



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 01:36 am BST

PDB ID : 3HBB
Title : Structures of dihydrofolate reductase-thymidylate synthase of *Trypanosoma cruzi* in the folate-free state and in complex with two antifolate drugs, trimetrexate and methotrexate
Authors : Schormann, N.; Senkovich, O.; Chattopadhyay, D.
Deposited on : 2009-05-04
Resolution : 3.00 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 : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

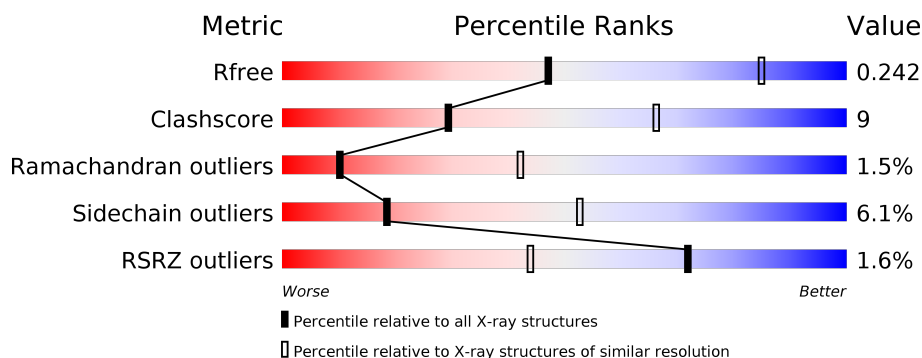
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	521	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>••</div> </div> </div>
1	B	521	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>17%</div> <div>5%</div> <div>••</div> </div> </div>
1	C	521	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>17%</div> <div>••</div> </div> </div>
1	D	521	<div> <div>%</div> <div> <div></div> <div>83%</div> <div>13%</div> <div>••</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	TMQ	A	611	-	-	X	-
3	TMQ	B	612	-	-	X	X
3	TMQ	C	613	-	-	X	-
3	TMQ	D	614	-	-	X	-
4	SO4	C	822	-	-	-	X

2 Entry composition ⓘ

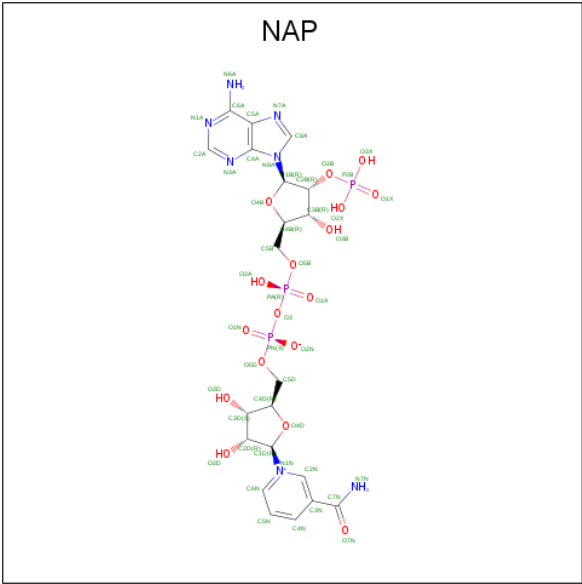
There are 6 unique types of molecules in this entry. The entry contains 16812 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydrofolate reductase-thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	510	Total	C	N	O	S	0	0	0
			4052	2571	714	749	18			
1	B	505	Total	C	N	O	S	0	0	0
			4010	2546	707	739	18			
1	C	510	Total	C	N	O	S	0	0	0
			4049	2570	712	748	19			
1	D	507	Total	C	N	O	S	0	0	0
			4023	2553	711	741	18			

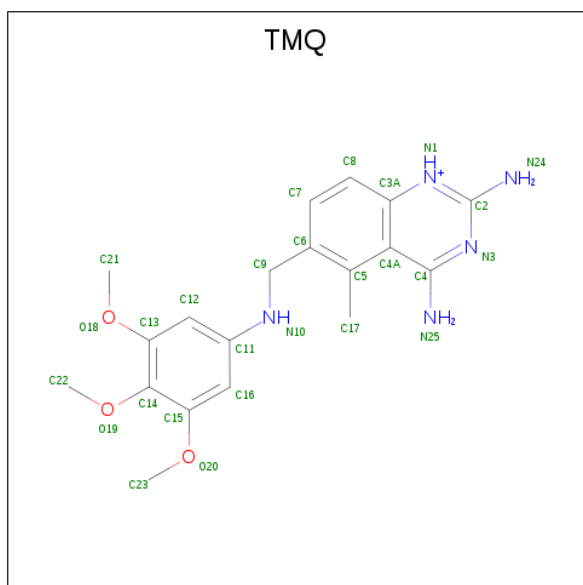
- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



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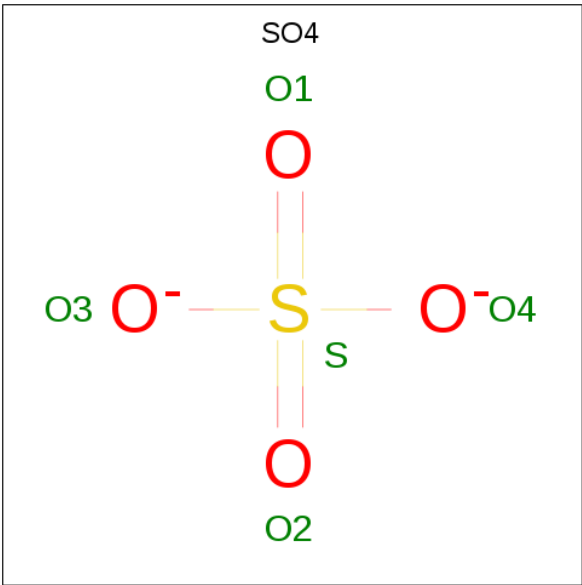
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	P	
			48	21	7	17	3	
2	D	1	Total	C	N	O	P	
			48	21	7	17	3	

- Molecule 3 is TRIMETREXATE (three-letter code: TMQ) (formula: $C_{19}H_{24}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O		
			27	19	5	3		
3	B	1	Total	C	N	O		
			27	19	5	3		
3	C	1	Total	C	N	O		
			27	19	5	3		
3	D	1	Total	C	N	O		
			27	19	5	3		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



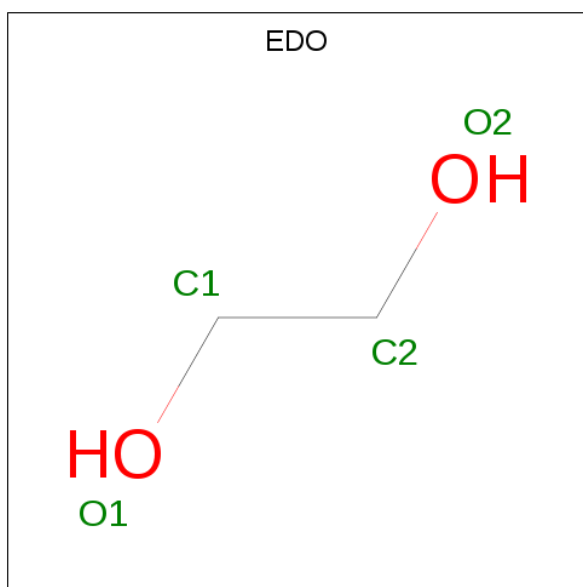
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		

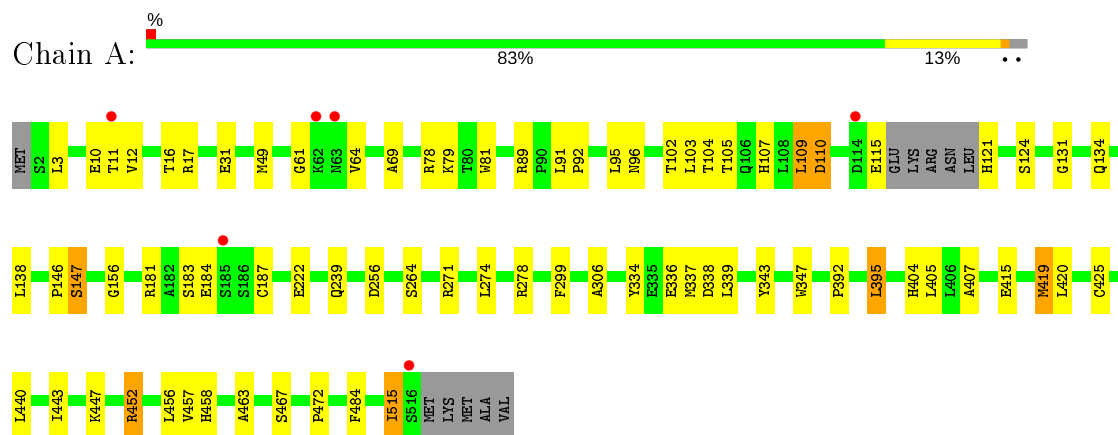
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	43	Total 43	O 43	0	0
6	B	45	Total 45	O 45	0	0
6	C	56	Total 56	O 56	0	0
6	D	46	Total 46	O 46	0	0

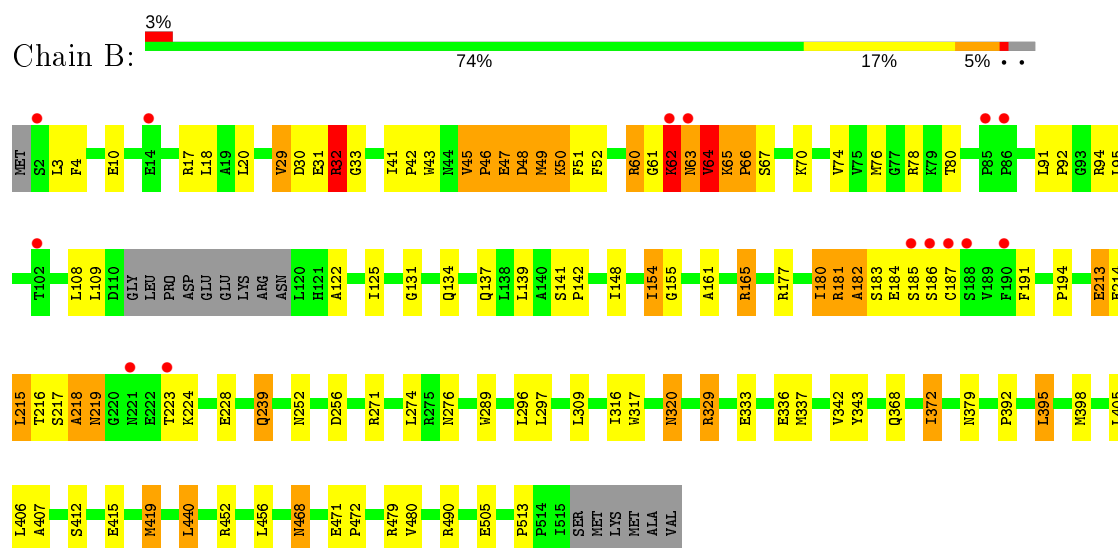
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

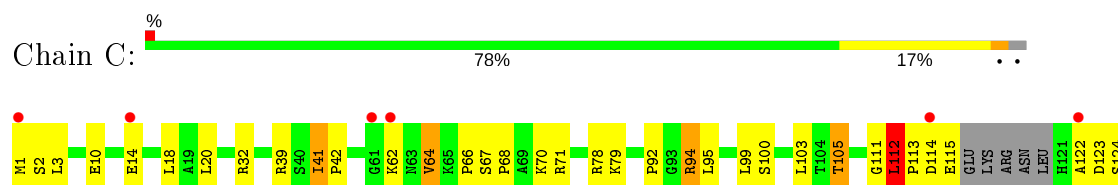
• Molecule 1: Dihydrofolate reductase-thymidylate synthase

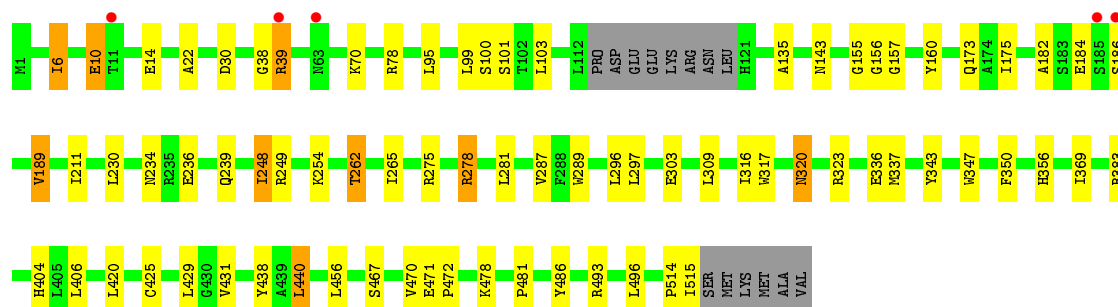


• Molecule 1: Dihydrofolate reductase-thymidylate synthase



• Molecule 1: Dihydrofolate reductase-thymidylate synthase





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	176.57Å 176.57Å 251.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 3.00 19.90 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.1 (19.94-3.00) 97.1 (19.90-3.00)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.16 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.5.0053	Depositor
R, R_{free}	0.201 , 0.239 0.204 , 0.242	Depositor DCC
R_{free} test set	3898 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	60.9	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 61.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	16812	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.99% 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: TMQ, NAP, SO4, EDO

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	0.38	0/4150	0.56	1/5632 (0.0%)
1	B	0.41	0/4106	0.68	1/5572 (0.0%)
1	C	0.42	1/4146 (0.0%)	0.56	0/5626
1	D	0.38	0/4120	0.53	0/5590
All	All	0.39	1/16522 (0.0%)	0.59	2/22420 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	153	CYS	CB-SG	-5.94	1.72	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	VAL	C-N-CD	-27.63	59.82	120.60
1	A	109	LEU	CA-CB-CG	5.30	127.48	115.30

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	4052	0	4000	47	4

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4010	0	3959	120	0
1	C	4049	0	4002	56	0
1	D	4023	0	3973	45	4
2	A	48	0	25	7	0
2	B	48	0	25	7	0
2	C	48	0	25	11	0
2	D	48	0	25	12	0
3	A	27	0	24	14	0
3	B	27	0	24	14	0
3	C	27	0	24	9	0
3	D	27	0	24	12	0
4	A	30	0	0	0	0
4	B	25	0	0	0	0
4	C	45	0	0	0	0
4	D	40	0	0	0	0
5	A	8	0	12	0	0
5	B	4	0	6	1	0
5	C	16	0	24	2	0
5	D	20	0	30	2	0
6	A	43	0	0	0	0
6	B	45	0	0	5	0
6	C	56	0	0	3	0
6	D	46	0	0	1	0
All	All	16812	0	16202	311	4

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 9.

