



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 12:56 PM BST

PDB ID : 6B0Z
Title : IDH1 R132H mutant in complex with IDH305
Authors : Xie, X.; Kulathila, R.
Deposited on : 2017-09-15
Resolution : 2.33 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 : 2.13.1
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) : 2.13.1

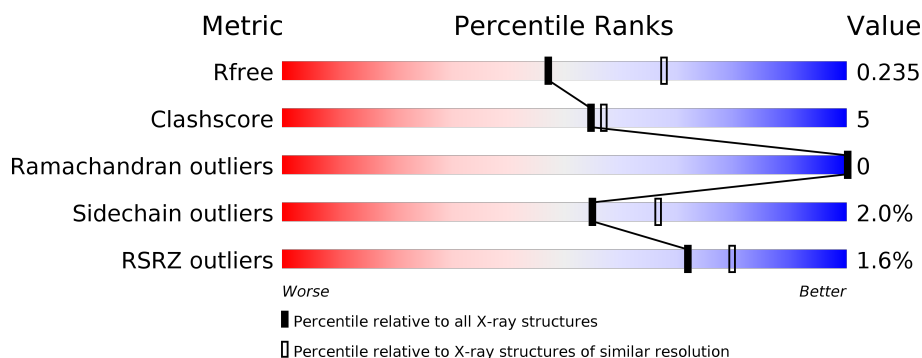
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	417	<div> <div style="width: 86%;"></div> <div style="width: 10%;"></div> <div style="width: 4%;"></div> </div> <div>86% 10% .</div>
1	B	417	<div> <div style="width: 88%;"></div> <div style="width: 10%;"></div> <div style="width: 2%;"></div> </div> <div>88% 10% .</div>
1	C	417	<div> <div style="width: 82%;"></div> <div style="width: 15%;"></div> <div style="width: 3%;"></div> </div> <div>82% 15% ..</div>
1	D	417	<div> <div style="width: 81%;"></div> <div style="width: 14%;"></div> <div style="width: 5%;"></div> </div> <div>81% 14% .</div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13796 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.

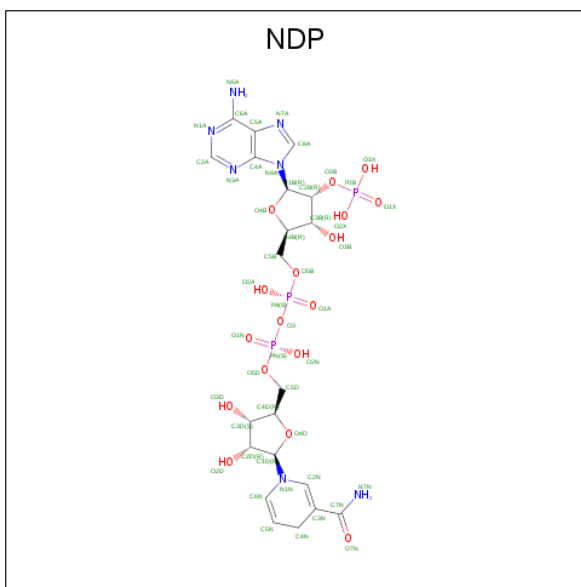
- Molecule 1 is a protein called Isocitrate dehydrogenase [NADP] cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	2	0
			3162	2012	535	596	19			
1	C	410	Total	C	N	O	S	0	0	0
			3249	2066	551	614	18			
1	B	409	Total	C	N	O	S	0	1	0
			3247	2064	551	614	18			
1	D	399	Total	C	N	O	S	0	3	0
			3173	2018	541	594	20			

There are 16 discrepancies between the modelled and reference sequences:

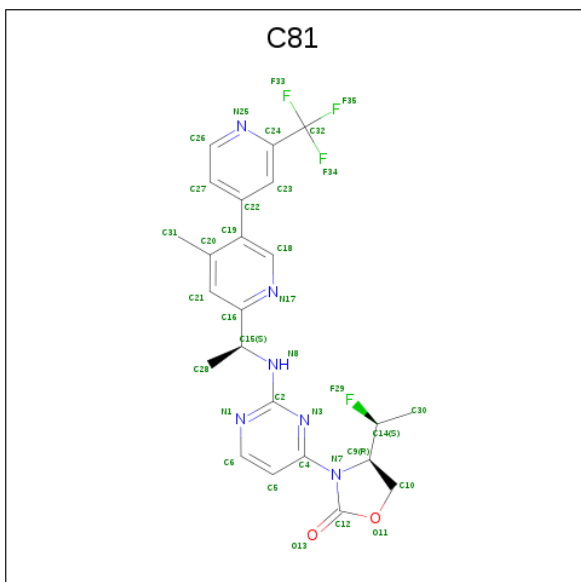
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP O75874
A	-1	PRO	-	expression tag	UNP O75874
A	0	GLY	-	expression tag	UNP O75874
A	132	HIS	ARG	engineered mutation	UNP O75874
C	-2	GLY	-	expression tag	UNP O75874
C	-1	PRO	-	expression tag	UNP O75874
C	0	GLY	-	expression tag	UNP O75874
C	132	HIS	ARG	engineered mutation	UNP O75874
B	-2	GLY	-	expression tag	UNP O75874
B	-1	PRO	-	expression tag	UNP O75874
B	0	GLY	-	expression tag	UNP O75874
B	132	HIS	ARG	engineered mutation	UNP O75874
D	-2	GLY	-	expression tag	UNP O75874
D	-1	PRO	-	expression tag	UNP O75874
D	0	GLY	-	expression tag	UNP O75874
D	132	HIS	ARG	engineered mutation	UNP O75874

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



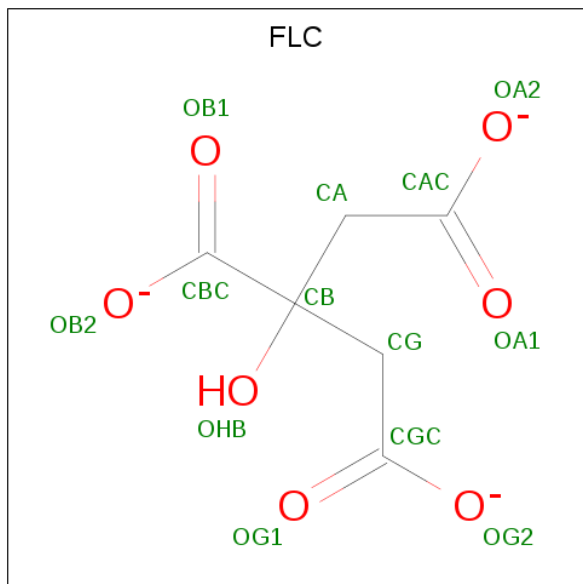
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	C	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	D	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 3 is (4R)-4-[(1S)-1-fluoroethyl]-3-[2-({(1S)-1-[4-methyl-2'-(trifluoromethyl)]3,4'-bipyridin]-6-yl)ethyl}amino)pyrimidin-4-yl]-1,3-oxazolidin-2-one (three-letter code: C81) (formula: C₂₃H₂₂F₄N₆O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			35	23	4	6	2		
3	A	1	Total	C	F	N	O	0	0
			35	23	4	6	2		
3	C	1	Total	C	F	N	O	0	0
			35	23	4	6	2		
3	B	1	Total	C	F	N	O	0	0
			35	23	4	6	2		
3	D	1	Total	C	F	N	O	0	0
			35	23	4	6	2		
3	D	1	Total	C	F	N	O	0	0
			35	23	4	6	2		

- Molecule 4 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7^-$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	6	7		
4	C	1	Total	C	O	0	0
			13	6	7		
4	B	1	Total	C	O	0	0
			13	6	7		
4	D	1	Total	C	O	0	0
			13	6	7		

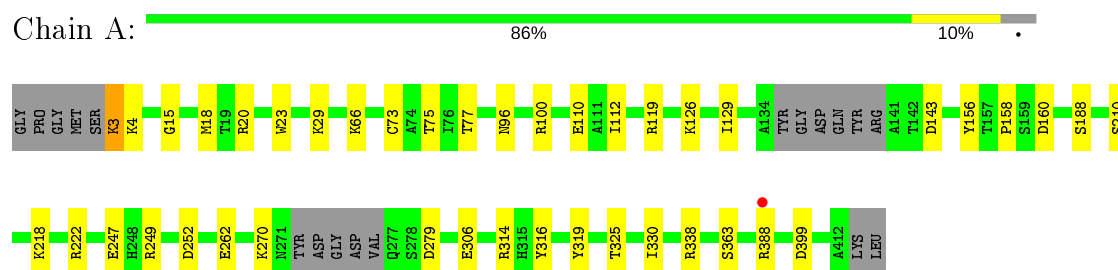
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	175	Total 175	O 175	0	0
5	C	75	Total 75	O 75	0	0
5	B	208	Total 208	O 208	0	0
5	D	53	Total 53	O 53	0	0

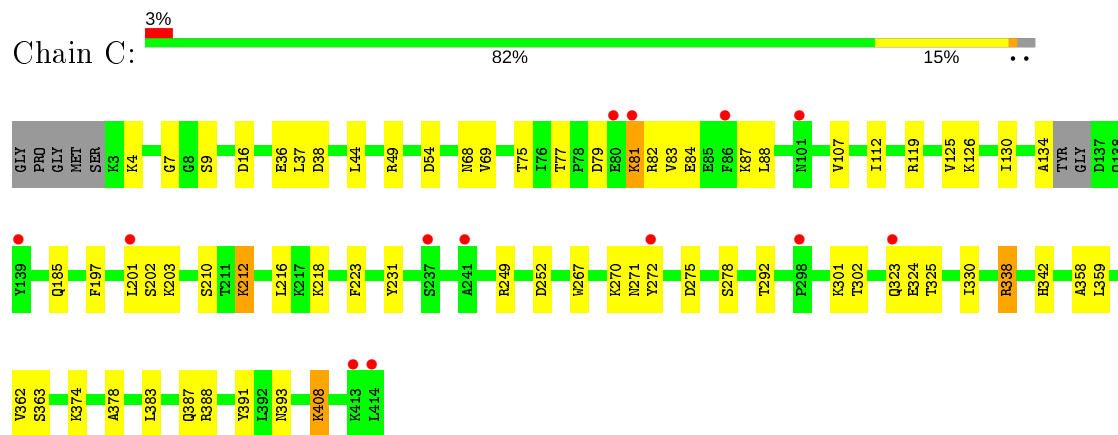
3 Residue-property plots [i](#)

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.

