



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:12 AM UTC

PDB ID : 6LHI / pdb_00006lhi
Title : Quadruple mutant (N51I+C59R+S108N+I164L) plasmodium falciparum dihydrofolate reductase-thymidylate synthase (PfDHFR-TS) complexed with C466 (compound 42) and NADPH
Authors : Vanichtanankul, J.; Vitsupakorn, D.
Deposited on : 2019-12-09
Resolution : 2.59 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

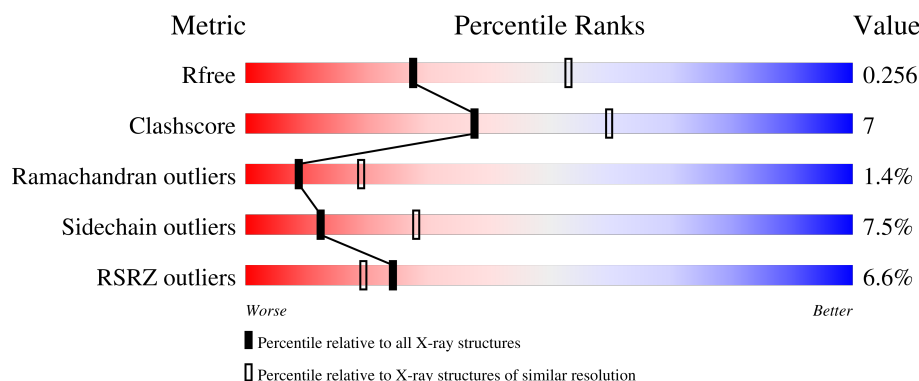
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	2% 69% 15% • 12%
1	B	608	9% 64% 19% • 13%

2 Entry composition [i](#)

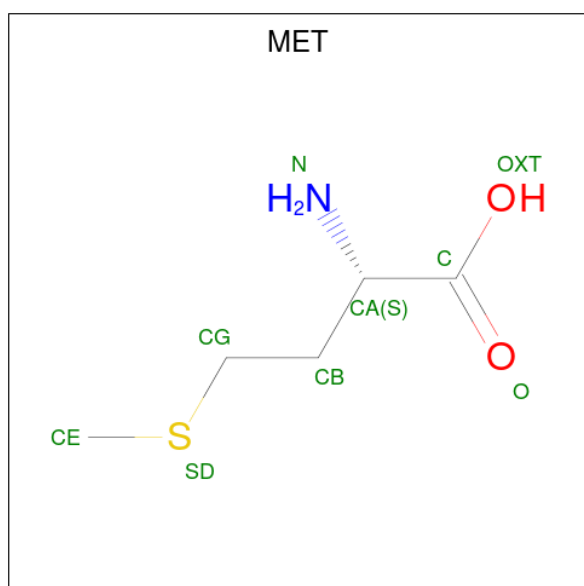
There are 6 unique types of molecules in this entry. The entry contains 18078 atoms, of which 8923 are hydrogens and 0 are deuteriums.

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.

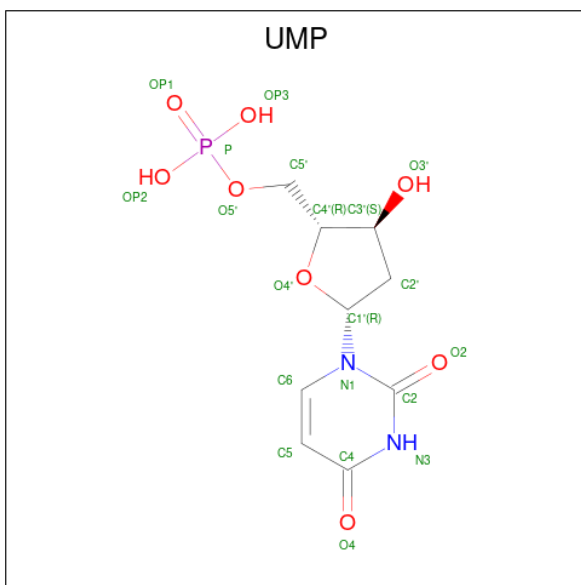
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	532	Total	C	H	N	O	S	256	0	0
			8832	2867	4399	733	809	24			
1	B	531	Total	C	H	N	O	S	254	0	0
			8813	2862	4389	730	807	25			

- Molecule 2 is METHIONINE (CCD ID: MET) (formula: $C_5H_{11}NO_2S$).



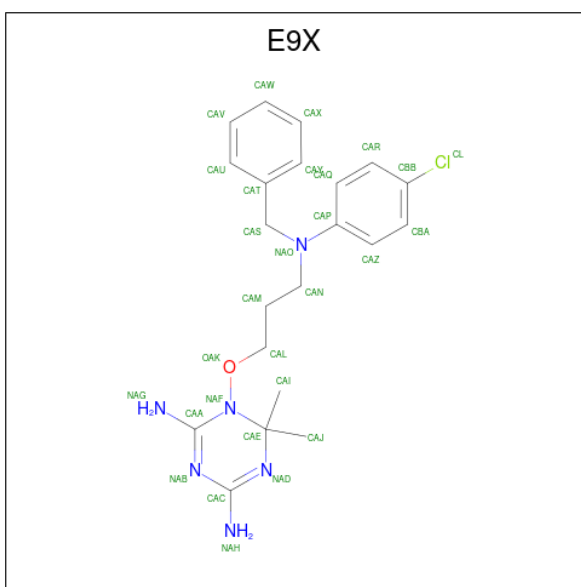
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	S	0	0
			19	5	11	1	1	1		

- Molecule 3 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (CCD ID: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 31	C 9	H 11	N 2	O 8	P 1	1	0
3	B	1	Total 31	C 9	H 11	N 2	O 8	P 1	1	0

- Molecule 4 is 1-[3-[(4-chlorophenyl)-(phenylmethyl)amino]propoxy]-6,6-dimethyl-1,3,5-triazine-2,4-diamine (CCD ID: E9X) (formula: $C_{21}H_{27}ClN_6O$) (labeled as "Ligand of Interest" by depositor).



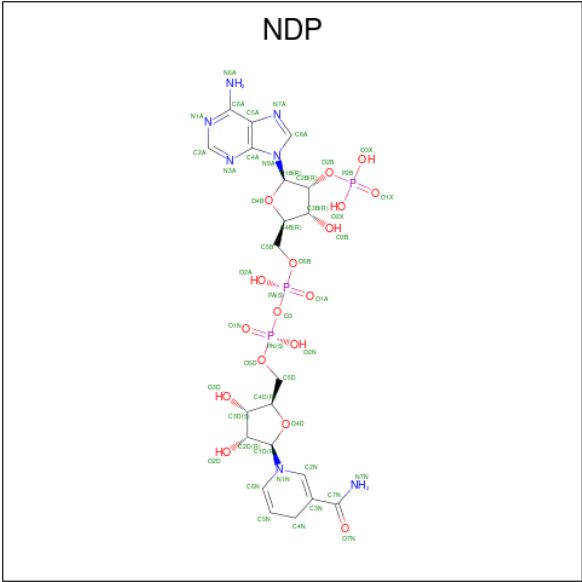
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Cl	H	N	O	4	0
			56	21	1	27	6	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total	C	Cl	H	N	O	4	1
			56	21	1	27	6	1		

- Molecule 5 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NDP) (formula: C₂₁H₃₀N₇O₁₇P₃).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	H	N	O	P	3	0
			72	21	24	7	17	3		
5	B	1	Total	C	H	N	O	P	3	0
			72	21	24	7	17	3		