All (311) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:GLU:CG	1:B:217:SER:OG	1.75	1.33
1:B:64:VAL:CG2	1:B:66:PRO:HD3	1.62	1.29
1:B:47:GLU:HG3	1:B:217:SER:OG	1.02	1.19
1:B:180:ILE:HG23	1:B:181:ARG:N	1.54	1.17
1:B:64:VAL:CG2	1:B:65:LYS:N	2.12	1.13
1:B:64:VAL:HG23	1:B:66:PRO:HD3	1.25	1.13
1:B:180:ILE:CG2	1:B:181:ARG:H	1.58	1.12
1:B:64:VAL:CG2	1:B:65:LYS:H	1.63	1.10
1:B:215:LEU:HD12	1:B:216:THR:N	1.67	1.09
1:B:64:VAL:HG23	1:B:65:LYS:N	1.65	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:612:TMQ:C17	3:B:612:TMQ:HN52	1.65	1.08
1:A:3:LEU:O	1:A:3:LEU:HD13	1.52	1.07
3:D:614:TMQ:HN52	3:D:614:TMQ:C17	1.65	1.05
3:B:612:TMQ:O20	3:B:612:TMQ:H222	1.56	1.05
3:C:613:TMQ:HN52	3:C:613:TMQ:C17	1.65	1.05
1:B:155:GLY:HA2	2:B:602:NAP:H5N	1.33	1.04
1:B:64:VAL:HG23	1:B:65:LYS:H	0.88	1.04
1:B:215:LEU:HD12	1:B:216:THR:H	0.89	1.04
1:B:45:VAL:HG21	1:B:180:ILE:HD11	1.37	1.04
3:A:611:TMQ:C17	3:A:611:TMQ:HN52	1.65	1.01
1:B:180:ILE:O	1:B:181:ARG:HG2	1.61	1.01
1:B:30:ASP:HB3	1:B:180:ILE:CG2	1.92	0.99
1:B:30:ASP:HA	1:B:180:ILE:HG22	1.39	0.99
1:B:46:PRO:O	1:B:47:GLU:HB2	1.64	0.98
1:B:30:ASP:CA	1:B:180:ILE:HG22	1.95	0.96
1:B:30:ASP:HB3	1:B:180:ILE:HG21	1.47	0.95
1:B:131:GLY:H	1:B:134:GLN:HE21	1.13	0.93
1:B:47:GLU:CG	1:B:217:SER:HG	1.70	0.93
1:B:31:GLU:O	1:B:32:ARG:HB2	1.67	0.92
1:B:61:GLY:O	1:B:62:LYS:HB2	1.66	0.92
1:D:78:ARG:NH1	2:D:604:NAP:O1X	2.03	0.90
1:C:192:ARG:HG3	6:C:568:HOH:O	1.72	0.89
1:C:100:SER:OG	2:C:603:NAP:O1X	1.90	0.89
1:B:215:LEU:CD1	1:B:216:THR:H	1.83	0.87
1:B:45:VAL:HG21	1:B:180:ILE:CD1	2.04	0.87
3:B:612:TMQ:HN52	3:B:612:TMQ:H172	1.41	0.86
3:D:614:TMQ:HN52	3:D:614:TMQ:H173	1.41	0.85
3:B:612:TMQ:HN52	3:B:612:TMQ:H173	1.41	0.85
3:D:614:TMQ:HN52	3:D:614:TMQ:H172	1.41	0.84
3:A:611:TMQ:H173	3:A:611:TMQ:HN52	1.41	0.84
3:C:613:TMQ:HN52	3:C:613:TMQ:H173	1.41	0.84
3:C:613:TMQ:HN52	3:C:613:TMQ:H172	1.41	0.83
1:A:79:LYS:HG3	2:A:601:NAP:H51A	1.58	0.83
3:C:613:TMQ:N25	3:C:613:TMQ:C17	2.38	0.83
3:A:611:TMQ:H172	3:A:611:TMQ:HN52	1.41	0.82
1:A:419:MET:HG2	1:A:457:VAL:HB	1.60	0.81
1:B:180:ILE:HG23	1:B:181:ARG:H	0.71	0.81
1:B:62:LYS:O	1:B:63:ASN:HB2	1.81	0.81
1:D:78:ARG:HD2	1:D:103:LEU:HD23	1.62	0.81
3:B:612:TMQ:C22	3:B:612:TMQ:O20	2.30	0.80
1:B:180:ILE:O	1:B:181:ARG:CG	2.30	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:156:GLY:O	1:C:157:GLY:C	2.20	0.79
1:B:46:PRO:O	1:B:47:GLU:CB	2.29	0.78
1:A:78:ARG:HD2	1:A:103:LEU:HD23	1.63	0.78
1:B:30:ASP:CB	1:B:180:ILE:CG2	2.61	0.78
1:B:218:ALA:HA	6:B:561:HOH:O	1.84	0.76
1:B:43:TRP:HD1	1:B:45:VAL:HG23	1.50	0.75
3:D:614:TMQ:H213	3:D:614:TMQ:H222	1.68	0.75
1:C:155:GLY:HA2	1:C:160:TYR:CZ	2.21	0.75
3:D:614:TMQ:N25	3:D:614:TMQ:C17	2.38	0.74
3:A:611:TMQ:C17	3:A:611:TMQ:N25	2.38	0.74
1:D:78:ARG:HD3	2:D:604:NAP:O1X	1.85	0.74
1:B:64:VAL:HG21	1:B:66:PRO:HD3	1.70	0.74
3:B:612:TMQ:H223	6:B:558:HOH:O	1.88	0.74
2:A:601:NAP:C4N	3:A:611:TMQ:H172	2.18	0.73
1:B:43:TRP:CD1	1:B:45:VAL:HG23	2.23	0.73
1:B:64:VAL:HG22	1:B:65:LYS:N	2.01	0.73
2:B:602:NAP:C4N	3:B:612:TMQ:H172	2.17	0.73
1:C:78:ARG:NH1	2:C:603:NAP:O3X	2.21	0.73
1:C:288:PHE:HA	5:C:909:EDO:H11	1.71	0.73
1:D:99:LEU:O	2:D:604:NAP:H1B	1.89	0.73
1:B:216:THR:HB	1:B:224:LYS:HA	1.71	0.71
1:B:180:ILE:CG2	1:B:181:ARG:N	2.30	0.70
3:B:612:TMQ:C17	3:B:612:TMQ:N25	2.38	0.70
1:A:31:GLU:OE2	1:A:181:ARG:HD2	1.92	0.69
3:D:614:TMQ:H232	3:D:614:TMQ:H223	1.74	0.69
1:B:30:ASP:CB	1:B:180:ILE:HG22	2.22	0.69
1:C:78:ARG:HH11	2:C:603:NAP:P2B	2.16	0.69
1:B:180:ILE:O	1:B:181:ARG:CB	2.40	0.68
1:B:64:VAL:HG23	1:B:66:PRO:CD	2.15	0.68
1:B:48:ASP:OD1	1:B:48:ASP:C	2.29	0.68
1:D:211:ILE:HD11	5:D:901:EDO:H11	1.76	0.67
1:A:131:GLY:H	1:A:134:GLN:HE21	1.43	0.67
1:A:3:LEU:O	1:A:3:LEU:CD1	2.36	0.67
3:A:611:TMQ:H232	3:A:611:TMQ:H223	1.77	0.66
2:B:602:NAP:C5N	3:B:612:TMQ:H172	2.26	0.66
1:B:45:VAL:CG2	1:B:180:ILE:HD11	2.19	0.65
1:D:78:ARG:HH11	2:D:604:NAP:P2B	2.19	0.64
2:A:601:NAP:H4N	3:A:611:TMQ:H172	1.78	0.64
1:A:79:LYS:HG3	2:A:601:NAP:C5B	2.27	0.64
1:B:47:GLU:HG3	1:B:217:SER:HG	0.75	0.64
1:C:39:ARG:HH12	1:C:79:LYS:HD3	1.63	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:184:GLU:HG3	1:D:186:SER:HB3	1.80	0.64
1:B:379:ASN:HB3	6:B:560:HOH:O	1.97	0.63
1:B:213:GLU:HG2	1:B:214:GLU:N	2.13	0.63
1:B:80:THR:HG21	1:B:155:GLY:HA3	1.81	0.62
1:B:181:ARG:HG2	1:B:224:LYS:O	1.98	0.62
1:D:347:TRP:HB3	1:D:369:ILE:HD13	1.80	0.62
1:B:20:LEU:HD23	1:B:137:GLN:HG3	1.81	0.62
3:C:613:TMQ:N25	3:C:613:TMQ:H173	2.12	0.62
1:B:76:MET:HB3	1:B:154:ILE:HD11	1.81	0.62
1:B:161:ALA:O	1:B:165:ARG:HG2	1.99	0.62
1:D:101:SER:N	2:D:604:NAP:O2X	2.29	0.61
1:B:154:ILE:O	3:B:612:TMQ:N25	2.33	0.61
3:A:611:TMQ:H172	3:A:611:TMQ:N25	2.12	0.61
1:C:112:LEU:HB2	1:C:113:PRO:CA	2.31	0.61
1:B:29:VAL:O	1:B:180:ILE:N	2.30	0.60
1:D:471:GLU:HB3	1:D:472:PRO:HD3	1.83	0.60
3:D:614:TMQ:H173	3:D:614:TMQ:N25	2.12	0.60
1:B:43:TRP:CD1	1:B:45:VAL:CG2	2.84	0.60
1:B:78:ARG:HB3	2:B:602:NAP:O3B	2.02	0.59
1:C:71:ARG:HH12	1:C:124:SER:HB3	1.67	0.59
1:B:181:ARG:CG	1:B:224:LYS:O	2.51	0.59
1:B:47:GLU:HG2	1:B:217:SER:OG	1.92	0.59
1:B:296:LEU:HD22	1:B:440:LEU:HB3	1.85	0.58
1:D:296:LEU:HD22	1:D:440:LEU:HB3	1.85	0.58
3:A:611:TMQ:H173	3:A:611:TMQ:N25	2.12	0.58
1:B:65:LYS:H	1:B:66:PRO:HD3	1.68	0.58
1:D:38:GLY:HA2	2:D:604:NAP:H4D	1.85	0.58
3:D:614:TMQ:H172	3:D:614:TMQ:N25	2.12	0.58
3:C:613:TMQ:N25	3:C:613:TMQ:H172	2.12	0.57
1:A:407:ALA:HA	1:A:419:MET:O	2.05	0.57
1:D:184:GLU:CG	1:D:186:SER:HB3	2.35	0.57
1:D:100:SER:HA	2:D:604:NAP:O2X	2.05	0.57
1:B:49:MET:O	1:B:51:PHE:N	2.39	0.56
1:A:10:GLU:H	1:A:10:GLU:CD	2.09	0.56
3:B:612:TMQ:N25	3:B:612:TMQ:H173	2.12	0.56
1:B:74:VAL:HB	1:B:154:ILE:HD12	1.88	0.56
1:C:422:GLN:NE2	1:C:458:HIS:NE2	2.54	0.55
1:A:79:LYS:CG	2:A:601:NAP:H51A	2.33	0.55
1:B:239:GLN:HE22	1:B:271:ARG:H	1.54	0.55
1:C:156:GLY:HA2	2:C:603:NAP:PA	2.47	0.55
1:A:515:ILE:C	1:A:515:ILE:HD13	2.27	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:LEU:O	1:A:110:ASP:HB2	2.05	0.55
1:B:329:ARG:NH1	1:B:398:MET:O	2.40	0.55
1:B:46:PRO:C	1:B:48:ASP:H	2.09	0.55
1:D:254:LYS:HB2	1:D:262:THR:HG22	1.88	0.55
1:A:306:ALA:HB2	1:A:339:LEU:HD11	1.89	0.54
1:C:156:GLY:O	1:C:158:SER:N	2.40	0.54
1:C:146:PRO:HG3	1:C:505:GLU:HG2	1.89	0.54
1:B:64:VAL:O	1:B:65:LYS:CB	2.54	0.54
1:B:64:VAL:C	1:B:65:LYS:HG2	2.27	0.53
1:C:99:LEU:O	2:C:603:NAP:H1B	2.08	0.53
2:C:603:NAP:H4N	3:C:613:TMQ:H172	1.90	0.53
1:C:156:GLY:CA	2:C:603:NAP:PA	2.96	0.53
2:A:601:NAP:C5N	3:A:611:TMQ:H172	2.39	0.53
1:B:182:ALA:H	1:B:223:THR:HG23	1.73	0.53
1:C:32:ARG:O	1:C:192:ARG:NH2	2.42	0.53
3:B:612:TMQ:N25	3:B:612:TMQ:H172	2.12	0.53
1:C:20:LEU:HB2	1:C:171:LEU:CD1	2.39	0.53
1:C:14:GLU:OE1	1:C:14:GLU:HA	2.08	0.52
1:B:471:GLU:HB2	1:B:472:PRO:HD3	1.90	0.52
1:B:131:GLY:N	1:B:134:GLN:HE21	1.96	0.52
1:B:392:PRO:HA	1:B:395:LEU:HD22	1.91	0.52
