



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:03 PM UTC

PDB ID : 8WGM / pdb_00008wgm
Title : Quadruple mutant Plasmodium falciparum dihydrofolate reductase-thymidylate synthase (PfDHFR-TS V1/S, N51I+C59R+S108N+I164L) complexed with LA1, NADPH and dUMP
Authors : Vanichtanankul, J.; Saeyang, T.; Vitsupakorn, D.; Arwon, U.; Decharuangsilp, S.; Yuthavong, Y.; Kamchonwongpaisan, S.
Deposited on : 2023-09-22
Resolution : 2.15 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](#) ①) 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

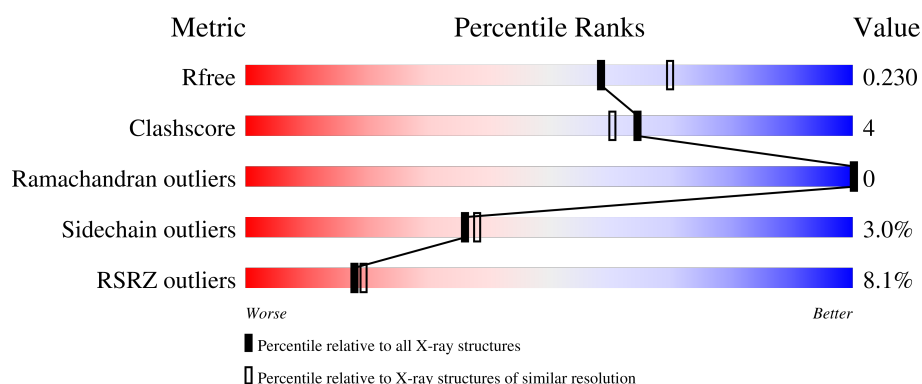
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	608	5% 79% 7% 12%
1	B	608	9% 75% 11% 13%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9780 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 Bifunctional dihydrofolate reductase-thymidylate synthase.

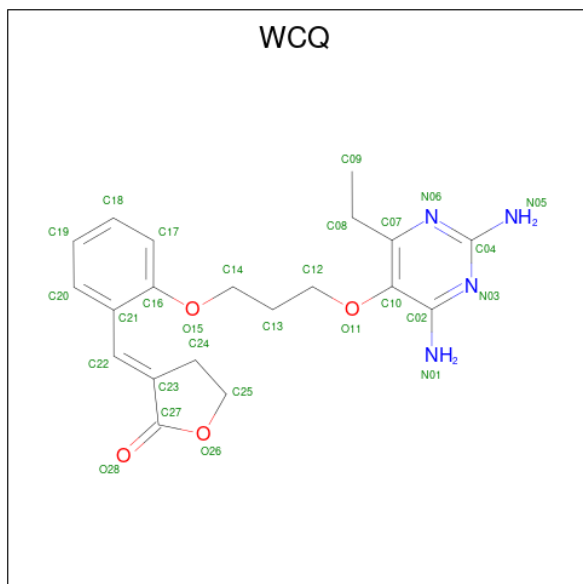
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	532	Total	C	N	O	S	0	0	0
			4422	2863	729	805	25			
1	B	528	Total	C	N	O	S	0	0	0
			4392	2845	725	798	24			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	26	ALA	GLY	conflict	UNP A7UD81
A	27	ALA	LYS	conflict	UNP A7UD81
A	28	ALA	LYS	conflict	UNP A7UD81
A	51	ILE	ASN	engineered mutation	UNP A7UD81
A	59	ARG	CYS	engineered mutation	UNP A7UD81
A	108	ASN	SER	engineered mutation	UNP A7UD81
A	164	LEU	ILE	engineered mutation	UNP A7UD81
B	26	ALA	GLY	conflict	UNP A7UD81
B	27	ALA	LYS	conflict	UNP A7UD81
B	28	ALA	LYS	conflict	UNP A7UD81
B	51	ILE	ASN	engineered mutation	UNP A7UD81
B	59	ARG	CYS	engineered mutation	UNP A7UD81
B	108	ASN	SER	engineered mutation	UNP A7UD81
B	164	LEU	ILE	engineered mutation	UNP A7UD81

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).

- Molecule 4 is (3 {E})-3-[[2-[3-[2,4-bis(azanyl)-6-ethyl-pyrimidin-5-yl]oxypropoxy]phenyl]methylidene]oxolan-2-one (CCD ID: WCQ) (formula: C₂₀H₂₄N₄O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			28	20	4	4		
4	B	1	Total	C	N	O	0	0
			28	20	4	4		

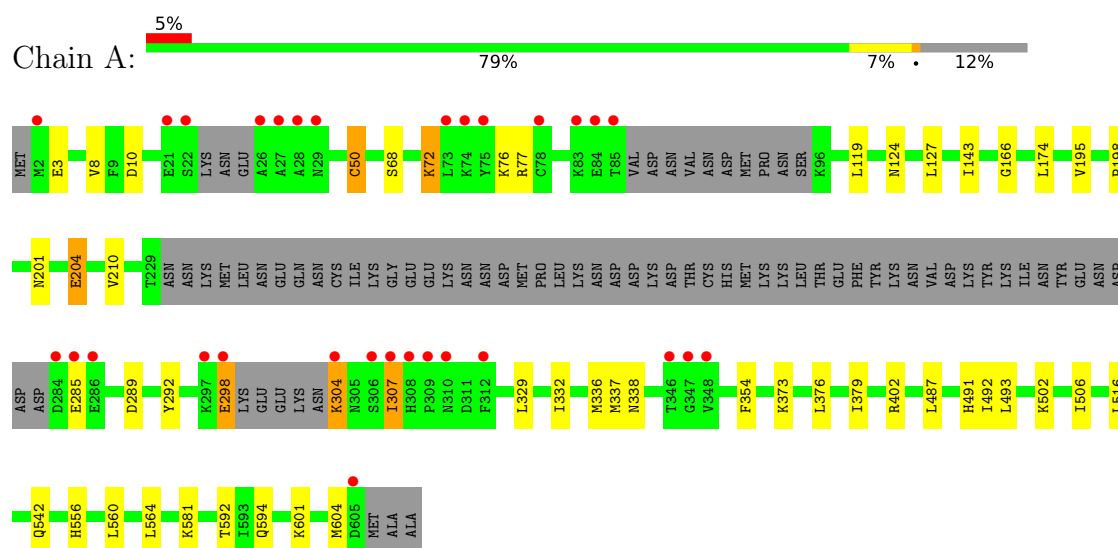
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	421	Total	O	0	0
			421	421		
5	B	353	Total	O	0	0
			353	353		

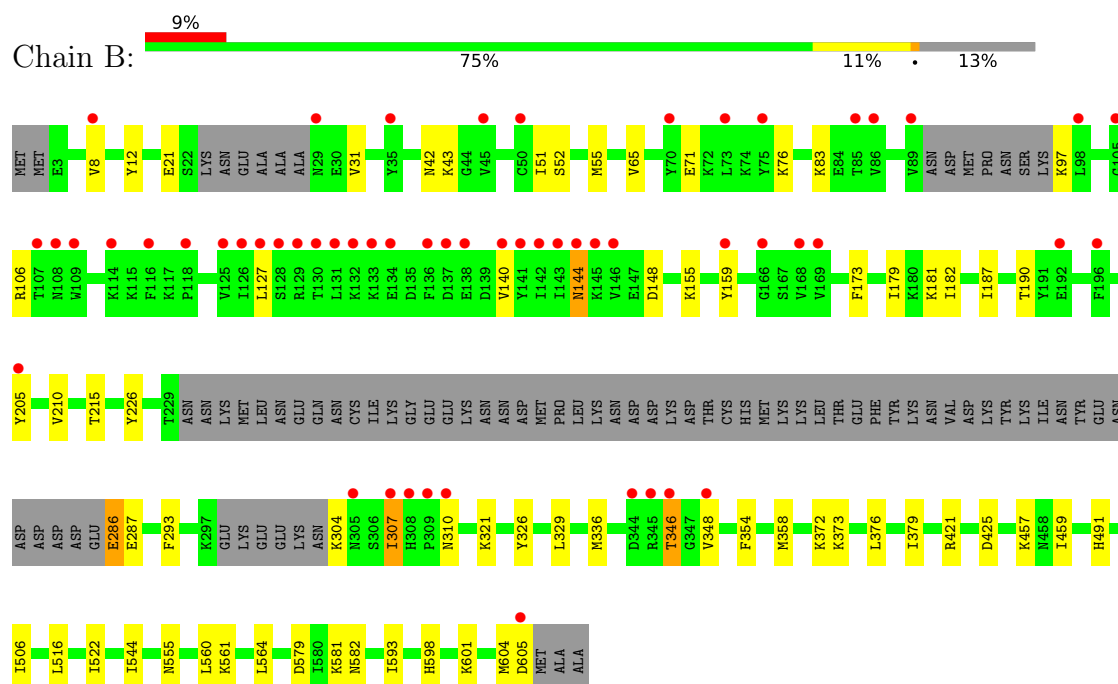
3 Residue-property plots [i](#)

