



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:49 AM UTC

PDB ID : 6X6V / pdb_00006x6v
Title : Crystal structure of inactive enzymatic binary toxin component from *Clostridium difficile* in complex with NADPH
Authors : Pozharski, E.
Deposited on : 2020-05-29
Resolution : 2.42 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 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

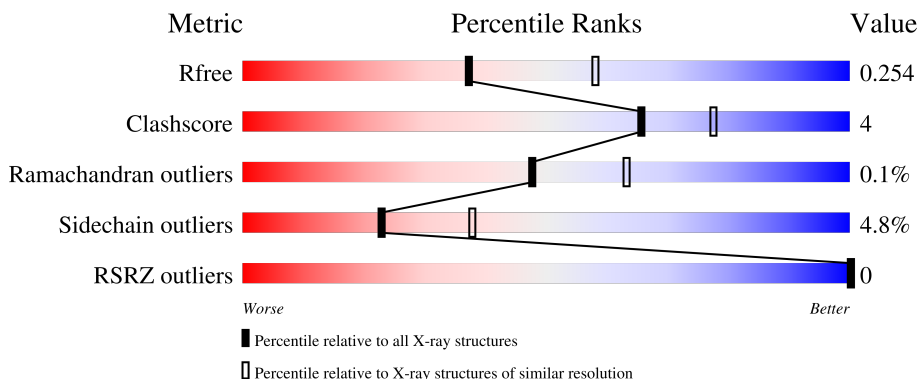
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	429	
1	B	429	
1	C	429	
1	D	429	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14229 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 CdtA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	416	Total	C	N	O	S	0	2	0
			3355	2141	556	651	7			
1	B	416	Total	C	N	O	S	0	2	0
			3345	2139	554	645	7			
1	C	416	Total	C	N	O	S	0	1	0
			3361	2145	560	650	6			
1	D	416	Total	C	N	O	S	0	2	0
			3339	2133	556	644	6			

There are 56 discrepancies between the modelled and reference sequences:

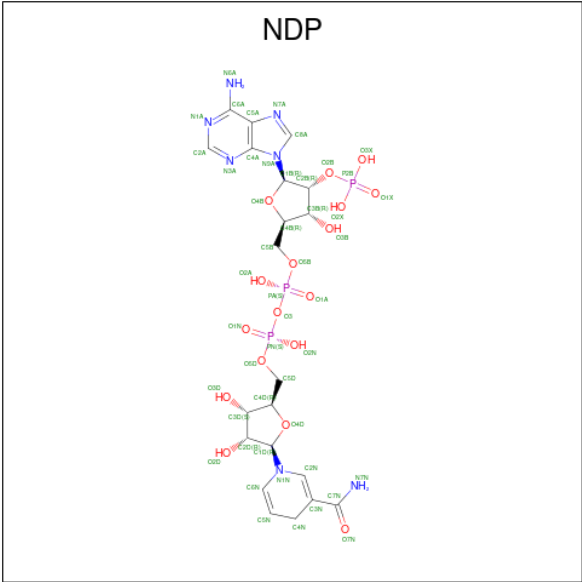
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP F5B5W8
A	1	VAL	-	expression tag	UNP F5B5W8
A	2	ALA	-	expression tag	UNP F5B5W8
A	345	PHE	SER	conflict	UNP F5B5W8
A	385	GLN	GLU	conflict	UNP F5B5W8
A	387	GLN	GLU	conflict	UNP F5B5W8
A	421	LEU	-	expression tag	UNP F5B5W8
A	422	GLU	-	expression tag	UNP F5B5W8
A	423	HIS	-	expression tag	UNP F5B5W8
A	424	HIS	-	expression tag	UNP F5B5W8
A	425	HIS	-	expression tag	UNP F5B5W8
A	426	HIS	-	expression tag	UNP F5B5W8
A	427	HIS	-	expression tag	UNP F5B5W8
A	428	HIS	-	expression tag	UNP F5B5W8
B	0	MET	-	initiating methionine	UNP F5B5W8
B	1	VAL	-	expression tag	UNP F5B5W8
B	2	ALA	-	expression tag	UNP F5B5W8
B	345	PHE	SER	conflict	UNP F5B5W8
B	385	GLN	GLU	conflict	UNP F5B5W8
B	387	GLN	GLU	conflict	UNP F5B5W8
B	421	LEU	-	expression tag	UNP F5B5W8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	422	GLU	-	expression tag	UNP F5B5W8
B	423	HIS	-	expression tag	UNP F5B5W8
B	424	HIS	-	expression tag	UNP F5B5W8
B	425	HIS	-	expression tag	UNP F5B5W8
B	426	HIS	-	expression tag	UNP F5B5W8
B	427	HIS	-	expression tag	UNP F5B5W8
B	428	HIS	-	expression tag	UNP F5B5W8
C	0	MET	-	initiating methionine	UNP F5B5W8
C	1	VAL	-	expression tag	UNP F5B5W8
C	2	ALA	-	expression tag	UNP F5B5W8
C	345	PHE	SER	conflict	UNP F5B5W8
C	385	GLN	GLU	conflict	UNP F5B5W8
C	387	GLN	GLU	conflict	UNP F5B5W8
C	421	LEU	-	expression tag	UNP F5B5W8
C	422	GLU	-	expression tag	UNP F5B5W8
C	423	HIS	-	expression tag	UNP F5B5W8
C	424	HIS	-	expression tag	UNP F5B5W8
C	425	HIS	-	expression tag	UNP F5B5W8
C	426	HIS	-	expression tag	UNP F5B5W8
C	427	HIS	-	expression tag	UNP F5B5W8
C	428	HIS	-	expression tag	UNP F5B5W8
D	0	MET	-	initiating methionine	UNP F5B5W8
D	1	VAL	-	expression tag	UNP F5B5W8
D	2	ALA	-	expression tag	UNP F5B5W8
D	345	PHE	SER	conflict	UNP F5B5W8
D	385	GLN	GLU	conflict	UNP F5B5W8
D	387	GLN	GLU	conflict	UNP F5B5W8
D	421	LEU	-	expression tag	UNP F5B5W8
D	422	GLU	-	expression tag	UNP F5B5W8
D	423	HIS	-	expression tag	UNP F5B5W8
D	424	HIS	-	expression tag	UNP F5B5W8
D	425	HIS	-	expression tag	UNP F5B5W8
D	426	HIS	-	expression tag	UNP F5B5W8
D	427	HIS	-	expression tag	UNP F5B5W8
D	428	HIS	-	expression tag	UNP F5B5W8

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: $C_{21}H_{30}N_7O_{17}P_3$) (labeled as "Ligand of Interest" by depositor).