1:A:472:PRO:HB3	1:A:515:ILE:HD12	1.91	0.52
1:C:156:GLY:O	1:C:159:VAL:N	2.43	0.52
1:B:320:ASN:N	1:B:320:ASN:HD22	2.07	0.52
1:B:64:VAL:CG2	1:B:66:PRO:CD	2.59	0.52
1:A:131:GLY:N	1:A:134:GLN:HE21	2.06	0.52
1:A:278:ARG:HH21	1:A:484:PHE:HD2	1.56	0.52
1:B:60:ARG:NH2	6:B:552:HOH:O	2.42	0.52
2:D:604:NAP:H4N	3:D:614:TMQ:H172	1.92	0.52
1:C:78:ARG:HD2	1:C:103:LEU:HD23	1.92	0.51
1:A:69:ALA:O	1:A:147:SER:HA	2.09	0.51
1:C:234:ASN:HB3	1:C:481:PRO:HB2	1.92	0.51
1:B:215:LEU:CD1	1:B:216:THR:N	2.54	0.51
1:B:47:GLU:O	1:B:47:GLU:OE1	2.29	0.51
1:C:105:THR:HB	1:C:127:ALA:H	1.74	0.51
1:B:65:LYS:O	1:B:67:SER:N	2.43	0.51
1:C:112:LEU:HB2	1:C:113:PRO:HA	1.91	0.51
1:D:248:ILE:HD13	1:D:470:VAL:HG22	1.93	0.51
1:C:297:LEU:HB3	1:C:301:ARG:HH21	1.76	0.51
1:D:6:ILE:HG13	1:D:496:LEU:HD23	1.93	0.51
1:A:104:THR:H	1:A:107:HIS:HB2	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:GLU:O	1:B:32:ARG:CB	2.51	0.50
1:B:64:VAL:O	1:B:65:LYS:HB3	2.11	0.50
1:B:252:ASN:HD22	5:B:906:EDO:H12	1.77	0.50
1:D:336:GLU:O	1:D:337:MET:HB2	2.12	0.50
1:D:514:PRO:O	1:D:515:ILE:HB	2.12	0.50
1:A:105:THR:O	1:A:109:LEU:HD23	2.11	0.50
1:B:180:ILE:O	1:B:224:LYS:O	2.29	0.50
1:B:47:GLU:OE1	1:B:47:GLU:CA	2.60	0.50
3:C:613:TMQ:O20	3:C:613:TMQ:H223	2.11	0.50
1:B:185:SER:O	1:B:187:CYS:N	2.39	0.50
1:C:64:VAL:HG12	1:C:66:PRO:HD3	1.93	0.49
1:D:157:GLY:HA3	1:D:189:VAL:HG11	1.93	0.49
1:D:404:HIS:HB2	1:D:420:LEU:HD11	1.93	0.49
1:A:49:MET:HE1	3:A:611:TMQ:C13	2.42	0.49
1:C:157:GLY:HA3	1:C:189:VAL:HG11	1.93	0.49
1:B:479:ARG:NH2	1:B:513:PRO:O	2.45	0.49
1:C:404:HIS:HB2	1:C:420:LEU:HD11	1.93	0.49
1:B:131:GLY:H	1:B:134:GLN:NE2	1.95	0.49
1:B:180:ILE:O	1:B:181:ARG:HB2	2.12	0.49
1:D:156:GLY:HA3	2:D:604:NAP:PA	2.53	0.48
1:A:61:GLY:HA3	1:A:64:VAL:CG2	2.43	0.48
1:B:30:ASP:CA	1:B:180:ILE:CG2	2.79	0.48
1:C:174:ALA:HB2	5:C:910:EDO:H21	1.95	0.48
1:D:425:CYS:HB3	1:D:431:VAL:CG2	2.43	0.48
1:D:234:ASN:HB3	1:D:481:PRO:HB2	1.94	0.48
2:D:604:NAP:O2N	2:D:604:NAP:O2A	2.30	0.48
1:A:61:GLY:HA3	1:A:64:VAL:HG23	1.96	0.48
1:C:92:PRO:O	1:C:94:ARG:NH1	2.46	0.48
1:A:256:ASP:HB3	1:D:383:ARG:NH2	2.29	0.47
1:B:17:ARG:HD2	1:B:276:ASN:HB3	1.95	0.47
1:A:420:LEU:HD23	1:A:458:HIS:CD2	2.48	0.47
1:D:10:GLU:H	1:D:10:GLU:CD	2.17	0.47
1:A:392:PRO:HA	1:A:395:LEU:HD22	1.95	0.47
1:C:207:GLN:HE22	1:D:249:ARG:HH11	1.61	0.47
1:C:20:LEU:HB2	1:C:171:LEU:HD13	1.96	0.47
1:A:443:ILE:O	1:A:447:LYS:HG3	2.15	0.47
1:A:425:CYS:HB2	1:A:463:ALA:HA	1.96	0.47
1:C:333:GLU:HA	1:C:333:GLU:OE1	2.14	0.47
1:A:334:TYR:HB3	1:A:338:ASP:HB3	1.96	0.47
1:A:415:GLU:HA	1:A:452:ARG:O	2.14	0.47
1:C:99:LEU:HD21	1:C:135:ALA:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LYS:HE3	1:B:70:LYS:HD3	1.96	0.47
1:C:303:GLU:HG2	1:C:308:LYS:HD2	1.97	0.47
1:D:30:ASP:HB2	1:D:182:ALA:O	2.15	0.47
1:C:358:ASP:HB2	6:C:555:HOH:O	2.13	0.47
1:B:139:LEU:HD22	1:B:148:ILE:HG21	1.97	0.46
1:D:236:GLU:O	1:D:239:GLN:HB2	2.13	0.46
1:D:320:ASN:N	1:D:320:ASN:HD22	2.13	0.46
1:A:156:GLY:HA3	2:A:601:NAP:PA	2.55	0.46
1:B:78:ARG:N	2:B:602:NAP:H4B	2.31	0.46
1:B:141:SER:HB2	1:B:142:PRO:HD2	1.97	0.46
1:C:342:VAL:HG12	1:C:398:MET:HB3	1.97	0.46
1:B:490:ARG:HD2	1:B:505:GLU:OE1	2.16	0.46
1:A:404:HIS:H	1:A:404:HIS:CD2	2.34	0.46
1:C:175:ILE:HB	1:C:230:LEU:HB2	1.98	0.46
1:C:100:SER:CB	2:C:603:NAP:O1X	2.63	0.46
1:A:239:GLN:HE22	1:A:271:ARG:H	1.62	0.46
1:B:48:ASP:O	1:B:49:MET:C	2.54	0.46
1:B:62:LYS:HD3	1:B:62:LYS:HA	1.49	0.46
1:C:67:SER:HB2	1:C:68:PRO:HD2	1.98	0.46
1:B:65:LYS:N	1:B:66:PRO:HD3	2.29	0.46
1:C:254:LYS:HB2	1:C:262:THR:HG22	1.97	0.46
1:B:217:SER:O	1:B:219:ASN:N	2.49	0.45
1:B:407:ALA:HA	1:B:419:MET:O	2.17	0.45
1:B:78:ARG:H	2:B:602:NAP:H4B	1.81	0.45
1:D:289:TRP:HH2	1:D:440:LEU:HG	1.81	0.45
3:A:611:TMQ:H232	3:A:611:TMQ:C22	2.45	0.45
1:C:10:GLU:CD	1:C:10:GLU:H	2.20	0.45
1:C:407:ALA:HA	1:C:419:MET:O	2.16	0.45
1:B:405:LEU:HD12	1:B:406:LEU:HB2	1.98	0.45
1:B:49:MET:O	1:B:50:LYS:C	2.55	0.45
1:A:299:PHE:HB3	1:A:347:TRP:CZ3	2.53	0.45
1:B:49:MET:C	1:B:51:PHE:N	2.70	0.44
1:C:281:LEU:HD22	1:C:287:VAL:HB	2.00	0.44
1:C:336:GLU:O	1:C:337:MET:HB2	2.16	0.44
1:B:91:LEU:HA	1:B:92:PRO:HD3	1.85	0.44
1:D:22:ALA:HB2	1:D:486:TYR:CE1	2.53	0.44
1:B:316:ILE:HG13	1:B:317:TRP:CD1	2.52	0.44
1:D:275:ARG:O	1:D:278:ARG:HB2	2.18	0.44
1:A:109:LEU:O	1:A:110:ASP:CB	2.66	0.44
1:A:16:THR:HA	1:A:452:ARG:HH21	1.83	0.44
3:B:612:TMQ:H92	3:B:612:TMQ:H171	1.72	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:603:NAP:C4N	3:C:613:TMQ:H172	2.48	0.44
1:D:175:ILE:HB	1:D:230:LEU:HB2	2.00	0.44
1:C:469:HIS:HD2	6:C:557:HOH:O	2.00	0.43
1:A:131:GLY:H	1:A:134:GLN:NE2	2.12	0.43
1:A:49:MET:HE2	3:A:611:TMQ:H222	2.00	0.43
1:A:81:TRP:CE2	1:A:89:ARG:HD3	2.53	0.43
1:B:48:ASP:O	1:B:51:PHE:HB3	2.18	0.43
1:B:76:MET:HB3	1:B:154:ILE:CD1	2.49	0.43
1:C:215:LEU:O	1:C:224:LYS:HA	2.18	0.43
1:C:473:CYS:O	1:C:477:LEU:HG	2.18	0.43
1:D:478:LYS:HA	1:D:478:LYS:HD3	1.70	0.43
1:C:305:TYR:OH	1:C:307:LYS:HE3	2.18	0.43
1:B:33:GLY:O	1:B:191:PHE:N	2.52	0.43
1:B:48:ASP:OD1	1:B:49:MET:N	2.52	0.43
1:B:336:GLU:O	1:B:337:MET:HB2	2.19	0.43
1:B:41:ILE:HA	1:B:42:PRO:HD3	1.83	0.42
2:C:603:NAP:O2A	2:C:603:NAP:O2N	2.30	0.42
1:B:406:LEU:HD11	1:C:405:LEU:HD11	2.01	0.42
1:D:316:ILE:HG13	1:D:317:TRP:CD1	2.54	0.42
1:D:404:HIS:CB	1:D:420:LEU:HD11	2.49	0.42
3:D:614:TMQ:H223	3:D:614:TMQ:C23	2.46	0.42
1:C:467:SER:OG	5:D:901:EDO:H21	2.18	0.42
1:A:96:ASN:HD22	1:A:96:ASN:N	2.17	0.42
1:B:468:ASN:HD22	1:B:468:ASN:C	2.23	0.42
2:B:602:NAP:H4N	3:B:612:TMQ:N25	2.35	0.42
3:D:614:TMQ:H92	3:D:614:TMQ:H171	1.70	0.42
1:D:99:LEU:HD21	1:D:135:ALA:HB2	2.00	0.42
1:B:289:TRP:HH2	1:B:440:LEU:HG	1.84	0.42
1:D:155:GLY:HA3	1:D:160:TYR:CZ	2.54	0.42
1:B:180:ILE:HA	1:B:180:ILE:HD12	1.89	0.42
1:C:425:CYS:HB3	1:C:431:VAL:HG22	2.01	0.42
2:D:604:NAP:C4N	3:D:614:TMQ:H172	2.49	0.41
1:B:177:ARG:HH21	1:B:228:GLU:CD	2.23	0.41
1:A:405:LEU:HD11	1:D:406:LEU:HD11	2.02	0.41
1:D:420:LEU:HD13	1:D:438:TYR:CZ	2.56	0.41
1:D:100:SER:CA	2:D:604:NAP:O2X	2.68	0.41
1:A:103:LEU:HD12	1:A:107:HIS:HB3	2.02	0.41
1:B:368:GLN:O	1:B:372:ILE:HD12	2.21	0.41
1:C:41:ILE:HA	1:C:42:PRO:HD3	1.92	0.41
1:A:115:GLU:O	1:A:121:HIS:HD2	2.04	0.41
1:B:415:GLU:HA	1:B:452:ARG:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:356:HIS:HB2	6:D:542:HOH:O	2.20	0.41
1:A:392:PRO:HD2	1:D:350:PHE:CZ	2.56	0.41
1:C:350:PHE:CE2	1:C:391:ASN:HB2	2.56	0.41
2:C:603:NAP:H8A	2:C:603:NAP:H52A	2.03	0.41
1:B:194:PRO:HG3	6:B:555:HOH:O	2.20	0.40
1:B:4:PHE:CD1	1:B:4:PHE:N	2.89	0.40
1:B:49:MET:O	1:B:52:PHE:N	2.54	0.40
1:A:91:LEU:HA	1:A:92:PRO:HD3	1.85	0.40
1:C:425:CYS:HB3	1:C:431:VAL:CG2	2.52	0.40
1:D:281:LEU:HD22	1:D:287:VAL:HB	2.03	0.40
1:A:336:GLU:O	1:A:337:MET:HB2	2.21	0.40
1:A:49:MET:HE1	3:A:611:TMQ:O18	2.21	0.40
1:B:342:VAL:HG12	1:B:398:MET:HB3	2.03	0.40
1:C:331:LEU:HD22	1:C:334:TYR:CE2	2.55	0.40