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: Bifunctional dihydrofolate reductase-thymidylate synthase



- Molecule 1: Bifunctional dihydrofolate reductase-thymidylate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.35Å 156.28Å 165.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.32 – 2.15 24.32 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.2 (24.32-2.15) 89.1 (24.32-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.30 (at 2.15Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.190 , 0.233 0.194 , 0.230	Depositor DCC
R_{free} test set	1982 reflections (2.39%)	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.360	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 51.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9780	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UMP, NDP, WCQ

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	0.34	0/4525	0.53	0/6111
1	B	0.33	0/4495	0.51	0/6073
All	All	0.34	0/9020	0.52	0/12184

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	4422	0	4379	34	0
1	B	4392	0	4347	38	0
2	A	48	0	26	1	0
2	B	48	0	26	2	0
3	A	20	0	11	0	0
3	B	20	0	11	0	0
4	A	28	0	0	0	0
4	B	28	0	0	1	0
5	A	421	0	0	5	0
5	B	353	0	0	6	0
All	All	9780	0	8800	69	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 (69) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:LEU:H	1:A:124:ASN:HD21	1.27	0.81
1:B:8:VAL:HA	1:B:76:LYS:HD3	1.68	0.76
1:A:491:HIS:HD2	5:A:979:HOH:O	1.77	0.68
1:A:201:ASN:HB3	1:A:204:GLU:HB2	1.81	0.63
1:A:373:LYS:HB3	1:A:601:LYS:HB3	1.84	0.60
1:B:376:LEU:HD22	1:B:379:ILE:HD11	1.83	0.59
1:B:421:ARG:NE	5:B:809:HOH:O	2.31	0.59
1:B:51:ILE:HD13	1:B:187:ILE:HD12	1.85	0.58
1:A:542:GLN:OE1	5:A:801:HOH:O	2.17	0.58
1:B:346:THR:HB	1:B:348:VAL:HG23	1.83	0.58
1:B:286:GLU:CD	1:B:287:GLU:H	2.11	0.58
1:A:502:LYS:NZ	5:A:803:HOH:O	2.36	0.58
1:B:210:VAL:HB	5:B:873:HOH:O	2.04	0.56
1:A:592:THR:HG22	1:A:594:GLN:HG3	1.88	0.55
1:A:332:ILE:HD13	1:A:560:LEU:HD22	1.89	0.55
1:A:354:PHE:CE2	1:B:506:ILE:HG13	2.45	0.52
1:A:8:VAL:HA	1:A:76:LYS:HD3	1.90	0.52
1:A:556:HIS:CD2	1:A:604:MET:HB3	2.46	0.50
1:A:373:LYS:HD3	1:A:601:LYS:HE2	1.94	0.50
1:A:329:LEU:HD22	1:A:564:LEU:HD12	1.94	0.49
1:B:506:ILE:HG12	1:B:544:ILE:HB	1.93	0.49
1:B:55:MET:HE1	4:B:703:WCQ:C21	2.43	0.49
1:B:336:MET:HE1	1:B:560:LEU:HB2	1.95	0.49
1:B:372:LYS:NZ	1:B:604:MET:HE3	2.28	0.48
1:B:21:GLU:HG2	1:B:190:THR:HG22	1.95	0.48
1:A:10:ASP:OD2	1:A:77:ARG:NH2	2.47	0.48
1:B:491:HIS:HD2	5:B:933:HOH:O	1.96	0.47
1:B:12:TYR:CD1	1:B:181:LYS:HB2	2.50	0.47
1:B:179:ILE:HB	1:B:205:TYR:OH	2.13	0.47
1:B:321:LYS:HD2	1:B:326:TYR:CE1	2.49	0.46
1:A:50:CYS:SG	5:A:1180:HOH:O	2.45	0.46
1:B:582:ASN:HB3	5:B:852:HOH:O	2.14	0.46
1:A:72:LYS:O	1:A:72:LYS:HD3	2.16	0.46
1:A:307:ILE:HB	1:A:337:MET:HE1	1.98	0.46
1:A:304:LYS:HG3	1:A:337:MET:O	2.16	0.45
1:B:329:LEU:HD22	1:B:564:LEU:HD12	1.97	0.45
1:A:119:LEU:H	1:A:124:ASN:ND2	2.04	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:307:ILE:HG23	1:B:561:LYS:HE2	1.98	0.44
1:B:555:ASN:ND2	1:B:605:ASP:O	2.50	0.44
1:B:376:LEU:HD12	1:B:593:ILE:HG13	1.99	0.44
1:B:293:PHE:O	5:B:801:HOH:O	2.21	0.44
1:B:579:ASP:O	1:B:581:LYS:HG2	2.17	0.43
1:B:358:MET:HE1	1:B:522:ILE:HD11	2.01	0.43
1:A:210:VAL:HB	5:A:980:HOH:O	2.18	0.43
1:A:166:GLY:HA3	2:A:701:NDP:O5D	2.18	0.43
1:B:173:PHE:HB3	1:B:179:ILE:HD11	1.99	0.43
1:B:373:LYS:HG3	1:B:598:HIS:CE1	2.54	0.43
1:A:127:LEU:HD23	1:A:143:ILE:HG13	2.01	0.43
1:A:492:ILE:HG22	1:A:493:LEU:HD23	2.00	0.43
1:B:106:ARG:NH2	2:B:701:NDP:O1X	2.49	0.43
1:A:289:ASP:OD2	1:B:181:LYS:NZ	2.46	0.42
1:B:182:ILE:HB	1:B:226:TYR:HB2	2.02	0.42
1:A:304:LYS:HE3	1:A:304:LYS:HB2	1.75	0.42
1:B:287:GLU:HB3	5:B:817:HOH:O	2.19	0.42
1:A:298:GLU:H	1:A:298:GLU:HG3	1.62	0.42
1:B:42:ASN:HD21	1:B:43:LYS:HE2	1.85	0.42
1:A:289:ASP:HA	1:A:292:TYR:CD2	2.55	0.41
1:B:144:ASN:ND2	1:B:148:ASP:OD2	2.49	0.41
1:B:65:VAL:HG22	1:B:159:TYR:CB	2.50	0.41
1:A:581:LYS:HA	1:A:581:LYS:HD3	1.80	0.41
1:B:83:LYS:HE3	1:B:83:LYS:HB2	1.81	0.41
1:B:106:ARG:HE	2:B:701:NDP:P2B	2.44	0.41
1:A:304:LYS:HD3	1:A:338:ASN:HA	2.03	0.41
1:A:376:LEU:HD22	1:A:379:ILE:HD11	2.02	0.41
1:A:506:ILE:HG13	1:B:354:PHE:CE2	2.56	0.40
1:A:516:LEU:HA	1:A:516:LEU:HD23	1.82	0.40
1:B:421:ARG:HD2	1:B:425:ASP:O	2.21	0.40
1:A:174:LEU:HD12	1:A:198:PRO:HG2	2.04	0.40
1:A:307:ILE:HD12	1:A:336:MET:HB3	2.02	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	522/608 (86%)	509 (98%)	13 (2%)	0	100	100
1	B	518/608 (85%)	502 (97%)	16 (3%)	0	100	100
All	All	1040/1216 (86%)	1011 (97%)	29 (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	495/568 (87%)	483 (98%)	12 (2%)	43	47
1	B	493/568 (87%)	475 (96%)	18 (4%)	30	30
All	All	988/1136 (87%)	958 (97%)	30 (3%)	36	38

