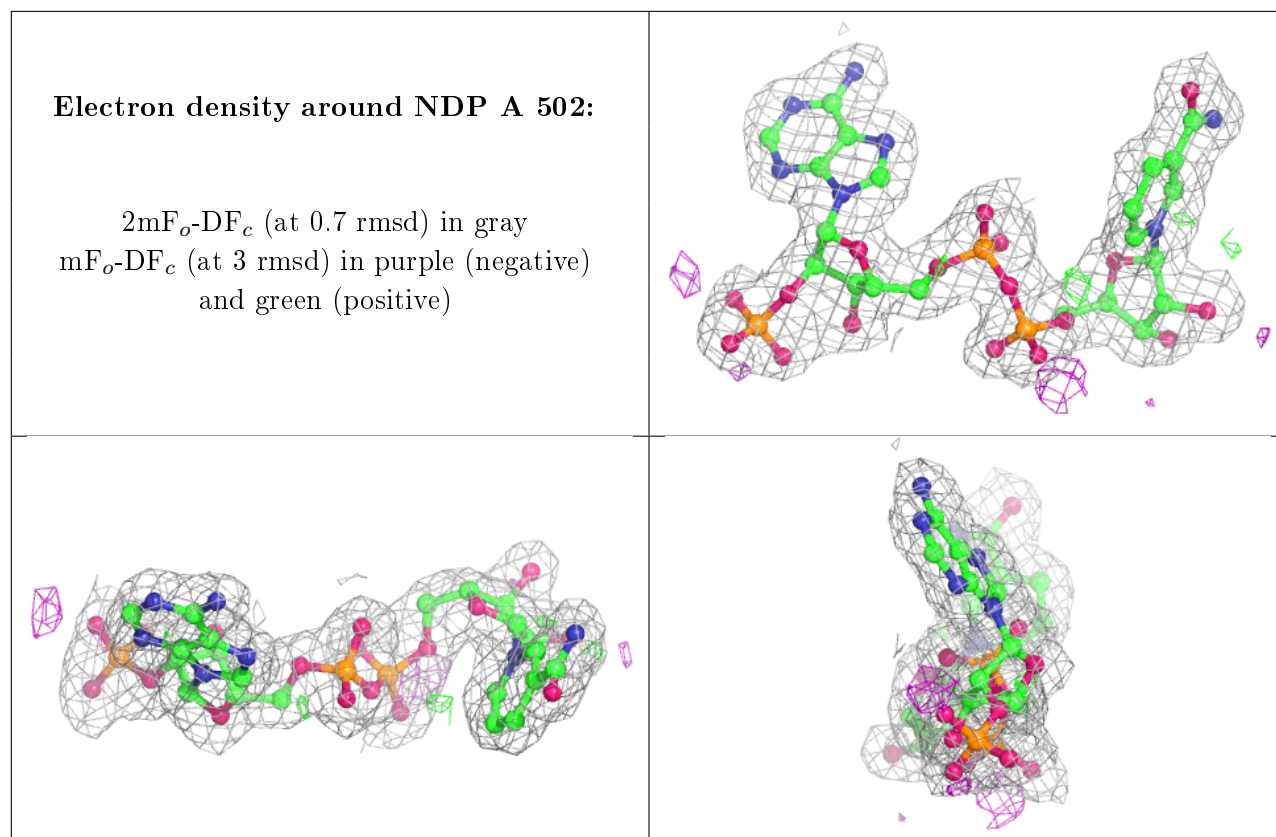




**Electron density around 9UO A 501 (B):**

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





## 6.5 Other polymers [i](#)

There are no such residues in this entry.