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

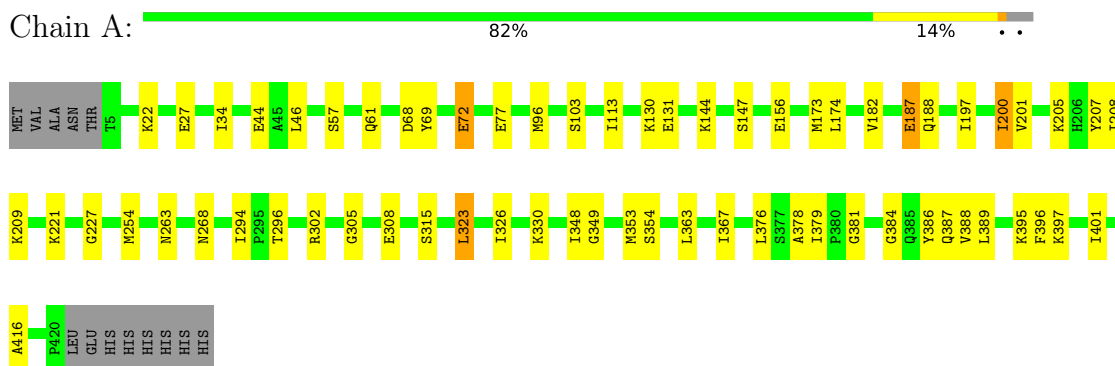
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	140	Total	O	0	0
			140	140		
3	B	177	Total	O	0	0
			177	177		
3	C	165	Total	O	0	0
			165	165		
3	D	155	Total	O	0	0
			155	155		

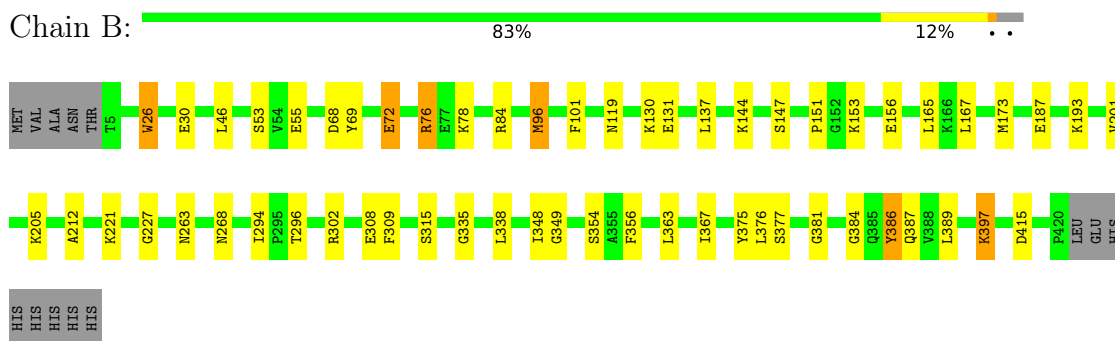
3 Residue-property plots [i](#)

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.

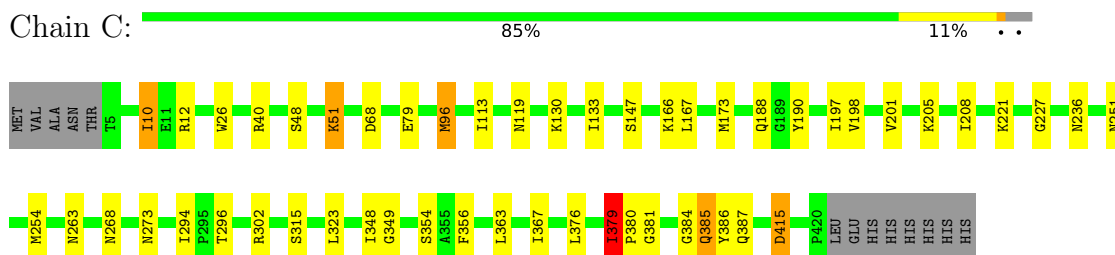
• Molecule 1: CdtA



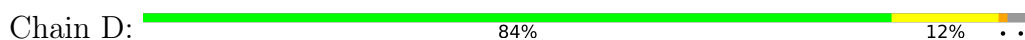
• Molecule 1: CdtA

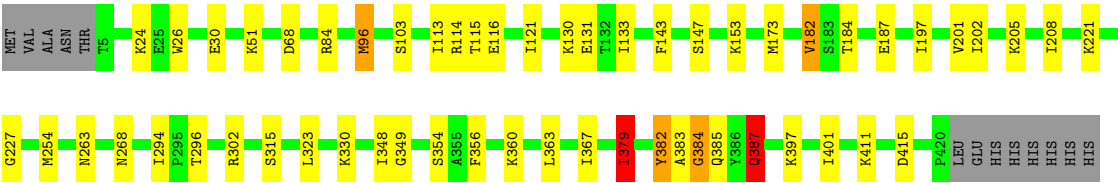


• Molecule 1: CdtA



• Molecule 1: CdtA





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	82.77Å 101.10Å 104.33Å 98.80° 112.36° 106.72°	Depositor
Resolution (Å)	38.91 – 2.42 38.91 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.4 (38.91-2.42) 80.3 (38.91-2.42)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.42Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.206 , 0.234 0.225 , 0.254	Depositor DCC
R_{free} test set	5428 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 15.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14229	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% 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: NDP

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	0.82	4/3441 (0.1%)	1.32	9/4650 (0.2%)
1	B	0.82	0/3428	1.32	12/4633 (0.3%)
1	C	0.83	3/3441 (0.1%)	1.35	14/4648 (0.3%)
1	D	0.81	1/3425 (0.0%)	1.32	11/4630 (0.2%)
All	All	0.82	8/13735 (0.1%)	1.33	46/18561 (0.2%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	379	ILE	CA-C	6.09	1.59	1.52
1	D	379	ILE	CA-C	5.94	1.58	1.53
1	C	379	ILE	CA-CB	5.85	1.59	1.54
1	A	188	GLN	CA-C	5.74	1.58	1.53
1	C	96	MET	SD-CE	-5.66	1.65	1.79
1	A	379	ILE	CA-C	5.22	1.57	1.53
1	A	378	ALA	C-N	5.14	1.38	1.33
1	A	254	MET	SD-CE	-5.09	1.66	1.79

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	415	ASP	CA-CB-CG	8.78	121.38	112.60
1	C	384	GLY	CA-C-N	8.34	136.71	121.70
1	C	384	GLY	C-N-CA	8.34	136.71	121.70
1	C	415	ASP	CA-CB-CG	8.20	120.80	112.60
1	B	268	ASN	CA-C-N	6.66	125.99	121.13
1	B	268	ASN	C-N-CA	6.66	125.99	121.13
1	D	268	ASN	CA-C-N	6.31	125.74	121.13
1	D	268	ASN	C-N-CA	6.31	125.74	121.13
1	A	268	ASN	CA-C-N	6.28	125.71	121.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	ASN	C-N-CA	6.28	125.71	121.13
1	C	10	ILE	CA-C-N	6.19	130.09	120.75
1	C	10	ILE	C-N-CA	6.19	130.09	120.75
1	D	379	ILE	CB-CA-C	6.18	118.33	111.04
1	B	415	ASP	CA-CB-CG	6.12	118.72	112.60
1	C	268	ASN	CA-C-N	6.03	125.53	121.13
1	C	268	ASN	C-N-CA	6.03	125.53	121.13
1	C	10	ILE	CB-CA-C	6.00	117.86	110.73
1	C	68	ASP	CA-CB-CG	5.93	118.53	112.60
1	B	386	TYR	CA-C-N	5.87	129.21	120.87
1	B	386	TYR	C-N-CA	5.87	129.21	120.87
1	C	379	ILE	CB-CA-C	5.82	116.61	111.08
1	A	68	ASP	CA-CB-CG	5.80	118.40	112.60
1	D	68	ASP	CA-CB-CG	5.62	118.22	112.60
1	C	356	PHE	CA-C-N	5.61	127.80	120.28
1	C	356	PHE	C-N-CA	5.61	127.80	120.28
1	B	26	TRP	CA-C-N	5.60	127.79	120.28
1	B	26	TRP	C-N-CA	5.60	127.79	120.28
1	C	263	ASN	CA-C-N	5.50	127.66	120.28
1	C	263	ASN	C-N-CA	5.50	127.66	120.28
1	D	263	ASN	CA-C-N	5.42	127.49	120.44
1	D	263	ASN	C-N-CA	5.42	127.49	120.44
1	A	77	GLU	CB-CG-CD	5.42	121.81	112.60
1	A	263	ASN	CA-C-N	5.39	127.51	120.28
1	A	263	ASN	C-N-CA	5.39	127.51	120.28
1	B	68	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	103	SER	N-CA-C	-5.28	103.41	110.07
1	A	131	GLU	CB-CG-CD	5.24	121.51	112.60
1	A	388	VAL	N-CA-C	-5.24	100.02	107.77
1	D	387	GLN	N-CA-C	5.22	118.45	110.20
1	D	356	PHE	CA-C-N	5.12	127.40	120.38
1	D	356	PHE	C-N-CA	5.12	127.40	120.38
1	B	263	ASN	CA-C-N	5.12	127.13	120.28
1	B	263	ASN	C-N-CA	5.12	127.13	120.28
1	D	103	SER	N-CA-C	-5.06	103.66	109.93
1	B	356	PHE	CA-C-N	5.05	127.30	120.38
1	B	356	PHE	C-N-CA	5.05	127.30	120.38