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	50	CYS
1	A	68	SER
1	A	72	LYS
1	A	195	VAL
1	A	204	GLU
1	A	285	GLU
1	A	298	GLU
1	A	304	LYS
1	A	307	ILE
1	A	402	ARG
1	A	487	LEU
1	B	31	VAL
1	B	52	SER

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Mol	Chain	Res	Type
1	B	71	GLU
1	B	97	LYS
1	B	127	LEU
1	B	140	VAL
1	B	144	ASN
1	B	155	LYS
1	B	215	THR
1	B	286	GLU
1	B	304	LYS
1	B	307	ILE
1	B	310	ASN
1	B	346	THR
1	B	457	LYS
1	B	459	ILE
1	B	516	LEU
1	B	601	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	206	GLN
1	A	294	ASN
1	A	407	ASN
1	A	415	ASN
1	A	491	HIS
1	A	534	GLN
1	A	539	GLN
1	A	542	GLN
1	A	554	ASN
1	B	42	ASN
1	B	121	ASN
1	B	206	GLN
1	B	294	ASN
1	B	340	ASN
1	B	407	ASN
1	B	450	ASN
1	B	491	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

6 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	B	701	-	51,52,52	2.54	10 (19%)	71,80,80	1.51	13 (18%)
3	UMP	A	702	-	21,21,21	0.88	0	30,31,31	1.66	7 (23%)
4	WCQ	B	703	-	30,30,30	2.33	6 (20%)	37,40,40	2.09	10 (27%)
3	UMP	B	702	-	21,21,21	0.92	0	30,31,31	1.46	6 (20%)
4	WCQ	A	703	-	30,30,30	2.27	7 (23%)	37,40,40	2.40	12 (32%)
2	NDP	A	701	-	51,52,52	2.50	8 (15%)	71,80,80	1.36	9 (12%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NDP	B	701	-	-	2/34/77/77	0/5/5/5
3	UMP	A	702	-	-	1/10/22/22	0/2/2/2
4	WCQ	B	703	-	-	0/14/24/24	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	UMP	B	702	-	-	2/10/22/22	0/2/2/2
4	WCQ	A	703	-	-	0/14/24/24	0/3/3/3
2	NDP	A	701	-	-	5/34/77/77	0/5/5/5

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	701	NDP	P2B-O2B	12.56	1.81	1.59
2	B	701	NDP	P2B-O2B	11.97	1.80	1.59
2	B	701	NDP	PA-O3	9.18	1.69	1.59
2	A	701	NDP	PA-O3	7.89	1.68	1.59
4	B	703	WCQ	O26-C27	6.58	1.46	1.36
4	A	703	WCQ	O26-C27	6.29	1.46	1.36
4	A	703	WCQ	C04-N05	5.63	1.45	1.33
4	B	703	WCQ	C04-N05	5.43	1.44	1.33
2	A	701	NDP	PN-O5D	4.85	1.78	1.59
4	A	703	WCQ	C24-C23	-4.65	1.41	1.50
2	B	701	NDP	PN-O5D	4.60	1.77	1.59
4	B	703	WCQ	C24-C23	-4.60	1.42	1.50
4	B	703	WCQ	C21-C22	4.52	1.54	1.46
4	B	703	WCQ	C02-N01	4.41	1.45	1.34
4	A	703	WCQ	C21-C22	4.03	1.53	1.46
4	A	703	WCQ	C02-N01	3.85	1.44	1.34
2	B	701	NDP	O2B-C2B	-3.28	1.32	1.44
2	A	701	NDP	C7N-N7N	3.10	1.42	1.33
2	A	701	NDP	O2B-C2B	-2.98	1.33	1.44
2	B	701	NDP	PN-O3	2.87	1.62	1.59
2	B	701	NDP	C2A-N1A	2.56	1.38	1.33
2	B	701	NDP	C5A-C4A	2.40	1.43	1.39
4	B	703	WCQ	C27-C23	2.39	1.52	1.47
2	B	701	NDP	C4A-N3A	2.34	1.38	1.34
4	A	703	WCQ	C22-C23	-2.34	1.32	1.34
2	A	701	NDP	C5A-C4A	2.31	1.43	1.39
2	B	701	NDP	C2B-C1B	2.29	1.58	1.53
2	B	701	NDP	C7N-N7N	2.21	1.39	1.33
2	A	701	NDP	C2A-N1A	2.21	1.37	1.33
2	A	701	NDP	C8A-N9A	2.18	1.41	1.37
4	A	703	WCQ	C21-C16	-2.14	1.37	1.41

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	703	WCQ	C04-N06-C07	9.02	122.84	116.26
4	B	703	WCQ	C04-N06-C07	6.57	121.06	116.26
4	A	703	WCQ	C25-O26-C27	-5.45	106.11	110.74
2	B	701	NDP	P2B-O2B-C2B	-4.43	111.60	123.43
3	A	702	UMP	C4-N3-C2	-4.22	121.37	126.61
4	A	703	WCQ	C10-C07-N06	-4.13	117.35	122.46
4	A	703	WCQ	O26-C27-O28	3.85	125.83	121.09
3	B	702	UMP	C4-N3-C2	-3.76	121.94	126.61
2	B	701	NDP	O2B-P2B-O1X	-3.69	96.19	109.33
3	A	702	UMP	C5-C4-N3	3.66	119.92	114.80
2	B	701	NDP	O2N-PN-O3	3.61	117.03	107.27
4	B	703	WCQ	C25-O26-C27	-3.57	107.71	110.74
3	B	702	UMP	C5-C4-N3	3.50	119.70	114.80
4	B	703	WCQ	C10-C07-N06	-3.43	118.21	122.46
4	B	703	WCQ	O26-C27-O28	3.41	125.29	121.09
4	B	703	WCQ	O15-C16-C21	3.38	119.16	115.78
2	A	701	NDP	O2N-PN-O3	3.31	116.22	107.27
2	A	701	NDP	O2B-P2B-O1X	-3.11	98.26	109.33
3	A	702	UMP	C1'-N1-C2	3.07	123.66	117.66
4	A	703	WCQ	C16-C21-C22	-2.97	114.77	118.73
4	B	703	WCQ	N01-C02-N03	2.96	121.02	117.03
3	A	702	UMP	N3-C2-N1	2.93	118.71	114.89
3	B	702	UMP	O4-C4-C5	-2.93	120.10	125.16
3	A	702	UMP	C1'-N1-C6	-2.92	115.79	121.53
2	A	701	NDP	C3B-C2B-C1B	-2.82	97.40	102.81
2	B	701	NDP	PA-O5B-C5B	-2.81	105.25	121.35
2	B	701	NDP	PN-O5D-C5D	-2.73	105.71	121.35
2	B	701	NDP	O3X-P2B-O2X	2.72	118.02	107.80
4	A	703	WCQ	N01-C02-N03	2.72	120.70	117.03
3	B	702	UMP	C1'-N1-C6	-2.68	116.25	121.53
2	A	701	NDP	O3X-P2B-O2X	2.68	117.85	107.80
3	B	702	UMP	C1'-N1-C2	2.62	122.78	117.66
4	A	703	WCQ	O28-C27-C23	-2.56	125.87	129.65
4	A	703	WCQ	N06-C04-N03	-2.53	121.60	125.48
3	A	702	UMP	O2-C2-N3	-2.49	116.89	121.49
4	A	703	WCQ	O26-C25-C24	-2.49	105.02	107.16
3	A	702	UMP	O4-C4-C5	-2.41	121.01	125.16
2	A	701	NDP	PA-O5B-C5B	-2.40	107.62	121.35
2	B	701	NDP	C2A-N1A-C6A	-2.40	114.80	118.73
2	A	701	NDP	C2A-N1A-C6A	-2.37	114.83	118.73
2	B	701	NDP	O5D-PN-O1N	-2.37	99.53	108.94
4	B	703	WCQ	C25-C24-C23	2.36	107.77	103.33
2	B	701	NDP	O2A-PA-O1A	2.34	123.33	112.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	703	WCQ	C14-O15-C16	-2.33	112.07	117.69
4	B	703	WCQ	C04-N03-C02	2.29	120.03	117.28
2	B	701	NDP	N3A-C4A-N9A	2.21	130.93	127.17
4	B	703	WCQ	O28-C27-C23	-2.20	126.39	129.65
2	A	701	NDP	C2B-C3B-C4B	-2.19	97.29	101.99
4	A	703	WCQ	N05-C04-N03	2.13	120.42	117.22
4	A	703	WCQ	O15-C16-C21	2.09	117.87	115.78
2	B	701	NDP	O3X-P2B-O2B	-2.08	97.75	105.85
2	B	701	NDP	C5A-C4A-N3A	-2.07	123.86	126.72
2	B	701	NDP	O3-PA-O1A	-2.06	104.50	110.70
4	A	703	WCQ	C08-C07-N06	2.06	121.03	116.61
3	B	702	UMP	N3-C2-N1	2.04	117.55	114.89
2	A	701	NDP	PN-O5D-C5D	-2.03	109.73	121.35
2	A	701	NDP	O2N-PN-O1N	2.02	121.86	112.44

