



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 05:37 AM BST

PDB ID : 6WEP  
Title : Crystal structure of TS-DHFR from *Cryptosporidium hominis* with Apo-TS site  
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Deposited on : 2020-04-02  
Resolution : 2.79 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](mailto: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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

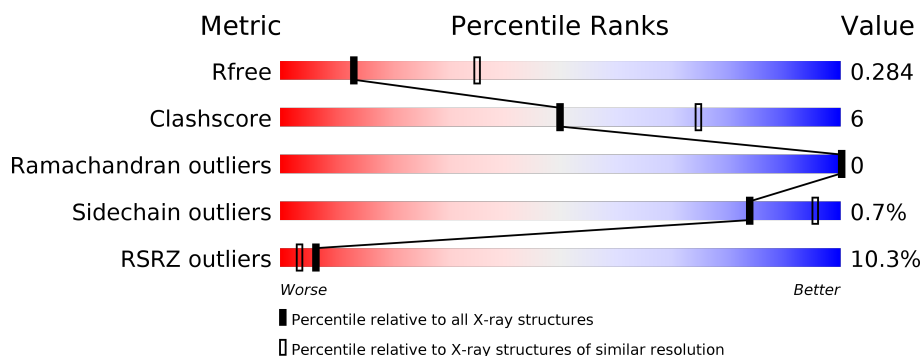
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.79 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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>8%</div> <div>84%</div> <div>13%</div> <div>.</div> </div>
1	B	521	<div> <div>12%</div> <div>82%</div> <div>14%</div> <div>.</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8009 atoms, of which 0 are hydrogens and 0 are deuteriums.

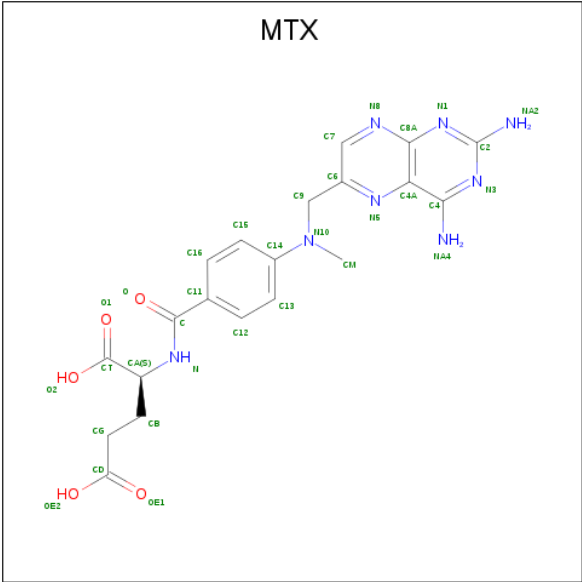
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Bifunctional dihydrofolate reductase-thymidylate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	503	Total	C	N	O	S	0	0	0
			3904	2510	660	714	20			
1	B	503	Total	C	N	O	S	0	0	0
			3926	2517	665	725	19			

- Molecule 2 is NADPH DIHYDRO-NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NDP) (formula: C<sub>21</sub>H<sub>30</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			33	20	8	5		
3	B	1	Total	C	N	O	0	0
			33	20	8	5		

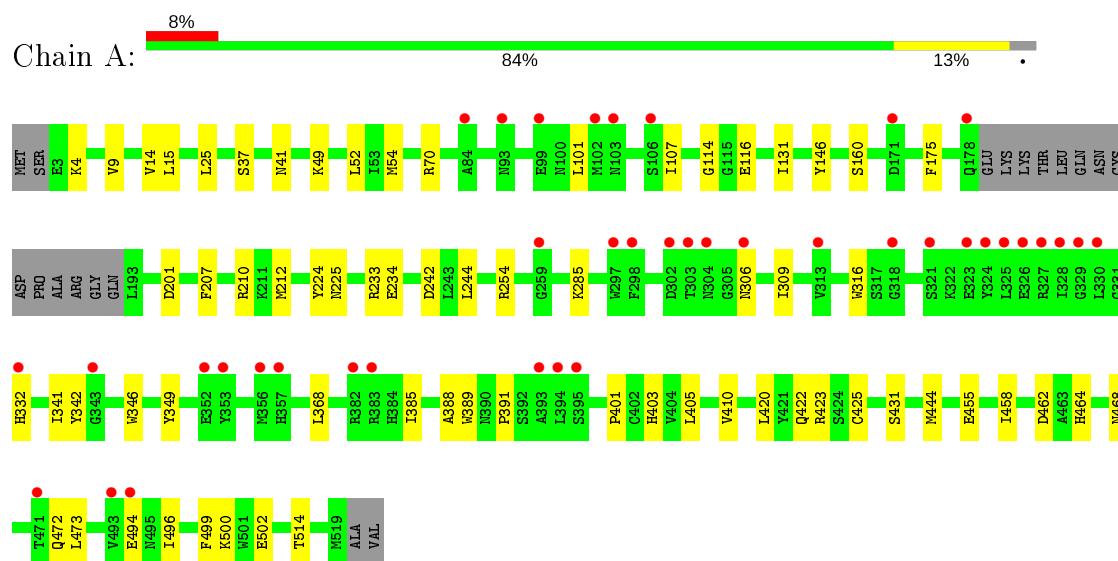
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		
4	B	8	Total	O	0	0
			8	8		