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	3355	0	3297	30	0
1	B	3345	0	3291	31	0
1	C	3361	0	3319	21	0
1	D	3339	0	3281	23	0
2	A	48	0	26	2	0
2	B	48	0	26	3	0
2	C	48	0	26	1	0
2	D	48	0	26	2	0
3	A	140	0	0	2	0
3	B	177	0	0	0	0
3	C	165	0	0	2	0
3	D	155	0	0	0	0
All	All	14229	0	13292	103	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.

All (103) 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:D:384:GLY:HA3	1:D:385:GLN:HG2	1.56	0.85
1:B:76:ARG:HG2	1:B:76:ARG:HH21	1.45	0.81
1:A:308:GLU:HG3	2:A:501:NDP:O3X	1.90	0.71
1:B:376:LEU:CD1	1:B:389:LEU:HB2	2.21	0.71
1:B:26:TRP:CZ2	1:B:30:GLU:HG3	2.25	0.71
1:A:376:LEU:CD1	1:A:389:LEU:HB2	2.22	0.69
1:C:119:ASN:HA	1:C:205:LYS:HG3	1.73	0.69
1:C:254:MET:SD	1:C:376:LEU:HD12	2.34	0.68
1:B:76:ARG:HH21	1:B:76:ARG:CG	2.09	0.65
1:C:48:SER:HA	1:C:51:LYS:HD2	1.84	0.59
1:D:227:GLY:HA3	1:D:348:ILE:HD13	1.85	0.59
1:A:353:MET:HE1	3:A:660:HOH:O	2.01	0.59
1:C:130:LYS:HE3	3:C:729:HOH:O	2.01	0.59
1:C:227:GLY:HA3	1:C:348:ILE:HD13	1.84	0.59
1:B:227:GLY:HA3	1:B:348:ILE:HD13	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:GLY:HA3	1:A:348:ILE:HD13	1.84	0.57
1:B:55:GLU:HB3	1:B:76:ARG:NH1	2.20	0.57
1:A:384:GLY:HA2	3:A:660:HOH:O	2.05	0.57
1:B:376:LEU:HD11	1:B:389:LEU:HB2	1.87	0.56
1:C:96:MET:HB3	1:C:173[A]:MET:HE1	1.87	0.56
1:A:376:LEU:HD11	1:A:389:LEU:HB2	1.88	0.56
1:A:200:ILE:HG12	1:A:209:LYS:HB2	1.88	0.55
1:B:76:ARG:HG2	1:B:76:ARG:NH2	2.18	0.55
1:C:113:ILE:HD13	1:C:208:ILE:HD12	1.89	0.55
1:D:113:ILE:HD13	1:D:208:ILE:HD12	1.90	0.54
1:C:380:PRO:HB2	1:C:381:GLY:HA2	1.89	0.54
1:B:96:MET:HB3	1:B:173[B]:MET:HE1	1.90	0.54
1:D:26:TRP:CZ2	1:D:30:GLU:HG3	2.41	0.54
1:C:130:LYS:HB2	1:C:197:ILE:HD11	1.88	0.54
1:D:387:GLN:HG2	2:D:501:NDP:H6N	1.90	0.53
1:D:254:MET:HE1	1:D:379:ILE:HD12	1.89	0.53
1:A:113:ILE:HD13	1:A:208:ILE:HD12	1.91	0.53
1:A:69:TYR:HA	1:A:72:GLU:HG3	1.92	0.52
1:C:385:GLN:HG2	1:C:387:GLN:HE21	1.74	0.52
1:A:376:LEU:HD11	1:A:389:LEU:CB	2.39	0.52
1:B:376:LEU:HD11	1:B:389:LEU:CB	2.40	0.51
1:A:323:LEU:HD22	1:A:326:ILE:HD12	1.93	0.51
1:B:69:TYR:HA	1:B:72:GLU:HG3	1.93	0.51
1:D:130:LYS:HB2	1:D:197:ILE:HD11	1.93	0.51
1:A:130:LYS:HB2	1:A:197:ILE:HD11	1.92	0.50
1:A:348:ILE:HG13	1:A:386:TYR:HB3	1.95	0.49
1:A:381:GLY:HA2	1:B:69:TYR:O	2.13	0.49
1:D:24:LYS:HD3	1:D:182:VAL:HG23	1.95	0.48
1:D:173[A]:MET:HG2	1:D:184:THR:CG2	2.43	0.48
1:D:173[B]:MET:HG2	1:D:184:THR:CG2	2.44	0.48
1:D:26:TRP:CE2	1:D:30:GLU:HG3	2.49	0.48
1:D:360:LYS:HA	1:D:411:LYS:HG2	1.95	0.48
1:A:46:LEU:HD13	1:A:174:LEU:HD21	1.97	0.46
1:B:221:LYS:O	1:B:349:GLY:HA2	2.15	0.46
1:C:167:LEU:HD12	1:C:173[A]:MET:HE2	1.97	0.46
1:B:119:ASN:HA	1:B:205:LYS:HG3	1.97	0.46
1:B:26:TRP:CE2	1:B:30:GLU:HG3	2.51	0.46
1:D:84:ARG:HG3	1:D:187:GLU:OE1	2.15	0.46
1:A:221:LYS:O	1:A:349:GLY:HA2	2.16	0.46
1:C:188:GLN:HA	1:C:190:TYR:CE2	2.51	0.46
1:B:375:TYR:CE2	1:B:377:SER:HB2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:MET:HB3	1:D:173[A]:MET:HE1	1.98	0.46
1:B:165:LEU:HA	1:B:212:ALA:O	2.16	0.45
1:B:302:ARG:HB3	1:B:363:LEU:HB3	1.97	0.45
1:B:294:ILE:HG13	1:B:367:ILE:HG21	1.98	0.45
1:A:187:GLU:O	1:A:187:GLU:HG3	2.14	0.45
1:B:335:GLY:C	1:B:397:LYS:HE2	2.42	0.45
1:C:221:LYS:O	1:C:349:GLY:HA2	2.17	0.45
1:A:302:ARG:HB3	1:A:363:LEU:HB3	1.98	0.44
1:D:382:TYR:CD1	1:D:383:ALA:N	2.86	0.44
1:C:348:ILE:HG13	1:C:386:TYR:HB3	1.99	0.44
1:D:254:MET:O	1:D:382:TYR:CE1	2.71	0.44
1:C:294:ILE:HG13	1:C:367:ILE:HG21	1.99	0.43
1:C:379:ILE:HG22	3:C:703:HOH:O	2.18	0.43
1:C:251:ASN:HA	1:C:254:MET:HE2	1.99	0.43
1:D:387:GLN:HG2	2:D:501:NDP:C6N	2.47	0.43
1:B:376:LEU:HD12	1:B:389:LEU:HB2	1.99	0.43
1:D:114:ARG:HG2	1:D:121:ILE:HG12	2.00	0.43
1:D:302:ARG:HB3	1:D:363:LEU:HB3	2.00	0.43
1:B:167:LEU:HD12	1:B:173[B]:MET:HE2	2.00	0.43
1:B:308:GLU:HG2	2:B:501:NDP:C5A	2.49	0.43
1:D:221:LYS:O	1:D:349:GLY:HA2	2.17	0.43
1:B:384:GLY:HA3	2:B:501:NDP:H6N	1.99	0.43
1:D:143:PHE:CE1	1:D:187:GLU:HG3	2.54	0.43
1:B:84:ARG:HG3	1:B:187:GLU:OE1	2.17	0.43
1:C:130:LYS:HD3	1:C:197:ILE:HG12	2.00	0.42
1:C:387:GLN:HG2	2:C:501:NDP:H6N	2.01	0.42
1:D:294:ILE:HG13	1:D:367:ILE:HG21	2.01	0.42
1:B:348:ILE:HG13	1:B:386:TYR:HB3	2.01	0.42
1:A:57:SER:O	1:A:61:GLN:HG3	2.20	0.42
1:B:309:PHE:CE2	1:B:338:LEU:HD22	2.55	0.42
1:A:27:GLU:HG2	1:A:182:VAL:HG22	2.02	0.41
1:C:302:ARG:HB3	1:C:363:LEU:HB3	2.01	0.41
1:A:330:LYS:HG2	1:A:401:ILE:HB	2.02	0.41
1:A:384:GLY:HA3	1:A:387:GLN:NE2	2.35	0.41
1:B:387:GLN:HG2	2:B:501:NDP:H5N	2.02	0.41
1:A:294:ILE:HG13	1:A:367:ILE:HG21	2.02	0.41
1:D:330:LYS:HG2	1:D:401:ILE:HB	2.02	0.41
1:B:137:LEU:HD23	1:B:193:LYS:HB3	2.03	0.41
1:C:12:ARG:HB3	1:C:26:TRP:CD1	2.55	0.41
1:A:384:GLY:HA3	1:A:387:GLN:CD	2.46	0.41
1:A:34:ILE:HD13	1:A:34:ILE:HA	1.95	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:305:GLY:HA3	2:A:501:NDP:O1X	2.21	0.40
1:B:101:PHE:CE2	1:B:151:PRO:HG2	2.56	0.40
1:A:205:LYS:HB3	1:A:207:TYR:CE1	2.57	0.40
1:A:396:PHE:HB3	1:A:416:ALA:HB1	2.03	0.40
1:A:69:TYR:O	1:B:381:GLY:HA2	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	416/429 (97%)	409 (98%)	7 (2%)	0	100	100
1	B	416/429 (97%)	404 (97%)	12 (3%)	0	100	100
1	C	415/429 (97%)	403 (97%)	12 (3%)	0	100	100
1	D	416/429 (97%)	403 (97%)	11 (3%)	2 (0%)	24	35
All	All	1663/1716 (97%)	1619 (97%)	42 (2%)	2 (0%)	48	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	382	TYR
1	D	384	GLY