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	702	UMP	C3'-C4'-C5'-O5'
3	B	702	UMP	O4'-C4'-C5'-O5'
2	A	701	NDP	PA-O3-PN-O5D
2	A	701	NDP	C5B-O5B-PA-O1A
2	A	701	NDP	O4D-C1D-N1N-C2N
2	B	701	NDP	O4D-C1D-N1N-C2N
2	B	701	NDP	C2D-C1D-N1N-C2N
2	A	701	NDP	C2B-O2B-P2B-O1X
2	A	701	NDP	C2D-C1D-N1N-C2N
3	A	702	UMP	O4'-C4'-C5'-O5'

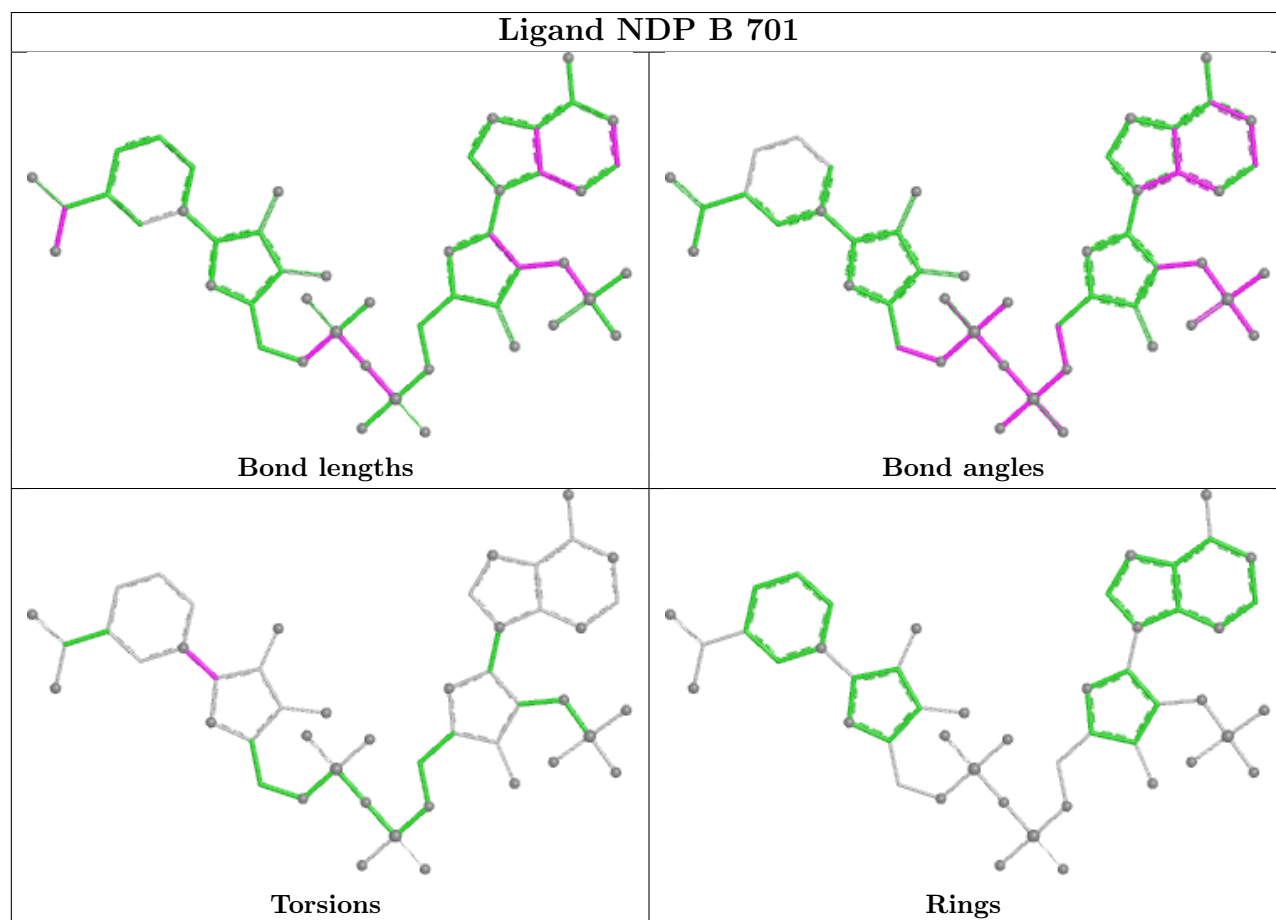
There are no ring outliers.