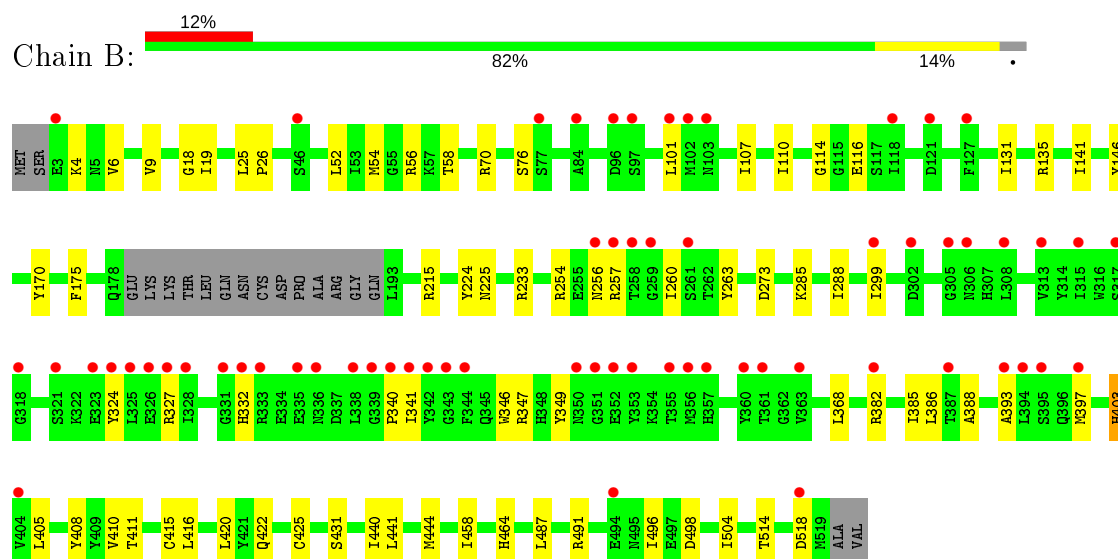
### 3 Residue-property plots

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, $\alpha$ , $\beta$ , $\gamma$	76.82Å 118.11Å 153.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.27 – 2.79 49.27 – 2.79	Depositor EDS
% Data completeness (in resolution range)	99.5 (49.27-2.79) 99.5 (49.27-2.79)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.20	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.93 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, $R_{free}$	0.239 , 0.285 0.239 , 0.284	Depositor DCC
$R_{free}$ test set	1765 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	73.8	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 41.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	8009	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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, MTX

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	0.25	0/3999	0.42	0/5438
1	B	0.25	0/4021	0.43	0/5466
All	All	0.25	0/8020	0.43	0/10904

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	3904	0	3693	44	0
1	B	3926	0	3713	53	0
2	A	48	0	26	1	0
2	B	48	0	26	2	0
3	A	33	0	20	4	0
3	B	33	0	20	2	0
4	A	9	0	0	1	0
4	B	8	0	0	0	0
All	All	8009	0	7498	89	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 6.