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	367/386 (95%)	350 (95%)	17 (5%)	24	39
1	B	363/386 (94%)	346 (95%)	17 (5%)	23	39
1	C	368/386 (95%)	350 (95%)	18 (5%)	22	37
1	D	363/386 (94%)	344 (95%)	19 (5%)	21	34
All	All	1461/1544 (95%)	1390 (95%)	71 (5%)	23	37

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	44	GLU
1	A	72	GLU
1	A	96[A]	MET
1	A	96[B]	MET
1	A	144	LYS
1	A	147	SER
1	A	156	GLU
1	A	187	GLU
1	A	200	ILE
1	A	201	VAL
1	A	296	THR
1	A	315	SER
1	A	323	LEU
1	A	354	SER
1	A	395	LYS
1	A	397	LYS
1	B	46	LEU
1	B	53	SER
1	B	72	GLU
1	B	76	ARG
1	B	78	LYS
1	B	96	MET
1	B	130	LYS
1	B	131	GLU
1	B	144	LYS
1	B	147	SER
1	B	153	LYS
1	B	156	GLU
1	B	201	VAL
1	B	296	THR

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Mol	Chain	Res	Type
1	B	315	SER
1	B	354	SER
1	B	397	LYS
1	C	10	ILE
1	C	40	ARG
1	C	51	LYS
1	C	79	GLU
1	C	133	ILE
1	C	147	SER
1	C	166	LYS
1	C	198	VAL
1	C	201	VAL
1	C	236	ASN
1	C	273	ASN
1	C	296	THR
1	C	315	SER
1	C	323	LEU
1	C	354	SER
1	C	379	ILE
1	C	385	GLN
1	C	415	ASP
1	D	51	LYS
1	D	96	MET
1	D	115	THR
1	D	116	GLU
1	D	131	GLU
1	D	133	ILE
1	D	147	SER
1	D	153	LYS
1	D	182	VAL
1	D	201	VAL
1	D	202	ILE
1	D	205	LYS
1	D	296	THR
1	D	315	SER
1	D	323	LEU
1	D	354	SER
1	D	379	ILE
1	D	387	GLN
1	D	397	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	135	ASN
1	A	188	GLN
1	A	262	ASN
1	A	387	GLN
1	A	399	ASN
1	B	36	GLN
1	B	119	ASN
1	B	127	ASN
1	B	134	GLN
1	B	262	ASN
1	B	385	GLN
1	B	387	GLN
1	B	399	ASN
1	C	307	GLN
1	C	392	HIS
1	C	399	ASN
1	D	85	ASN
1	D	134	GLN
1	D	188	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	D	501	-	51,52,52	0.31	0	71,80,80	0.55	0
2	NDP	A	501	-	51,52,52	0.40	0	71,80,80	0.69	1 (1%)
2	NDP	B	501	-	51,52,52	0.41	0	71,80,80	0.59	1 (1%)
2	NDP	C	501	-	51,52,52	0.39	0	71,80,80	0.58	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	D	501	-	-	9/34/77/77	0/5/5/5
2	NDP	A	501	-	-	8/34/77/77	0/5/5/5
2	NDP	B	501	-	-	5/34/77/77	0/5/5/5
2	NDP	C	501	-	-	12/34/77/77	0/5/5/5

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NDP	O5D-PN-O1N	2.69	119.61	108.94
2	B	501	NDP	O5D-PN-O1N	2.11	117.31	108.94