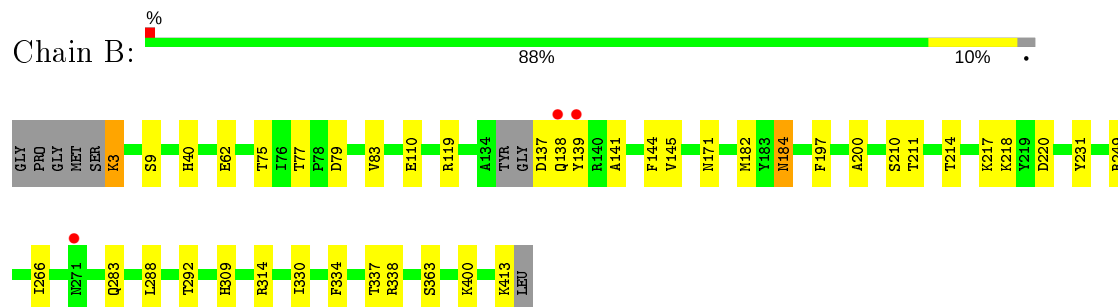
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



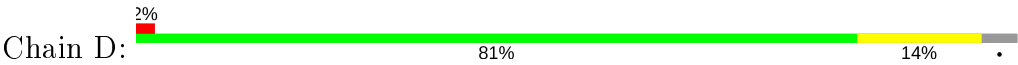
- Molecule 1: Isocitrate dehydrogenase [NADP] cytoplasmic



- 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 21	Depositor
Cell constants a, b, c, α , β , γ	81.50Å 155.31Å 163.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.70 – 2.33 112.46 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.70-2.33) 97.0 (112.46-2.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.85 (at 2.34Å)	Xtriage
Refinement program	PHENIX 1.12 _2829	Depositor
R, R_{free}	0.189 , 0.235 0.189 , 0.235	Depositor DCC
R_{free} test set	4444 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	41.0	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.018 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13796	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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, FLC, C81

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	0.26	0/3232	0.44	0/4355
1	B	0.26	0/3317	0.44	0/4471
1	C	0.26	0/3316	0.45	0/4470
1	D	0.27	0/3246	0.47	1/4372 (0.0%)
All	All	0.26	0/13111	0.45	1/17668 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	314	ARG	NE-CZ-NH1	7.37	123.98	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3162	0	3143	25	0
1	B	3247	0	3217	23	0
1	C	3249	0	3220	42	0
1	D	3173	0	3163	46	0
2	A	48	0	26	3	0
2	B	48	0	26	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	48	0	26	1	0
2	D	48	0	26	0	0
3	A	70	0	0	0	0
3	B	35	0	0	0	0
3	C	35	0	0	0	0
3	D	70	0	0	1	0
4	A	13	0	5	2	0
4	B	13	0	5	3	0
4	C	13	0	5	2	0
4	D	13	0	5	0	0
5	A	175	0	0	3	0
5	B	208	0	0	1	0
5	C	75	0	0	2	0
5	D	53	0	0	1	0
All	All	13796	0	12867	136	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 (136) 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:212:LYS:NZ	1:C:271:ASN:OD1	1.97	0.98
1:D:55:GLN:HE22	1:D:58:LYS:HE2	1.32	0.91
1:D:7:GLY:HA3	1:D:37:LEU:HD23	1.58	0.86
1:D:52:THR:HG21	1:D:56:VAL:HG12	1.67	0.76
1:D:55:GLN:NE2	1:D:59:ASP:OD1	2.19	0.74
1:D:19:THR:HG21	1:D:74:ALA:HB3	1.71	0.72
1:A:218:LYS:NZ	5:A:603:HOH:O	2.21	0.72
1:D:83:VAL:HG23	1:D:88:LEU:HB2	1.75	0.68
2:B:501:NDP:H42N	4:B:503:FLC:HA2	1.81	0.63
1:B:77:THR:OG1	4:B:503:FLC:OB2	2.16	0.63
1:D:52:THR:HG23	1:D:54:ASP:N	2.13	0.62
1:C:77:THR:OG1	4:C:503:FLC:OB2	2.17	0.61
1:A:96:ASN:O	1:A:100:ARG:HG3	2.00	0.61
1:C:323:GLN:OE1	1:C:323:GLN:N	2.34	0.61
1:B:283:GLN:HB2	1:B:288:LEU:HD22	1.81	0.60
1:C:16:ASP:OD2	1:C:49:ARG:NH2	2.35	0.60
1:D:47:GLU:HG2	1:D:48:ASN:N	2.17	0.60
1:C:330:ILE:HD12	1:C:363:SER:HB3	1.84	0.59
1:B:144:PHE:HA	1:D:218:LYS:HD3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:LEU:HG	1:B:309:HIS:HB3	1.84	0.59
1:C:130:ILE:HG21	1:C:272:TYR:CZ	2.38	0.58
1:B:110:GLU:O	1:B:292:THR:HG22	2.04	0.58
1:B:9:SER:OG	1:B:40:HIS:NE2	2.36	0.58
1:D:52:THR:OG1	1:D:55:GLN:HB3	2.05	0.57
1:D:155:THR:HG21	1:D:164:LYS:HE2	1.85	0.57
1:D:48:ASN:O	1:D:52:THR:HG22	2.05	0.57
1:A:77:THR:OG1	4:A:503:FLC:OB1	2.17	0.56
1:C:112:ILE:HD13	1:C:330:ILE:HG22	1.87	0.56
1:A:15:GLY:O	1:A:20:ARG:NH2	2.36	0.55
1:B:400:LYS:NZ	5:B:606:HOH:O	2.36	0.55
1:C:185:GLN:NE2	5:C:601:HOH:O	2.32	0.55
1:C:378:ALA:HB1	1:C:383:LEU:HA	1.88	0.54
1:A:218:LYS:NZ	5:A:602:HOH:O	2.21	0.54
1:C:49:ARG:O	1:C:54:ASP:N	2.37	0.54
1:B:200:ALA:HA	1:B:266:ILE:HD12	1.89	0.53
1:C:201:LEU:HD12	1:C:202:SER:N	2.23	0.53
1:A:388:ARG:NH2	1:A:388:ARG:O	2.40	0.53
1:C:44:LEU:O	1:C:49:ARG:NE	2.41	0.53
1:A:210:SER:HA	1:A:249:ARG:O	2.09	0.53
1:C:69:VAL:HG22	1:C:342:HIS:HD2	1.74	0.52
1:C:75:THR:O	2:C:501:NDP:H2N	2.09	0.52
1:D:67:HIS:O	1:D:343:ARG:NH1	2.42	0.51
1:D:8:GLY:O	1:D:38:ASP:N	2.37	0.51
1:C:107:VAL:HG23	1:C:134:ALA:HB2	1.92	0.51
1:B:314:ARG:NE	2:B:501:NDP:O2X	2.36	0.51
1:D:55:GLN:HE22	1:D:58:LYS:CE	2.15	0.51
1:C:387:GLN:NE2	5:C:607:HOH:O	2.44	0.51
1:C:84:GLU:OE1	1:C:87:LYS:HE2	2.10	0.51
1:A:112:ILE:HD13	1:A:330:ILE:HG22	1.93	0.50
1:A:330:ILE:HD12	1:A:363:SER:HB3	1.92	0.50
1:C:130:ILE:HD13	1:C:267:TRP:HB3	1.92	0.50
1:B:330:ILE:HD12	1:B:363:SER:HB3	1.92	0.50
1:A:319:TYR:HB2	1:A:325:THR:HG21	1.94	0.50
1:C:7:GLY:HA3	1:C:37:LEU:HD23	1.93	0.50
1:D:52:THR:HG23	1:D:54:ASP:H	1.76	0.50
1:D:174:GLU:O	5:D:601:HOH:O	2.20	0.49
1:B:145:VAL:HG23	1:D:218:LYS:HB3	1.95	0.49
1:C:374:LYS:HD2	1:C:391:TYR:CZ	2.47	0.49
1:C:270:LYS:HE3	1:C:271:ASN:OD1	2.12	0.49
1:D:47:GLU:N	1:D:47:GLU:OE2	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:23:TRP:CD2	1:A:73:CYS:HB2	2.47	0.49
1:D:330:ILE:HD12	1:D:363:SER:HB3	1.94	0.49
1:C:252:ASP:OD1	1:C:252:ASP:N	2.42	0.48
1:C:223:PHE:CE2	1:C:270:LYS:HG3	2.48	0.48
1:D:13:MET:HG2	1:D:44:LEU:HD13	1.95	0.48
1:C:358:ALA:O	1:C:362:VAL:HG12	2.13	0.48
4:C:503:FLC:OHB	4:C:503:FLC:OG1	2.29	0.48
1:D:279:ASP:N	1:D:279:ASP:OD1	2.32	0.48
1:D:113:ILE:HG23	1:D:119:ARG:HD3	1.96	0.47
1:A:126:LYS:NZ	1:A:262:GLU:OE1	2.39	0.47
1:D:22:ILE:HD11	1:D:327:THR:HB	1.95	0.47
1:D:44:LEU:O	1:D:49:ARG:NE	2.43	0.47
1:D:49:ARG:HA	1:D:52:THR:HG22	1.97	0.47
1:D:12:GLU:OE1	1:D:27:LYS:NZ	2.48	0.47
1:B:137:ASP:OD2	1:B:184:ASN:ND2	2.49	0.46
1:B:197:PHE:CZ	1:B:231:TYR:HB2	2.49	0.46
1:C:359:LEU:HA	1:C:362:VAL:HG12	1.96	0.46
1:C:79:ASP:OD1	1:C:82:ARG:HG2	2.15	0.46
1:C:83:VAL:HA	1:C:88:LEU:HD12	1.97	0.46
1:A:247:GLU:OE2	1:A:249:ARG:NH1	2.46	0.46
1:A:29:LYS:HE3	1:A:399:ASP:OD1	2.15	0.46
1:C:210:SER:HA	1:C:249:ARG:O	2.15	0.46
1:D:54:ASP:OD1	1:D:57:THR:OG1	2.23	0.46
1:C:4:LYS:HD2	1:C:36:GLU:HG3	1.96	0.46
1:B:214:THR:O	1:B:217:LYS:NZ	2.38	0.46
1:B:211:THR:HB	1:B:220:ASP:HB3	1.97	0.46
1:C:197:PHE:CZ	1:C:231:TYR:HB2	2.50	0.46
1:B:210:SER:HA	1:B:249:ARG:O	2.16	0.45
1:D:211:THR:HB	1:D:220:ASP:HB3	1.98	0.45
1:D:34:TYR:OH	1:D:406:LYS:HD3	2.16	0.45
1:D:9:SER:HA	1:D:38:ASP:HB3	1.99	0.45
1:A:156:TYR:CE2	1:A:158:PRO:HG3	2.52	0.45
1:A:306:GLU:OE1	5:A:601:HOH:O	2.21	0.45
1:C:81:LYS:HE2	1:C:81:LYS:HB2	1.89	0.44
1:D:77:THR:HA	1:D:78:PRO:HD3	1.89	0.44
1:A:314:ARG:NE	2:A:501:NDP:O2X	2.51	0.44
2:A:501:NDP:H42N	4:A:503:FLC:HA2	1.99	0.43
1:A:75:THR:O	2:A:501:NDP:H2N	2.18	0.43
1:D:47:GLU:CG	1:D:48:ASN:N	2.80	0.43
1:B:75:THR:O	2:B:501:NDP:H2N	2.18	0.43
1:D:44:LEU:HD21	1:D:57:THR:HA	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:270:LYS:HD3	1:A:270:LYS:C	2.38	0.43
1:D:392:LEU:HB3	1:D:396:GLU:HB2	2.00	0.43
1:B:413:LYS:HA	1:B:413:LYS:HD3	1.83	0.43
1:D:345:LYS:C	1:D:345:LYS:HD3	2.38	0.43
1:A:3:LYS:HD2	1:A:4:LYS:O	2.19	0.43
1:C:324:GLU:OE2	1:C:388:ARG:NH1	2.52	0.43
1:D:133:HIS:HE1	1:D:188:SER:HB3	1.84	0.42
1:A:110:GLU:HB2	1:A:129:ILE:HG12	2.01	0.42
1:C:125:VAL:HG23	1:C:126:LYS:HG3	2.02	0.42
1:B:3:LYS:HA	1:B:3:LYS:HD3	1.85	0.42
1:C:270:LYS:CE	1:C:271:ASN:OD1	2.68	0.42
1:D:210:SER:HA	1:D:249:ARG:O	2.19	0.42
1:A:18:MET:HB2	1:A:316:TYR:HB2	2.01	0.42
1:C:212:LYS:HE2	1:C:216:LEU:HD12	2.02	0.42
1:D:198:GLN:HE22	1:D:235:TYR:HE1	1.68	0.42
2:B:501:NDP:H42N	4:B:503:FLC:CA	2.50	0.41
1:B:79:ASP:O	1:B:83:VAL:HG23	2.20	0.41
1:C:325:THR:O	1:C:393:ASN:HB2	2.20	0.41
1:D:299:ASP:OD1	1:D:300:GLY:N	2.53	0.41
1:D:361:GLU:O	1:D:365:GLU:HG3	2.19	0.41
1:D:44:LEU:HD12	1:D:44:LEU:N	2.35	0.41
1:D:277:GLN:HB3	3:D:504:C81:C10	2.50	0.41
1:C:301:LYS:O	1:C:342:HIS:NE2	2.32	0.41
1:C:292:THR:HG21	1:C:338:ARG:HH11	1.85	0.41
1:D:133:HIS:HB2	1:D:192:PHE:CE2	2.56	0.41
1:D:13:MET:HG2	1:D:44:LEU:CD1	2.50	0.41
1:A:279:ASP:N	1:A:279:ASP:OD2	2.53	0.41
1:C:408:LYS:HB2	1:C:408:LYS:HE2	1.81	0.40
1:C:68:ASN:HA	1:C:302:THR:HG23	2.02	0.40
1:B:141:ALA:HB2	1:B:182:MET:HG2	2.03	0.40
1:B:334:PHE:HA	1:B:337:THR:OG1	2.22	0.40
1:A:143:ASP:O	1:C:218:LYS:HD2	2.21	0.40
1:C:9:SER:HA	1:C:38:ASP:HB3	2.03	0.40
1:A:66:LYS:HE2	1:B:62:GLU:OE2	2.22	0.40
1:D:160:ASP:CG	1:D:162:THR:HG1	2.25	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	395/417 (95%)	384 (97%)	11 (3%)	0	100	100
1	B	406/417 (97%)	394 (97%)	12 (3%)	0	100	100
1	C	406/417 (97%)	393 (97%)	13 (3%)	0	100	100
1	D	396/417 (95%)	385 (97%)	11 (3%)	0	100	100
All	All	1603/1668 (96%)	1556 (97%)	47 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	339/351 (97%)	332 (98%)	7 (2%)	53	65
1	B	347/351 (99%)	339 (98%)	8 (2%)	50	61
1	C	347/351 (99%)	339 (98%)	8 (2%)	50	61
1	D	341/351 (97%)	337 (99%)	4 (1%)	71	82
All	All	1374/1404 (98%)	1347 (98%)	27 (2%)	55	66