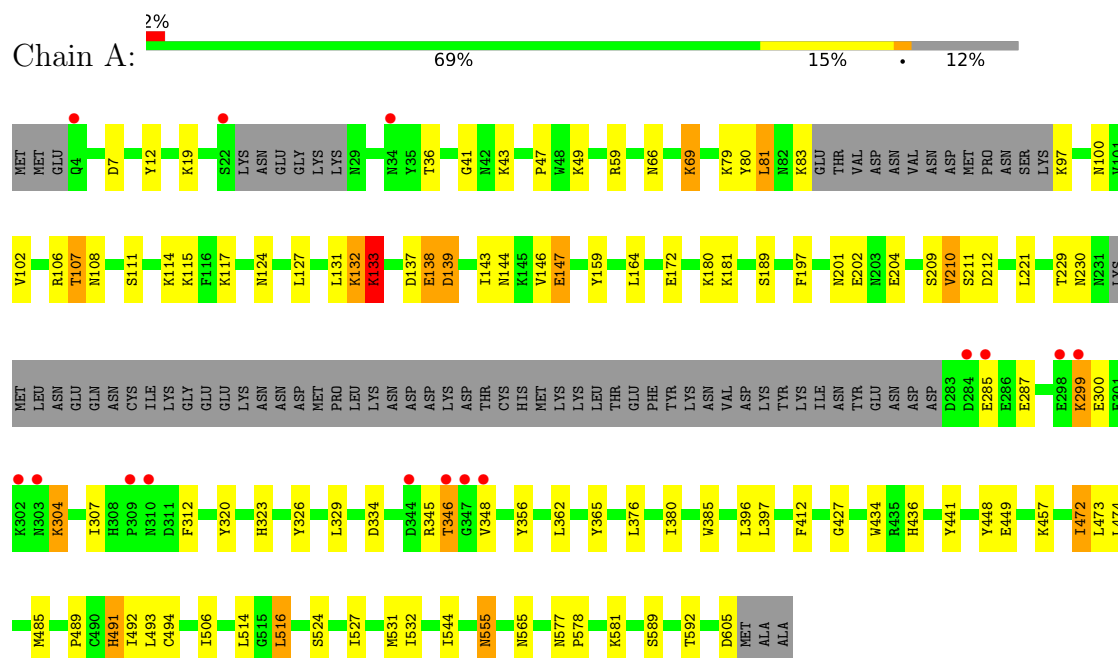
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	47	Total	O	0	0
			47	47		
6	B	49	Total	O	0	0
			49	49		

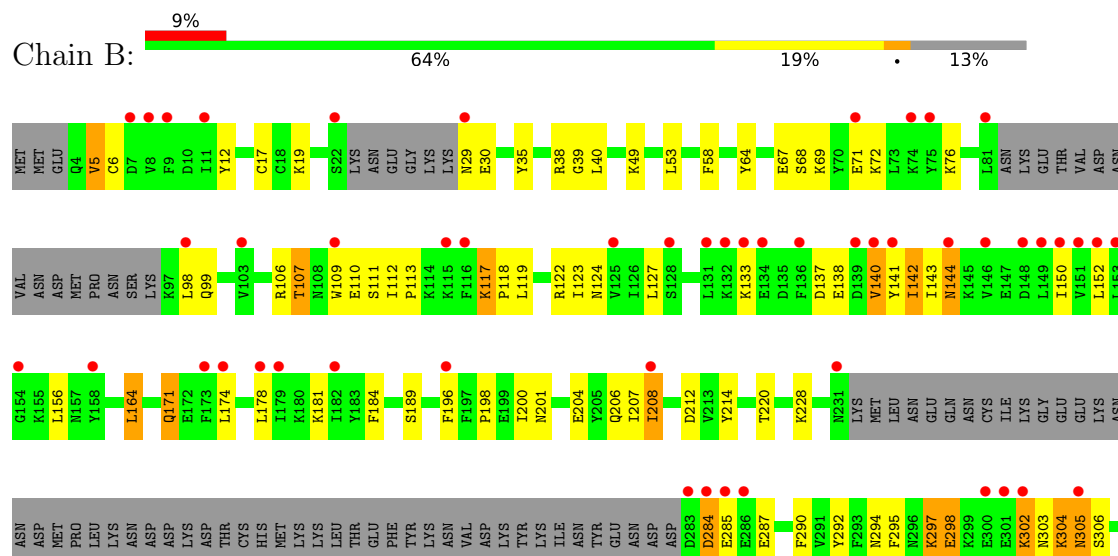
3 Residue-property plots [i](#)

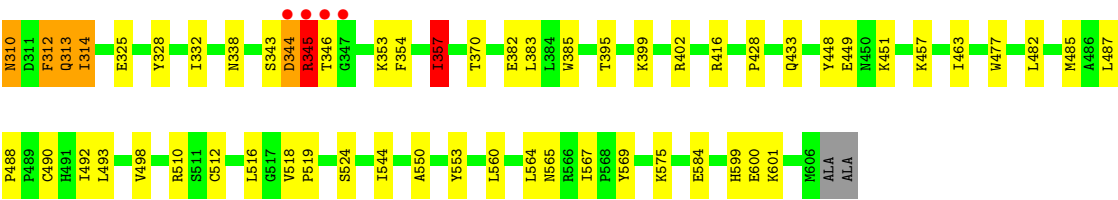
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.31Å 154.49Å 163.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.41 – 2.59 46.41 – 2.59	Depositor EDS
% Data completeness (in resolution range)	90.1 (46.41-2.59) 84.3 (46.41-2.59)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.243 , 0.311 0.189 , 0.256	Depositor DCC
R_{free} test set	2081 reflections (4.44%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.448	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18078	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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: NDP, UMP, E9X

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	1.03	2/4537 (0.0%)	1.43	9/6128 (0.1%)
1	B	1.08	0/4528	1.44	14/6116 (0.2%)
All	All	1.05	2/9065 (0.0%)	1.44	23/12244 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	6
All	All	0	7

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	138	GLU	C-O	5.87	1.31	1.24
1	A	491	HIS	CE1-NE2	5.25	1.37	1.32

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	553	TYR	CB-CA-C	6.55	120.12	109.90
1	B	357	ILE	CB-CA-C	6.38	118.79	110.12
1	A	472	ILE	CA-C-O	-6.36	117.10	122.63
1	A	229	THR	CA-C-N	6.22	128.52	120.44
1	A	229	THR	C-N-CA	6.22	128.52	120.44
1	B	284	ASP	CA-CB-CG	6.16	118.76	112.60
1	A	144	ASN	CB-CA-C	-6.15	97.46	110.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	402	ARG	CA-C-N	5.74	128.62	120.53
1	B	402	ARG	C-N-CA	5.74	128.62	120.53
1	B	212	ASP	CA-CB-CG	5.74	118.34	112.60
1	A	107	THR	N-CA-C	-5.65	104.74	111.69
1	A	578	PRO	CB-CA-C	-5.58	103.93	111.85
1	B	338	ASN	CA-C-N	5.58	125.27	119.92
1	B	338	ASN	C-N-CA	5.58	125.27	119.92
1	A	532	ILE	N-CA-C	-5.28	105.39	110.72
1	B	344	ASP	CA-CB-CG	5.28	117.88	112.60
1	B	5	VAL	CA-C-N	5.16	127.46	120.44
1	B	5	VAL	C-N-CA	5.16	127.46	120.44
1	A	197	PHE	CB-CA-C	5.08	116.73	109.26
1	B	353	LYS	CA-C-O	-5.08	115.82	121.51
1	B	19	LYS	CA-C-O	-5.07	115.73	121.72
1	A	555	ASN	CB-CA-C	5.03	119.06	110.01
1	B	312	PHE	CB-CA-C	-5.01	102.71	111.23

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	137	ASP	Peptide
1	B	113	PRO	Peptide
1	B	117	LYS	Peptide
1	B	297	LYS	Peptide
1	B	302	LYS	Peptide
1	B	306	SER	Peptide
1	B	345	ARG	Peptide

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	4433	4399	4384	56	0
1	B	4424	4389	4374	72	0
2	A	8	11	8	1	0
3	A	20	11	11	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	20	11	11	2	0
4	A	29	27	0	2	0
4	B	29	27	0	0	0
5	A	48	24	26	2	0
5	B	48	24	26	4	0
6	A	47	0	0	1	1
6	B	49	0	0	0	1
All	All	9155	8923	8840	125	1

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 7.