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NDP	C5D-O5D-PN-O3
2	A	501	NDP	C5D-O5D-PN-O1N
2	B	501	NDP	PN-O3-PA-O5B
2	B	501	NDP	C5D-O5D-PN-O3
2	B	501	NDP	C5D-O5D-PN-O1N
2	B	501	NDP	O4D-C1D-N1N-C6N
2	C	501	NDP	C5B-O5B-PA-O1A
2	C	501	NDP	C5B-O5B-PA-O3
2	C	501	NDP	PN-O3-PA-O5B
2	C	501	NDP	C2N-C3N-C7N-N7N

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Mol	Chain	Res	Type	Atoms
2	D	501	NDP	O4B-C4B-C5B-O5B
2	D	501	NDP	C5D-O5D-PN-O3
2	D	501	NDP	C5D-O5D-PN-O1N
2	D	501	NDP	C5D-O5D-PN-O2N
2	D	501	NDP	O4D-C1D-N1N-C6N
2	A	501	NDP	O4D-C1D-N1N-C6N
2	C	501	NDP	O4D-C1D-N1N-C6N
2	D	501	NDP	C3B-C4B-C5B-O5B
2	C	501	NDP	O4D-C4D-C5D-O5D
2	A	501	NDP	O4D-C4D-C5D-O5D
2	D	501	NDP	PN-O3-PA-O1A
2	A	501	NDP	C3D-C4D-C5D-O5D
2	C	501	NDP	O4B-C4B-C5B-O5B
2	A	501	NDP	PN-O3-PA-O5B
2	A	501	NDP	C5D-O5D-PN-O2N
2	B	501	NDP	C5D-O5D-PN-O2N
2	C	501	NDP	C5B-O5B-PA-O2A
2	C	501	NDP	C3B-C4B-C5B-O5B
2	C	501	NDP	C2D-C1D-N1N-C2N
2	D	501	NDP	C4B-C5B-O5B-PA
2	C	501	NDP	C3D-C4D-C5D-O5D
2	C	501	NDP	PN-O3-PA-O1A
2	A	501	NDP	C2B-O2B-P2B-O3X
2	D	501	NDP	PN-O3-PA-O2A

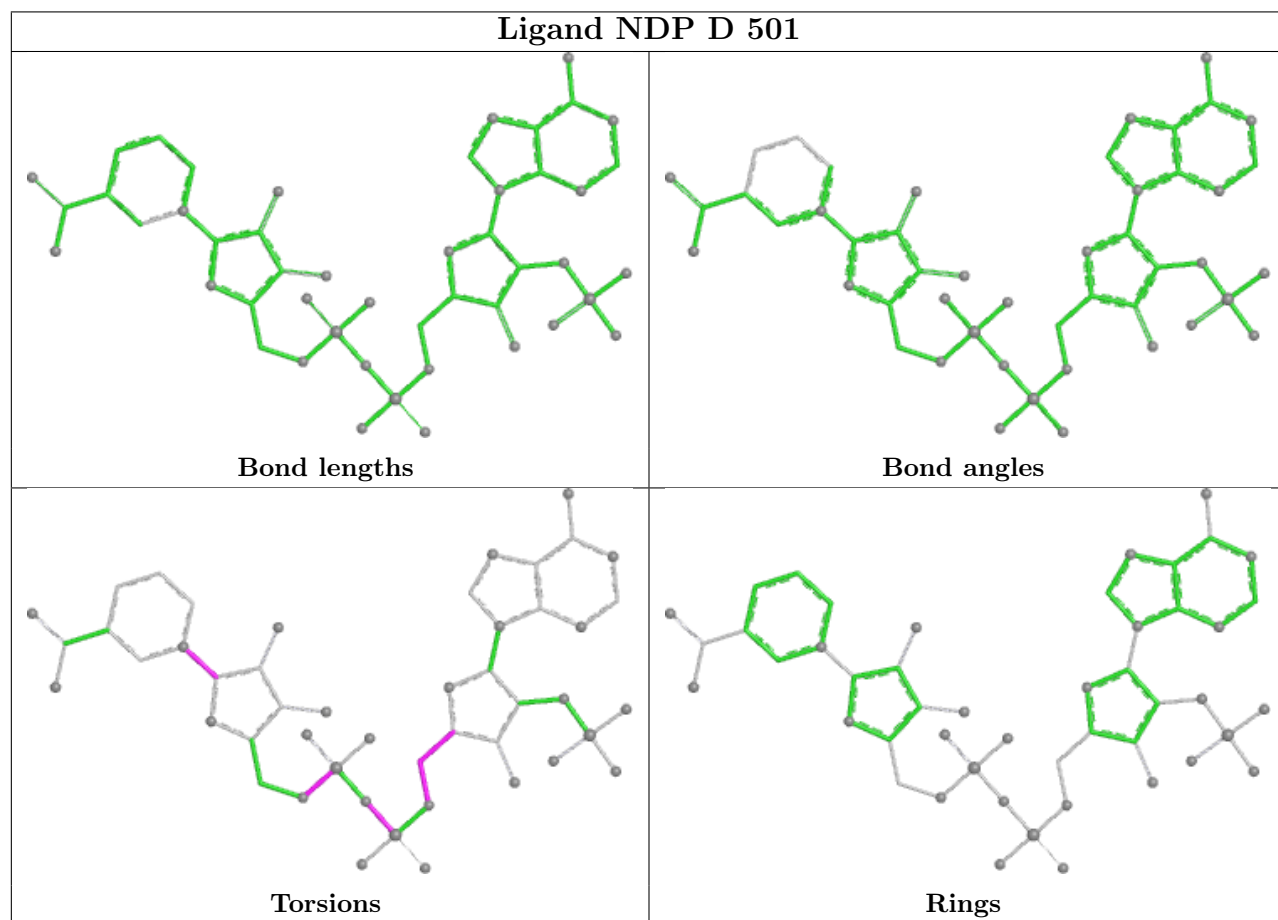
There are no ring outliers.