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	119	ARG

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Mol	Chain	Res	Type
1	A	160	ASP
1	A	188	SER
1	A	222	ARG
1	A	252	ASP
1	A	338	ARG
1	C	81	LYS
1	C	119	ARG
1	C	203	LYS
1	C	212	LYS
1	C	275	ASP
1	C	278	SER
1	C	338	ARG
1	C	408	LYS
1	B	3	LYS
1	B	119	ARG
1	B	138	GLN
1	B	139	TYR
1	B	171	ASN
1	B	184	ASN
1	B	218	LYS
1	B	338	ARG
1	D	12	GLU
1	D	140	ARG
1	D	252	ASP
1	D	277	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	90	GLN
1	D	171	ASN
1	D	198	GLN
1	D	277	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 monosaccharides 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$
4	FLC	C	503	-	3,12,12	1.18	0	3,17,17	2.52	2 (66%)
3	C81	D	504	-	35,38,38	0.98	2 (5%)	44,56,56	2.62	20 (45%)
3	C81	B	502	-	35,38,38	0.86	0	44,56,56	2.43	18 (40%)
2	NDP	B	501	-	45,52,52	1.32	3 (6%)	53,80,80	1.19	3 (5%)
3	C81	D	502	-	35,38,38	0.87	1 (2%)	44,56,56	2.56	19 (43%)
2	NDP	D	501	-	45,52,52	1.35	4 (8%)	53,80,80	1.59	8 (15%)
3	C81	A	504	-	35,38,38	0.95	1 (2%)	44,56,56	2.58	18 (40%)
4	FLC	D	503	-	3,12,12	1.12	0	3,17,17	1.87	1 (33%)
2	NDP	A	501	-	45,52,52	1.31	4 (8%)	53,80,80	1.22	3 (5%)
2	NDP	C	501	-	45,52,52	1.31	3 (6%)	53,80,80	1.23	4 (7%)
3	C81	C	502	-	35,38,38	0.86	0	44,56,56	2.38	16 (36%)
3	C81	A	502	-	35,38,38	0.84	0	44,56,56	2.46	18 (40%)
4	FLC	B	503	-	3,12,12	1.21	0	3,17,17	2.22	2 (66%)
4	FLC	A	503	-	3,12,12	1.11	0	3,17,17	1.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FLC	C	503	-	-	0/6/16/16	-
3	C81	D	504	-	-	4/26/39/39	0/4/4/4
3	C81	B	502	-	-	0/26/39/39	0/4/4/4
2	NDP	B	501	-	-	8/30/77/77	0/5/5/5
3	C81	D	502	-	-	0/26/39/39	0/4/4/4
2	NDP	D	501	-	-	8/30/77/77	0/5/5/5
3	C81	A	504	-	-	1/26/39/39	0/4/4/4
4	FLC	D	503	-	-	0/6/16/16	-
2	NDP	A	501	-	-	9/30/77/77	0/5/5/5
2	NDP	C	501	-	-	8/30/77/77	0/5/5/5
3	C81	C	502	-	-	0/26/39/39	0/4/4/4
3	C81	A	502	-	-	0/26/39/39	0/4/4/4
4	FLC	B	503	-	-	0/6/16/16	-
4	FLC	A	503	-	-	0/6/16/16	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	NDP	C4N-C3N	-5.39	1.39	1.49
2	B	501	NDP	C4N-C3N	-5.37	1.39	1.49
2	C	501	NDP	C4N-C3N	-5.37	1.39	1.49
2	A	501	NDP	C4N-C3N	-5.25	1.39	1.49
2	A	501	NDP	C6N-C5N	3.78	1.40	1.33
2	B	501	NDP	C6N-C5N	3.69	1.39	1.33
2	C	501	NDP	C6N-C5N	3.63	1.39	1.33
2	D	501	NDP	C6N-C5N	3.60	1.39	1.33
2	C	501	NDP	C4N-C5N	-3.56	1.39	1.48
2	D	501	NDP	C4N-C5N	-3.55	1.39	1.48
2	B	501	NDP	C4N-C5N	-3.52	1.39	1.48
2	A	501	NDP	C4N-C5N	-3.52	1.39	1.48
3	D	504	C81	C16-C15	2.61	1.55	1.52
3	A	504	C81	C16-C15	2.54	1.55	1.52
2	D	501	NDP	P2B-O2B	-2.48	1.54	1.59
3	D	502	C81	C9-N7	-2.12	1.45	1.48
3	D	504	C81	C9-N7	-2.02	1.45	1.48
2	A	501	NDP	O4B-C1B	2.01	1.43	1.41