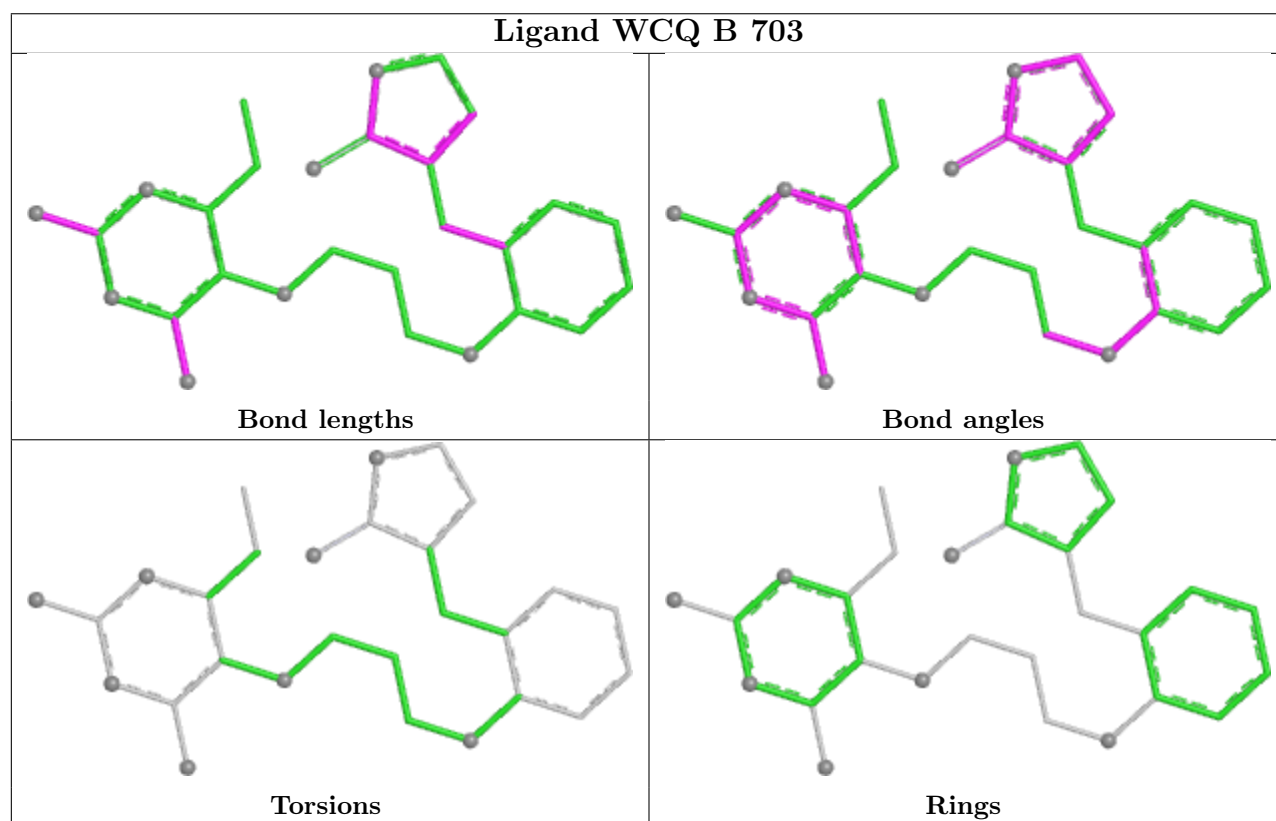
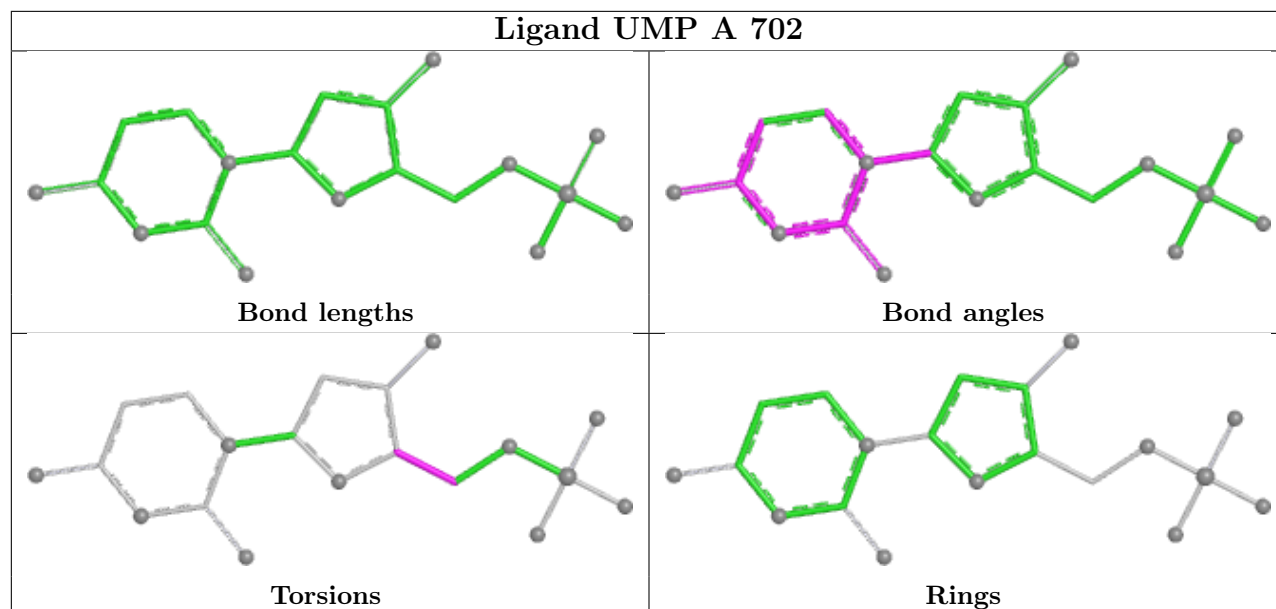
3 monomers are involved in 4 short contacts:

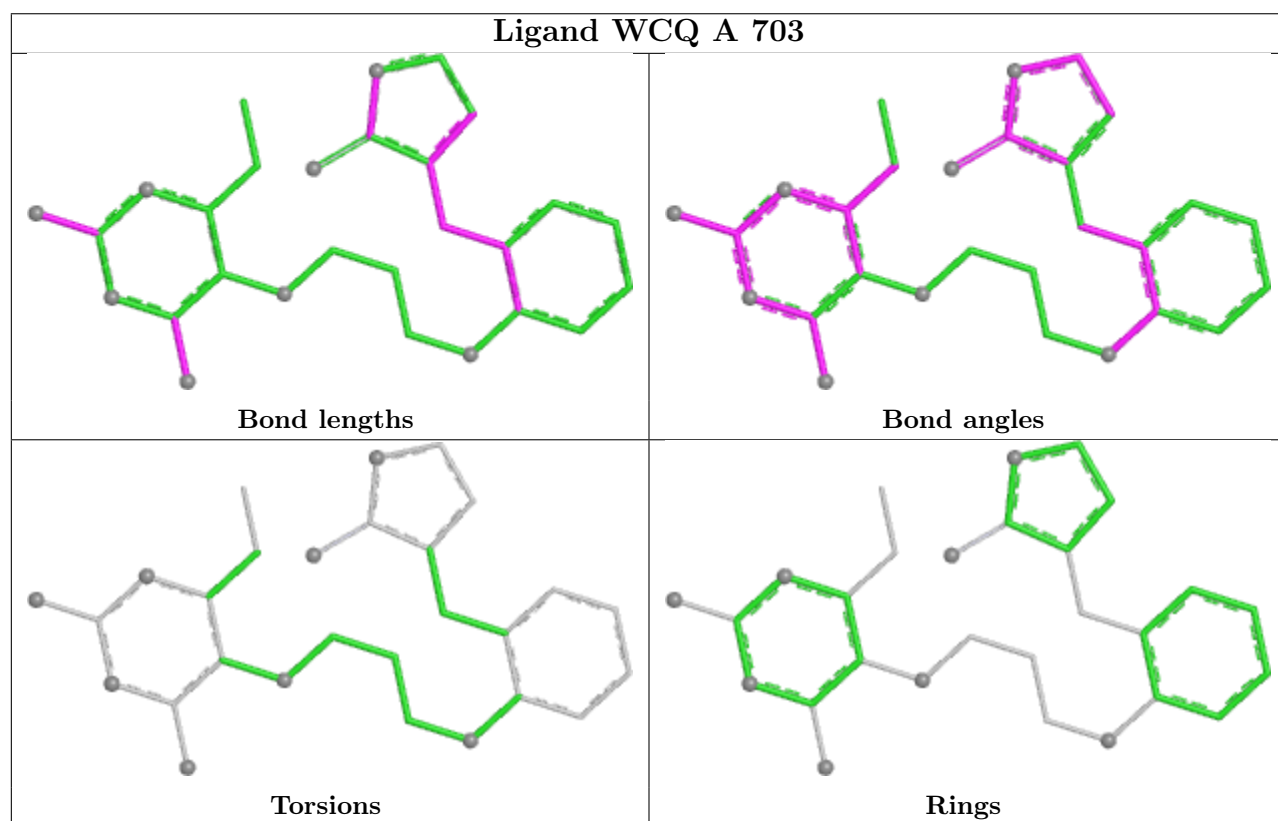
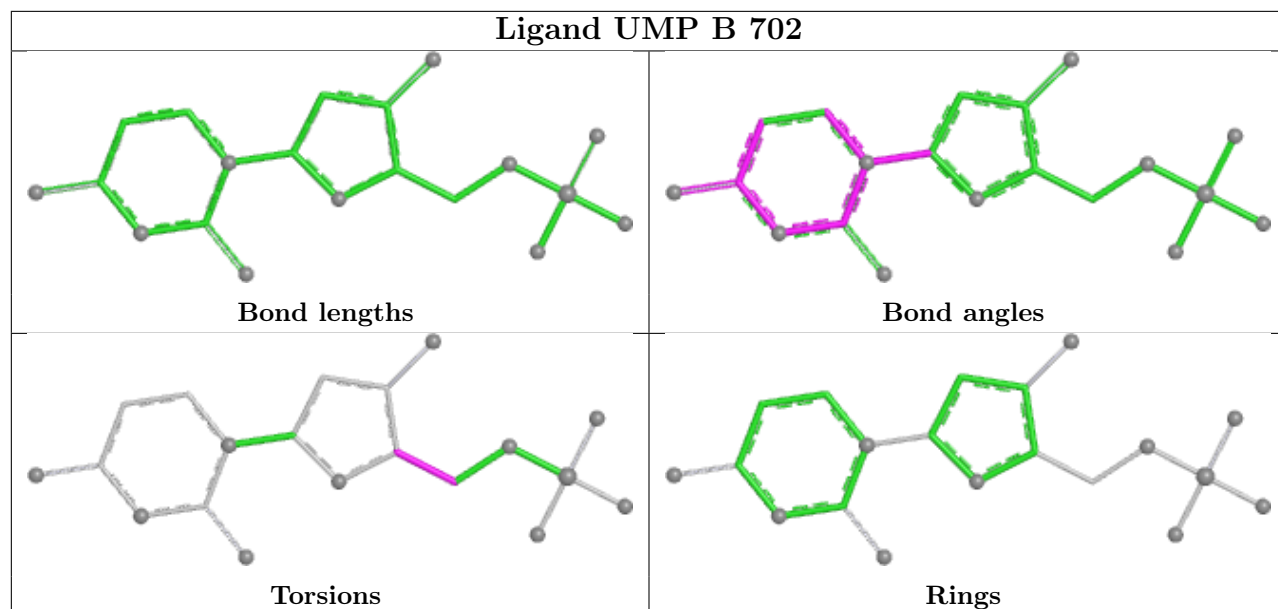
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	NDP	2	0
4	B	703	WCQ	1	0
2	A	701	NDP	1	0