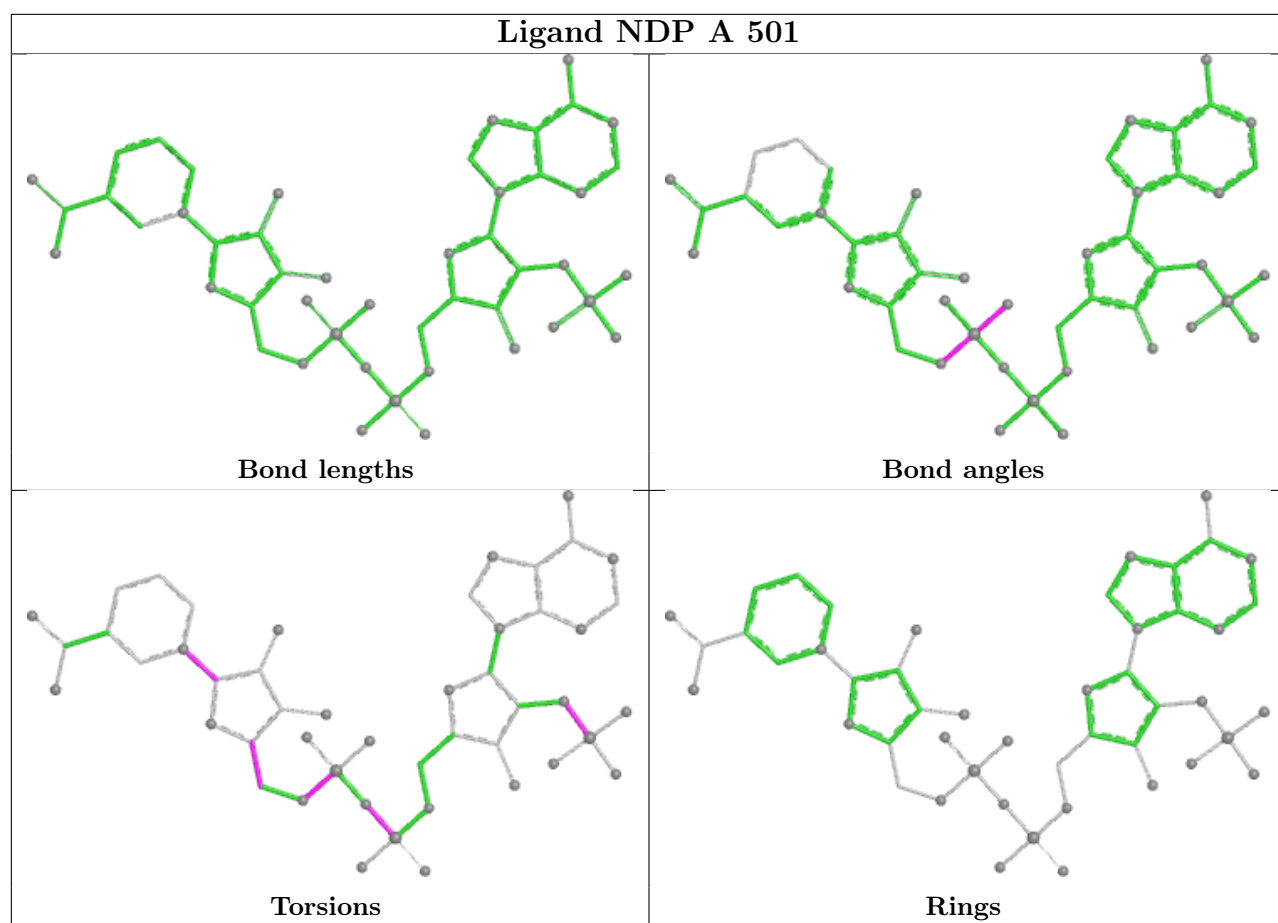
4 monomers are involved in 8 short contacts:

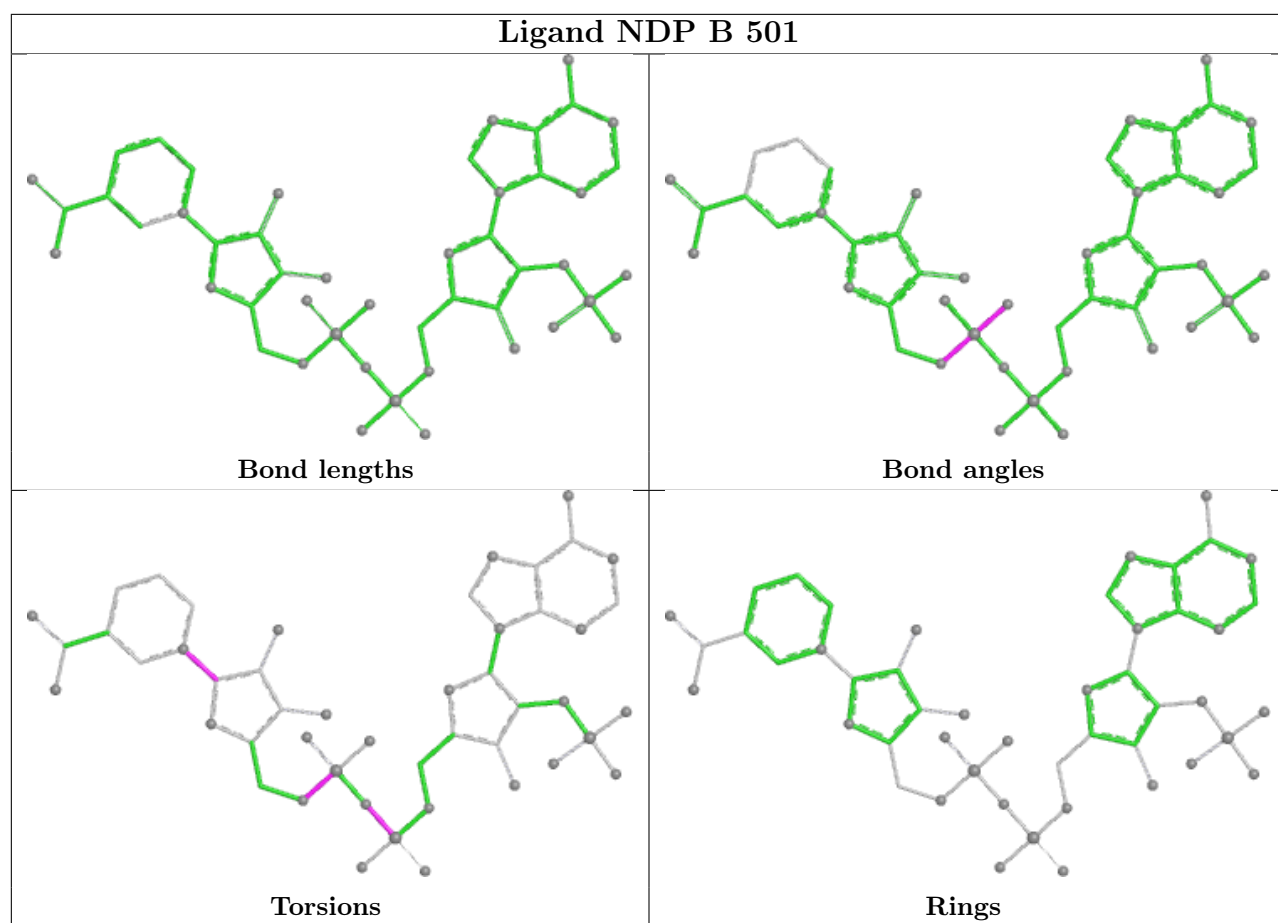
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	501	NDP	2	0
2	A	501	NDP	2	0
2	B	501	NDP	3	0
2	C	501	NDP	1	0