All (4) 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)
1:A:3:LEU:CG	1:D:39:ARG:CB[4_555]	1.85	0.35
1:A:3:LEU:CD1	1:D:39:ARG:CB[4_555]	2.05	0.15
1:A:3:LEU:CD2	1:D:39:ARG:CB[4_555]	2.14	0.06
1:A:3:LEU:CD2	1:D:39:ARG:CG[4_555]	2.18	0.02

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	506/521 (97%)	490 (97%)	13 (3%)	3 (1%)	25	64
1	B	501/521 (96%)	458 (91%)	24 (5%)	19 (4%)	3	18
1	C	506/521 (97%)	473 (94%)	25 (5%)	8 (2%)	9	40

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	D	503/521 (96%)	484 (96%)	18 (4%)	1 (0%)	47 82
All	All	2016/2084 (97%)	1905 (94%)	80 (4%)	31 (2%)	10 42

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	343	TYR
1	B	46	PRO
1	B	60	ARG
1	B	62	LYS
1	B	65	LYS
1	B	180	ILE
1	B	181	ARG
1	B	182	ALA
1	B	183	SER
1	B	186	SER
1	B	218	ALA
1	C	343	TYR
1	D	343	TYR
1	B	32	ARG
1	B	64	VAL
1	B	122	ALA
1	B	343	TYR
1	C	122	ALA
1	C	186	SER
1	B	49	MET
1	B	50	LYS
1	C	2	SER
1	A	110	ASP
1	B	66	PRO
1	C	112	LEU
1	B	63	ASN
1	B	219	ASN
1	C	62	LYS
1	C	64	VAL
1	C	111	GLY
1	A	146	PRO

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	436/446 (98%)	415 (95%)	21 (5%)	25	62
1	B	430/446 (96%)	395 (92%)	35 (8%)	11	40
1	C	435/446 (98%)	408 (94%)	27 (6%)	18	52
1	D	431/446 (97%)	408 (95%)	23 (5%)	22	58
All	All	1732/1784 (97%)	1626 (94%)	106 (6%)	18	53

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	THR
1	A	12	VAL
1	A	17	ARG
1	A	95	LEU
1	A	102	THR
1	A	124	SER
1	A	138	LEU
1	A	147	SER
1	A	183	SER
1	A	184	GLU
1	A	187	CYS
1	A	222	GLU
1	A	264	SER
1	A	274	LEU
1	A	395	LEU
1	A	419	MET
1	A	440	LEU
1	A	452	ARG
1	A	456	LEU
1	A	467	SER
1	A	515	ILE
1	B	3	LEU
1	B	10	GLU
1	B	18	LEU

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Mol	Chain	Res	Type
1	B	29	VAL
1	B	32	ARG
1	B	47	GLU
1	B	48	ASP
1	B	62	LYS
1	B	64	VAL
1	B	94	ARG
1	B	95	LEU
1	B	108	LEU
1	B	109	LEU
1	B	125	ILE
1	B	154	ILE
1	B	165	ARG
1	B	184	GLU
1	B	213	GLU
1	B	215	LEU
1	B	239	GLN
1	B	256	ASP
1	B	274	LEU
1	B	297	LEU
1	B	309	LEU
1	B	320	ASN
1	B	329	ARG
1	B	333	GLU
1	B	372	ILE
1	B	395	LEU
1	B	412	SER
1	B	419	MET
1	B	440	LEU
1	B	456	LEU
1	B	468	ASN
1	B	480	VAL
1	C	1	MET
1	C	3	LEU
1	C	18	LEU
1	C	41	ILE
1	C	70	LYS
1	C	94	ARG
1	C	95	LEU
1	C	105	THR
1	C	112	LEU
1	C	114	ASP

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Mol	Chain	Res	Type
1	C	115	GLU
1	C	123	ASP
1	C	125	ILE
1	C	173	GLN
1	C	189	VAL
1	C	196	SER
1	C	262	THR
1	C	274	LEU
1	C	286	ARG
1	C	297	LEU
1	C	303	GLU
1	C	309	LEU
1	C	358	ASP
1	C	386	LEU
1	C	397	ARG
1	C	419	MET
1	C	507	ILE
1	D	6	ILE
1	D	10	GLU
1	D	14	GLU
1	D	39	ARG
1	D	70	LYS
1	D	95	LEU
1	D	143	ASN
1	D	173	GLN
1	D	189	VAL
1	D	248	ILE
1	D	262	THR
1	D	265	ILE
1	D	278	ARG
1	D	297	LEU
1	D	303	GLU
1	D	309	LEU
1	D	320	ASN
1	D	323	ARG
1	D	429	LEU
1	D	440	LEU
1	D	456	LEU
1	D	467	SER
1	D	493	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	107	HIS
1	A	121	HIS
1	A	134	GLN
1	A	143	ASN
1	A	173	GLN
1	A	239	GLN
1	A	356	HIS
1	A	379	ASN
1	A	422	GLN
1	A	464	HIS
1	B	134	GLN
1	B	173	GLN
1	B	207	GLN
1	B	239	GLN
1	B	252	ASN
1	B	320	ASN
1	B	379	ASN
1	B	468	ASN
1	B	469	HIS
1	C	107	HIS
1	C	137	GLN
1	C	207	GLN
1	C	239	GLN
1	C	255	HIS
1	C	320	ASN
1	C	422	GLN
1	C	434	ASN
1	C	464	HIS
1	D	121	HIS
1	D	143	ASN
1	D	360	ASN
1	D	422	GLN
1	D	434	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry

48 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
5	EDO	A	904	-	3,3,3	0.43	0	2,2,2	0.37	0
4	SO4	D	810	-	4,4,4	0.13	0	6,6,6	0.06	0
5	EDO	C	910	-	3,3,3	0.48	0	2,2,2	0.36	0
2	NAP	A	601	-	45,52,52	1.19	5 (11%)	56,80,80	1.77	9 (16%)
5	EDO	D	901	-	3,3,3	0.44	0	2,2,2	0.34	0
5	EDO	C	908	-	3,3,3	0.49	0	2,2,2	0.32	0
5	EDO	D	905	-	3,3,3	0.49	0	2,2,2	0.24	0
4	SO4	B	809	-	4,4,4	0.12	0	6,6,6	0.09	0
5	EDO	D	911	-	3,3,3	0.41	0	2,2,2	0.43	0
4	SO4	B	823	-	4,4,4	0.13	0	6,6,6	0.15	0
5	EDO	A	902	-	3,3,3	0.52	0	2,2,2	0.21	0
4	SO4	D	811	-	4,4,4	0.14	0	6,6,6	0.10	0
5	EDO	B	906	-	3,3,3	0.49	0	2,2,2	0.36	0
4	SO4	A	805	-	4,4,4	0.16	0	6,6,6	0.12	0
4	SO4	C	812	-	4,4,4	0.14	0	6,6,6	0.11	0
3	TMQ	B	612	-	29,29,29	1.48	4 (13%)	39,41,41	2.02	7 (17%)
4	SO4	A	701	-	4,4,4	0.13	0	6,6,6	0.27	0
4	SO4	B	821	-	4,4,4	0.14	0	6,6,6	0.06	0
2	NAP	B	602	-	45,52,52	1.20	5 (11%)	56,80,80	1.77	10 (17%)
4	SO4	D	819	-	4,4,4	0.12	0	6,6,6	0.13	0
4	SO4	D	801	-	4,4,4	0.14	0	6,6,6	0.06	0
4	SO4	A	817	-	4,4,4	0.13	0	6,6,6	0.12	0
4	SO4	C	808	-	4,4,4	0.14	0	6,6,6	0.12	0
4	SO4	C	802	-	4,4,4	0.13	0	6,6,6	0.10	0
2	NAP	D	604	-	45,52,52	1.40	6 (13%)	56,80,80	2.21	14 (25%)
4	SO4	D	803	-	4,4,4	0.14	0	6,6,6	0.08	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	A	813	-	4,4,4	0.16	0	6,6,6	0.14	0
4	SO4	C	822	-	4,4,4	0.13	0	6,6,6	0.12	0
4	SO4	C	824	-	4,4,4	0.12	0	6,6,6	0.16	0
4	SO4	A	814	-	4,4,4	0.15	0	6,6,6	0.09	0
4	SO4	C	807	-	4,4,4	0.13	0	6,6,6	0.07	0
3	TMQ	D	614	-	29,29,29	1.48	4 (13%)	39,41,41	2.01	7 (17%)
2	NAP	C	603	-	45,52,52	1.36	7 (15%)	56,80,80	2.09	11 (19%)
4	SO4	A	815	-	4,4,4	0.13	0	6,6,6	0.08	0
4	SO4	D	820	-	4,4,4	0.15	0	6,6,6	0.15	0
3	TMQ	C	613	-	29,29,29	1.47	4 (13%)	39,41,41	2.02	7 (17%)
5	EDO	C	912	-	3,3,3	0.56	0	2,2,2	0.12	0
4	SO4	D	704	-	4,4,4	0.11	0	6,6,6	0.18	0
3	TMQ	A	611	-	29,29,29	1.48	4 (13%)	39,41,41	2.01	7 (17%)
4	SO4	D	804	-	4,4,4	0.12	0	6,6,6	0.12	0
4	SO4	C	702	-	4,4,4	0.19	0	6,6,6	0.19	0
5	EDO	C	909	-	3,3,3	0.49	0	2,2,2	0.28	0
5	EDO	D	903	-	3,3,3	0.45	0	2,2,2	0.31	0
4	SO4	B	818	-	4,4,4	0.14	0	6,6,6	0.07	0
4	SO4	B	806	-	4,4,4	0.15	0	6,6,6	0.16	0
4	SO4	C	703	-	4,4,4	0.16	0	6,6,6	0.21	0
5	EDO	D	907	-	3,3,3	0.40	0	2,2,2	0.52	0
4	SO4	C	816	-	4,4,4	0.14	0	6,6,6	0.20	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
5	EDO	A	904	-	-	1/1/1/1	-
5	EDO	C	912	-	-	1/1/1/1	-
3	TMQ	B	612	-	-	4/11/11/11	0/3/3/3
2	NAP	A	601	-	-	10/31/67/67	0/5/5/5
5	EDO	D	901	-	-	0/1/1/1	-
5	EDO	C	908	-	-	1/1/1/1	-
5	EDO	C	910	-	-	1/1/1/1	-
5	EDO	D	905	-	-	1/1/1/1	-
5	EDO	D	911	-	-	1/1/1/1	-
2	NAP	B	602	-	-	3/31/67/67	0/5/5/5
5	EDO	A	902	-	-	0/1/1/1	-
3	TMQ	D	614	-	-	7/11/11/11	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAP	C	603	-	-	4/31/67/67	0/5/5/5
5	EDO	D	903	-	-	0/1/1/1	-
5	EDO	B	906	-	-	0/1/1/1	-
3	TMQ	A	611	-	-	4/11/11/11	0/3/3/3
5	EDO	D	907	-	-	1/1/1/1	-
3	TMQ	C	613	-	-	2/11/11/11	0/3/3/3
2	NAP	D	604	-	-	6/31/67/67	0/5/5/5
5	EDO	C	909	-	-	1/1/1/1	-