All (125) 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:605:ASP:C	2:A:701:MET:N	2.05	1.13
1:B:312:PHE:HA	1:B:565:ASN:HD21	1.21	1.00
1:A:312:PHE:HA	1:A:565:ASN:HD21	1.29	0.96
1:B:312:PHE:HA	1:B:565:ASN:ND2	1.94	0.81
1:B:310:ASN:ND2	1:B:310:ASN:H	1.83	0.75
1:A:506:ILE:HD12	1:B:354:PHE:CZ	2.26	0.71
1:A:127:LEU:HD23	1:A:143:ILE:HG13	1.72	0.70
1:B:127:LEU:HD12	1:B:143:ILE:HD11	1.74	0.70
1:B:140:VAL:O	1:B:141:TYR:CD2	2.46	0.69
1:B:328:TYR:CZ	1:B:332:ILE:HD11	2.28	0.69
1:B:207:ILE:HB	1:B:567:ILE:HD13	1.78	0.66
4:A:703:E9X:CAU	4:A:703:E9X:CAZ	2.74	0.66
1:B:40:LEU:HD12	1:B:196:PHE:C	2.22	0.65
1:B:107:THR:HG23	5:B:803:NDP:PA	2.38	0.63
1:A:346:THR:HG23	1:A:348:VAL:HG23	1.80	0.62
1:B:140:VAL:C	1:B:141:TYR:CD2	2.78	0.62
1:B:99:GLN:HE21	1:B:123:ILE:HG13	1.66	0.60
1:B:305:ASN:O	1:B:305:ASN:ND2	2.35	0.59
1:A:493:LEU:C	1:A:493:LEU:HD12	2.28	0.59
1:B:482:LEU:HD22	1:B:488:PRO:HB3	1.84	0.59
1:B:428:PRO:O	1:B:485:MET:HE2	2.04	0.58
1:B:490:CYS:SG	3:B:801:UMP:C6	2.96	0.58
1:A:312:PHE:HA	1:A:565:ASN:ND2	2.10	0.57
1:B:310:ASN:ND2	1:B:310:ASN:N	2.52	0.57
1:B:109:TRP:CE2	1:B:117:LYS:HE3	2.40	0.57
1:A:107:THR:HG21	6:A:818:HOH:O	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:485:MET:SD	1:A:489:PRO:HD3	2.45	0.56
1:B:171:GLN:HE21	1:B:171:GLN:HA	1.71	0.55
1:B:127:LEU:HD23	5:B:803:NDP:C4A	2.37	0.55
1:B:490:CYS:SG	3:B:801:UMP:C5	3.00	0.55
1:A:66:ASN:HD22	1:A:69:LYS:HE2	1.70	0.55
1:A:427:GLY:HA2	1:A:441:TYR:CE2	2.42	0.55
1:B:122:ARG:O	1:B:124:ASN:ND2	2.41	0.54
1:B:297:LYS:C	1:B:298:GLU:HG3	2.33	0.54
1:B:292:TYR:O	1:B:295:PHE:HB3	2.07	0.54
1:A:506:ILE:HD12	1:B:354:PHE:CE1	2.43	0.54
1:A:12:TYR:HD1	1:A:181:LYS:HB2	1.71	0.54
1:A:19:LYS:HG2	1:A:36:THR:HG22	1.89	0.53
1:B:344:ASP:O	1:B:345:ARG:C	2.51	0.53
1:B:310:ASN:O	1:B:313:GLN:NE2	2.42	0.53
1:A:376:LEU:O	1:A:380:ILE:HG13	2.09	0.53
1:B:107:THR:O	1:B:111:SER:HB3	2.08	0.52
1:B:448:TYR:O	1:B:451:LYS:HB2	2.10	0.52
1:B:144:ASN:N	1:B:144:ASN:HD22	2.07	0.52
1:A:436:HIS:CE1	1:A:448:TYR:O	2.62	0.52
1:A:221:LEU:HD23	1:A:221:LEU:N	2.25	0.51
1:B:332:ILE:HD13	1:B:560:LEU:HD22	1.93	0.51
1:A:326:TYR:HA	1:A:329:LEU:HB2	1.92	0.51
1:A:172:GLU:OE2	5:A:704:NDP:N7A	2.45	0.49
1:A:527:ILE:O	1:A:531:MET:HG3	2.12	0.49
1:B:492:ILE:HD11	1:B:510:ARG:HD3	1.94	0.49
1:A:66:ASN:ND2	1:A:69:LYS:HE2	2.27	0.48
1:A:299:LYS:HD3	1:A:304:LYS:HB2	1.94	0.48
1:A:577:ASN:OD1	1:A:577:ASN:C	2.55	0.48
1:B:357:ILE:HD11	1:B:544:ILE:HG23	1.94	0.48
1:A:492:ILE:HG21	1:B:493:LEU:CD2	2.44	0.48
1:B:290:PHE:CE2	1:B:294:ASN:ND2	2.81	0.48
1:B:493:LEU:HD12	1:B:493:LEU:C	2.39	0.48
1:B:512:CYS:HB2	1:B:550:ALA:HA	1.96	0.47
1:A:108:ASN:O	1:A:111:SER:OG	2.31	0.47
1:A:7:ASP:HA	1:A:180:LYS:HD2	1.96	0.47
1:B:17:CYS:HB2	1:B:184:PHE:CZ	2.50	0.47
1:A:211:SER:OG	1:A:212:ASP:O	2.31	0.47
1:B:395:THR:O	1:B:399:LYS:HG3	2.15	0.47
1:A:506:ILE:HA	1:A:544:ILE:O	2.15	0.46
1:B:107:THR:HA	1:B:110:GLU:HG2	1.95	0.46
1:A:138:GLU:O	1:A:139:ASP:CB	2.63	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:THR:HG23	5:B:803:NDP:O5B	2.16	0.46
1:B:58:PHE:HZ	1:B:164:LEU:HD13	1.81	0.46
1:B:127:LEU:CD2	5:B:803:NDP:C4A	2.93	0.46
1:B:152:LEU:HD12	1:B:152:LEU:O	2.16	0.46
1:B:357:ILE:HD13	1:B:357:ILE:C	2.41	0.46
1:B:118:PRO:O	1:B:119:LEU:C	2.59	0.45
1:B:144:ASN:HD22	1:B:144:ASN:H	1.63	0.45
1:B:518:VAL:N	1:B:519:PRO:CD	2.80	0.45
1:B:382:GLU:O	1:B:385:TRP:HB3	2.17	0.45
1:A:41:GLY:HA2	1:A:47:PRO:HD3	1.98	0.45
1:A:201:ASN:HD22	1:A:204:GLU:HG3	1.82	0.45
1:B:35:TYR:O	1:B:38:ARG:NH1	2.48	0.45
1:A:493:LEU:CD2	1:B:492:ILE:HG21	2.47	0.45
1:B:285:GLU:C	1:B:287:GLU:H	2.24	0.45
1:A:287:GLU:OE1	1:B:69:LYS:HD3	2.16	0.45
1:B:208:ILE:O	1:B:208:ILE:HG12	2.17	0.44
1:A:210:VAL:CG2	1:A:323:HIS:HB2	2.47	0.44
1:B:6:CYS:HB3	1:B:178:LEU:C	2.43	0.44
1:B:17:CYS:HA	1:B:39:GLY:O	2.16	0.44
1:B:142:ILE:CG2	1:B:143:ILE:N	2.80	0.44
1:A:516:LEU:HD23	1:A:516:LEU:HA	1.83	0.44
1:B:383:LEU:HD12	1:B:383:LEU:HA	1.90	0.44
1:B:207:ILE:HB	1:B:567:ILE:CD1	2.47	0.44
1:B:477:TRP:HB2	1:B:492:ILE:HG23	2.00	0.44
1:B:12:TYR:CD1	1:B:181:LYS:HB2	2.53	0.43
1:A:131:LEU:C	1:A:132:LYS:HG2	2.43	0.43
1:A:106:ARG:NE	5:A:704:NDP:O1X	2.47	0.43
1:B:310:ASN:N	1:B:310:ASN:HD22	2.16	0.43
1:B:325:GLU:CD	1:B:370:THR:H	2.26	0.43
1:A:472:ILE:C	1:A:473:LEU:HD12	2.44	0.43
4:A:703:E9X:CAU	4:A:703:E9X:CAP	2.97	0.43
1:B:457:LYS:NZ	1:B:584:GLU:OE1	2.51	0.43
1:A:81:LEU:HB3	1:A:83:LYS:HD3	2.00	0.43
1:A:80:TYR:CD2	1:A:80:TYR:O	2.72	0.43
1:A:412:PHE:C	1:A:412:PHE:CD1	2.96	0.43
1:A:514:LEU:HD13	1:A:514:LEU:HA	1.90	0.43
1:B:569:TYR:CE2	1:B:599:HIS:CD2	3.07	0.42
1:A:132:LYS:O	1:A:133:LYS:C	2.63	0.42
1:A:146:VAL:O	1:A:147:GLU:C	2.63	0.41
1:A:434:TRP:CE2	1:A:474:LEU:HD21	2.55	0.41
1:B:142:ILE:HG23	1:B:143:ILE:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:TRP:CE3	1:A:396:LEU:HD11	2.55	0.41
1:B:174:LEU:HD12	1:B:198:PRO:HG2	2.02	0.41
1:B:201:ASN:HB3	1:B:204:GLU:HB3	2.02	0.41
1:A:473:LEU:HD12	1:A:473:LEU:N	2.34	0.41
1:B:428:PRO:HB2	1:B:433:GLN:NE2	2.35	0.41
1:A:334:ASP:OD2	1:A:356:TYR:OH	2.31	0.41
1:B:463:ILE:HG23	1:B:498:VAL:HG21	2.03	0.41
1:A:397:LEU:HD23	1:A:397:LEU:HA	1.92	0.41
1:B:214:TYR:O	1:B:220:THR:HA	2.21	0.41
1:A:102:VAL:HB	1:A:164:LEU:HD11	2.03	0.41
1:A:80:TYR:HD2	1:A:81:LEU:HD13	1.86	0.40
1:A:362:LEU:HA	1:A:365:TYR:O	2.22	0.40
1:A:493:LEU:HD12	1:A:493:LEU:O	2.22	0.40
1:A:494:CYS:HA	1:A:506:ILE:O	2.21	0.40
1:B:314:ILE:HB	1:B:565:ASN:OD1	2.20	0.40
1:A:209:SER:HB3	1:A:320:TYR:HB2	2.03	0.40
1:A:100:ASN:OD1	1:A:159:TYR:HB3	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:846:HOH:O	6:B:946:HOH:O[1_455]	1.83	0.37