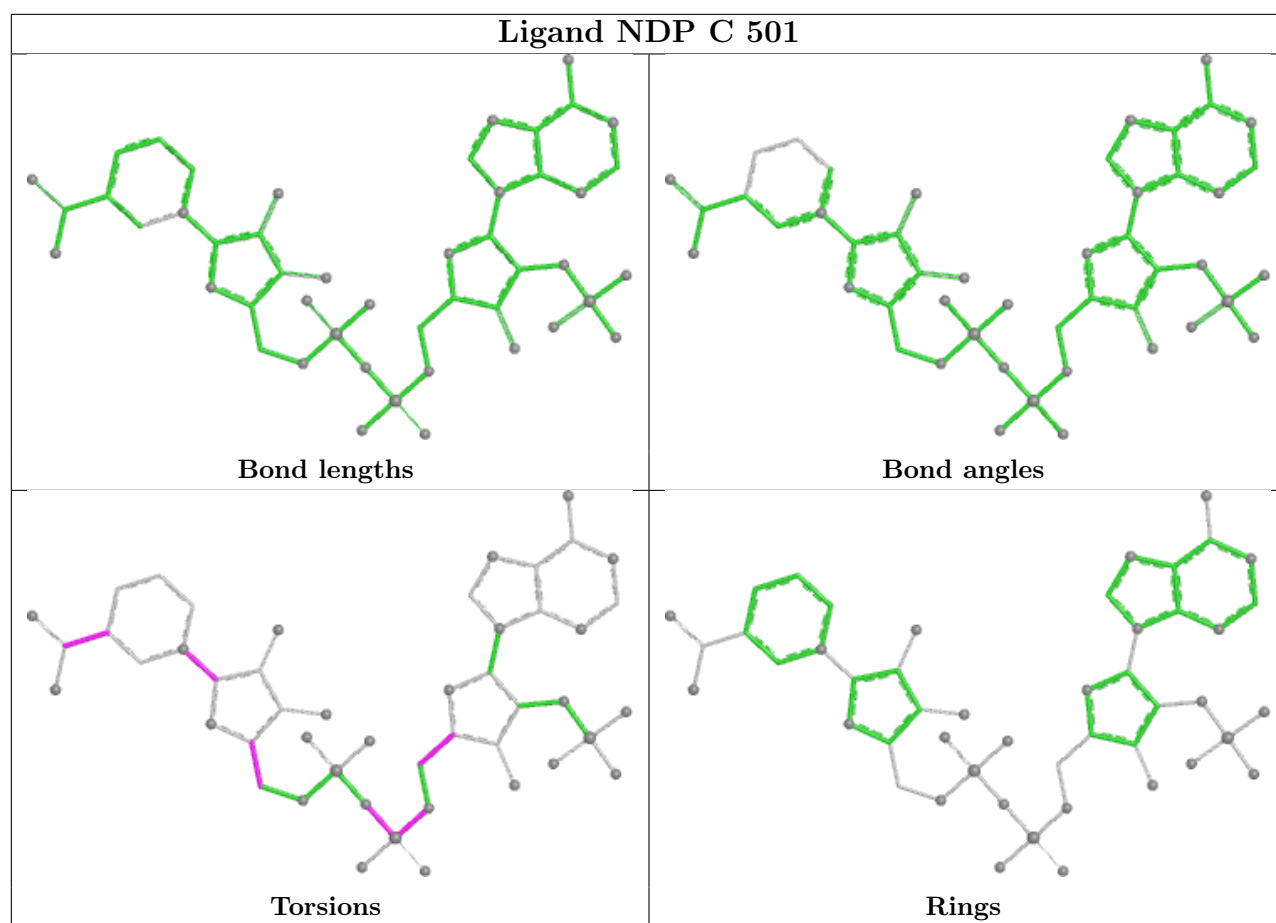
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	416/429 (96%)	-1.66	0 100 100	13, 29, 52, 70	2 (0%)
1	B	416/429 (96%)	-1.69	0 100 100	13, 27, 46, 69	2 (0%)
1	C	416/429 (96%)	-1.64	0 100 100	15, 29, 55, 88	1 (0%)
1	D	416/429 (96%)	-1.61	0 100 100	15, 31, 57, 79	2 (0%)
All	All	1664/1716 (96%)	-1.65	0 100 100	13, 29, 53, 88	7 (0%)

There are no RSRZ outliers to report.

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](#)

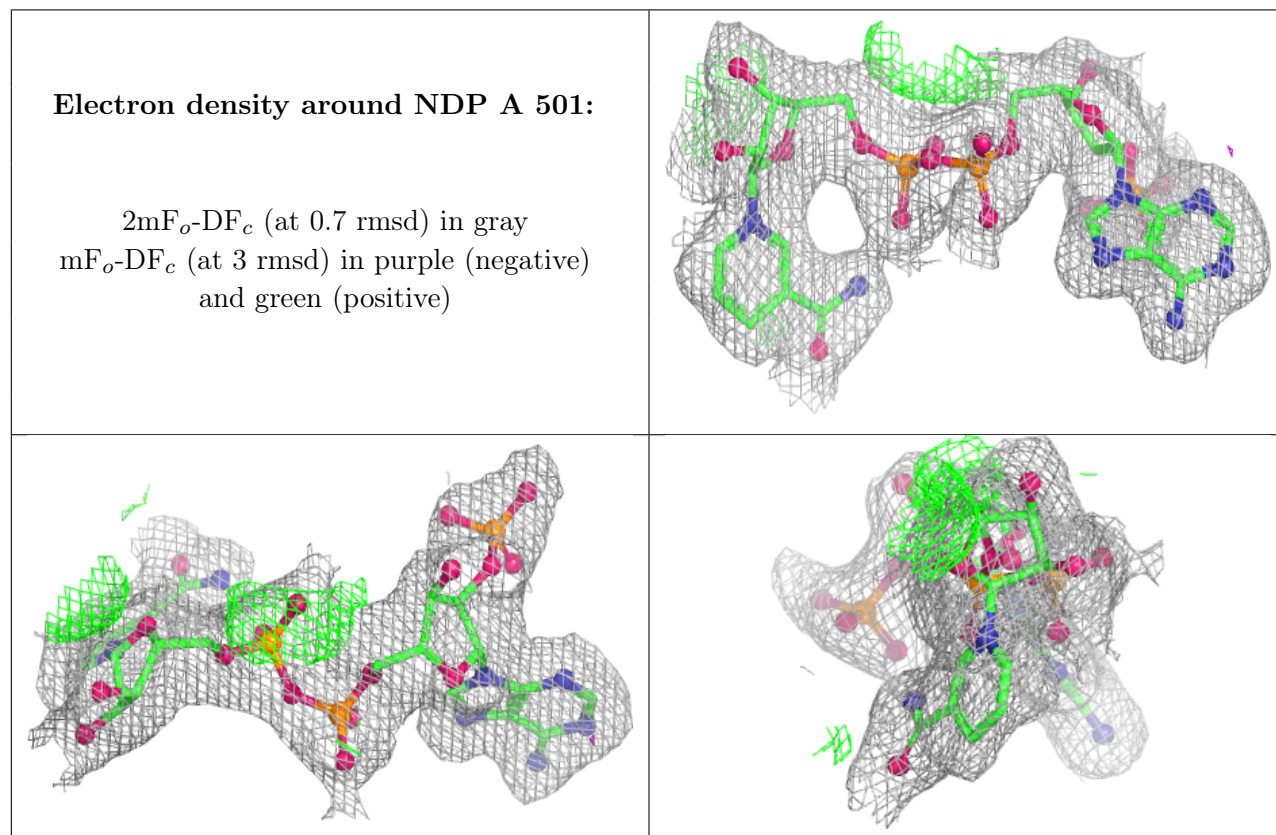
There are no oligosaccharides in this entry.

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.