All (89) 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:4:LYS:HE2	1:A:101:LEU:HA	1.64	0.78
1:A:25:LEU:HD11	3:A:602:MTX:H7	1.68	0.75
1:A:368:LEU:HD21	1:A:496:ILE:HG21	1.71	0.73
1:A:410:VAL:O	1:B:254:ARG:NH2	2.21	0.73
1:A:254:ARG:NH2	1:B:410:VAL:O	2.20	0.73
1:B:285:LYS:HB3	1:B:514:THR:HG22	1.71	0.72
1:A:425:CYS:HG	1:A:431:SER:HG	1.36	0.71
1:B:25:LEU:HD11	3:B:602:MTX:H7	1.73	0.71
1:B:441:LEU:HA	1:B:444:MET:HE2	1.72	0.70
1:A:225:ASN:O	1:A:233:ARG:NH2	2.24	0.69
1:B:256:ASN:ND2	1:B:260:ILE:O	2.28	0.67
1:A:52:LEU:HD11	1:A:70:ARG:HD2	1.76	0.66
1:B:52:LEU:HD11	1:B:70:ARG:HD2	1.81	0.63
1:B:4:LYS:HE3	1:B:101:LEU:HA	1.81	0.62
1:B:4:LYS:NZ	1:B:107:ILE:O	2.29	0.61
1:B:225:ASN:O	1:B:233:ARG:NH2	2.28	0.61
1:B:288:ILE:HD11	1:B:440:ILE:HD11	1.83	0.60
1:B:257:ARG:NH1	1:B:327:ARG:HH12	2.01	0.59
1:A:285:LYS:HB3	1:A:514:THR:HG22	1.86	0.58
1:A:349:TYR:HH	1:B:349:TYR:HH	1.53	0.56
1:B:116:GLU:HB3	1:B:146:TYR:O	2.06	0.55
1:A:468:ASN:O	1:A:472:GLN:NE2	2.40	0.54
1:B:131:ILE:HB	1:B:175:PHE:HB2	1.90	0.54
1:B:299:ILE:O	1:B:347:ARG:NH1	2.39	0.54
1:B:403:HIS:HB3	1:B:420:LEU:HD11	1.91	0.53
1:A:37:SER:OG	3:A:602:MTX:O1	2.27	0.52
1:A:14:VAL:HG13	1:A:15:LEU:HD13	1.91	0.52
1:A:455:GLU:OE2	1:B:215:ARG:NH1	2.43	0.52
1:A:423:ARG:HB3	1:B:385:ILE:HD11	1.91	0.52
2:A:601:NDP:H42N	3:A:602:MTX:N5	2.24	0.52
1:A:464:HIS:CD2	1:B:382:ARG:HH22	2.28	0.52
1:A:346:TRP:HE1	1:A:388:ALA:HB2	1.73	0.52
1:A:420:LEU:HD23	1:A:458:ILE:HD12	1.91	0.52
1:B:420:LEU:HD23	1:B:458:ILE:HD12	1.92	0.51
1:A:444:MET:HE3	1:A:499:PHE:CD2	2.47	0.50
1:A:131:ILE:HB	1:A:175:PHE:HB2	1.93	0.50
1:A:306:ASN:HA	1:A:309:ILE:HG12	1.93	0.50
1:A:201:ASP:OD1	4:A:701:HOH:O	2.20	0.49
1:B:257:ARG:HH12	1:B:327:ARG:HH12	1.60	0.49
1:A:341:ILE:HG12	1:A:342:TYR:H	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:GLN:NE2	1:B:425:CYS:SG	2.85	0.49
2:B:601:NDP:H52A	2:B:601:NDP:H8A	1.94	0.48
1:B:9:VAL:HG12	3:B:602:MTX:N3	2.29	0.48
1:B:18:GLY:HA2	1:B:146:TYR:HA	1.95	0.48
1:A:116:GLU:HB3	1:A:146:TYR:O	2.14	0.48
1:A:494:GLU:N	1:A:494:GLU:OE1	2.39	0.48
1:B:6:VAL:HG22	1:B:110:ILE:HB	1.96	0.47
1:A:207:PHE:HB3	1:A:210:ARG:HB2	1.94	0.47
1:B:518:ASP:N	1:B:518:ASP:OD1	2.48	0.47
1:A:212:MET:SD	1:B:273:ASP:HB2	2.55	0.47
1:B:26:PRO:HB3	1:B:141:ILE:HD11	1.97	0.47
1:B:425:CYS:SG	1:B:431:SER:OG	2.56	0.47
1:A:425:CYS:SG	1:A:431:SER:OG	2.54	0.47
1:A:160:SER:HA	1:A:234:GLU:HB3	1.97	0.46
1:B:487:LEU:HD11	1:B:504:ILE:HG23	1.97	0.46
1:A:500:LYS:HE2	1:A:502:GLU:HB2	1.98	0.46
1:A:9:VAL:HG12	3:A:602:MTX:N3	2.31	0.45
1:B:491:ARG:NH2	1:B:498:ASP:O	2.49	0.45
1:B:346:TRP:HE1	1:B:388:ALA:HB2	1.81	0.45
1:B:393:ALA:O	1:B:397:MET:HG3	2.16	0.45
1:A:244:LEU:HD11	1:A:473:LEU:HD13	1.97	0.45
1:A:462:ASP:O	1:B:382:ARG:NH2	2.50	0.45
1:B:346:TRP:CG	1:B:386:LEU:HD11	2.51	0.45
1:A:391:PRO:HD2	1:B:349:TYR:CE2	2.52	0.45
1:A:385:ILE:HG23	1:A:405:LEU:HD11	1.98	0.44
1:B:385:ILE:HG23	1:B:405:LEU:HD11	2.00	0.44
1:A:464:HIS:CD2	1:B:382:ARG:NH2	2.86	0.44
1:A:346:TRP:NE1	1:A:388:ALA:HB2	2.32	0.44
1:B:256:ASN:HD21	1:B:260:ILE:HG13	1.82	0.43
1:B:341:ILE:HG22	1:B:397:MET:HE3	1.99	0.43
1:A:422:GLN:NE2	1:A:425:CYS:SG	2.92	0.43
1:A:233:ARG:NH1	1:A:242:ASP:OD1	2.52	0.43
1:B:56:ARG:HB2	1:B:76:SER:HB2	2.01	0.43
1:A:54:MET:HA	1:A:114:GLY:HA3	2.00	0.43
1:B:324:TYR:CE1	1:B:327:ARG:NH2	2.87	0.43
1:A:316:TRP:CH2	1:A:341:ILE:HD11	2.54	0.43
1:A:37:SER:O	1:A:41:ASN:ND2	2.52	0.42
1:B:54:MET:HA	1:B:114:GLY:HA3	2.01	0.42
1:B:408:TYR:HB3	1:B:416:LEU:HD11	2.01	0.42
1:B:368:LEU:HD21	1:B:496:ILE:HG21	2.01	0.42
1:B:19:ILE:O	2:B:601:NDP:H2N	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:TYR:O	1:B:464:HIS:HA	2.20	0.41
1:B:58:THR:HG21	1:B:114:GLY:HA2	2.03	0.41
1:B:411:THR:OG1	1:B:415:CYS:HB2	2.21	0.41
1:B:340:PRO:O	1:B:397:MET:HG2	2.20	0.41
1:B:135:ARG:O	1:B:170:TYR:HA	2.21	0.41
1:A:389:TRP:HE3	1:A:401:PRO:HG2	1.86	0.40
1:B:346:TRP:CD2	1:B:386:LEU:HD21	2.55	0.40
1:A:49:LYS:O	1:A:107:ILE:HA	2.21	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	499/521 (96%)	465 (93%)	34 (7%)	0	100	100
1	B	499/521 (96%)	468 (94%)	31 (6%)	0	100	100
All	All	998/1042 (96%)	933 (94%)	65 (6%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	402/470 (86%)	399 (99%)	3 (1%)	84	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	408/470 (87%)	405 (99%)	3 (1%)	84	95
All	All	810/940 (86%)	804 (99%)	6 (1%)	84	95