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	524/608 (86%)	492 (94%)	27 (5%)	5 (1%)	12	28
1	B	523/608 (86%)	470 (90%)	43 (8%)	10 (2%)	6	13
All	All	1047/1216 (86%)	962 (92%)	70 (7%)	15 (1%)	9	19

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	139	ASP
1	B	49	LYS
1	B	303	ASN
1	B	304	LYS
1	B	345	ARG
1	A	133	LYS
1	A	345	ARG
1	B	76	LYS
1	B	137	ASP
1	B	140	VAL
1	A	299	LYS
1	B	64	TYR
1	B	98	LEU
1	B	150	ILE

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	498/570 (87%)	467 (94%)	31 (6%)	16	36
1	B	497/570 (87%)	454 (91%)	43 (9%)	9	21
All	All	995/1140 (87%)	921 (93%)	74 (7%)	12	29

All (74) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LYS
1	A	59	ARG
1	A	69	LYS
1	A	79	LYS
1	A	81	LEU
1	A	97	LYS
1	A	114	LYS
1	A	115	LYS

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Mol	Chain	Res	Type
1	A	117	LYS
1	A	124	ASN
1	A	132	LYS
1	A	133	LYS
1	A	147	GLU
1	A	189	SER
1	A	202	GLU
1	A	210	VAL
1	A	230	ASN
1	A	285	GLU
1	A	300	GLU
1	A	304	LYS
1	A	307	ILE
1	A	346	THR
1	A	449	GLU
1	A	457	LYS
1	A	491	HIS
1	A	516	LEU
1	A	524	SER
1	A	555	ASN
1	A	581	LYS
1	A	589	SER
1	A	592	THR
1	B	5	VAL
1	B	29	ASN
1	B	30	GLU
1	B	53	LEU
1	B	67	GLU
1	B	68	SER
1	B	71	GLU
1	B	72	LYS
1	B	106	ARG
1	B	107	THR
1	B	112	ILE
1	B	133	LYS
1	B	138	GLU
1	B	142	ILE
1	B	144	ASN
1	B	156	LEU
1	B	164	LEU
1	B	171	GLN
1	B	189	SER

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Mol	Chain	Res	Type
1	B	200	ILE
1	B	206	GLN
1	B	208	ILE
1	B	228	LYS
1	B	284	ASP
1	B	298	GLU
1	B	302	LYS
1	B	304	LYS
1	B	305	ASN
1	B	310	ASN
1	B	313	GLN
1	B	314	ILE
1	B	343	SER
1	B	346	THR
1	B	357	ILE
1	B	416	ARG
1	B	449	GLU
1	B	487	LEU
1	B	516	LEU
1	B	524	SER
1	B	564	LEU
1	B	575	LYS
1	B	600	GLU
1	B	601	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	ASN
1	A	201	ASN
1	A	316	ASN
1	A	338	ASN
1	A	394	ASN
1	A	415	ASN
1	A	424	ASN
1	A	436	HIS
1	A	465	ASN
1	A	555	ASN
1	A	556	HIS
1	B	29	ASN
1	B	99	GLN
1	B	144	ASN

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Mol	Chain	Res	Type
1	B	157	ASN
1	B	171	GLN
1	B	206	GLN
1	B	231	ASN
1	B	303	ASN
1	B	310	ASN
1	B	313	GLN
1	B	394	ASN
1	B	407	ASN
1	B	424	ASN
1	B	458	ASN
1	B	539	GLN
1	B	554	ASN
1	B	556	HIS
1	B	563	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](#)