All (132) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	504	C81	N1-C2-N3	-7.17	119.76	126.55
3	A	504	C81	N1-C2-N3	-6.93	119.99	126.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	C81	N1-C2-N3	-6.80	120.10	126.55
3	D	502	C81	N1-C2-N3	-6.71	120.19	126.55
3	C	502	C81	N1-C2-N3	-6.70	120.20	126.55
3	A	502	C81	N1-C2-N3	-6.67	120.23	126.55
3	D	504	C81	C6-N1-C2	6.12	120.88	115.45
3	A	504	C81	C6-N1-C2	5.99	120.76	115.45
3	B	502	C81	C6-N1-C2	5.90	120.68	115.45
3	C	502	C81	C6-N1-C2	5.82	120.61	115.45
3	A	502	C81	C6-N1-C2	5.71	120.52	115.45
3	D	502	C81	C6-N1-C2	5.71	120.51	115.45
3	A	504	C81	C18-N17-C16	5.64	122.06	117.95
3	D	504	C81	C18-N17-C16	5.51	121.97	117.95
3	A	504	C81	C32-C24-N25	5.33	121.01	114.61
3	D	502	C81	C10-C9-N7	-5.21	95.56	100.11
3	D	504	C81	C21-C16-N17	-5.16	118.65	123.09
2	D	501	NDP	N3A-C2A-N1A	-5.09	120.72	128.68
3	D	504	C81	C32-C24-N25	5.07	120.70	114.61
3	A	504	C81	C21-C16-N17	-5.05	118.74	123.09
3	D	502	C81	C15-C16-N17	5.04	120.88	116.01
3	A	502	C81	C18-N17-C16	5.03	121.62	117.95
3	D	504	C81	C10-C9-N7	-4.89	95.84	100.11
2	D	501	NDP	O2B-C2B-C1B	-4.84	92.66	110.10
2	B	501	NDP	N3A-C2A-N1A	-4.81	121.15	128.68
2	C	501	NDP	N3A-C2A-N1A	-4.80	121.18	128.68
3	C	502	C81	C18-N17-C16	4.79	121.44	117.95
2	A	501	NDP	N3A-C2A-N1A	-4.70	121.33	128.68
3	D	502	C81	C18-N17-C16	4.68	121.36	117.95
3	B	502	C81	C18-N17-C16	4.49	121.23	117.95
3	A	504	C81	C10-C9-N7	-4.47	96.20	100.11
3	A	502	C81	C21-C16-N17	-4.14	119.53	123.09
3	A	502	C81	C32-C24-N25	4.10	119.54	114.61
3	C	502	C81	C21-C16-N17	-4.09	119.57	123.09
3	B	502	C81	C10-C9-N7	-4.05	96.57	100.11
3	D	502	C81	C21-C16-N17	-4.04	119.61	123.09
3	A	502	C81	C10-C9-N7	-4.00	96.61	100.11
3	B	502	C81	C21-C16-N17	-3.94	119.70	123.09
3	C	502	C81	C32-C24-N25	3.74	119.10	114.61
3	B	502	C81	C15-C16-N17	3.67	119.56	116.01
3	A	502	C81	N8-C2-N1	3.54	122.68	116.65
3	A	502	C81	C15-C16-N17	3.42	119.31	116.01
3	D	502	C81	N8-C2-N1	3.38	122.42	116.65
3	C	502	C81	C15-C16-N17	3.38	119.27	116.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	C81	C32-C24-N25	3.37	118.66	114.61
2	C	501	NDP	PN-O3-PA	-3.34	121.38	132.83
4	C	503	FLC	CB-CA-CAC	-3.31	109.69	114.98
3	D	502	C81	C26-N25-C24	3.25	121.94	117.49
3	B	502	C81	C5-C6-N1	-3.19	120.00	123.96
2	D	501	NDP	PN-O3-PA	-3.18	121.93	132.83
3	B	502	C81	C26-N25-C24	3.17	121.82	117.49
3	C	502	C81	C10-C9-N7	-3.15	97.35	100.11
3	C	502	C81	C5-C6-N1	-3.14	120.06	123.96
3	A	502	C81	C5-C6-N1	-3.10	120.11	123.96
2	D	501	NDP	O4B-C1B-C2B	-3.10	101.21	106.59
3	B	502	C81	N8-C2-N1	3.10	121.93	116.65
3	D	502	C81	C27-C26-N25	-3.09	120.11	123.96
3	D	504	C81	N8-C2-N1	3.08	121.90	116.65
2	A	501	NDP	PN-O3-PA	-3.07	122.31	132.83
3	D	502	C81	C5-C6-N1	-3.06	120.15	123.96
3	C	502	C81	C26-N25-C24	3.05	121.66	117.49
3	D	504	C81	C5-C6-N1	-3.04	120.18	123.96
3	B	502	C81	C27-C26-N25	-3.04	120.18	123.96
3	A	504	C81	C5-C6-N1	-3.04	120.19	123.96
3	C	502	C81	C27-C26-N25	-3.01	120.22	123.96
3	C	502	C81	N8-C2-N1	2.99	121.75	116.65
3	D	504	C81	C26-N25-C24	2.99	121.58	117.49
3	C	502	C81	C5-C4-N3	-2.98	118.66	123.53
4	B	503	FLC	CB-CG-CGC	-2.98	110.22	114.98
3	A	504	C81	N8-C2-N1	2.93	121.65	116.65
3	A	504	C81	C15-C16-N17	2.93	118.84	116.01
3	D	502	C81	C32-C24-N25	2.89	118.08	114.61
3	A	502	C81	C27-C26-N25	-2.89	120.37	123.96
3	D	502	C81	C10-O11-C12	-2.88	106.71	109.94
3	D	504	C81	C15-C16-N17	2.88	118.79	116.01
2	D	501	NDP	O4D-C1D-N1N	2.87	113.67	108.06
3	A	502	C81	C26-N25-C24	2.86	121.41	117.49
3	A	504	C81	C26-N25-C24	2.86	121.40	117.49
3	B	502	C81	C5-C4-N3	-2.84	118.89	123.53
2	B	501	NDP	PN-O3-PA	-2.84	123.08	132.83
3	A	502	C81	C5-C4-N7	2.84	124.56	122.42
3	D	502	C81	C5-C4-N3	-2.80	118.95	123.53
3	A	502	C81	C5-C4-N3	-2.80	118.96	123.53
4	C	503	FLC	CB-CG-CGC	-2.79	110.51	114.98
2	C	501	NDP	O4D-C1D-N1N	2.72	113.36	108.06
2	D	501	NDP	O3X-P2B-O2B	-2.69	93.95	105.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	504	C81	C23-C24-N25	-2.59	119.10	123.31
3	A	504	C81	C5-C4-N3	-2.59	119.31	123.53
4	D	503	FLC	CB-CA-CAC	-2.57	110.87	114.98
3	D	504	C81	C9-N7-C12	2.52	114.21	111.03
2	B	501	NDP	C1D-N1N-C2N	-2.51	116.94	121.11
3	A	504	C81	C9-N7-C12	2.50	114.19	111.03
3	C	502	C81	C9-N7-C12	2.49	114.18	111.03
3	D	502	C81	C5-C4-N7	2.48	124.29	122.42
3	D	504	C81	C5-C4-N3	-2.45	119.53	123.53
3	A	504	C81	C23-C24-N25	-2.43	119.35	123.31
3	A	502	C81	C28-C15-C16	-2.40	108.18	110.95
4	B	503	FLC	CB-CA-CAC	-2.39	111.15	114.98
2	A	501	NDP	C1D-N1N-C2N	-2.38	117.15	121.11
3	B	502	C81	C9-N7-C12	2.37	114.03	111.03
3	D	502	C81	F33-C32-C24	-2.36	108.43	112.47
3	A	502	C81	C9-N7-C12	2.36	114.01	111.03
3	D	502	C81	C19-C18-N17	-2.29	120.88	124.49
3	D	504	C81	C5-C4-N7	2.28	124.14	122.42
3	A	502	C81	C2-N3-C4	2.27	122.13	114.64
3	A	504	C81	C5-C4-N7	2.27	124.13	122.42
3	B	502	C81	C5-C4-N7	2.26	124.13	122.42
3	D	502	C81	C2-N3-C4	2.26	122.10	114.64
3	A	504	C81	C27-C26-N25	-2.25	121.16	123.96
3	D	504	C81	C10-O11-C12	-2.25	107.42	109.94
3	C	502	C81	C2-N3-C4	2.24	122.03	114.64
3	B	502	C81	C2-N3-C4	2.24	122.03	114.64
3	A	502	C81	C19-C18-N17	-2.21	121.01	124.49
2	C	501	NDP	C1D-N1N-C2N	-2.21	117.44	121.11
3	D	504	C81	C2-N3-C4	2.20	121.89	114.64
3	D	504	C81	C27-C26-N25	-2.20	121.23	123.96
3	A	504	C81	C2-N3-C4	2.18	121.84	114.64
3	A	504	C81	C19-C18-N17	-2.18	121.06	124.49
3	B	502	C81	F33-C32-C24	-2.17	108.75	112.47
3	C	502	C81	C5-C4-N7	2.16	124.05	122.42
3	D	504	C81	C23-C22-C19	-2.16	117.04	120.61
3	D	502	C81	O11-C10-C9	-2.13	103.22	105.66
3	D	504	C81	C19-C18-N17	-2.11	121.17	124.49
3	A	504	C81	C10-O11-C12	-2.10	107.59	109.94
3	B	502	C81	F34-C32-C24	-2.10	108.89	112.47
2	D	501	NDP	O2X-P2B-O1X	2.07	118.80	110.68
3	D	502	C81	C9-N7-C12	2.07	113.65	111.03
3	A	502	C81	F34-C32-C24	-2.06	108.96	112.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	504	C81	C27-C22-C19	2.05	124.24	120.91
3	B	502	C81	C19-C18-N17	-2.03	121.30	124.49
3	C	502	C81	F34-C32-C24	-2.01	109.03	112.47
2	D	501	NDP	C1D-N1N-C2N	-2.00	117.78	121.11

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	NDP	C5D-O5D-PN-O1N
2	B	501	NDP	C5D-O5D-PN-O2N
2	D	501	NDP	C2B-O2B-P2B-O2X
2	D	501	NDP	C5D-O5D-PN-O1N
2	D	501	NDP	C5D-O5D-PN-O2N
2	A	501	NDP	C2B-O2B-P2B-O1X
2	A	501	NDP	C5D-O5D-PN-O1N
2	A	501	NDP	C5D-O5D-PN-O2N
2	C	501	NDP	C5D-O5D-PN-O1N
2	C	501	NDP	C5D-O5D-PN-O2N
2	D	501	NDP	O4B-C4B-C5B-O5B
2	D	501	NDP	C3B-C4B-C5B-O5B
2	C	501	NDP	C3B-C4B-C5B-O5B
2	C	501	NDP	O4B-C4B-C5B-O5B
2	B	501	NDP	O4B-C4B-C5B-O5B
2	A	501	NDP	O4B-C4B-C5B-O5B
2	B	501	NDP	C3B-C4B-C5B-O5B
2	A	501	NDP	C3B-C4B-C5B-O5B
2	D	501	NDP	C2B-O2B-P2B-O1X
2	A	501	NDP	C2B-O2B-P2B-O2X
2	C	501	NDP	C2B-O2B-P2B-O2X
2	B	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	O4D-C1D-N1N-C2N
2	C	501	NDP	O4D-C1D-N1N-C2N
3	D	504	C81	C18-C19-C22-C23
3	D	504	C81	C18-C19-C22-C27
2	B	501	NDP	C2D-C1D-N1N-C2N
2	C	501	NDP	C2D-C1D-N1N-C2N
2	D	501	NDP	O4D-C1D-N1N-C2N
2	A	501	NDP	C2D-C1D-N1N-C2N
2	B	501	NDP	C5D-O5D-PN-O3
2	D	501	NDP	C5D-O5D-PN-O3
2	A	501	NDP	C5D-O5D-PN-O3