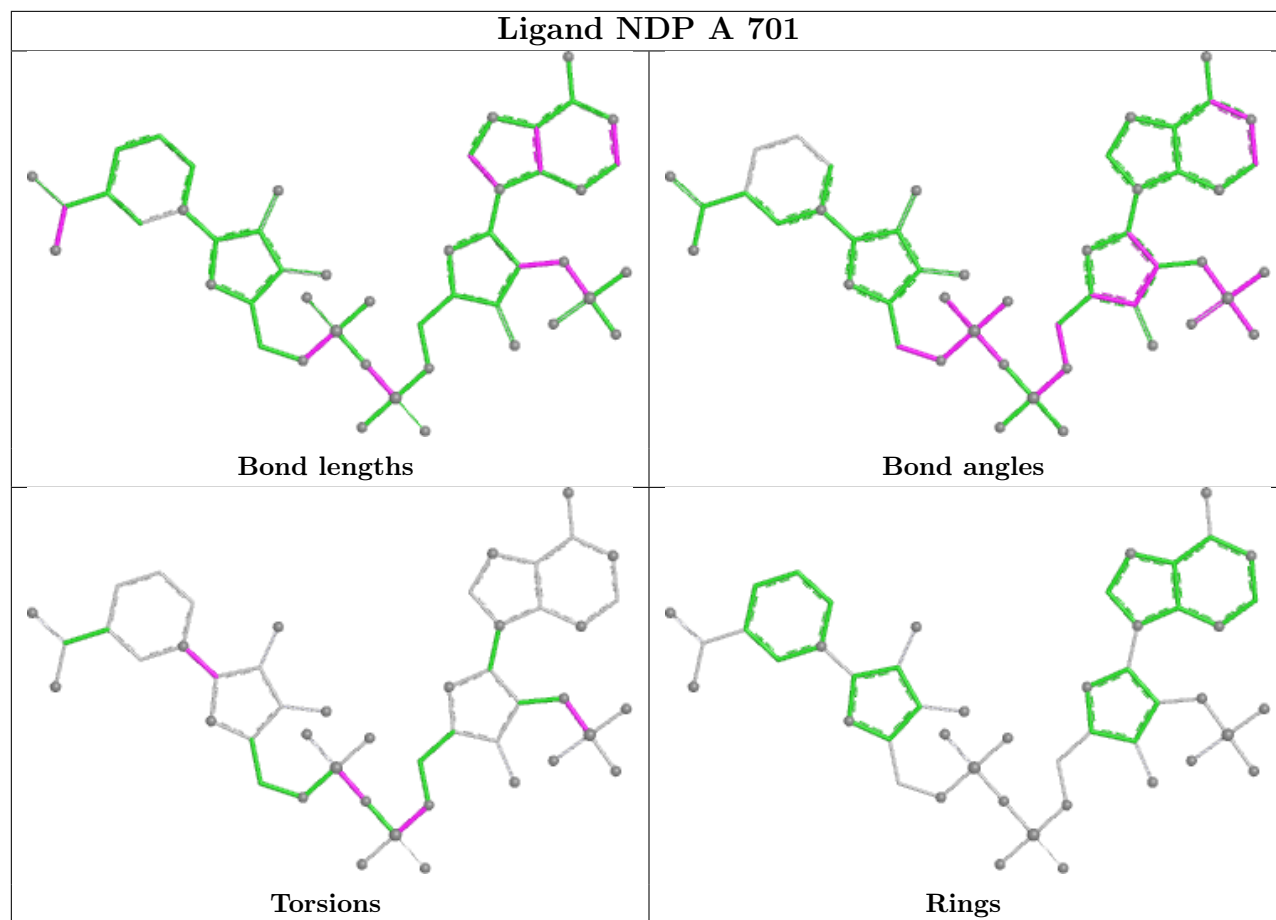
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

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	532/608 (87%)	0.05	30 (5%) 30 34	14, 27, 62, 87	0
1	B	528/608 (86%)	0.39	56 (10%) 11 12	14, 32, 70, 96	0
All	All	1060/1216 (87%)	0.22	86 (8%) 18 19	14, 29, 67, 96	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	307	ILE	4.4
1	B	308	HIS	4.3
1	B	131	LEU	4.1
1	A	85	THR	4.1
1	B	142	ILE	4.1
1	B	136	PHE	4.1
1	B	168	VAL	4.0
1	B	89	VAL	4.0
1	B	75	TYR	3.8
1	A	75	TYR	3.6
1	B	309	PRO	3.6
1	A	26	ALA	3.6
1	B	144	ASN	3.6
1	B	310	ASN	3.5
1	A	2	MET	3.5
1	B	86	VAL	3.3
1	A	307	ILE	3.3
1	B	118	PRO	3.3
1	B	605	ASP	3.3
1	B	126	ILE	3.2
1	A	27	ALA	3.2
1	B	143	ILE	3.2
1	B	130	THR	3.2
1	A	284	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	605	ASP	3.1
1	B	140	VAL	3.1
1	A	285	GLU	3.0
1	A	28	ALA	3.0
1	A	308	HIS	3.0
1	A	309	PRO	2.9
1	A	286	GLU	2.9
1	B	85	THR	2.9
1	B	205	TYR	2.9
1	B	132	LYS	2.8
1	B	107	THR	2.8
1	B	45	VAL	2.8
1	B	345	ARG	2.7
1	B	108	ASN	2.7
1	B	192	GLU	2.7
1	B	73	LEU	2.7
1	B	109	TRP	2.6
1	B	138	GLU	2.5
1	B	146	VAL	2.5
1	B	105	GLY	2.5
1	B	169	VAL	2.5
1	B	346	THR	2.5
1	B	29	ASN	2.5
1	B	141	TYR	2.5
1	A	306	SER	2.5
1	B	98	LEU	2.5
1	A	29	ASN	2.4
1	B	134	GLU	2.4
1	B	129	ARG	2.4
1	B	348	VAL	2.4
1	B	114	LYS	2.4
1	B	133	LYS	2.4
1	B	137	ASP	2.4
1	A	312	PHE	2.4
1	B	196	PHE	2.4
1	B	35	TYR	2.3
1	A	74	LYS	2.3
1	B	125	VAL	2.3
1	A	84	GLU	2.3
1	A	346	THR	2.3
1	B	166	GLY	2.3
1	B	127	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	22	SER	2.2
1	B	8	VAL	2.2
1	A	21	GLU	2.2
1	B	128	SER	2.2
1	A	304	LYS	2.2
1	A	310	ASN	2.2
1	A	348	VAL	2.1
1	B	116	PHE	2.1
1	A	83	LYS	2.1
1	A	297	LYS	2.1
1	A	298	GLU	2.1
1	B	70	TYR	2.1
1	B	305	ASN	2.0
1	B	344	ASP	2.0
1	A	347	GLY	2.0
1	B	159	TYR	2.0
1	A	73	LEU	2.0
1	A	78	CYS	2.0
1	B	50	CYS	2.0
1	B	145	LYS	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 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.