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	224	TYR
1	A	332	HIS
1	A	403	HIS
1	B	224	TYR
1	B	332	HIS
1	B	403	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	MTX	B	602	-	29,35,35	1.81	6 (20%)	38,49,49	1.92	9 (23%)
2	NDP	B	601	-	45,52,52	1.93	11 (24%)	53,80,80	1.13	6 (11%)
3	MTX	A	602	-	29,35,35	1.80	6 (20%)	38,49,49	2.03	11 (28%)
2	NDP	A	601	-	45,52,52	1.94	11 (24%)	53,80,80	1.10	6 (11%)

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
3	MTX	B	602	-	-	7/19/25/25	0/3/3/3
2	NDP	B	601	-	-	10/30/77/77	0/5/5/5
3	MTX	A	602	-	-	7/19/25/25	0/3/3/3
2	NDP	A	601	-	-	8/30/77/77	0/5/5/5

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	MTX	C2-NA2	5.47	1.44	1.33
3	B	602	MTX	C2-NA2	5.35	1.44	1.33
3	B	602	MTX	C-N	5.09	1.45	1.34
3	A	602	MTX	C-N	4.97	1.45	1.34
2	B	601	NDP	C4N-C3N	-4.51	1.41	1.49
2	A	601	NDP	C4N-C3N	-4.50	1.41	1.49
2	A	601	NDP	C6A-N6A	4.08	1.48	1.34
2	B	601	NDP	C6A-N6A	4.07	1.48	1.34
2	A	601	NDP	C7N-N7N	4.06	1.44	1.33
2	B	601	NDP	C7N-N7N	4.02	1.44	1.33
2	A	601	NDP	C3B-C2B	-3.82	1.44	1.52
2	B	601	NDP	C2D-C3D	-3.81	1.42	1.53
2	A	601	NDP	C2D-C3D	-3.76	1.43	1.53
2	B	601	NDP	C3B-C2B	-3.68	1.44	1.52
2	A	601	NDP	C6N-C5N	3.62	1.39	1.33
2	B	601	NDP	C6N-C5N	3.60	1.39	1.33
2	B	601	NDP	C4N-C5N	-3.22	1.40	1.48
2	A	601	NDP	C4N-C5N	-3.21	1.40	1.48
3	A	602	MTX	CB-CA	-3.20	1.49	1.53
3	B	602	MTX	CB-CA	-3.13	1.49	1.53
3	B	602	MTX	C4-NA4	2.52	1.43	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	MTX	C4-NA4	2.45	1.43	1.34
3	B	602	MTX	C7-N8	2.45	1.35	1.31
3	A	602	MTX	C7-N8	2.42	1.35	1.31
3	B	602	MTX	C14-N10	2.40	1.46	1.39
3	A	602	MTX	C14-N10	2.33	1.45	1.39
2	B	601	NDP	C3B-C4B	-2.29	1.47	1.53
2	A	601	NDP	C3B-C4B	-2.29	1.47	1.53
2	A	601	NDP	C4A-N3A	2.29	1.38	1.35
2	B	601	NDP	C4A-N3A	2.23	1.38	1.35
2	A	601	NDP	C3D-C4D	-2.22	1.47	1.53
2	B	601	NDP	C3D-C4D	-2.19	1.47	1.53
2	B	601	NDP	C2N-C3N	2.07	1.40	1.34
2	A	601	NDP	C2N-C3N	2.02	1.40	1.34

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	602	MTX	C6-C7-N8	-5.15	118.08	123.13
3	A	602	MTX	C6-C7-N8	-5.05	118.18	123.13
3	A	602	MTX	N8-C8A-N1	4.24	120.66	115.82
3	A	602	MTX	C7-N8-C8A	4.07	120.78	116.69
3	B	602	MTX	C7-N8-C8A	3.98	120.69	116.69
3	A	602	MTX	N1-C2-N3	-3.94	121.97	127.22
3	B	602	MTX	N8-C8A-N1	3.91	120.29	115.82
3	B	602	MTX	N1-C2-N3	-3.84	122.09	127.22
2	B	601	NDP	N3A-C2A-N1A	-3.81	122.73	128.68
2	A	601	NDP	N3A-C2A-N1A	-3.69	122.92	128.68
3	B	602	MTX	C2-N1-C8A	3.62	119.49	115.36
3	A	602	MTX	C2-N1-C8A	3.50	119.35	115.36
3	A	602	MTX	CG-CB-CA	3.36	119.82	113.04
3	A	602	MTX	C6-C9-N10	-3.09	108.30	113.60
3	B	602	MTX	CG-CB-CA	3.03	119.17	113.04
2	A	601	NDP	PN-O3-PA	-2.77	123.32	132.83
2	B	601	NDP	PN-O3-PA	-2.77	123.33	132.83
2	B	601	NDP	O5D-C5D-C4D	2.74	118.42	108.99
3	A	602	MTX	C4A-C4-N3	-2.52	119.35	121.01
3	B	602	MTX	C4A-C4-N3	-2.49	119.38	121.01
3	B	602	MTX	C6-C9-N10	-2.47	109.37	113.60
2	A	601	NDP	O5D-C5D-C4D	2.44	117.38	108.99
2	B	601	NDP	O5B-C5B-C4B	2.23	116.66	108.99
2	B	601	NDP	C3N-C7N-N7N	2.21	121.59	117.67
2	A	601	NDP	C4A-C5A-N7A	-2.16	107.14	109.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	NDP	C4A-C5A-N7A	-2.15	107.16	109.40
3	A	602	MTX	CM-N10-C9	2.14	120.51	114.84
3	A	602	MTX	CA-N-C	-2.12	119.61	122.34
2	A	601	NDP	C3N-C7N-N7N	2.05	121.31	117.67
2	A	601	NDP	O5B-C5B-C4B	2.05	116.04	108.99
3	A	602	MTX	C11-C-N	2.04	120.98	117.06
3	B	602	MTX	C7-C6-N5	2.02	122.17	120.85