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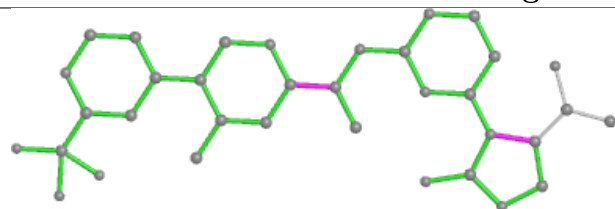
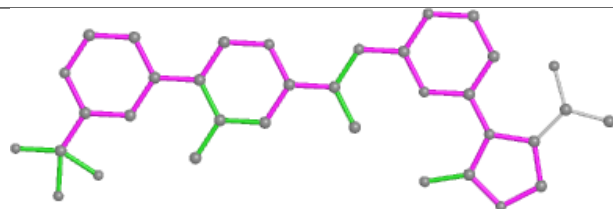
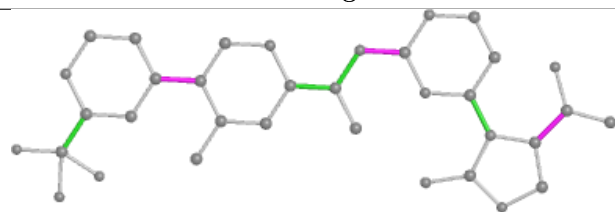
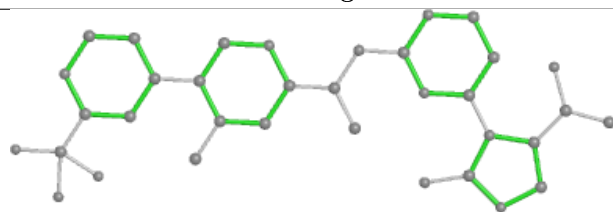
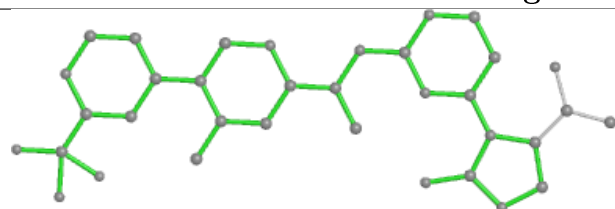
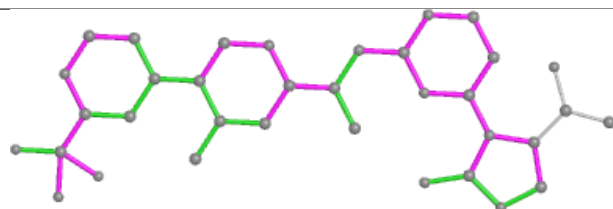
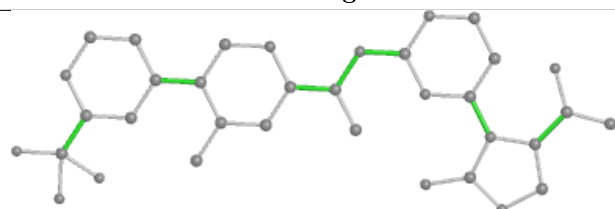
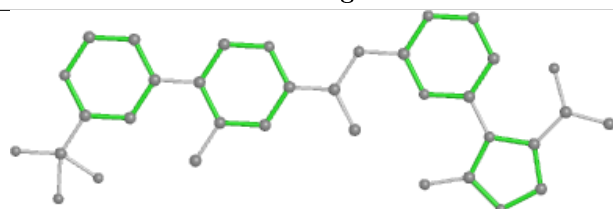
Mol	Chain	Res	Type	Atoms
2	C	501	NDP	C5D-O5D-PN-O3
3	D	504	C81	N1-C2-N8-C15
2	B	501	NDP	O4D-C1D-N1N-C6N
3	D	504	C81	C30-C14-C9-N7
3	A	504	C81	C30-C14-C9-N7

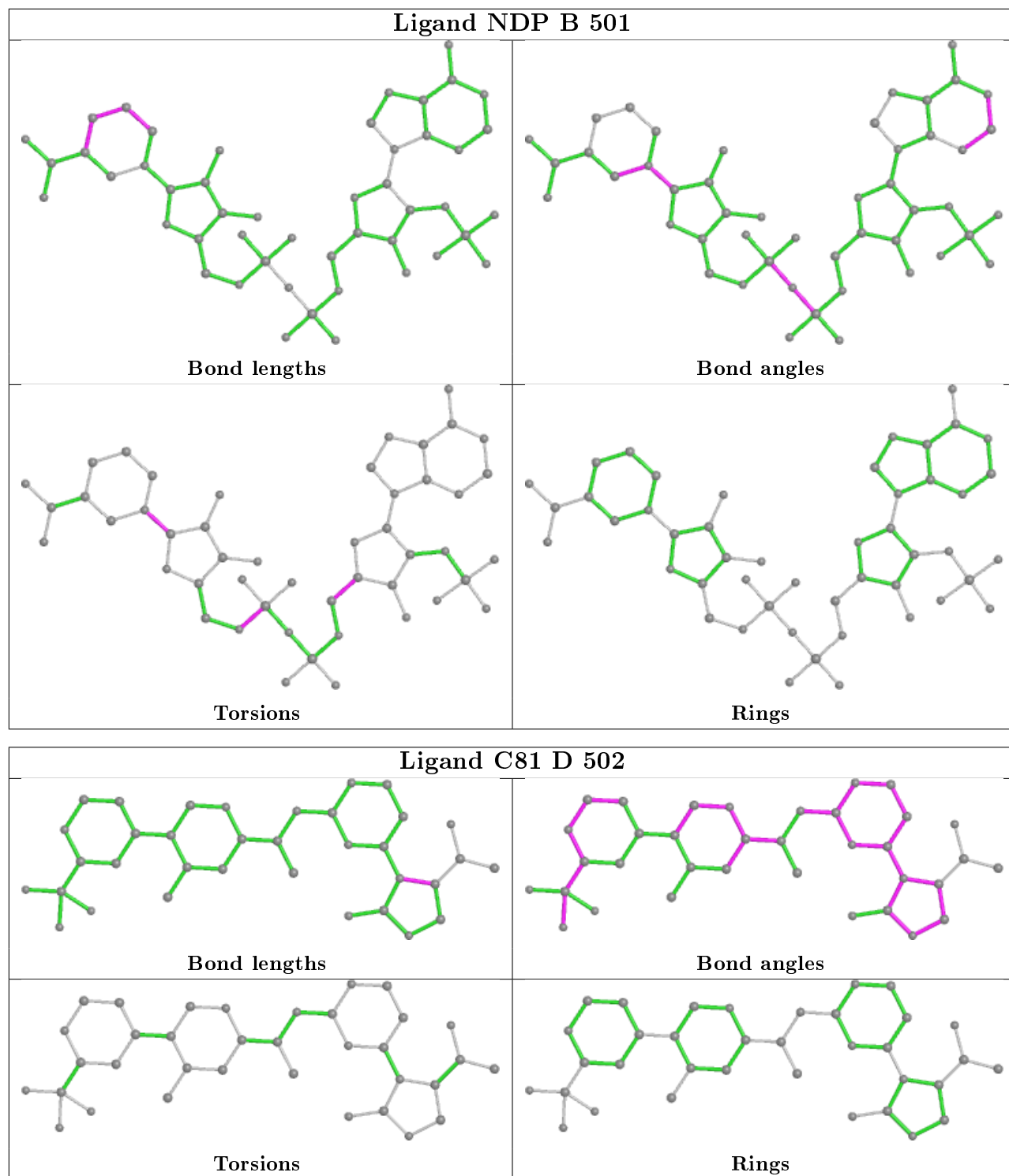
There are no ring outliers.

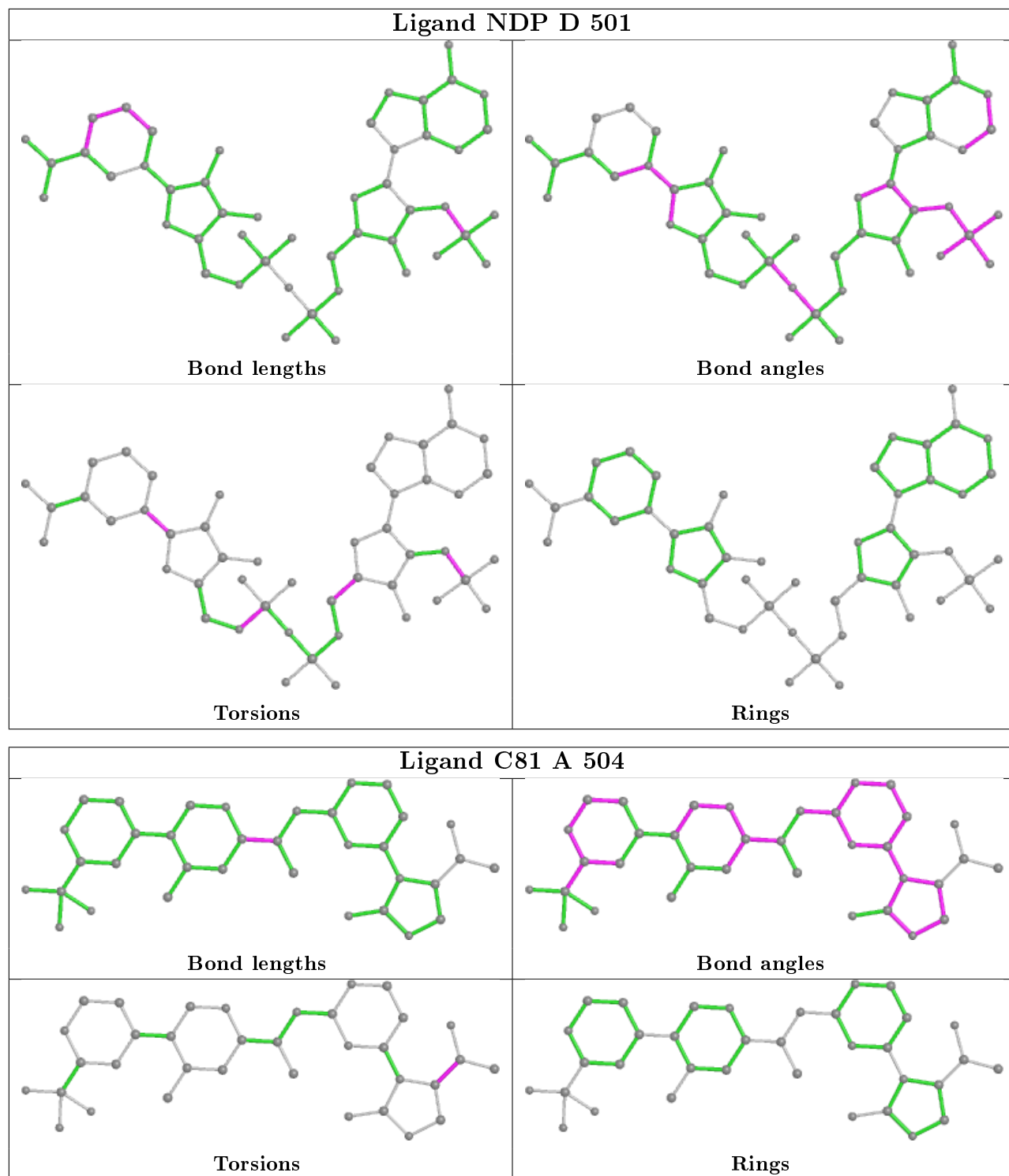
7 monomers are involved in 13 short contacts:

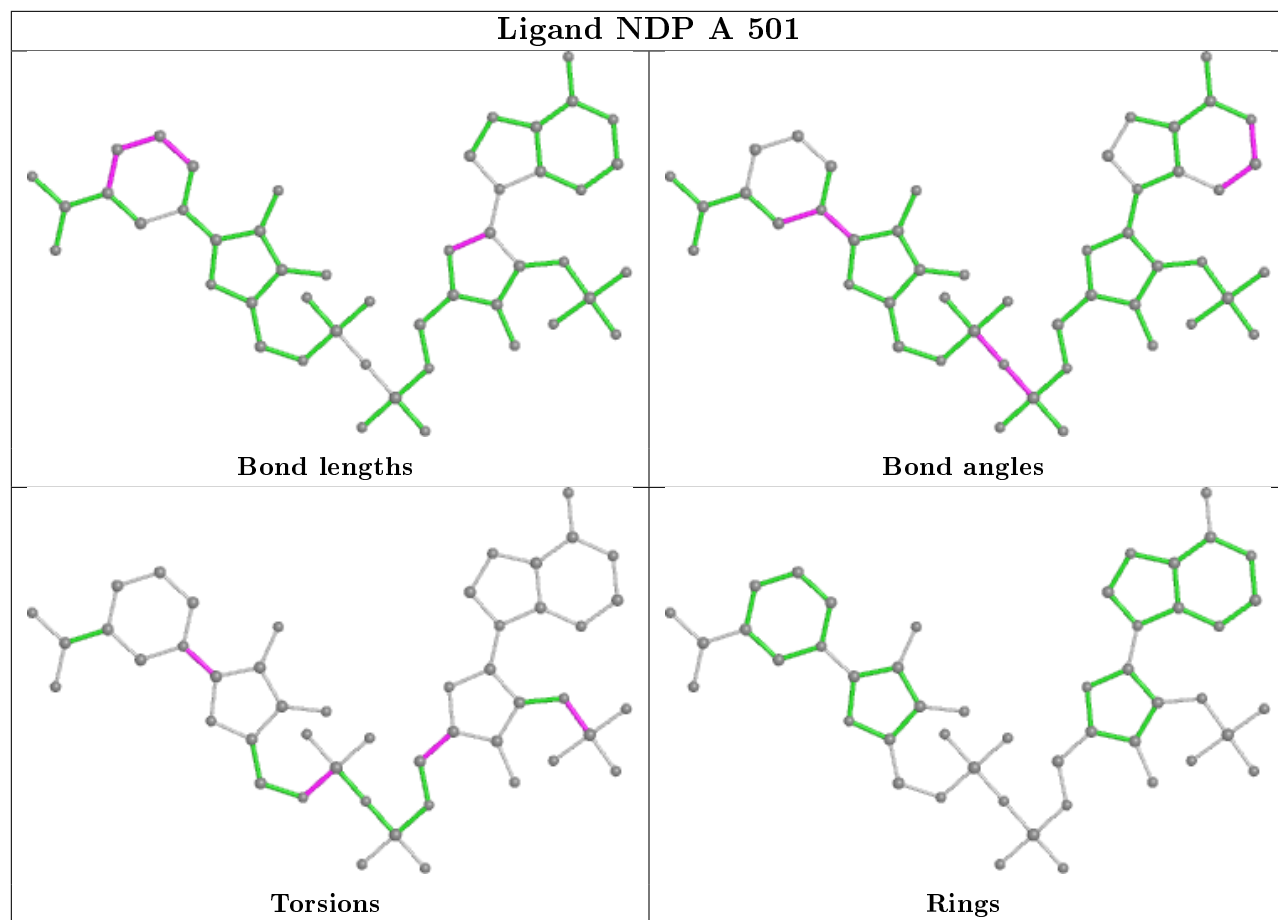
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	503	FLC	2	0
3	D	504	C81	1	0
2	B	501	NDP	4	0
2	A	501	NDP	3	0
2	C	501	NDP	1	0
4	B	503	FLC	3	0
4	A	503	FLC	2	0

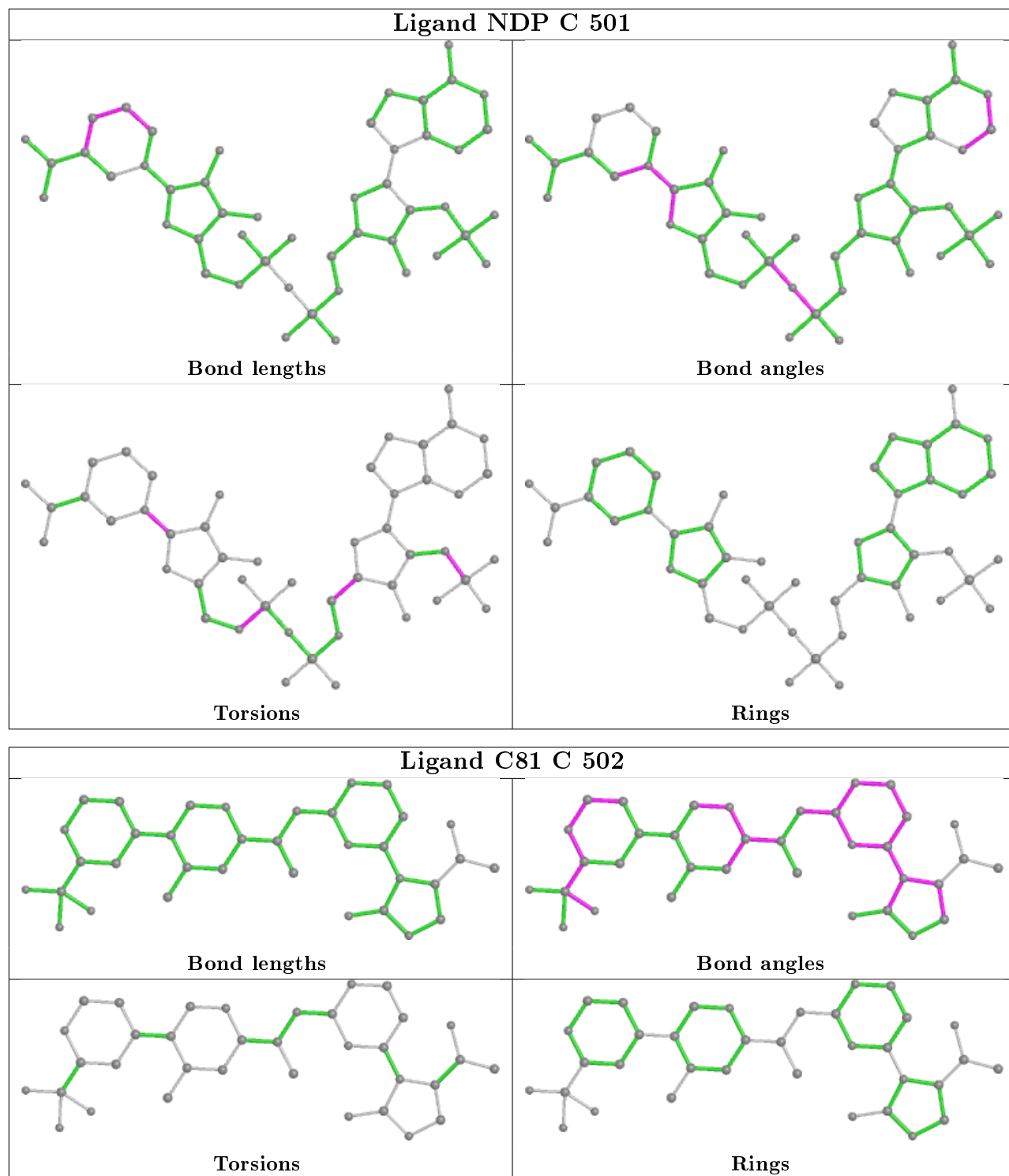
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

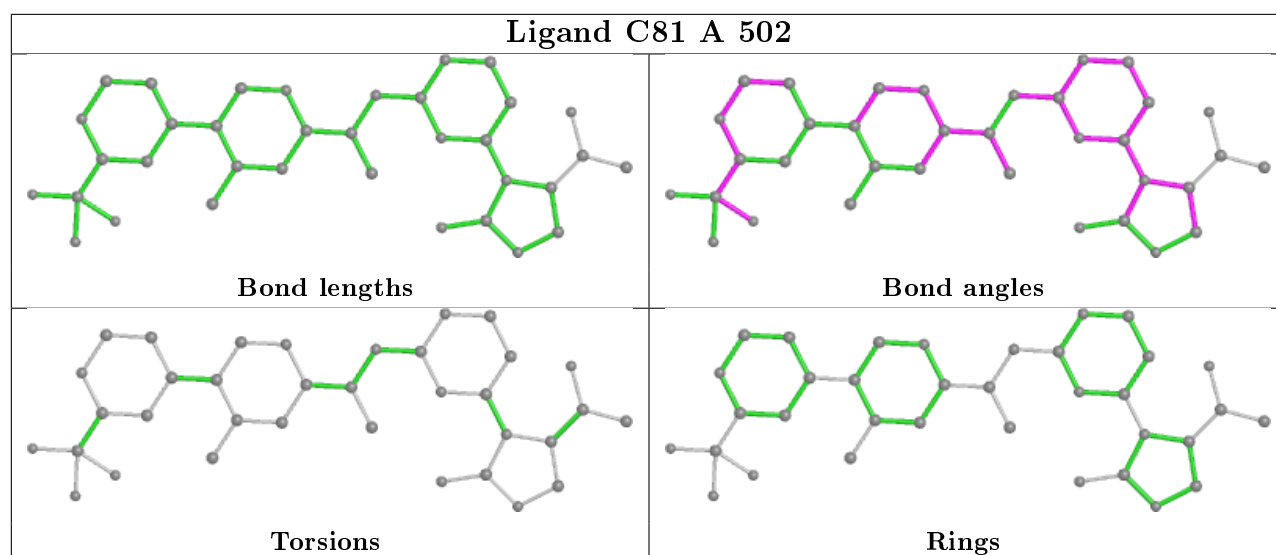
Ligand C81 D 504**Bond lengths****Bond angles****Torsions****Rings****Ligand C81 B 502****Bond lengths****Bond angles****Torsions****Rings**











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, 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	399/417 (95%)	-0.11	1 (0%) 94 97	28, 39, 61, 73	0
1	B	409/417 (98%)	-0.13	3 (0%) 87 92	26, 38, 55, 83	0
1	C	410/417 (98%)	0.13	13 (3%) 47 58	31, 51, 82, 98	0
1	D	399/417 (95%)	0.16	9 (2%) 60 69	32, 56, 81, 94	0
All	All	1617/1668 (96%)	0.01	26 (1%) 72 80	26, 45, 75, 98	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	86	PHE	5.1
1	C	139	TYR	4.9
1	B	139	TYR	3.7
1	D	10	VAL	3.6
1	D	86	PHE	3.2
1	B	271	ASN	3.0
1	C	80	GLU	3.0
1	D	88	LEU	3.0
1	B	138	GLN	2.7
1	D	411	GLN	2.7
1	C	272	TYR	2.5
1	C	414	LEU	2.5
1	C	201	LEU	2.4
1	D	37	LEU	2.4
1	C	241	ALA	2.4
1	C	323	GLN	2.3
1	A	388	ARG	2.3
1	D	125	VAL	2.3
1	D	383	LEU	2.2
1	D	89	LYS	2.1
1	D	271	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	81	LYS	2.1
1	C	298	PRO	2.1
1	C	413	LYS	2.1
1	C	101	ASN	2.1
1	C	237	SER	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 monosaccharides in this entry.