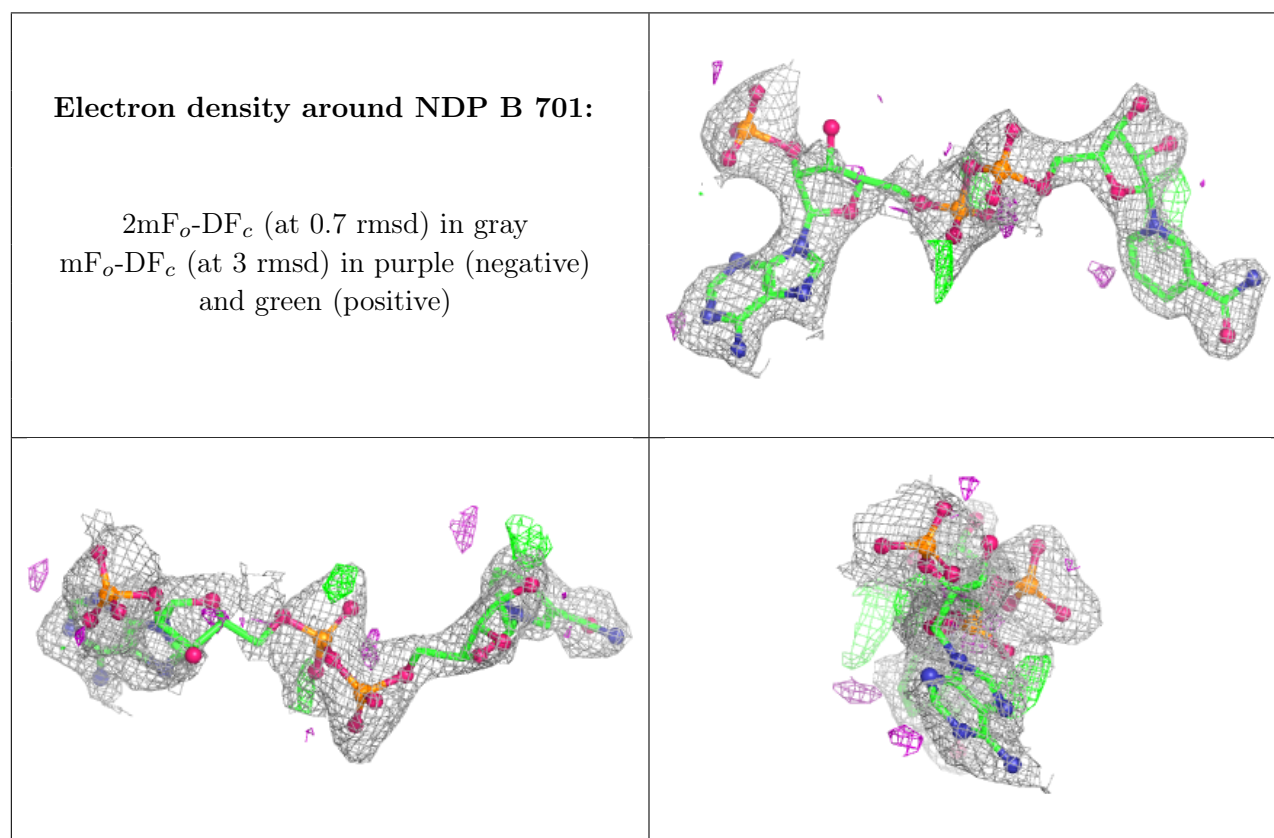
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NDP	B	701	48/48	0.83	0.14	39,63,76,78	0
4	WCQ	B	703	28/28	0.88	0.12	38,48,58,60	0
3	UMP	B	702	20/20	0.90	0.12	26,38,45,45	0

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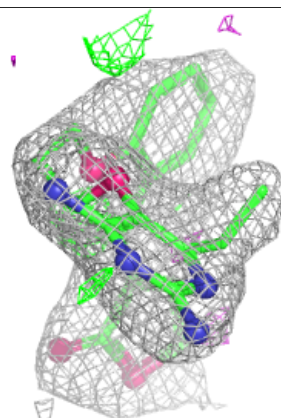
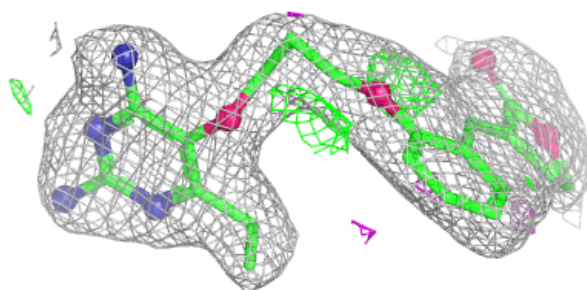
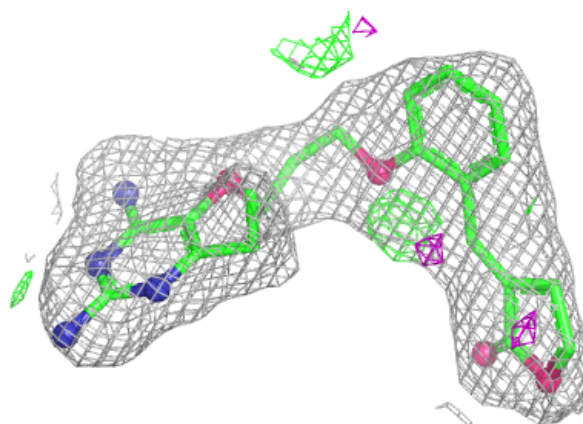
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	UMP	A	702	20/20	0.94	0.09	27,33,38,38	0
4	WCQ	A	703	28/28	0.95	0.07	17,24,35,36	0
2	NDP	A	701	48/48	0.96	0.08	22,32,39,41	0

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.