8 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	E9X	A	703	-	28,31,31	0.64	1 (3%)	34,43,43	1.16	3 (8%)
3	UMP	A	702	-	21,21,21	0.50	0	30,31,31	0.65	0
5	NDP	B	803	-	51,52,52	0.56	0	71,80,80	0.73	1 (1%)
2	MET	A	701	-	6,7,8	0.77	0	2,7,9	0.65	0
5	NDP	A	704	-	51,52,52	0.60	0	71,80,80	0.84	2 (2%)
3	UMP	B	801	-	21,21,21	0.44	0	30,31,31	0.85	1 (3%)

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	E9X	A	703	-	-	10/15/34/34	0/3/3/3
3	UMP	A	702	-	-	2/10/22/22	0/2/2/2
5	NDP	B	803	-	-	8/34/77/77	0/5/5/5
2	MET	A	701	-	-	4/5/6/8	-
5	NDP	A	704	-	-	4/34/77/77	0/5/5/5
3	UMP	B	801	-	-	2/10/22/22	0/2/2/2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	703	E9X	OAK-NAF	-2.09	1.38	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	703	E9X	CAT-CAS-NAO	4.34	121.16	114.13
4	A	703	E9X	NAG-CAA-NAF	2.61	119.61	117.46
4	A	703	E9X	CAS-NAO-CAP	2.41	124.89	120.72
5	A	704	NDP	P2B-O2B-C2B	-2.34	117.18	123.43
5	B	803	NDP	P2B-O2B-C2B	-2.29	117.31	123.43
5	A	704	NDP	O4D-C1D-C2D	-2.02	102.30	106.62
3	B	801	UMP	C2'-C3'-C4'	-2.01	98.72	102.80

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	MET	O-C-CA-CB
2	A	701	MET	N-CA-CB-CG
2	A	701	MET	C-CA-CB-CG
4	A	703	E9X	OAK-CAL-CAM-CAN
4	A	703	E9X	CAT-CAS-NAO-CAP
5	B	803	NDP	O4B-C4B-C5B-O5B
5	B	803	NDP	C3B-C4B-C5B-O5B
5	B	803	NDP	C2D-C1D-N1N-C6N
5	B	803	NDP	C2N-C3N-C7N-N7N
4	A	703	E9X	CAZ-CAP-NAO-CAN
5	B	803	NDP	C2D-C1D-N1N-C2N
4	A	703	E9X	CAQ-CAP-NAO-CAN
3	B	801	UMP	C3'-C4'-C5'-O5'
3	B	801	UMP	O4'-C4'-C5'-O5'
4	A	703	E9X	CAZ-CAP-NAO-CAS
3	A	702	UMP	C3'-C4'-C5'-O5'
2	A	701	MET	CA-CB-CG-SD
3	A	702	UMP	O4'-C4'-C5'-O5'
4	A	703	E9X	CAM-CAN-NAO-CAP
5	A	704	NDP	PA-O3-PN-O5D
5	B	803	NDP	PA-O3-PN-O5D
4	A	703	E9X	CAL-CAM-CAN-NAO
4	A	703	E9X	CAT-CAS-NAO-CAN
5	B	803	NDP	O4D-C1D-N1N-C6N
5	A	704	NDP	O4D-C1D-N1N-C2N
5	A	704	NDP	C2N-C3N-C7N-N7N
5	B	803	NDP	O4D-C1D-N1N-C2N
4	A	703	E9X	CAM-CAN-NAO-CAS
5	A	704	NDP	C2D-C1D-N1N-C2N
4	A	703	E9X	CAQ-CAP-NAO-CAS

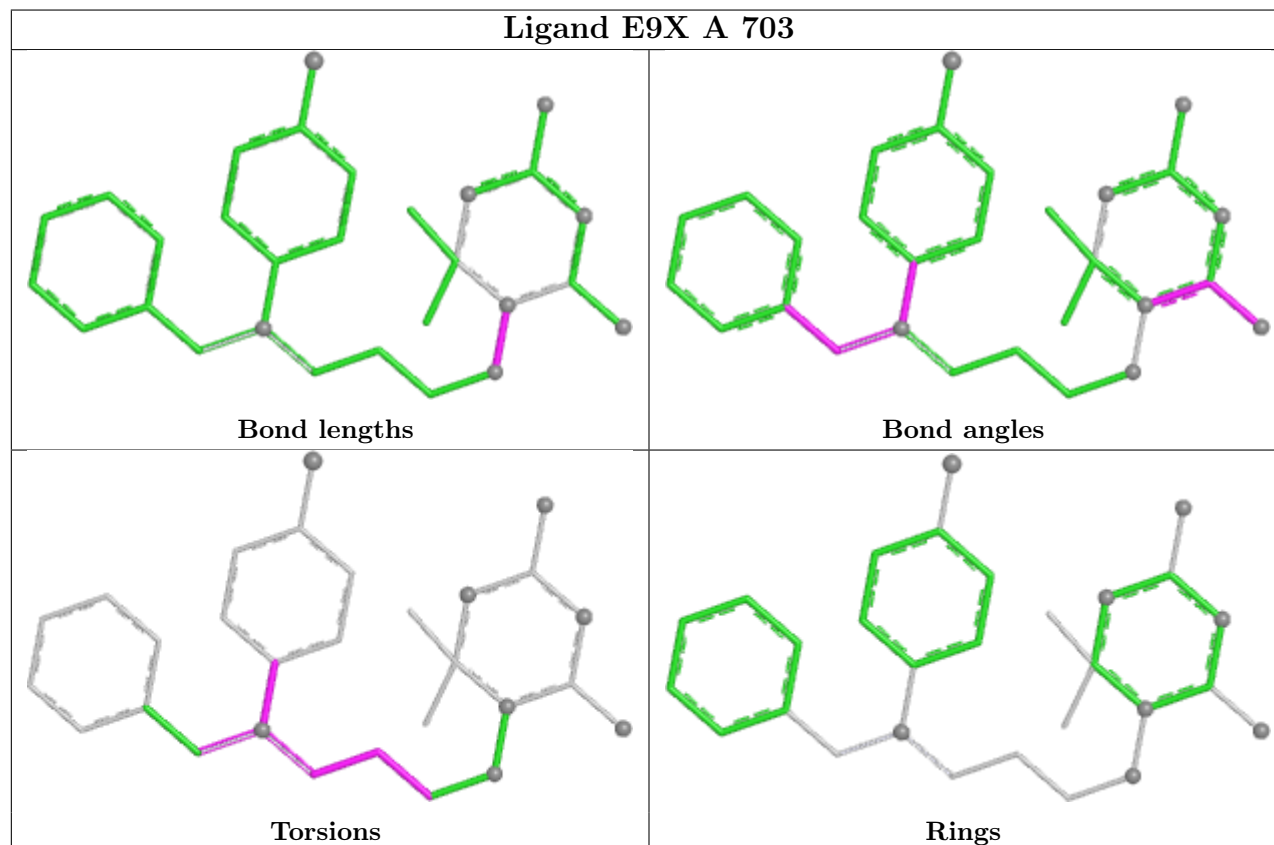
There are no ring outliers.