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	611	TMQ	C17-C5	-5.11	1.39	1.51
3	B	612	TMQ	C17-C5	-5.09	1.39	1.51
3	D	614	TMQ	C17-C5	-5.08	1.39	1.51
3	C	613	TMQ	C17-C5	-5.08	1.39	1.51
2	A	601	NAP	C2N-N1N	4.51	1.40	1.35
2	B	602	NAP	C2N-N1N	4.50	1.40	1.35
3	B	612	TMQ	C9-C6	-3.51	1.39	1.50
3	C	613	TMQ	C9-C6	-3.50	1.39	1.50
3	A	611	TMQ	C9-C6	-3.50	1.39	1.50
3	D	614	TMQ	C9-C6	-3.49	1.39	1.50
2	D	604	NAP	C2D-C1D	-3.48	1.48	1.53
2	D	604	NAP	C5B-C4B	2.85	1.60	1.51
2	C	603	NAP	C2D-C1D	-2.75	1.49	1.53
2	D	604	NAP	O4B-C1B	2.63	1.44	1.41
2	C	603	NAP	O4D-C4D	-2.61	1.39	1.45
2	D	604	NAP	C2D-C3D	-2.50	1.46	1.53
2	C	603	NAP	O4B-C1B	2.42	1.44	1.41
2	C	603	NAP	PN-O2N	-2.30	1.44	1.55
2	C	603	NAP	C5A-N7A	-2.24	1.31	1.39
2	D	604	NAP	C5A-N7A	-2.23	1.31	1.39
3	A	611	TMQ	C4A-C3A	-2.20	1.39	1.42
2	C	603	NAP	C4A-N3A	-2.20	1.32	1.35
3	C	613	TMQ	C4A-C3A	-2.20	1.39	1.42
3	D	614	TMQ	C5-C4A	-2.20	1.39	1.44
3	B	612	TMQ	C5-C4A	-2.20	1.39	1.44
3	B	612	TMQ	C4A-C3A	-2.19	1.39	1.42
3	C	613	TMQ	C5-C4A	-2.18	1.39	1.44
3	D	614	TMQ	C4A-C3A	-2.18	1.39	1.42
2	A	601	NAP	PA-O2A	2.16	1.65	1.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	602	NAP	PA-O2A	2.16	1.65	1.55
3	A	611	TMQ	C5-C4A	-2.16	1.39	1.44
2	B	602	NAP	O4B-C1B	2.15	1.44	1.41
2	A	601	NAP	O4B-C1B	2.12	1.44	1.41
2	D	604	NAP	PN-O2N	-2.11	1.45	1.55
2	A	601	NAP	O4D-C1D	2.10	1.44	1.41
2	B	602	NAP	O4D-C1D	2.10	1.44	1.41
2	C	603	NAP	C5B-C4B	2.07	1.58	1.51
2	B	602	NAP	C6N-N1N	2.06	1.40	1.35
2	A	601	NAP	C6N-N1N	2.03	1.40	1.35

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	603	NAP	O5B-PA-O1A	-7.64	79.20	109.07
2	D	604	NAP	O2A-PA-O5B	-7.27	74.00	107.75
2	D	604	NAP	O5B-PA-O1A	-7.26	80.70	109.07
2	C	603	NAP	O2A-PA-O5B	-6.97	75.39	107.75
2	B	602	NAP	O5B-PA-O1A	-6.41	84.01	109.07
2	A	601	NAP	O5B-PA-O1A	-6.41	84.02	109.07
2	A	601	NAP	O2A-PA-O5B	-6.34	78.28	107.75
2	B	602	NAP	O2A-PA-O5B	-6.34	78.29	107.75
3	B	612	TMQ	C23-O20-C15	-5.35	109.45	117.53
3	C	613	TMQ	C21-O18-C13	-5.35	109.45	117.53
3	B	612	TMQ	C21-O18-C13	-5.34	109.47	117.53
3	D	614	TMQ	C23-O20-C15	-5.33	109.48	117.53
3	C	613	TMQ	C23-O20-C15	-5.33	109.49	117.53
3	A	611	TMQ	C21-O18-C13	-5.32	109.50	117.53
3	D	614	TMQ	C21-O18-C13	-5.31	109.51	117.53
3	A	611	TMQ	C23-O20-C15	-5.30	109.53	117.53
3	C	613	TMQ	C4-C4A-C3A	5.16	118.37	114.86
3	B	612	TMQ	C4-C4A-C3A	5.13	118.35	114.86
3	A	611	TMQ	C4-C4A-C3A	5.12	118.35	114.86
3	D	614	TMQ	C4-C4A-C3A	5.09	118.32	114.86
2	D	604	NAP	O4D-C1D-C2D	-4.82	99.89	106.93
3	D	614	TMQ	N1-C2-N3	-4.70	120.96	127.22
3	B	612	TMQ	N1-C2-N3	-4.68	120.98	127.22
3	C	613	TMQ	N1-C2-N3	-4.68	120.98	127.22
2	D	604	NAP	N3A-C2A-N1A	-4.68	121.37	128.68
3	A	611	TMQ	N1-C2-N3	-4.67	121.00	127.22
2	D	604	NAP	C3D-C2D-C1D	-4.66	93.96	100.98
2	D	604	NAP	PN-O3-PA	-4.65	116.86	132.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	603	NAP	N3A-C2A-N1A	-4.62	121.46	128.68
2	B	602	NAP	N3A-C2A-N1A	-4.46	121.71	128.68
2	A	601	NAP	N3A-C2A-N1A	-4.45	121.73	128.68
2	C	603	NAP	O4D-C1D-C2D	-4.06	101.00	106.93
2	B	602	NAP	PN-O3-PA	-3.75	119.95	132.83
2	A	601	NAP	PN-O3-PA	-3.75	119.96	132.83
2	C	603	NAP	PN-O3-PA	-3.64	120.34	132.83
3	A	611	TMQ	C4A-C4-N25	-3.22	116.86	122.67
3	B	612	TMQ	C4A-C4-N25	-3.21	116.88	122.67
3	D	614	TMQ	C4A-C4-N25	-3.20	116.89	122.67
3	C	613	TMQ	C4A-C4-N25	-3.20	116.89	122.67
2	D	604	NAP	O3D-C3D-C2D	-3.05	101.95	111.82
3	C	613	TMQ	C5-C4A-C4	-2.72	121.18	127.18
3	A	611	TMQ	C5-C4A-C4	-2.72	121.19	127.18
3	D	614	TMQ	C5-C4A-C4	-2.71	121.20	127.18
3	B	612	TMQ	C5-C4A-C4	-2.70	121.22	127.18
2	C	603	NAP	C2N-C3N-C4N	2.65	121.26	118.26
2	D	604	NAP	O5B-C5B-C4B	2.65	118.10	108.99
2	C	603	NAP	O5B-C5B-C4B	2.43	117.34	108.99
2	B	602	NAP	C6N-N1N-C2N	-2.38	119.80	121.97
2	C	603	NAP	O2N-PN-O5D	2.38	118.78	107.75
2	A	601	NAP	C6N-N1N-C2N	-2.37	119.81	121.97
2	A	601	NAP	C3N-C7N-N7N	-2.31	114.98	117.75
2	B	602	NAP	C3N-C7N-N7N	-2.31	114.98	117.75
2	D	604	NAP	C3N-C7N-N7N	-2.27	115.03	117.75
2	C	603	NAP	C3B-C2B-C1B	-2.24	98.68	102.89
2	D	604	NAP	O2N-PN-O1N	2.21	123.19	112.24
2	D	604	NAP	O2N-PN-O5D	2.20	117.96	107.75
2	A	601	NAP	O2N-PN-O1N	2.18	123.01	112.24
2	D	604	NAP	C2N-C3N-C4N	2.18	120.72	118.26
2	C	603	NAP	O2N-PN-O1N	2.17	122.99	112.24
2	B	602	NAP	O2N-PN-O1N	2.17	122.97	112.24
2	B	602	NAP	O5D-PN-O1N	2.17	117.54	109.07
2	A	601	NAP	O5D-PN-O1N	2.17	117.53	109.07
2	C	603	NAP	C3N-C7N-N7N	-2.12	115.21	117.75
2	B	602	NAP	O2N-PN-O5D	2.11	117.55	107.75
2	A	601	NAP	O2N-PN-O5D	2.10	117.50	107.75
3	C	613	TMQ	N25-C4-N3	2.07	122.71	117.07
3	A	611	TMQ	N25-C4-N3	2.06	122.68	117.07
3	D	614	TMQ	N25-C4-N3	2.05	122.65	117.07
2	D	604	NAP	O4B-C4B-C5B	2.05	116.12	109.37
3	B	612	TMQ	N25-C4-N3	2.05	122.64	117.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	604	NAP	O5D-PN-O1N	2.04	117.03	109.07
2	B	602	NAP	O4D-C1D-C2D	-2.01	103.98	106.93

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	NAP	C2B-O2B-P2B-O1X
2	A	601	NAP	C5D-O5D-PN-O1N
2	D	604	NAP	C5D-O5D-PN-O1N
2	D	604	NAP	O4D-C1D-N1N-C6N
3	D	614	TMQ	C14-C15-O20-C23
3	B	612	TMQ	C12-C11-N10-C9
3	A	611	TMQ	C12-C11-N10-C9
3	A	611	TMQ	C16-C11-N10-C9
2	A	601	NAP	C3D-C4D-C5D-O5D
2	B	602	NAP	C3B-C4B-C5B-O5B
3	D	614	TMQ	C12-C13-O18-C21
3	D	614	TMQ	C16-C15-O20-C23
3	B	612	TMQ	C16-C11-N10-C9
3	C	613	TMQ	C12-C11-N10-C9
3	D	614	TMQ	C14-C13-O18-C21
3	A	611	TMQ	C16-C15-O20-C23
3	D	614	TMQ	C12-C11-N10-C9
3	C	613	TMQ	C16-C11-N10-C9
3	B	612	TMQ	C13-C14-O19-C22
2	A	601	NAP	C3B-C4B-C5B-O5B
3	D	614	TMQ	C16-C11-N10-C9
3	A	611	TMQ	C14-C15-O20-C23
5	C	908	EDO	O1-C1-C2-O2
5	C	909	EDO	O1-C1-C2-O2
2	A	601	NAP	O4B-C4B-C5B-O5B
2	A	601	NAP	O4D-C4D-C5D-O5D
2	B	602	NAP	O4B-C4B-C5B-O5B
3	B	612	TMQ	C15-C14-O19-C22
2	D	604	NAP	PN-O3-PA-O1A
2	C	603	NAP	PN-O3-PA-O1A
5	C	910	EDO	O1-C1-C2-O2
2	C	603	NAP	C2B-O2B-P2B-O1X
2	B	602	NAP	C5D-O5D-PN-O2N
2	C	603	NAP	C5D-O5D-PN-O2N
5	D	905	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	C	912	EDO	O1-C1-C2-O2
2	A	601	NAP	C3B-C2B-O2B-P2B
5	A	904	EDO	O1-C1-C2-O2
5	D	907	EDO	O1-C1-C2-O2
2	A	601	NAP	C1B-C2B-O2B-P2B
2	D	604	NAP	C2D-C1D-N1N-C2N
2	D	604	NAP	C2D-C1D-N1N-C6N
2	C	603	NAP	C2B-O2B-P2B-O2X
2	A	601	NAP	PN-O3-PA-O1A
3	D	614	TMQ	C6-C9-N10-C11
2	D	604	NAP	C5D-O5D-PN-O2N
5	D	911	EDO	O1-C1-C2-O2
2	A	601	NAP	C4B-C5B-O5B-PA