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	602	MTX	N-CA-CB-CG
3	B	602	MTX	CT-CA-CB-CG
3	A	602	MTX	N-CA-CB-CG
3	A	602	MTX	CT-CA-CB-CG
3	B	602	MTX	C13-C14-N10-CM
3	A	602	MTX	C13-C14-N10-CM
3	B	602	MTX	C15-C14-N10-CM
3	A	602	MTX	C15-C14-N10-CM
2	B	601	NDP	C3D-C4D-C5D-O5D
2	B	601	NDP	PA-O3-PN-O5D
2	A	601	NDP	PA-O3-PN-O5D
2	B	601	NDP	C2B-O2B-P2B-O1X
2	A	601	NDP	C2B-O2B-P2B-O1X
2	B	601	NDP	C5D-O5D-PN-O3
2	B	601	NDP	C5D-O5D-PN-O2N
3	B	602	MTX	C6-C9-N10-CM
3	A	602	MTX	C6-C9-N10-CM
2	B	601	NDP	C4D-C5D-O5D-PN
2	B	601	NDP	O4D-C4D-C5D-O5D
2	A	601	NDP	O4D-C1D-N1N-C2N
2	A	601	NDP	C4D-C5D-O5D-PN
2	A	601	NDP	C3D-C4D-C5D-O5D
2	B	601	NDP	O4D-C1D-N1N-C2N
3	B	602	MTX	C15-C14-N10-C9
3	A	602	MTX	C15-C14-N10-C9
3	B	602	MTX	CT-CA-N-C
3	A	602	MTX	CT-CA-N-C
2	B	601	NDP	C2B-O2B-P2B-O3X
2	A	601	NDP	C5D-O5D-PN-O3
2	B	601	NDP	C5B-O5B-PA-O1A

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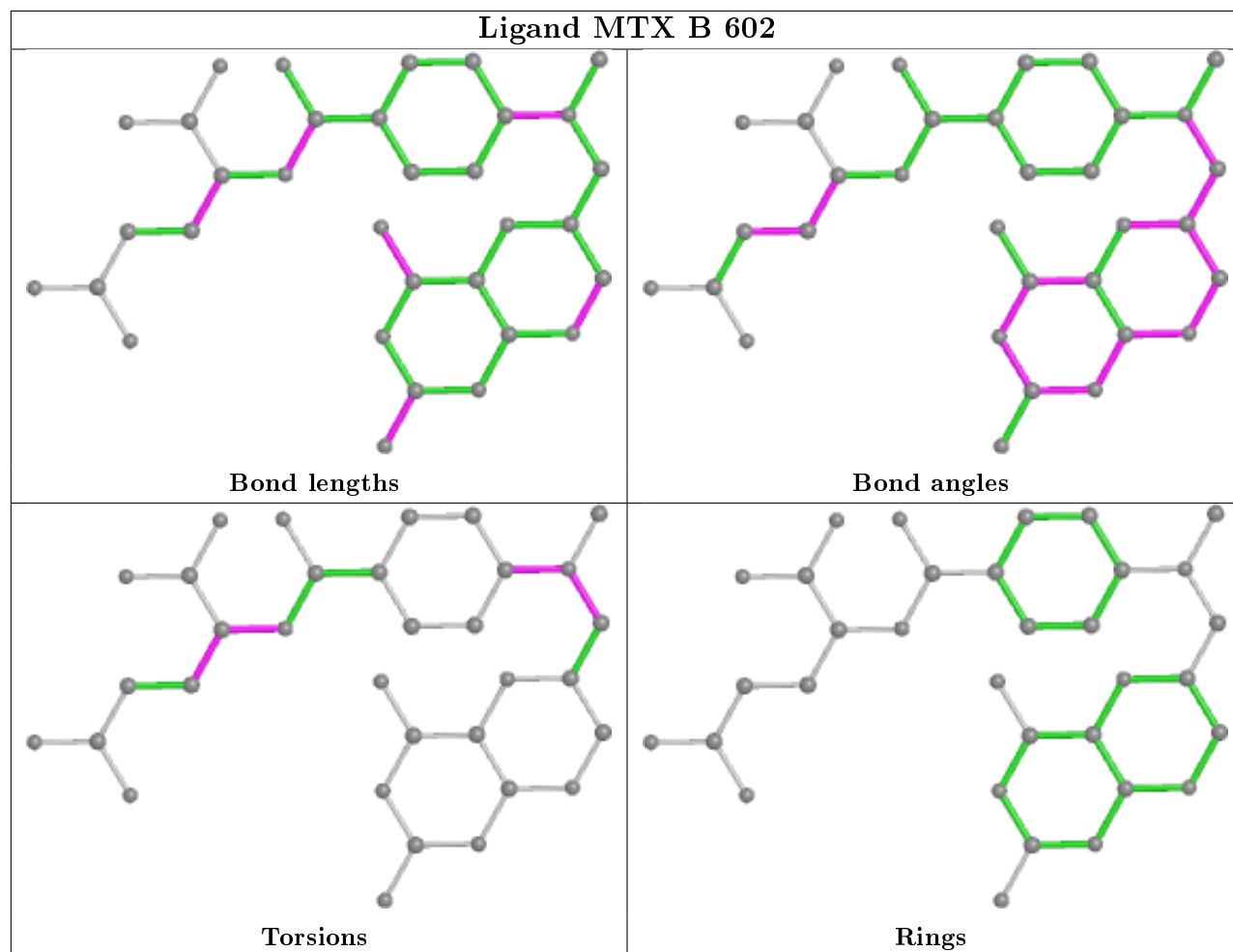
Mol	Chain	Res	Type	Atoms
2	A	601	NDP	C5B-O5B-PA-O1A
2	A	601	NDP	C5D-O5D-PN-O2N