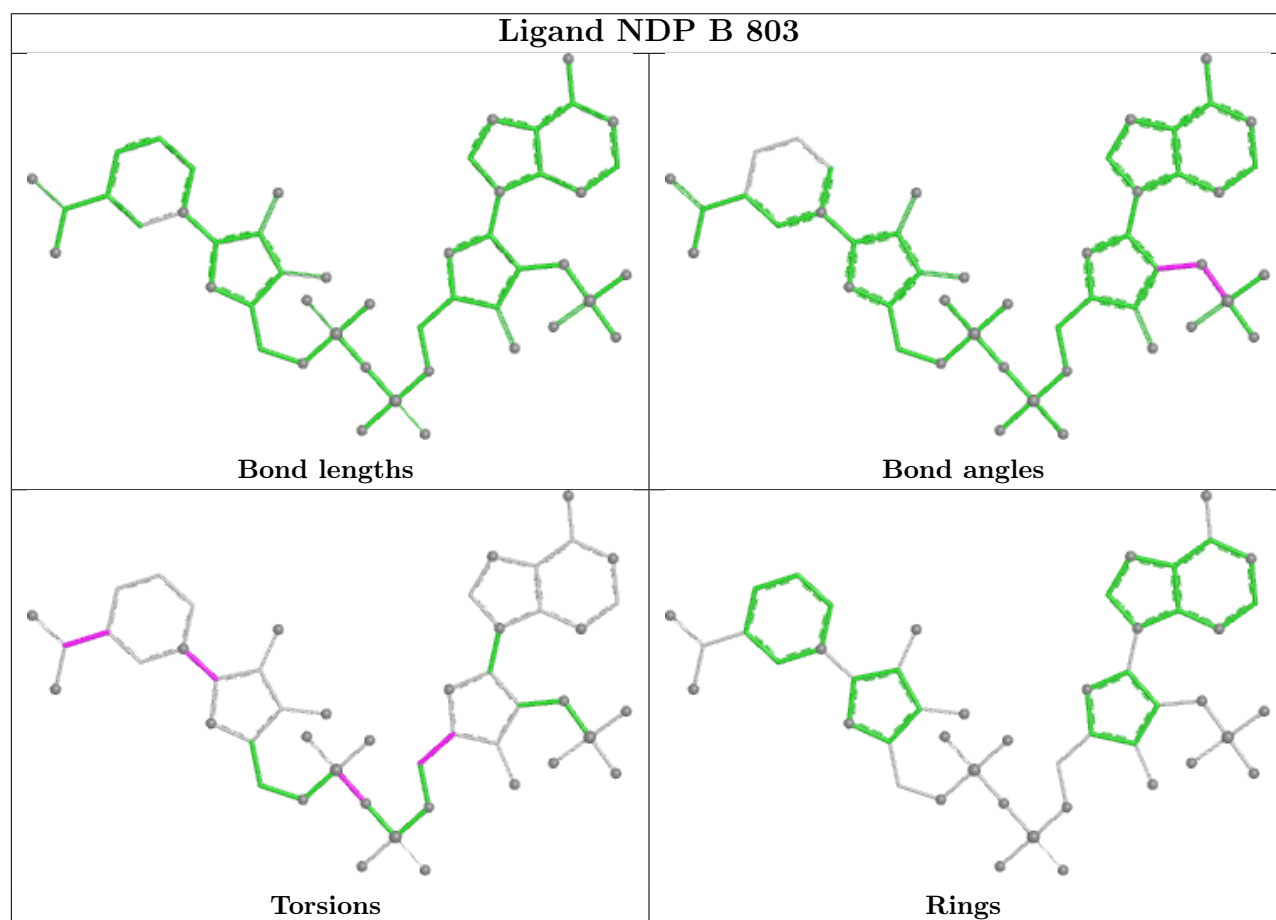
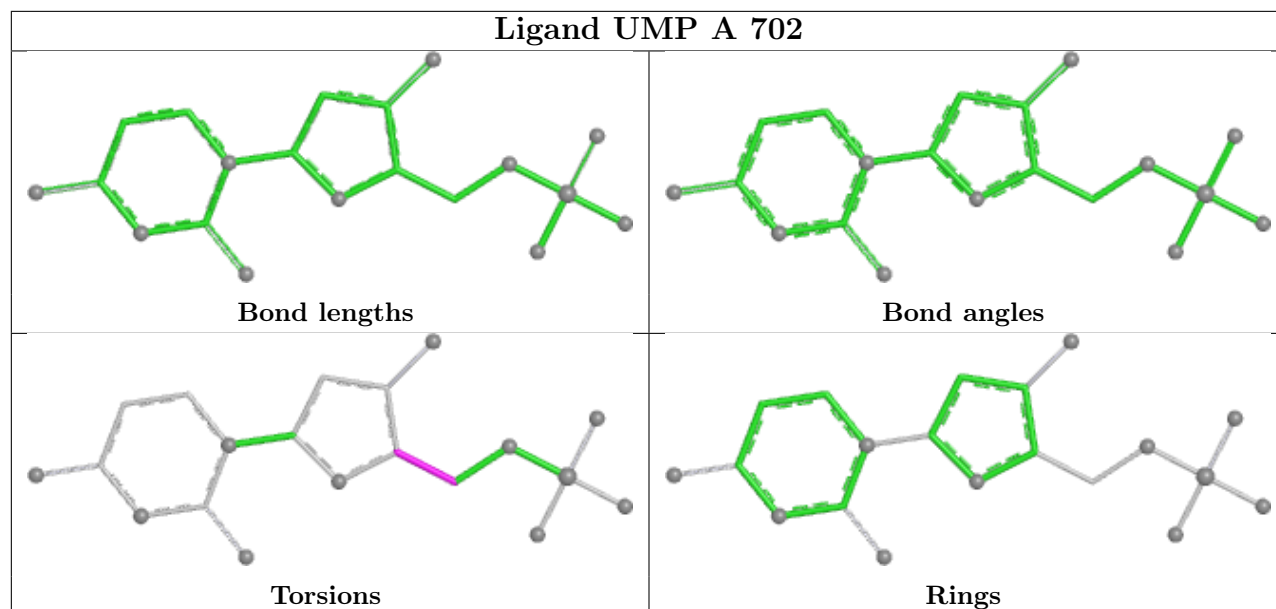
5 monomers are involved in 11 short contacts:

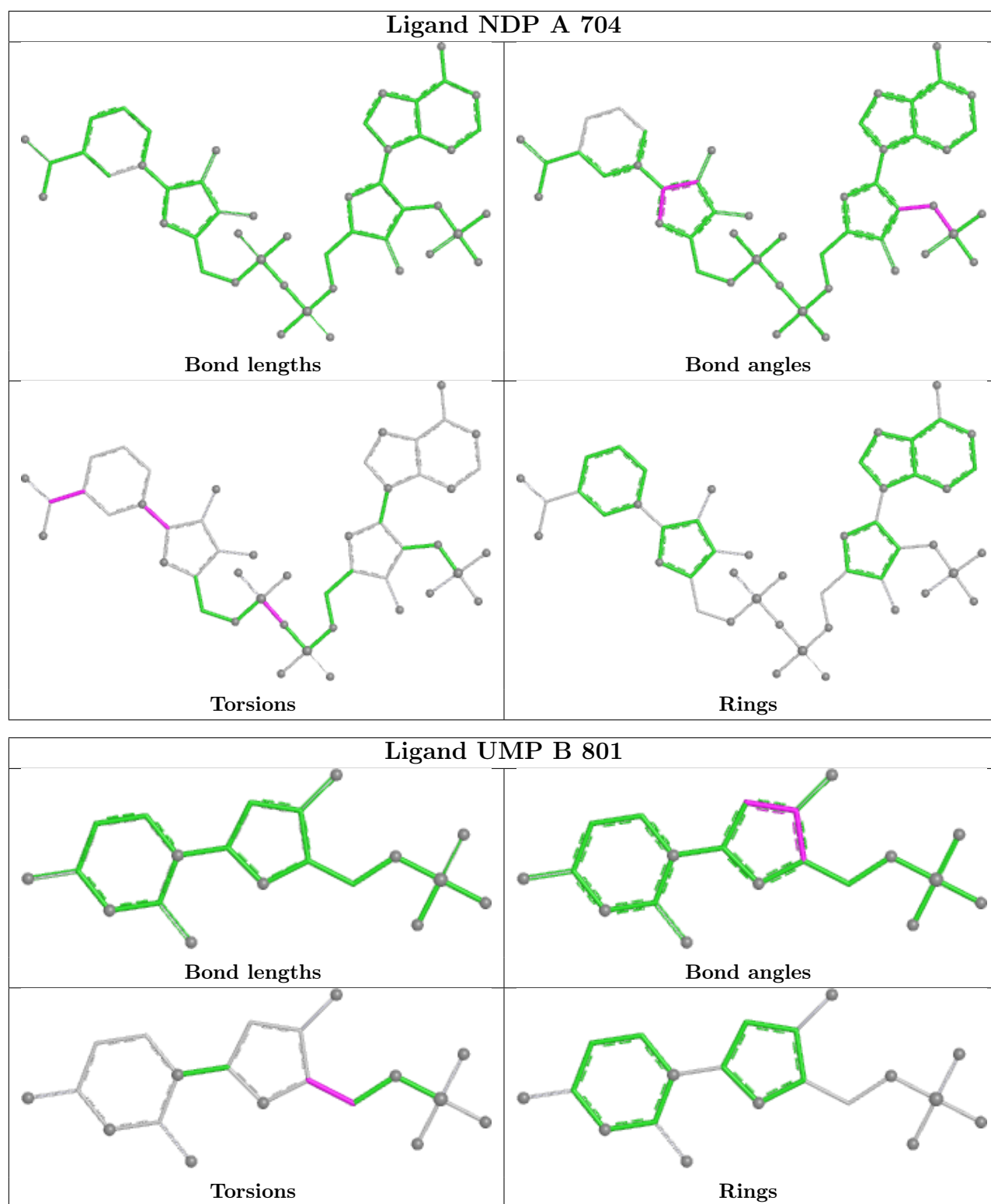
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	703	E9X	2	0
5	B	803	NDP	4	0
2	A	701	MET	1	0
5	A	704	NDP	2	0
3	B	801	UMP	2	0

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.







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	532/608 (87%)	-0.31	15 (2%) 55 49	14, 28, 70, 142	0
1	B	531/608 (87%)	0.19	55 (10%) 11 9	14, 32, 111, 136	0
All	All	1063/1216 (87%)	-0.06	70 (6%) 24 19	14, 29, 102, 142	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	136	PHE	5.2
1	A	347	GLY	5.0
1	B	150	ILE	4.5
1	B	140	VAL	4.4
1	B	116	PHE	4.2
1	B	285	GLU	4.0
1	A	348	VAL	4.0
1	B	115	LYS	3.8
1	A	346	THR	3.8
1	B	154	GLY	3.7
1	B	81	LEU	3.7
1	B	75	TYR	3.6
1	B	141	TYR	3.5
1	A	22	SER	3.3
1	B	302	LYS	3.3
1	B	125	VAL	3.2
1	B	139	ASP	3.2
1	B	11	ILE	3.1
1	A	303	ASN	3.1
1	A	299	LYS	3.0
1	A	344	ASP	3.0
1	B	149	LEU	3.0
1	B	109	TRP	3.0
1	B	103	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	148	ASP	2.9
1	B	283	ASP	2.9
1	B	8	VAL	2.7
1	B	71	GLU	2.7
1	B	301	GLU	2.6
1	B	182	ILE	2.6
1	B	131	LEU	2.6
1	B	286	GLU	2.6
1	B	153	LEU	2.5
1	B	344	ASP	2.5
1	A	34	ASN	2.5
1	B	146	VAL	2.5
1	B	151	VAL	2.5
1	B	144	ASN	2.5
1	B	231	ASN	2.5
1	A	285	GLU	2.4
1	B	347	GLY	2.4
1	B	208	ILE	2.4
1	B	132	LYS	2.4
1	B	173	PHE	2.4
1	B	133	LYS	2.4
1	A	298	GLU	2.4
1	B	22	SER	2.3
1	B	346	THR	2.3
1	B	179	ILE	2.3
1	B	7	ASP	2.3
1	B	174	LEU	2.3
1	A	302	LYS	2.3
1	B	284	ASP	2.3
1	B	196	PHE	2.3
1	B	74	LYS	2.2
1	A	284	ASP	2.2
1	A	4	GLN	2.1
1	B	305	ASN	2.1
1	B	134	GLU	2.1
1	B	9	PHE	2.1
1	B	29	ASN	2.1
1	B	152	LEU	2.1
1	B	128	SER	2.1
1	A	310	ASN	2.1
1	B	98	LEU	2.1
1	B	345	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	309	PRO	2.1
1	B	178	LEU	2.0
1	B	158	TYR	2.0
1	B	300	GLU	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](#)

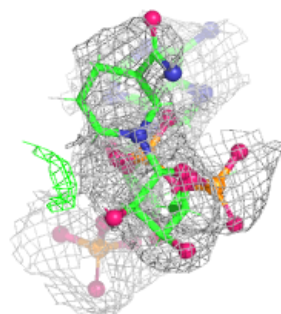
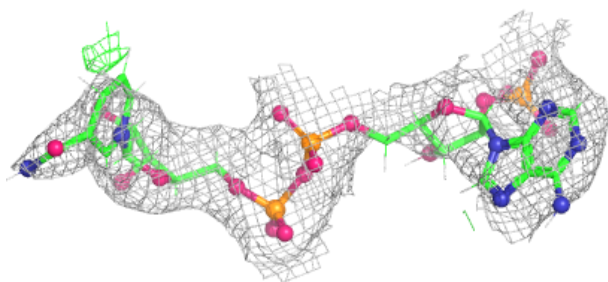
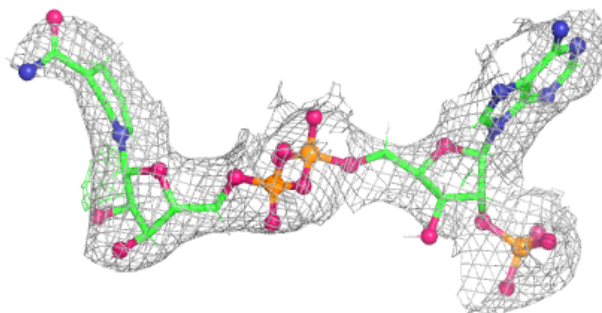
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NDP	B	803	48/48	0.84	0.13	57,91,122,125	3
4	E9X	B	802[B]	29/29	0.87	0.15	57,73,80,85	3
4	E9X	B	802[A]	29/29	0.87	0.15	57,73,80,85	3
2	MET	A	701	8/9	0.88	0.17	53,57,60,74	0
4	E9X	A	703	29/29	0.94	0.10	22,35,56,60	4
3	UMP	B	801	20/20	0.96	0.09	30,43,48,51	1
5	NDP	A	704	48/48	0.98	0.05	21,25,29,30	3
3	UMP	A	702	20/20	0.98	0.06	21,32,42,47	1