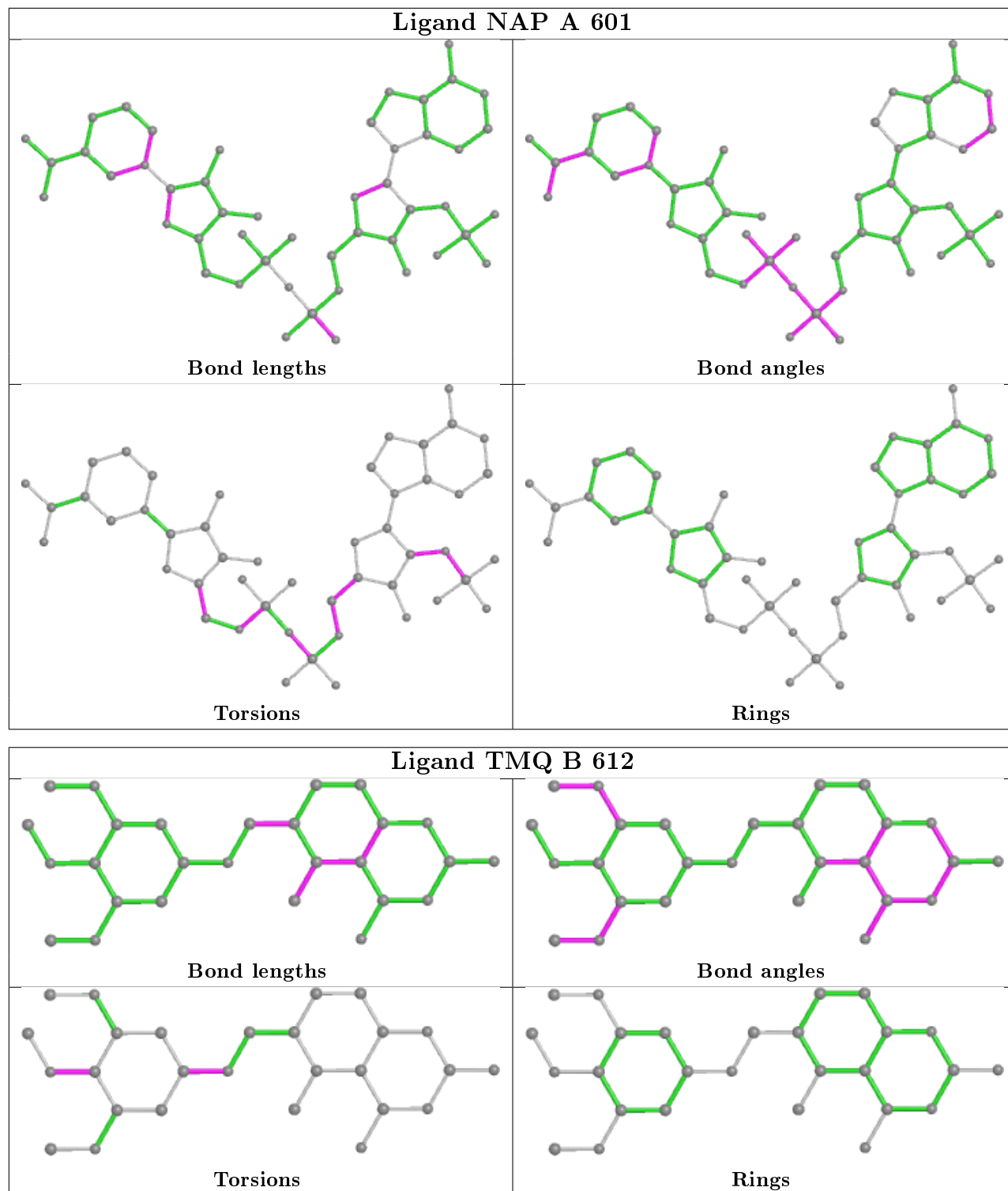
There are no ring outliers.

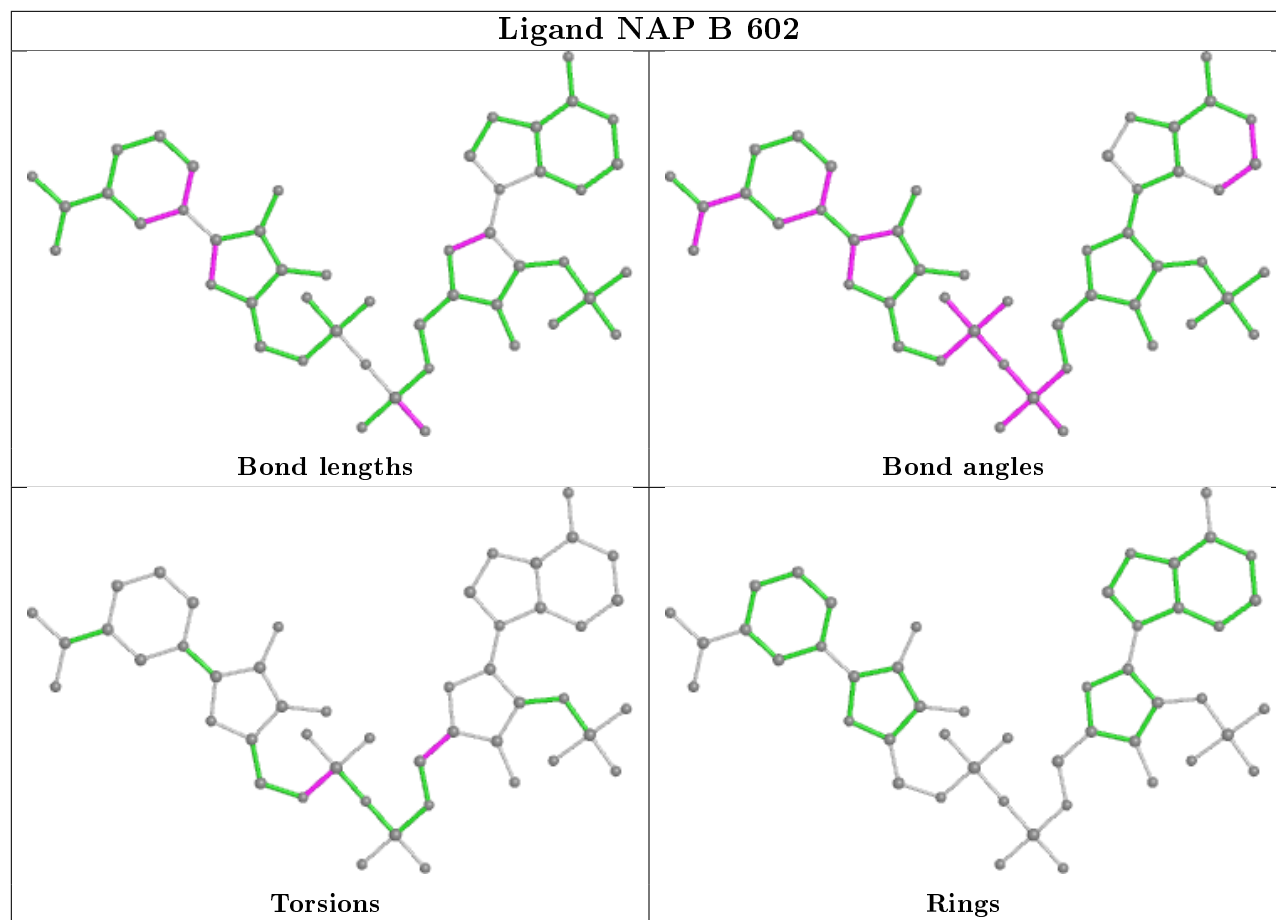
12 monomers are involved in 81 short contacts:

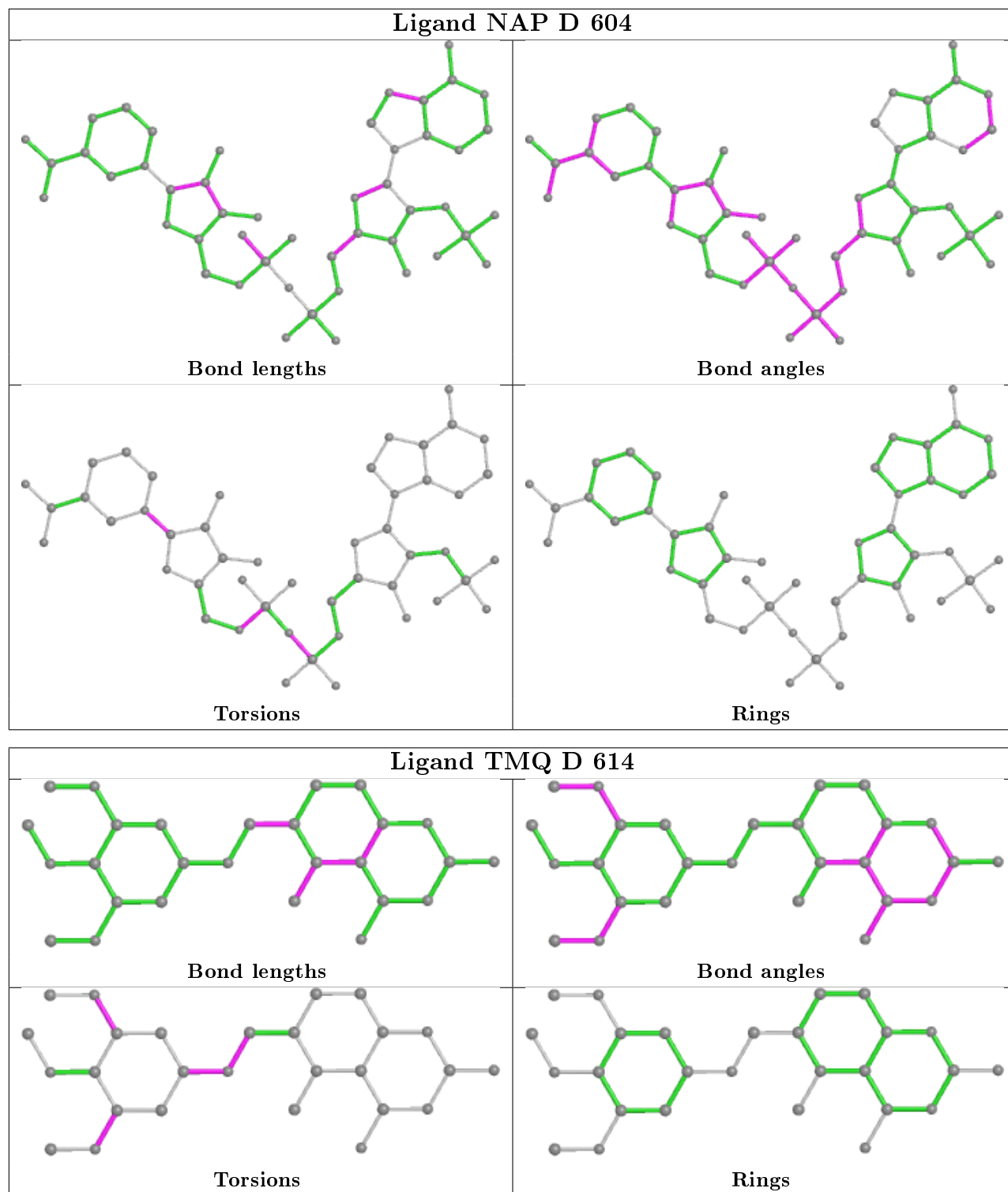
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	910	EDO	1	0
2	A	601	NAP	7	0
5	D	901	EDO	2	0
5	B	906	EDO	1	0
3	B	612	TMQ	14	0
2	B	602	NAP	7	0
2	D	604	NAP	12	0
3	D	614	TMQ	12	0
2	C	603	NAP	11	0
3	C	613	TMQ	9	0
3	A	611	TMQ	14	0
5	C	909	EDO	1	0

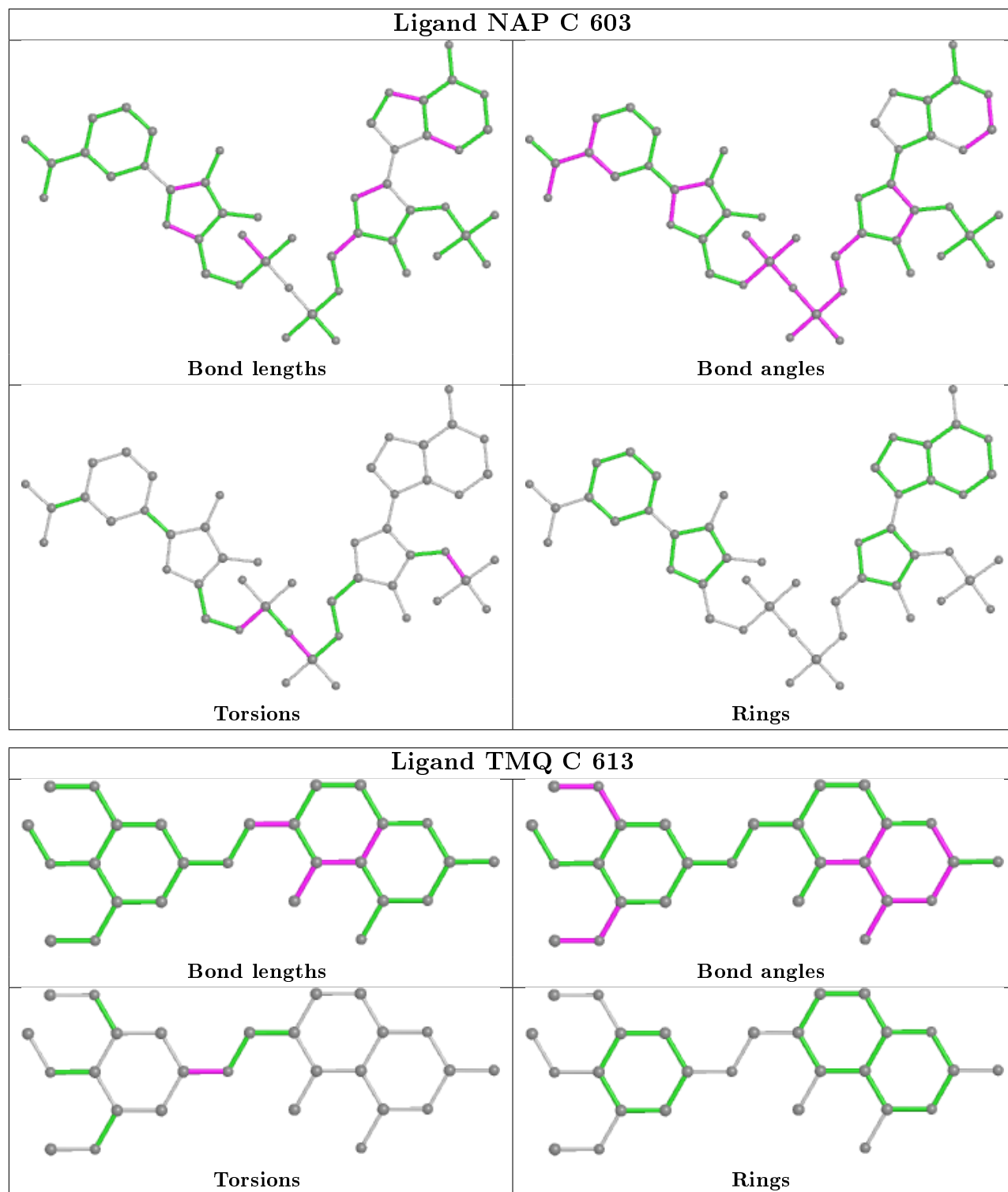
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