There are no ring outliers.

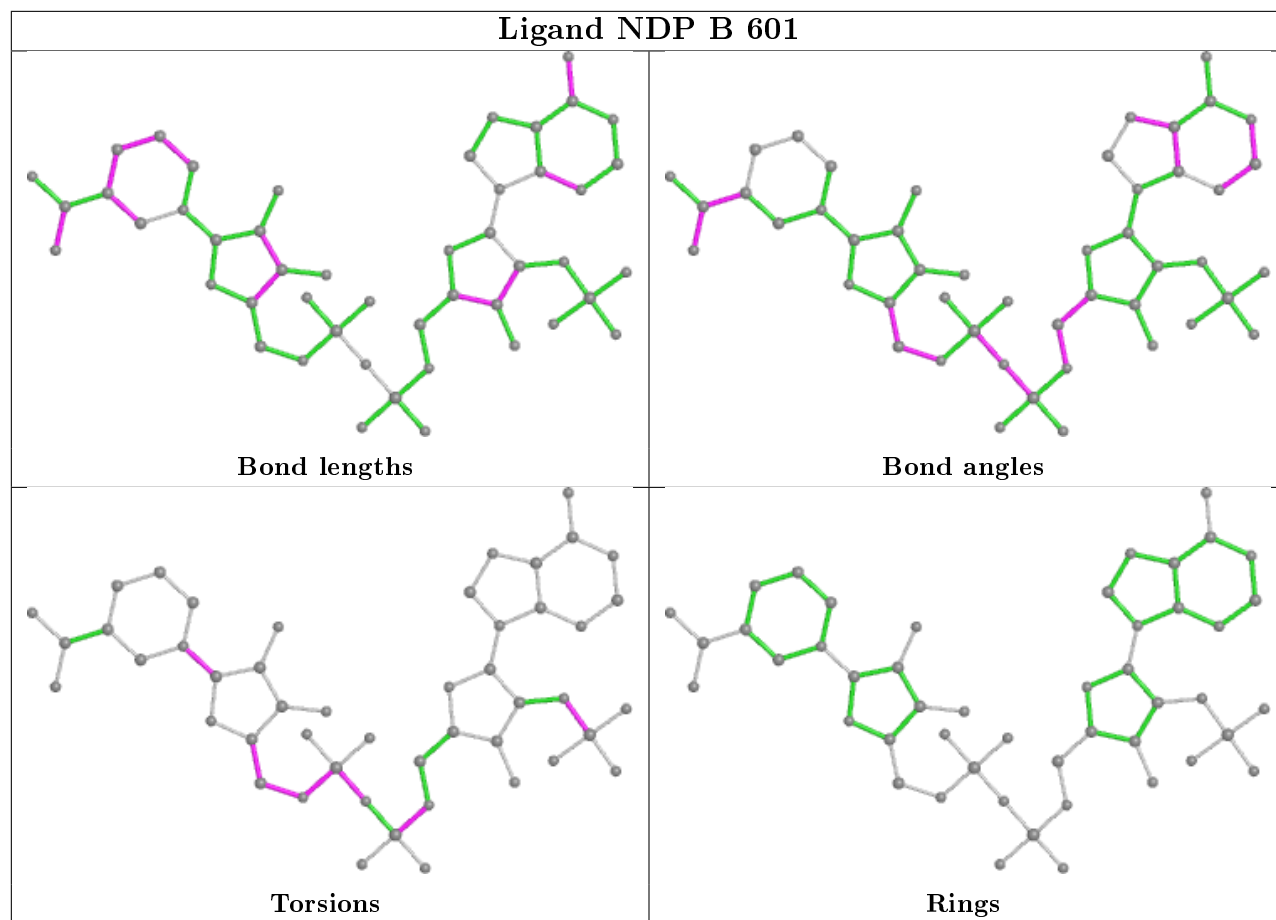
4 monomers are involved in 8 short contacts:

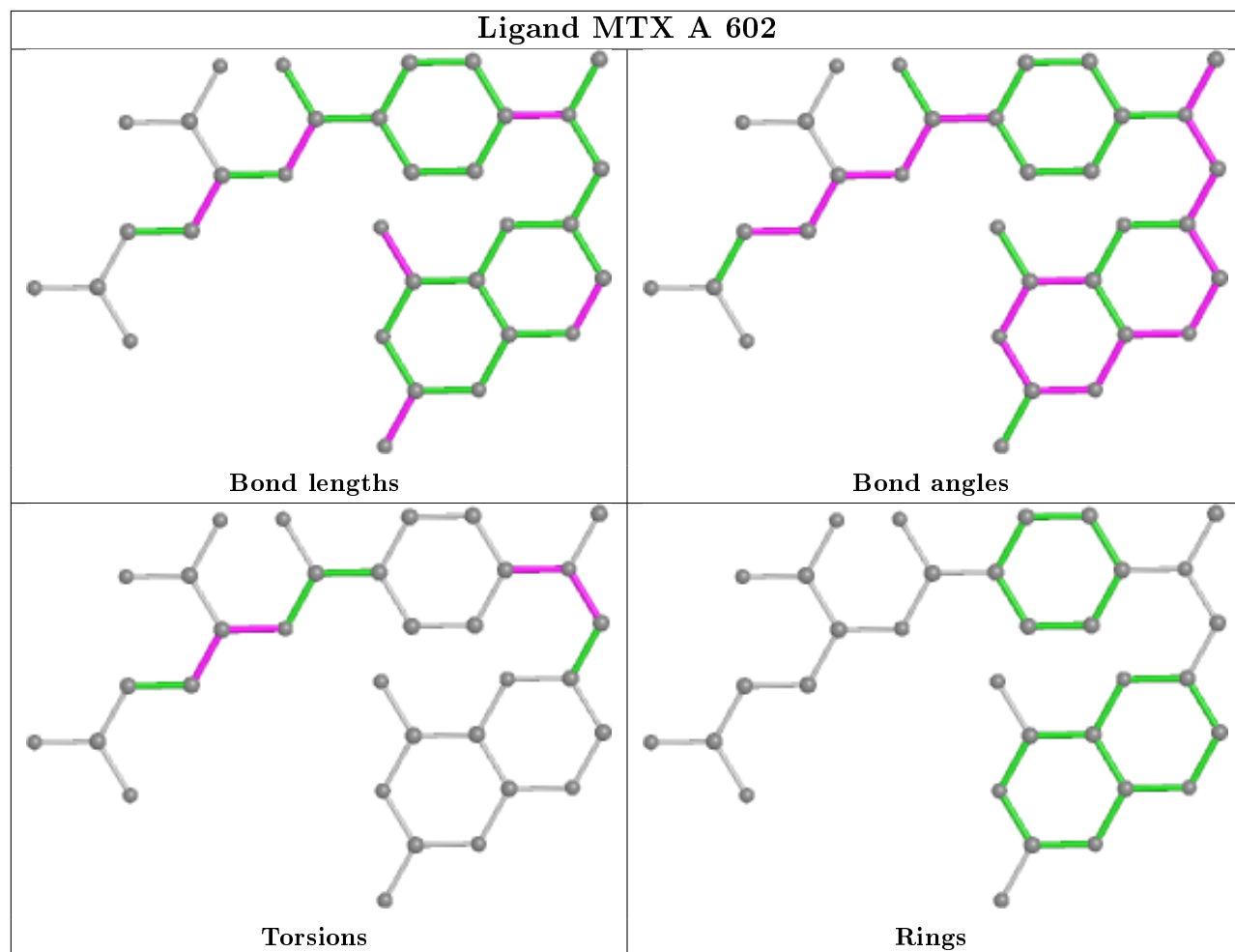
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	602	MTX	2	0
2	B	601	NDP	2	0
3	A	602	MTX	4	0
2	A	601	NDP	1	0