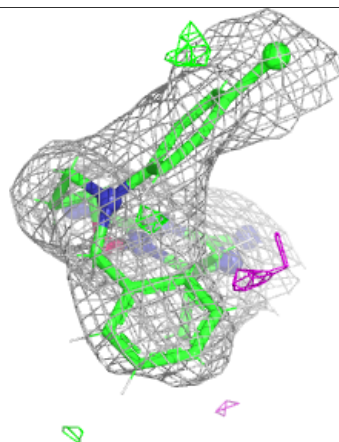
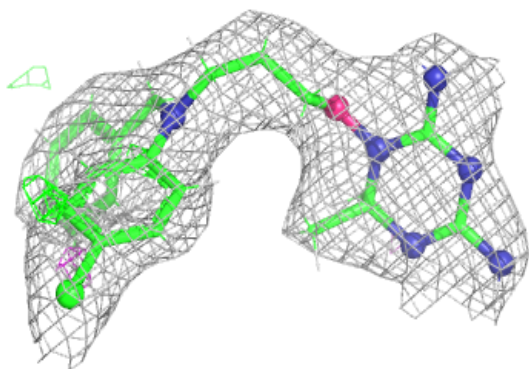
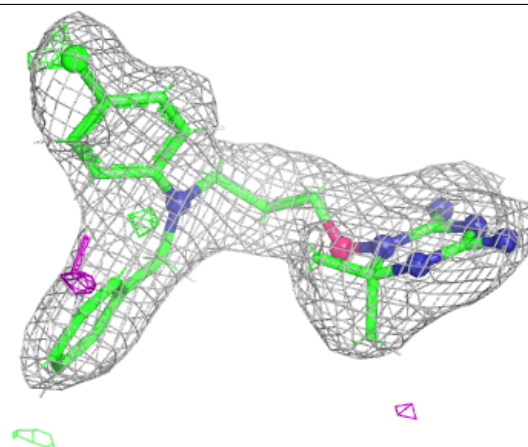
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around NDP B 803:

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

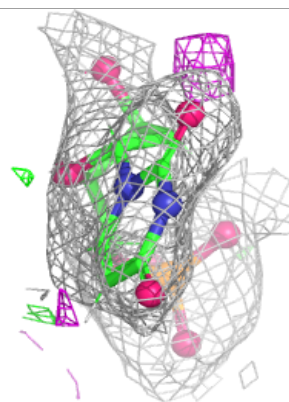
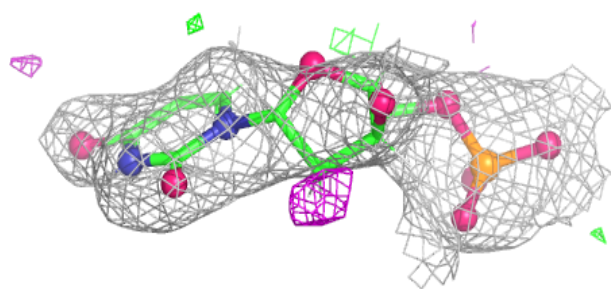
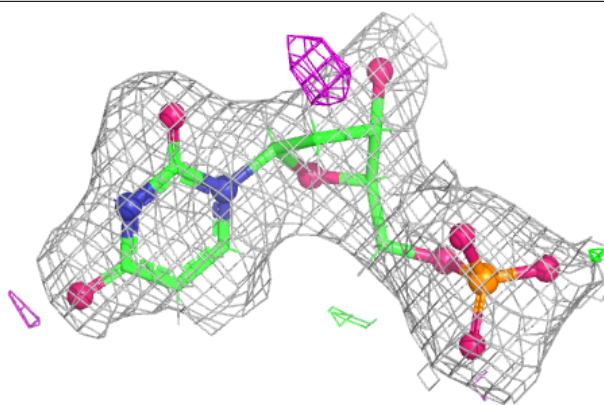
**Electron density around E9X A 703:**

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

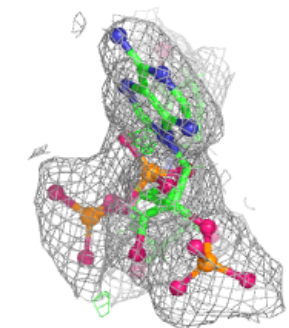
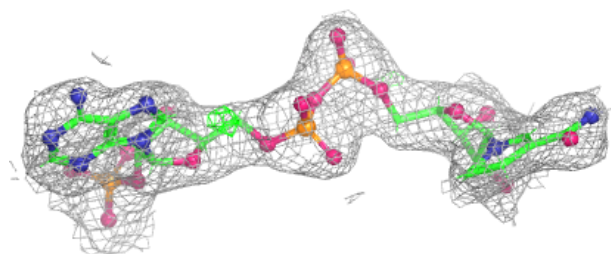
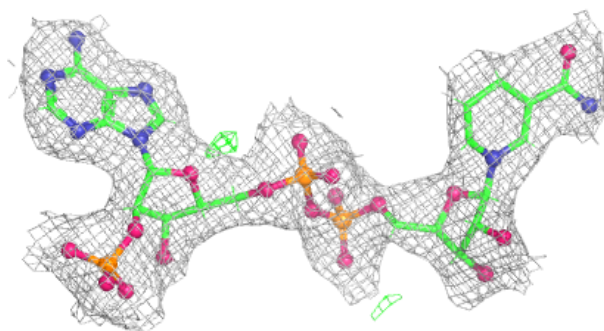


Electron density around UMP B 801:

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

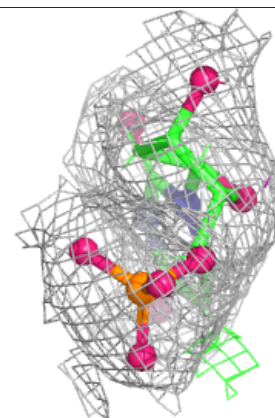
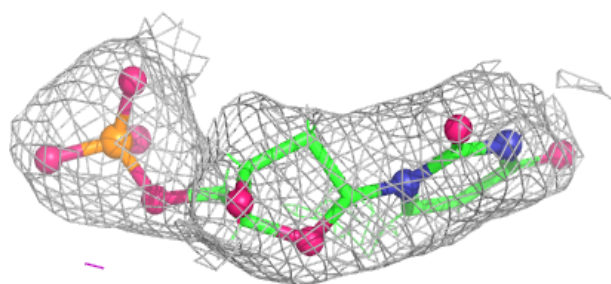
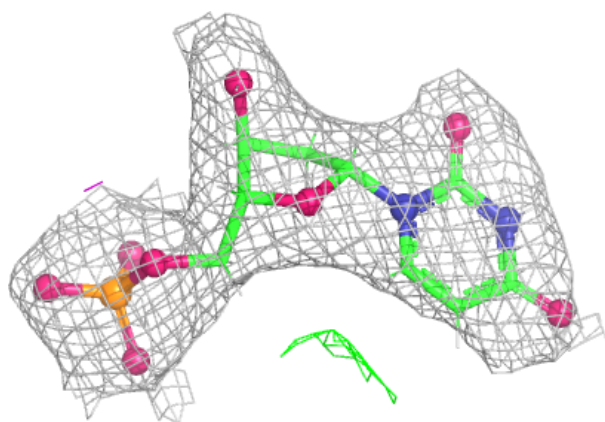
**Electron density around NDP A 704:**

$2mF_o-DF_c$ (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)



6.5 Other polymers [i](#)

There are no such residues in this entry.