6.4 Ligands [i](#)

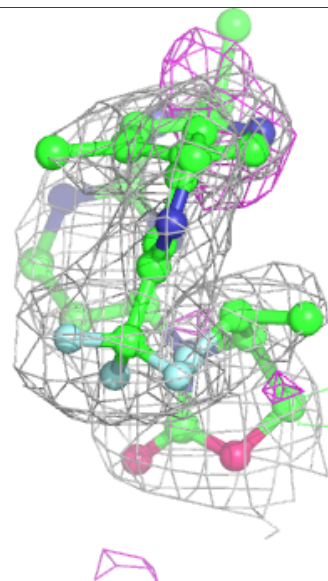
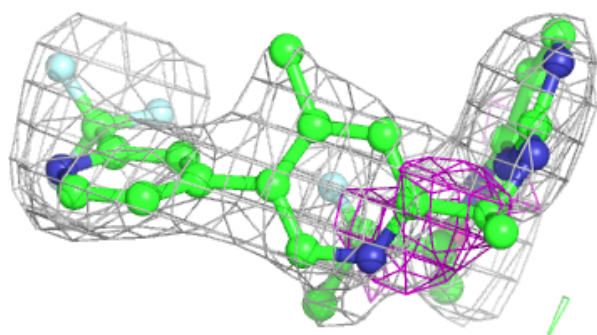
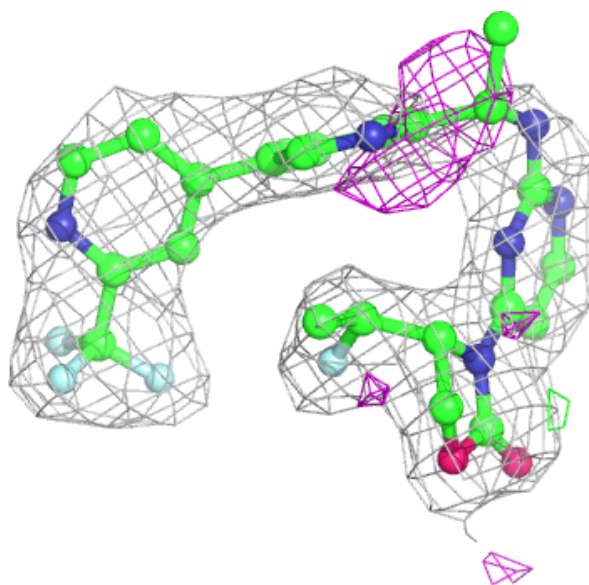
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	FLC	D	503	13/13	0.72	0.25	73,77,87,89	0
4	FLC	C	503	13/13	0.82	0.29	86,92,96,97	0
4	FLC	A	503	13/13	0.85	0.24	56,59,68,69	0
4	FLC	B	503	13/13	0.89	0.21	51,60,68,69	0
3	C81	A	504	35/35	0.92	0.24	53,59,65,66	0
3	C81	D	502	35/35	0.93	0.19	41,51,70,72	0
2	NDP	C	501	48/48	0.94	0.13	49,60,71,73	0
3	C81	D	504	35/35	0.94	0.26	50,63,69,70	0
2	NDP	D	501	48/48	0.94	0.13	42,59,69,81	0
3	C81	C	502	35/35	0.96	0.14	38,44,54,60	0
3	C81	A	502	35/35	0.96	0.15	30,39,54,56	0
2	NDP	A	501	48/48	0.98	0.12	28,39,47,49	0
3	C81	B	502	35/35	0.98	0.14	30,37,53,60	0
2	NDP	B	501	48/48	0.98	0.12	24,34,41,59	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.

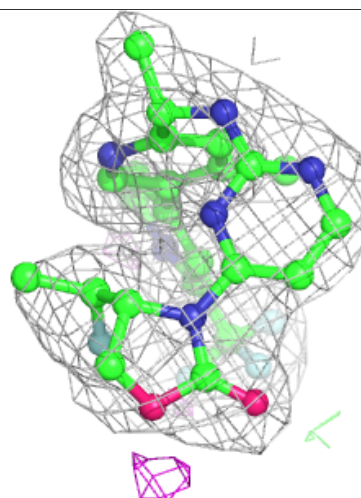
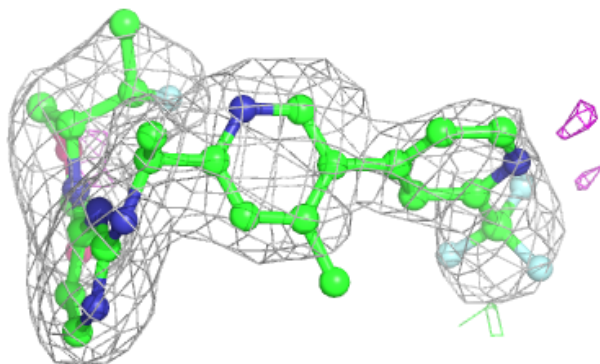
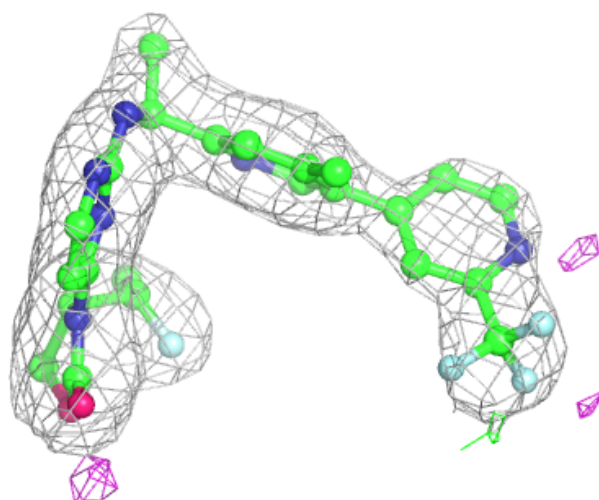
Electron density around C81 A 504:

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



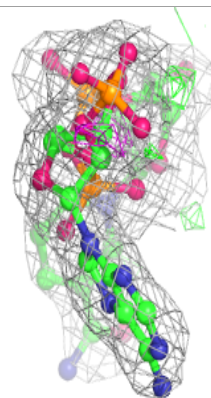
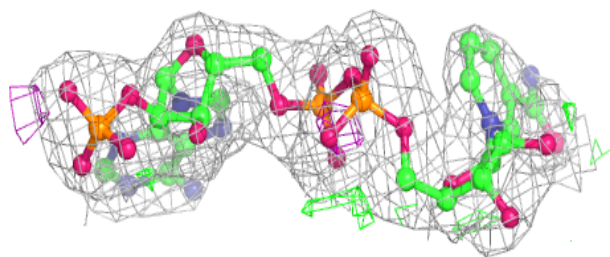
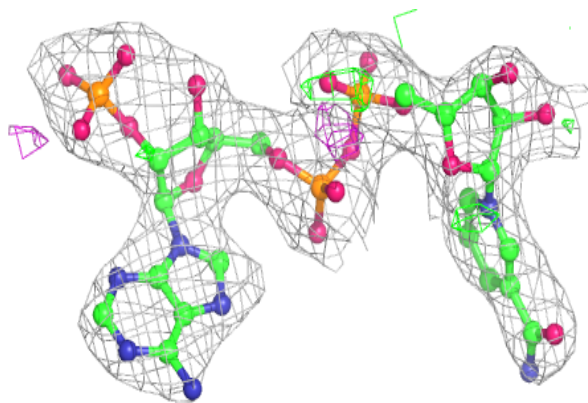
Electron density around C81 D 502:

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



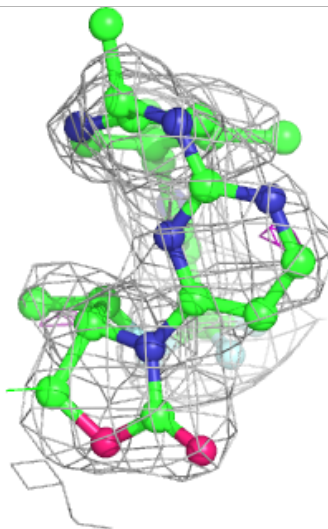
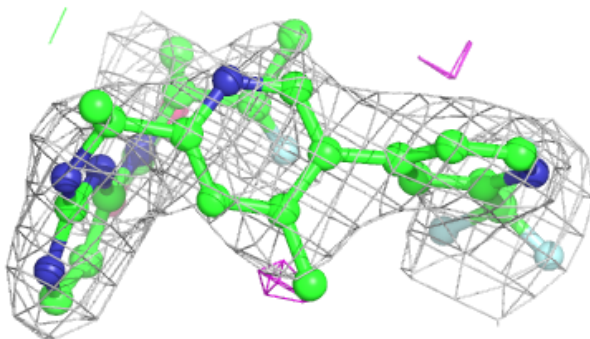
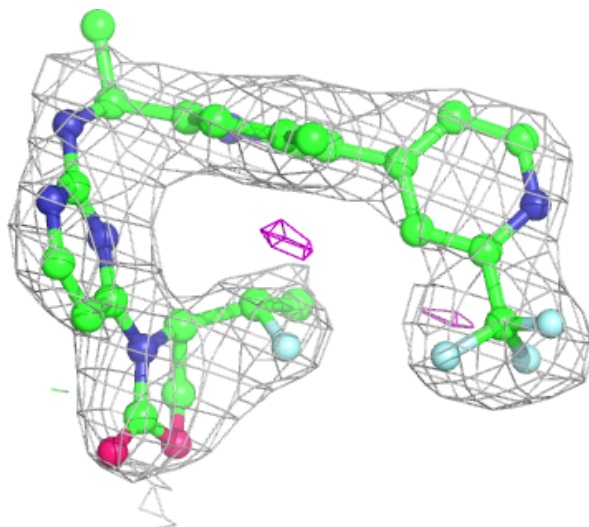
Electron density around NDP C 501:

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



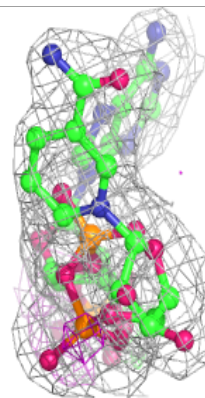
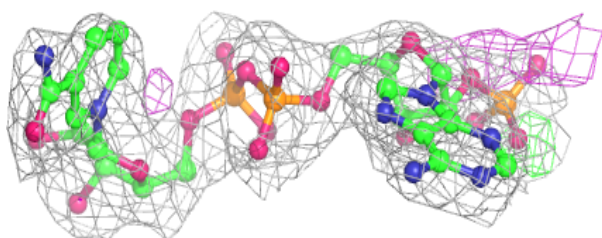
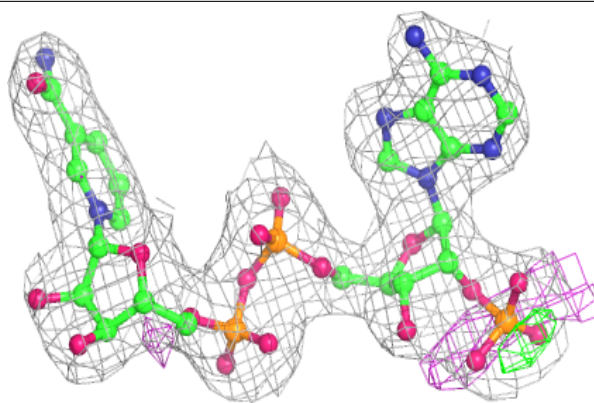
Electron density around C81 D 504:

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and green (positive)

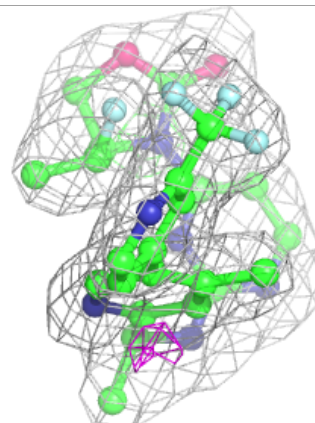
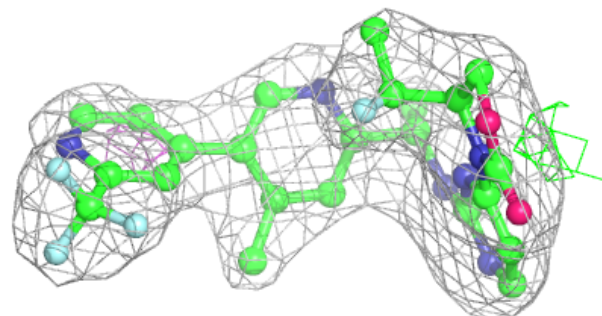
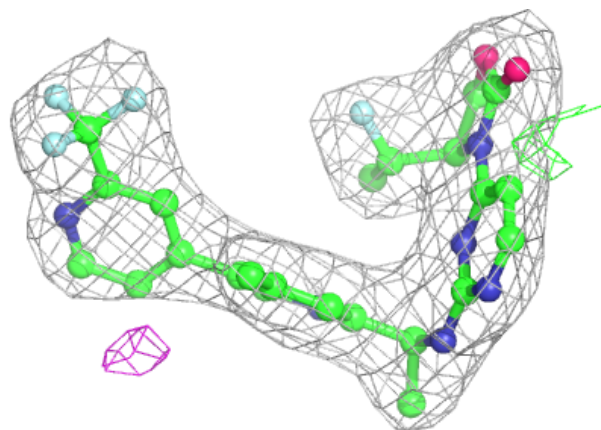


Electron density around NDP D 501:

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

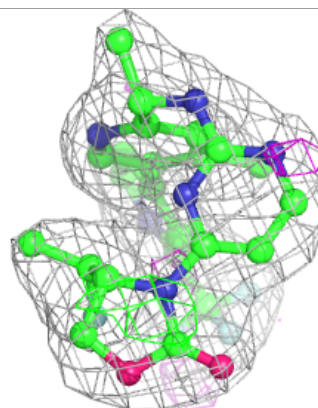
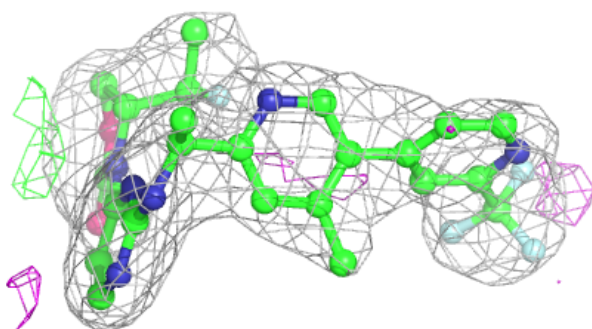
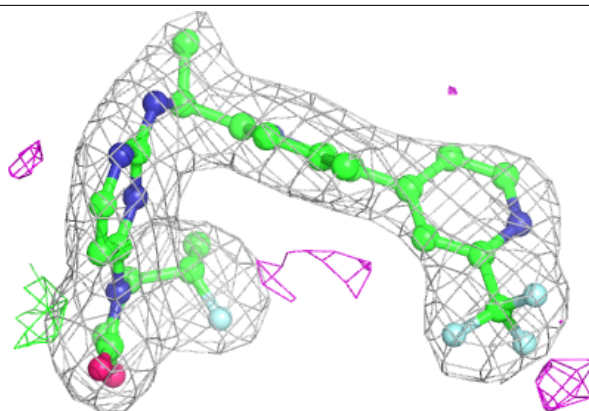
**Electron density around C81 C 502:**

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and green (positive)

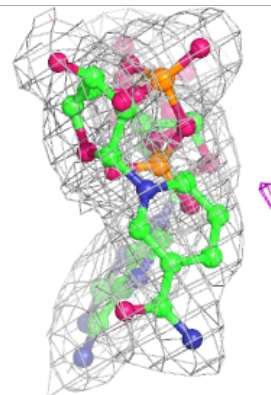
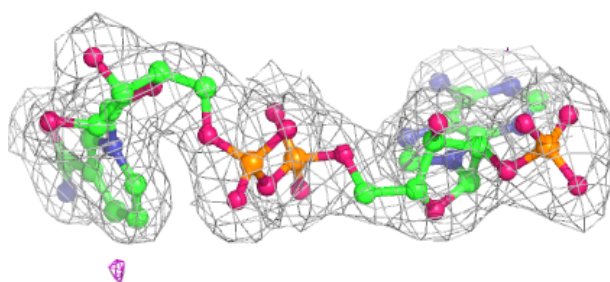
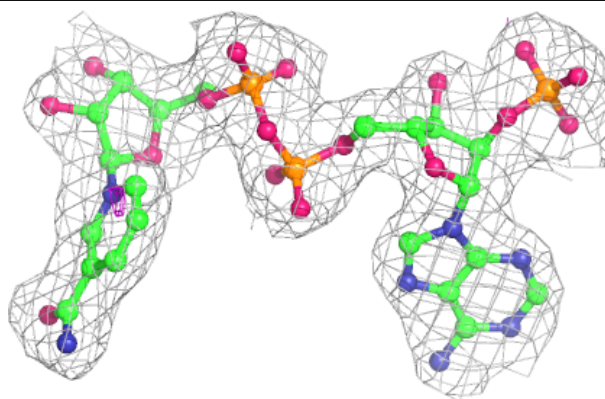


Electron density around C81 A 502:

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

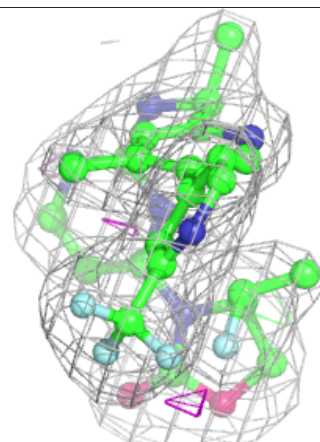
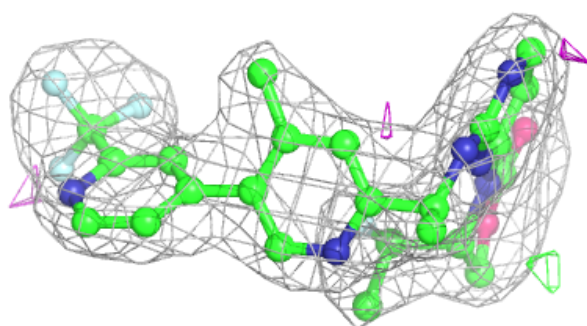
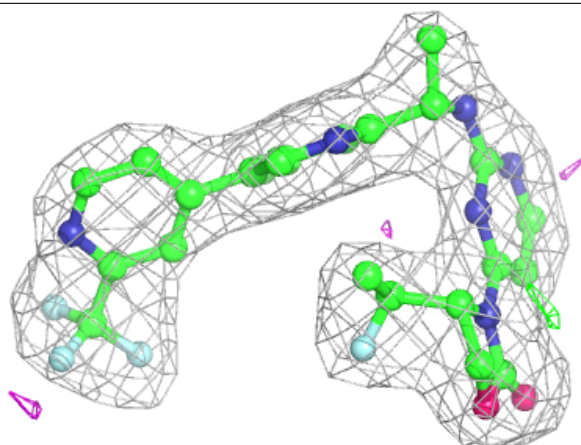
**Electron density around NDP A 501:**

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

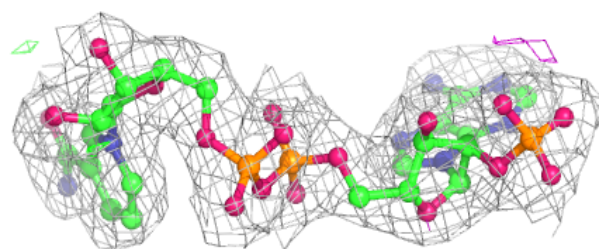
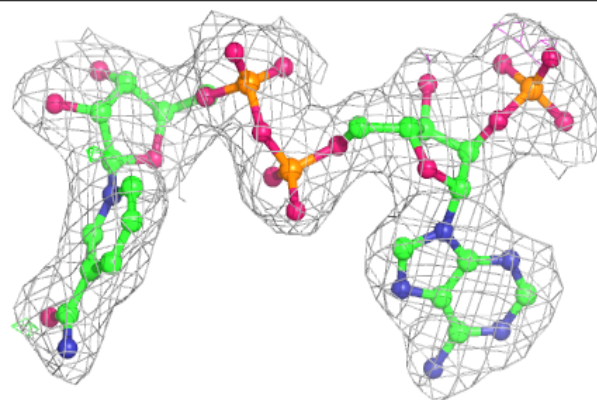


Electron density around C81 B 502:

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

**Electron density around NDP 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.