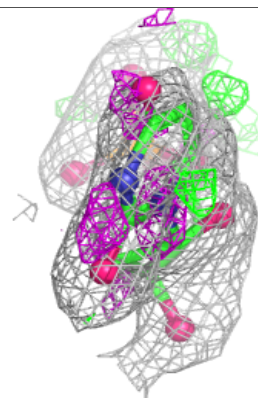
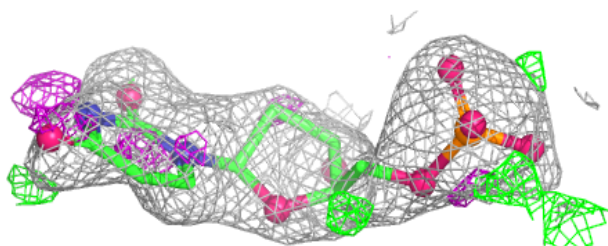
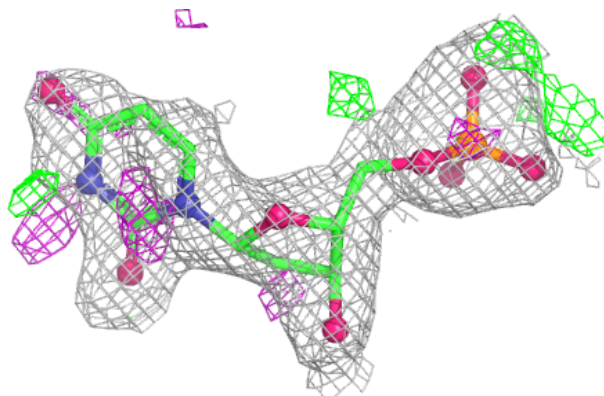


Electron density around WCQ B 703:

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

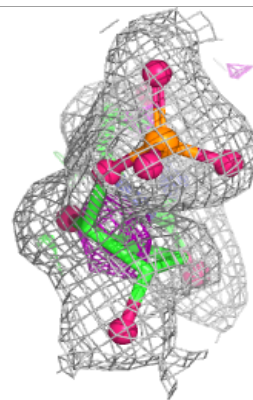
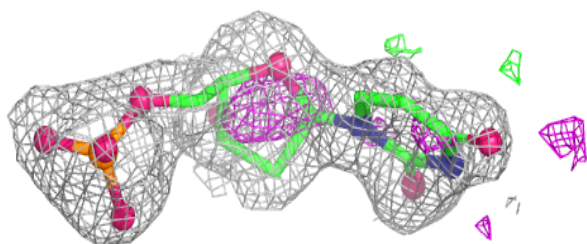
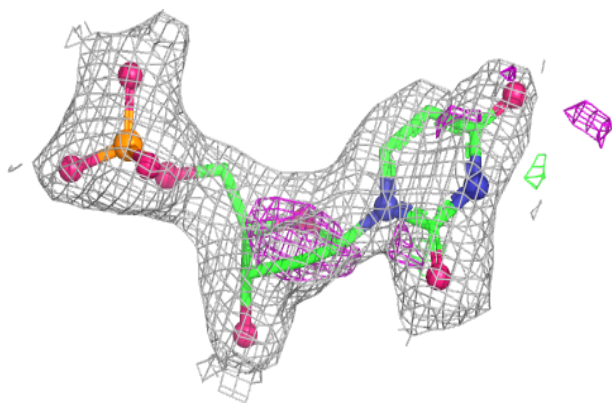
**Electron density around UMP B 702:**

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

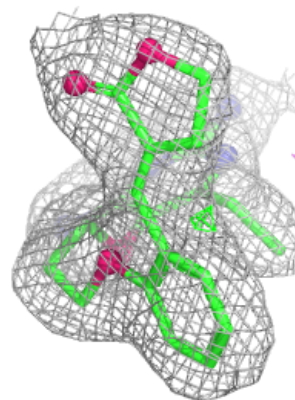
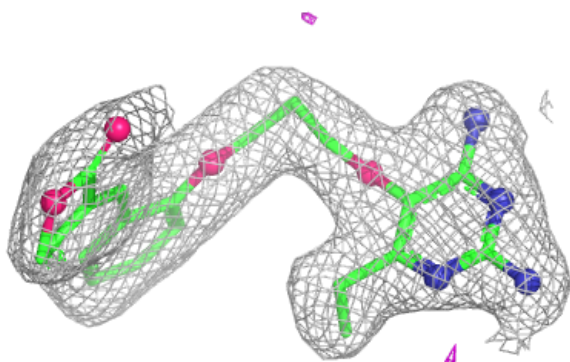
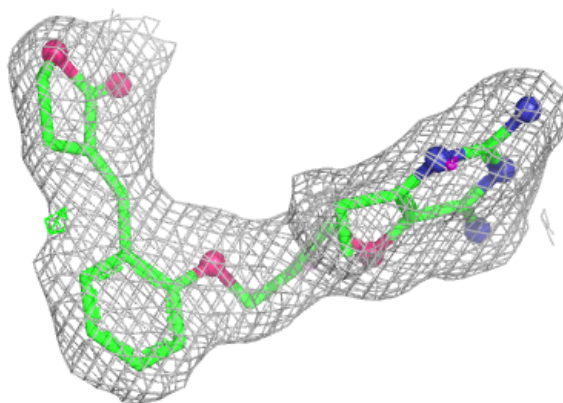


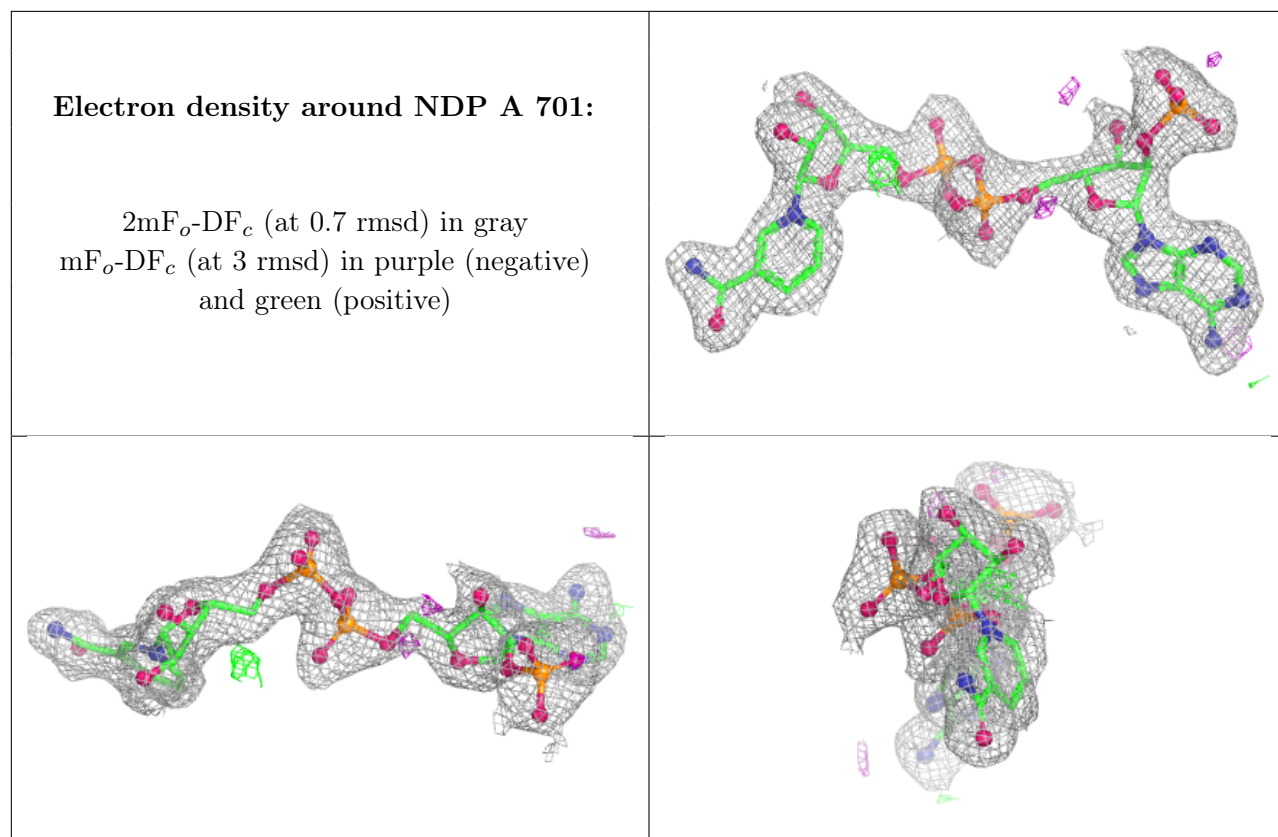
Electron density around UMP A 702:

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

**Electron density around WCQ A 703:**

$2mF_o-DF_c$ (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.