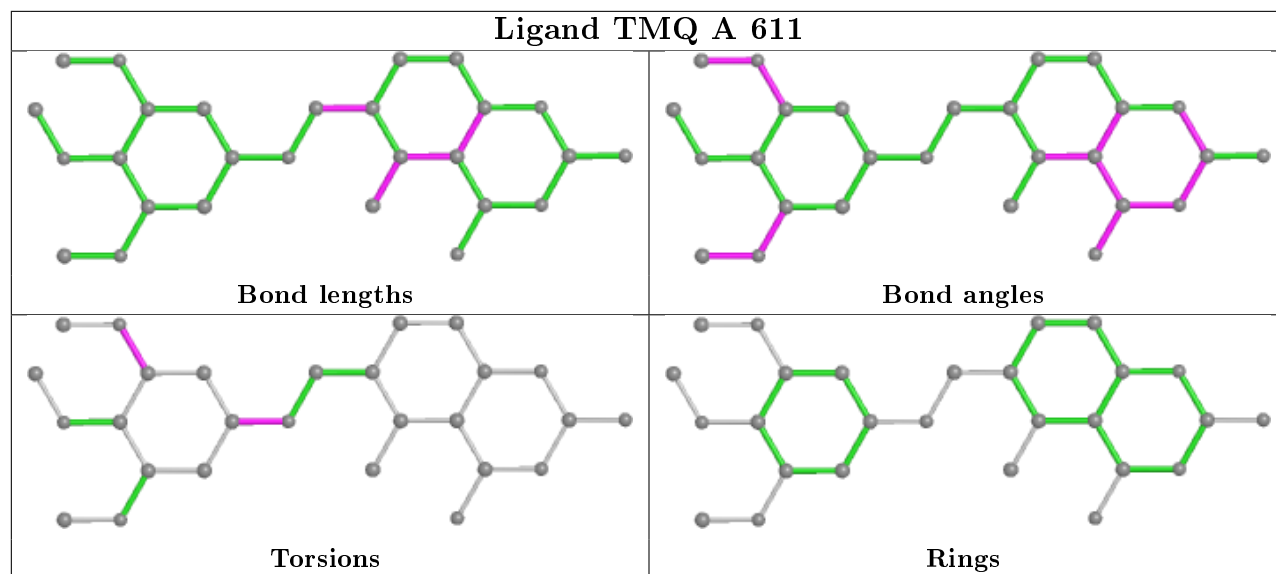
equivalents in the CSD to analyse the geometry.











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 ⓘ

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	510/521 (97%)	-0.43	6 (1%) 79 54	25, 38, 50, 57	0
1	B	505/521 (96%)	-0.27	14 (2%) 53 25	27, 37, 54, 62	0
1	C	510/521 (97%)	-0.46	7 (1%) 75 49	22, 35, 49, 61	0
1	D	507/521 (97%)	-0.43	5 (0%) 82 59	25, 37, 51, 56	0
All	All	2032/2084 (97%)	-0.40	32 (1%) 72 44	22, 37, 51, 62	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	185	SER	4.1
1	B	62	LYS	3.7
1	B	186	SER	3.5
1	B	188	SER	3.3
1	C	185	SER	3.2
1	C	114	ASP	3.2
1	A	516	SER	3.1
1	A	114	ASP	3.1
1	B	102	THR	3.1
1	A	185	SER	2.9
1	C	1	MET	2.8
1	B	63	ASN	2.8
1	D	39	ARG	2.8
1	C	122	ALA	2.7
1	B	223	THR	2.6
1	C	62	LYS	2.6
1	A	62	LYS	2.6
1	B	2	SER	2.6
1	A	63	ASN	2.5
1	C	14	GLU	2.5
1	D	186	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	63	ASN	2.4
1	A	11	THR	2.4
1	D	11	THR	2.4
1	B	187	CYS	2.3
1	B	190	PHE	2.3
1	B	221	ASN	2.3
1	B	185	SER	2.3
1	B	86	PRO	2.2
1	B	85	PRO	2.1
1	C	61	GLY	2.1
1	B	14	GLU	2.1

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 carbohydrates 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
3	TMQ	B	612	27/27	0.55	0.51	60,61,61,62	27
5	EDO	C	909	4/4	0.73	0.35	74,75,75,75	0
2	NAP	B	602	48/48	0.75	0.34	74,75,76,76	48
5	EDO	C	912	4/4	0.79	0.32	65,66,67,67	0
4	SO4	C	822	5/5	0.80	0.45	139,139,140,140	0
4	SO4	C	812	5/5	0.81	0.28	131,131,131,131	0
4	SO4	D	820	5/5	0.81	0.25	106,106,107,107	0
4	SO4	C	807	5/5	0.82	0.32	149,149,149,149	0
4	SO4	B	809	5/5	0.84	0.42	123,123,124,124	0
4	SO4	C	816	5/5	0.84	0.30	106,106,107,107	0
5	EDO	A	902	4/4	0.85	0.27	63,64,65,65	0
4	SO4	A	815	5/5	0.85	0.40	128,128,128,128	0

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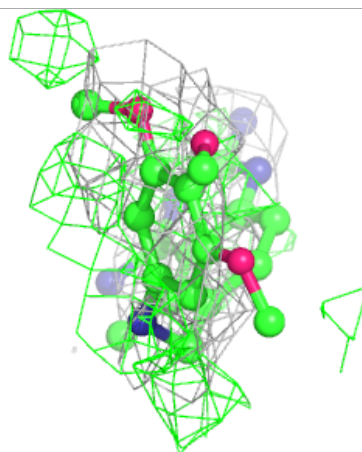
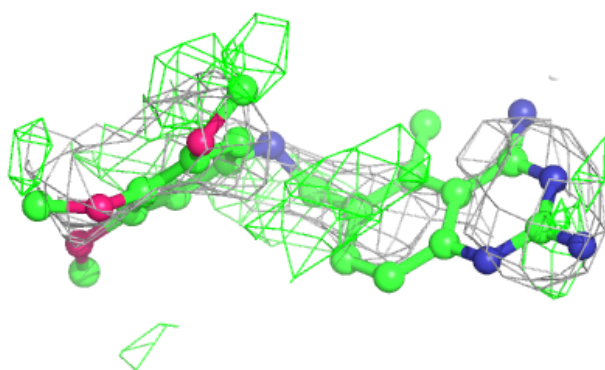
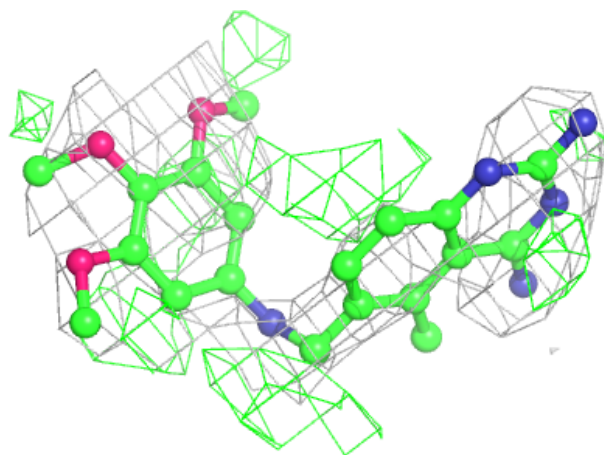
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	SO4	B	821	5/5	0.86	0.41	121,121,121,121	0
4	SO4	B	823	5/5	0.86	0.32	111,111,111,111	0
2	NAP	A	601	48/48	0.86	0.24	69,71,74,74	0
3	TMQ	D	614	27/27	0.87	0.22	68,70,75,76	0
4	SO4	D	810	5/5	0.87	0.44	134,134,134,134	0
2	NAP	C	603	48/48	0.89	0.23	73,79,84,84	0
3	TMQ	A	611	27/27	0.89	0.21	53,56,62,63	0
2	NAP	D	604	48/48	0.89	0.21	89,91,96,97	0
4	SO4	D	811	5/5	0.89	0.30	137,137,137,137	0
4	SO4	A	814	5/5	0.90	0.47	122,122,122,123	0
4	SO4	D	804	5/5	0.90	0.34	92,92,92,92	0
4	SO4	D	819	5/5	0.91	0.38	125,125,125,125	0
4	SO4	D	801	5/5	0.91	0.19	109,109,109,109	0
5	EDO	B	906	4/4	0.91	0.22	67,67,68,68	0
4	SO4	A	817	5/5	0.92	0.30	100,100,100,100	0
4	SO4	C	824	5/5	0.92	0.32	107,107,107,107	0
5	EDO	D	905	4/4	0.92	0.20	68,68,68,68	0
4	SO4	D	803	5/5	0.92	0.38	119,119,119,120	0
5	EDO	D	903	4/4	0.92	0.24	51,51,52,53	0
4	SO4	B	806	5/5	0.92	0.31	108,109,109,109	0
3	TMQ	C	613	27/27	0.92	0.16	53,57,62,64	0
5	EDO	C	908	4/4	0.93	0.27	55,56,56,57	0
5	EDO	C	910	4/4	0.93	0.18	46,48,49,52	0
4	SO4	A	805	5/5	0.94	0.29	106,106,106,106	0
5	EDO	D	907	4/4	0.94	0.22	57,58,58,60	0
5	EDO	D	911	4/4	0.94	0.19	64,64,64,64	0
4	SO4	B	818	5/5	0.95	0.30	121,121,121,121	0
4	SO4	A	813	5/5	0.95	0.16	104,104,104,104	0
4	SO4	C	802	5/5	0.96	0.13	88,89,89,89	0
5	EDO	D	901	4/4	0.96	0.16	45,46,46,48	0
4	SO4	D	704	5/5	0.97	0.10	71,72,72,72	0
5	EDO	A	904	4/4	0.98	0.12	50,50,50,51	0
4	SO4	C	702	5/5	0.98	0.11	58,59,60,60	0
4	SO4	C	808	5/5	0.98	0.35	106,106,106,106	0
4	SO4	A	701	5/5	0.98	0.09	64,64,64,65	0
4	SO4	C	703	5/5	0.99	0.08	62,63,63,63	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.