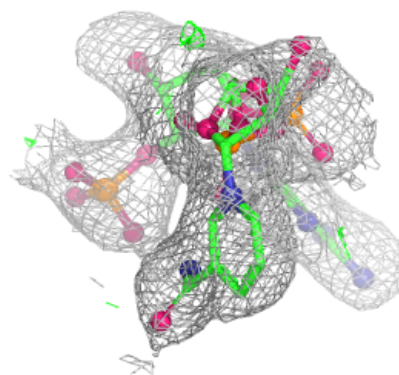
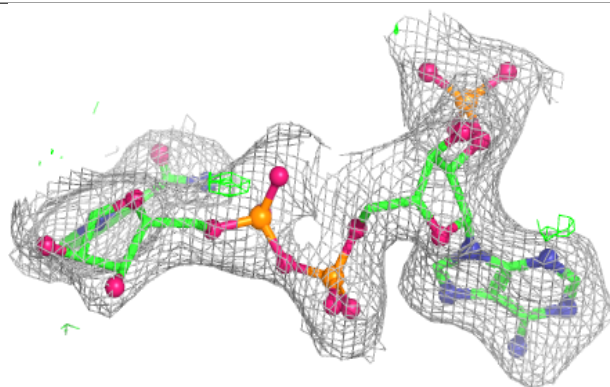
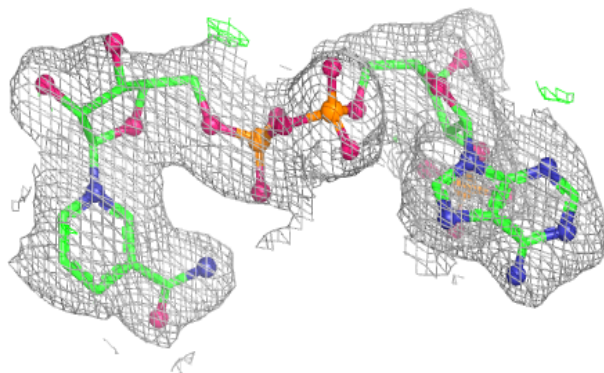
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NDP	A	501	48/48	0.99	0.03	15,36,51,53	0
2	NDP	C	501	48/48	0.99	0.03	35,56,72,73	0
2	NDP	D	501	48/48	0.99	0.03	33,58,66,70	0
2	NDP	B	501	48/48	1.00	0.02	13,39,47,49	0

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.

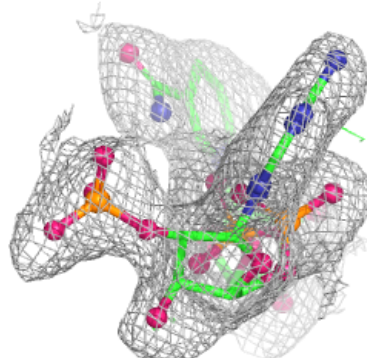
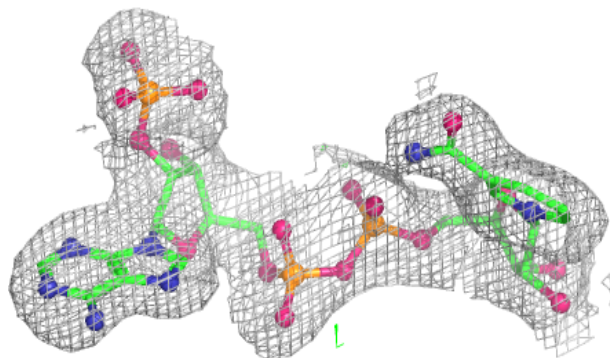
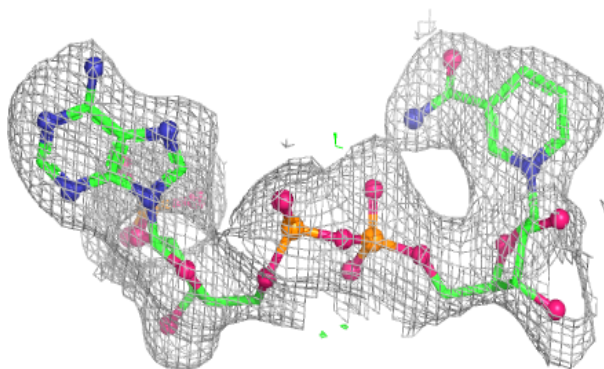


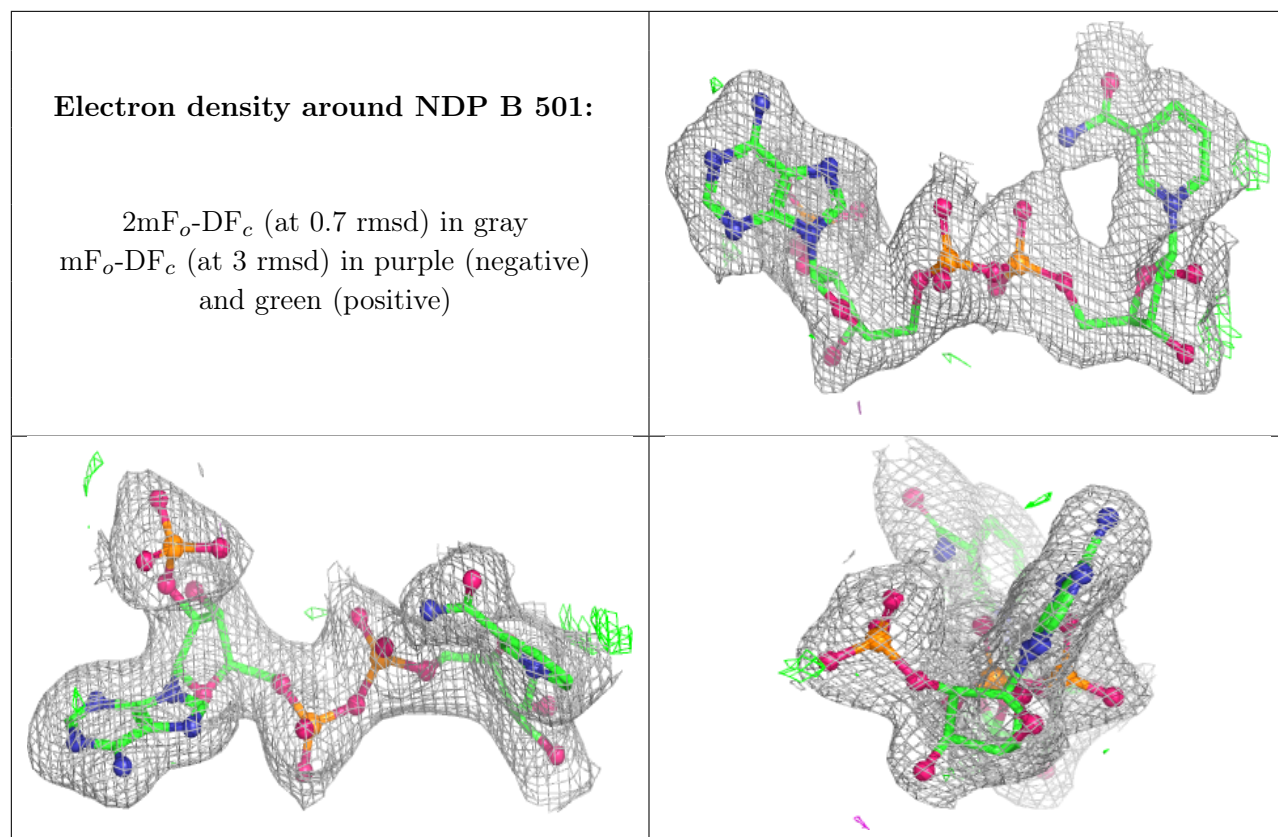
Electron density around NDP C 501:

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

**Electron density around NDP D 501:**

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





6.5 Other polymers [i](#)

There are no such residues in this entry.