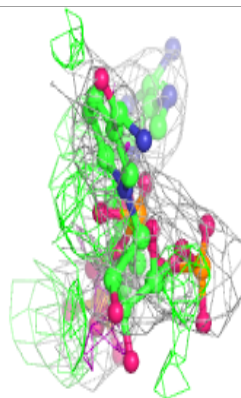
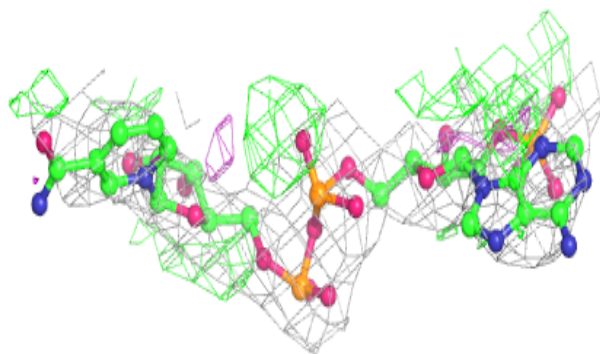
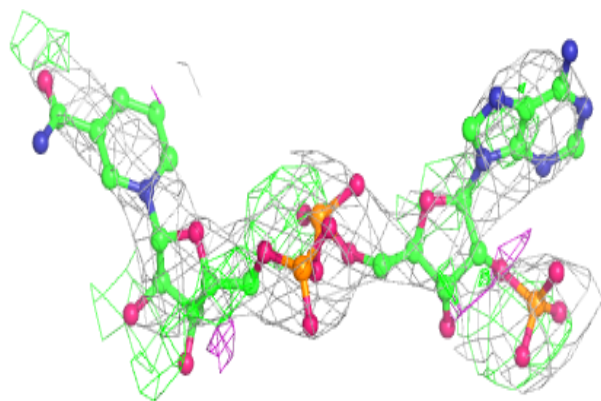
Electron density around TMQ B 612:

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

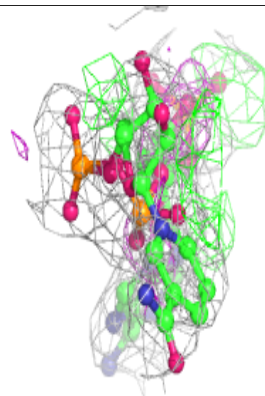
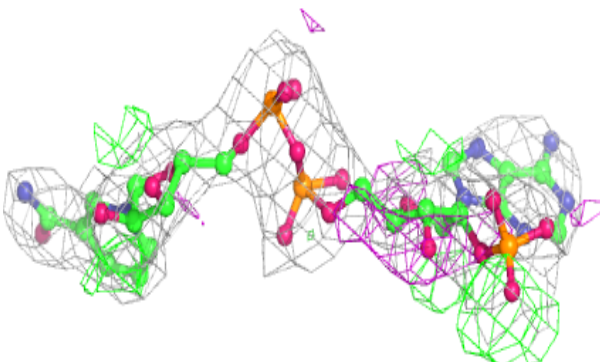
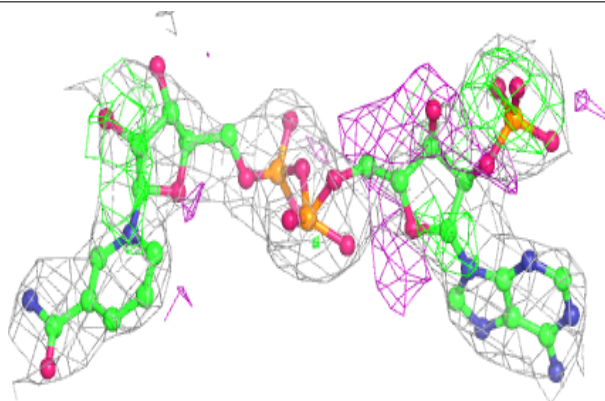


Electron density around NAP B 602:

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

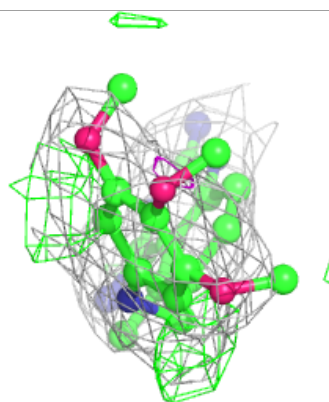
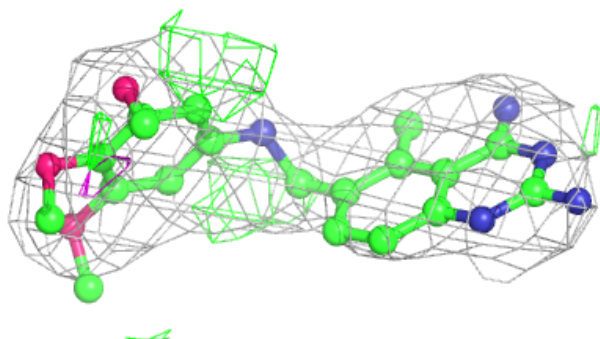
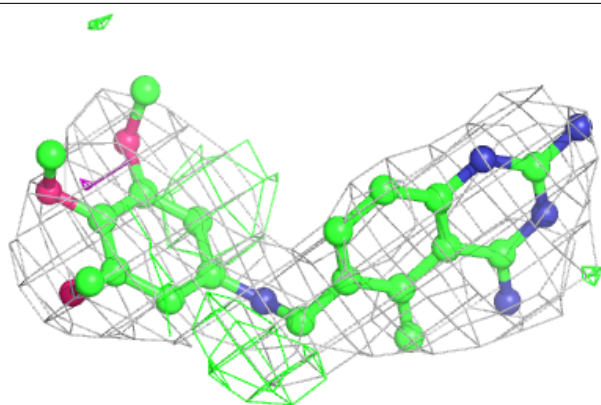
**Electron density around NAP A 601:**

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

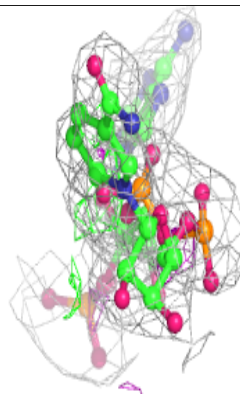
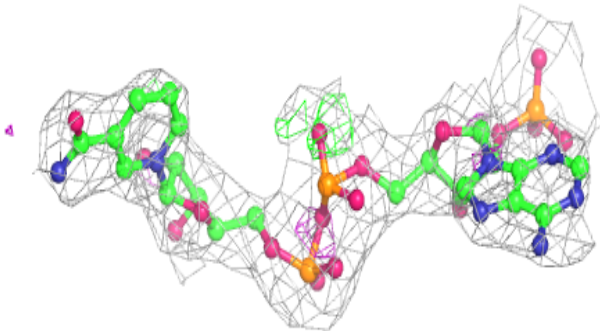
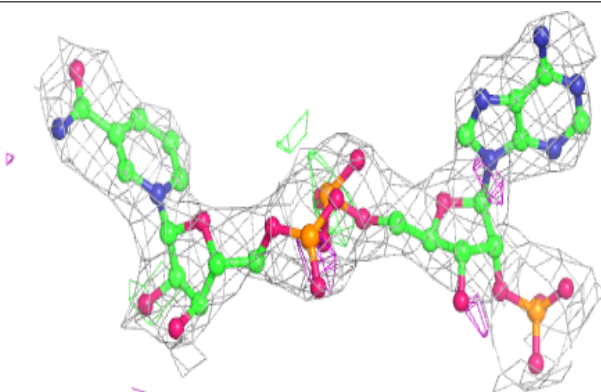


Electron density around TMQ D 614:

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

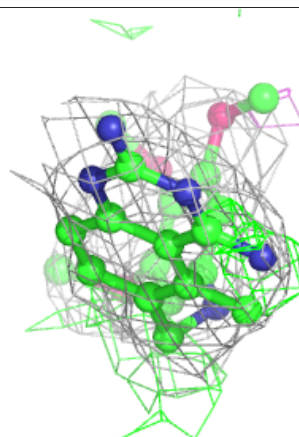
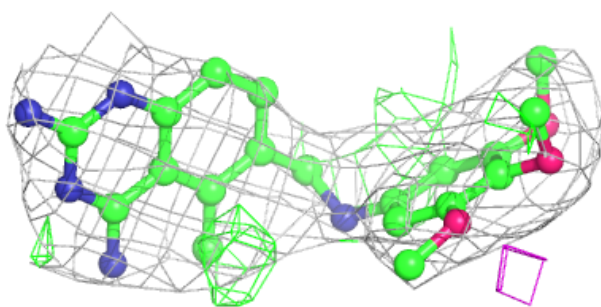
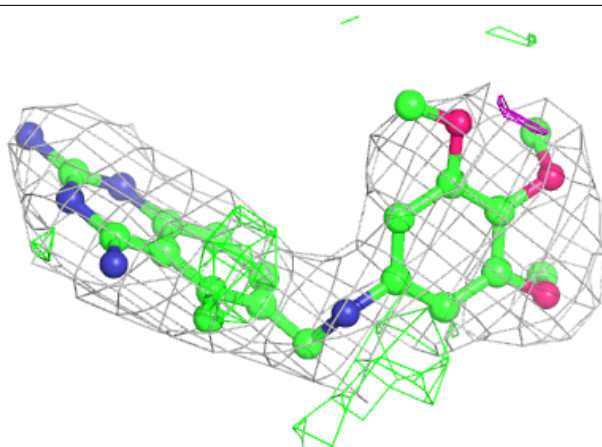
**Electron density around NAP C 603:**

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

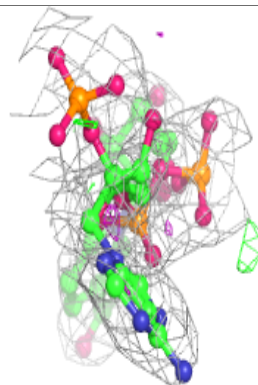
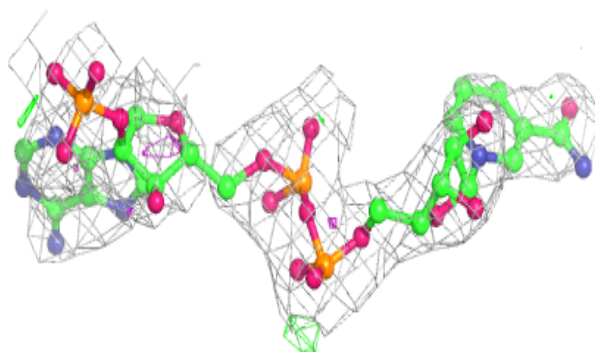
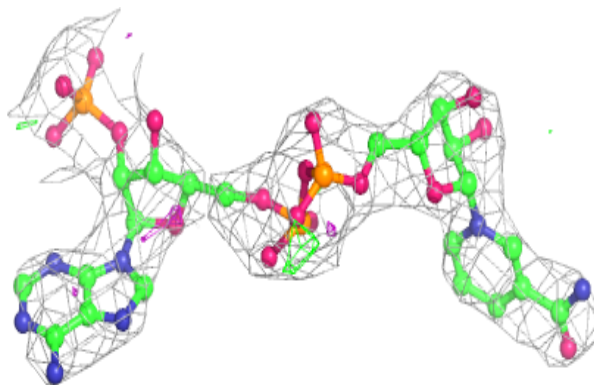


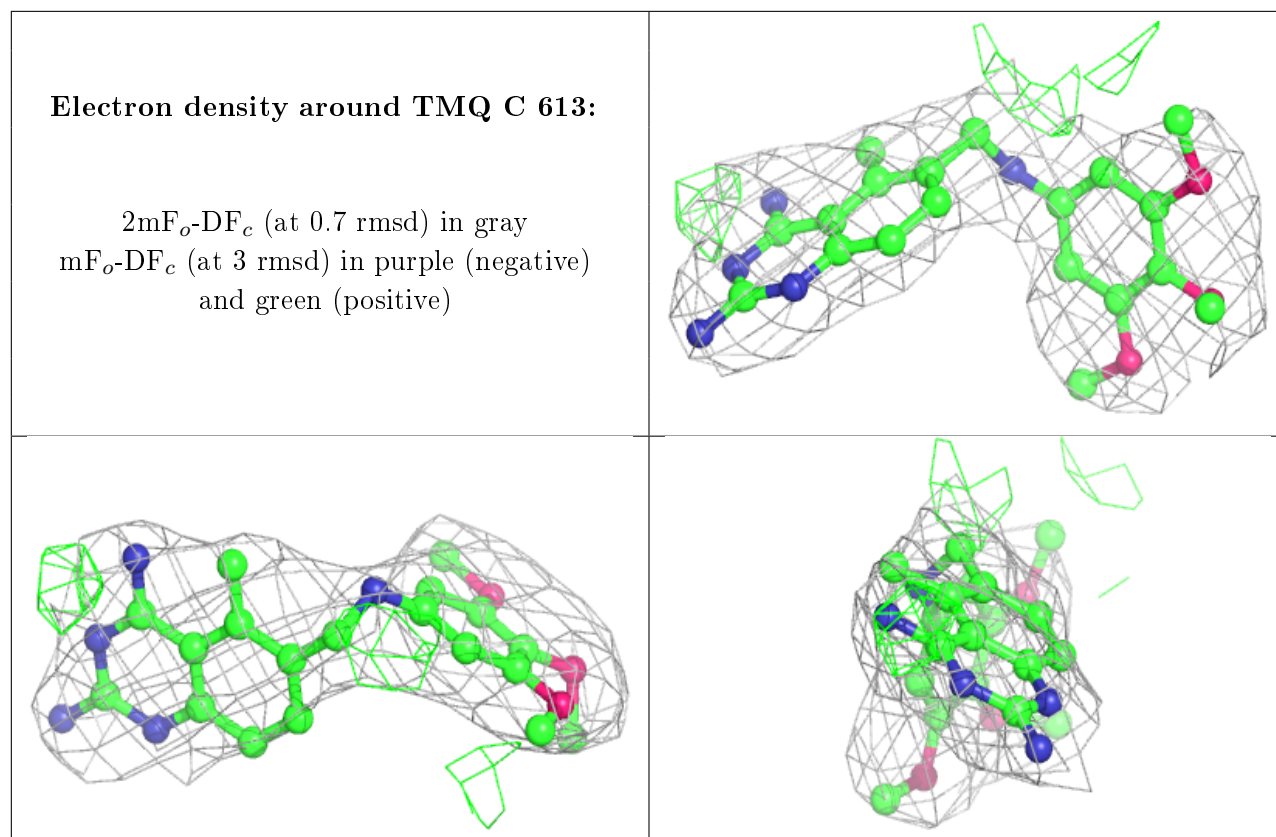
Electron density around TMQ A 611:

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

**Electron density around NAP D 604:**

$2mF_o-DF_c$ (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.