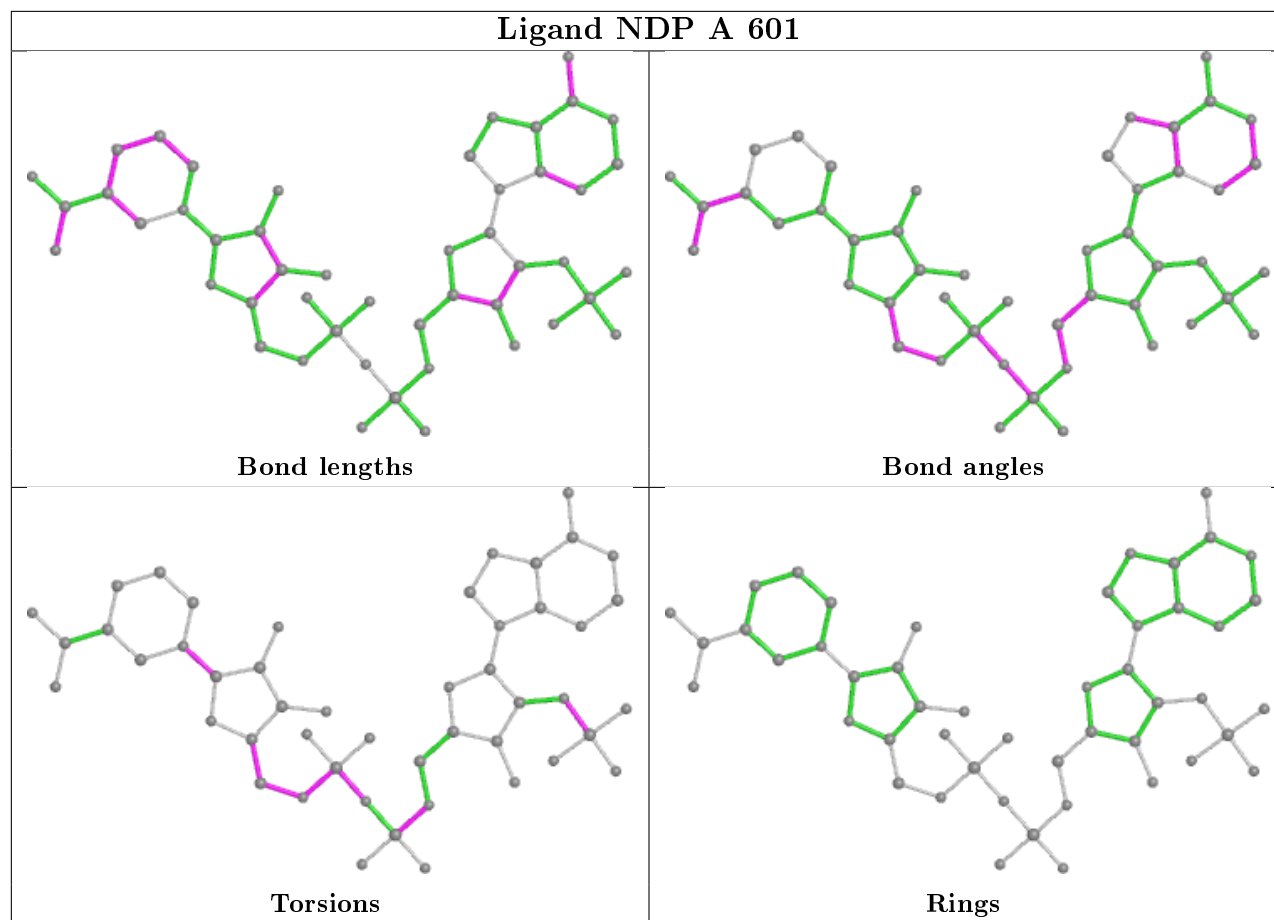
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	503/521 (96%)	0.70	40 (7%) <b>12</b> <b>6</b>	43, 67, 109, 135	0
1	B	503/521 (96%)	0.81	64 (12%) <b>3</b> <b>2</b>	42, 67, 117, 145	0
All	All	1006/1042 (96%)	0.75	104 (10%) <b>6</b> <b>3</b>	42, 67, 115, 145	0

All (104) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	363	VAL	6.3
1	B	324	TYR	6.3
1	A	321	SER	6.2
1	B	321	SER	6.0
1	B	357	HIS	5.9
1	A	103	ASN	5.9
1	A	318	GLY	5.8
1	B	351	GLY	5.8
1	B	323	GLU	5.6
1	B	102	MET	5.3
1	A	302	ASP	5.1
1	A	297	TRP	5.1
1	B	305	GLY	5.0
1	B	327	ARG	4.9
1	B	343	GLY	4.9
1	A	306	ASN	4.4
1	B	257	ARG	4.2
1	A	327	ARG	4.1
1	A	357	HIS	4.1
1	A	329	GLY	4.0
1	B	103	ASN	4.0
1	A	303	THR	3.8
1	B	258	THR	3.7
1	A	324	TYR	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	397	MET	3.5
1	B	494	GLU	3.4
1	A	325	LEU	3.4
1	B	332	HIS	3.3
1	A	326	GLU	3.3
1	B	518	ASP	3.3
1	B	256	ASN	3.2
1	A	304	ASN	3.2
1	A	102	MET	3.2
1	A	106	SER	3.2
1	B	328	ILE	3.1
1	A	356	MET	3.0
1	B	393	ALA	3.0
1	B	318	GLY	3.0
1	B	308	LEU	3.0
1	B	333	ARG	2.9
1	B	96	ASP	2.9
1	B	382	ARG	2.9
1	A	328	ILE	2.9
1	A	178	GLN	2.9
1	A	393	ALA	2.9
1	B	127	PHE	2.8
1	B	342	TYR	2.8
1	A	259	GLY	2.8
1	B	356	MET	2.8
1	B	353	TYR	2.8
1	B	331	GLY	2.8
1	A	383	ARG	2.7
1	A	313	VAL	2.7
1	B	302	ASP	2.7
1	A	99	GLU	2.7
1	A	323	GLU	2.6
1	B	361	THR	2.6
1	A	352	GLU	2.6
1	B	350	ASN	2.6
1	A	332	HIS	2.6
1	B	335	GLU	2.6
1	A	395	SER	2.6
1	B	315	ILE	2.5
1	B	355	THR	2.5
1	A	394	LEU	2.5
1	A	493	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	306	ASN	2.4
1	B	299	ILE	2.4
1	B	339	GLY	2.4
1	B	118	ILE	2.4
1	B	341	ILE	2.4
1	A	471	THR	2.4
1	A	171	ASP	2.3
1	B	101	LEU	2.3
1	A	84	ALA	2.3
1	B	97	SER	2.3
1	B	259	GLY	2.3
1	B	395	SER	2.3
1	A	93	ASN	2.3
1	B	336	ASN	2.3
1	B	313	VAL	2.3
1	B	261	SER	2.3
1	A	353	TYR	2.3
1	B	84	ALA	2.3
1	B	325	LEU	2.3
1	B	387	THR	2.2
1	B	326	GLU	2.2
1	B	121	ASP	2.2
1	B	360	TYR	2.2
1	A	330	LEU	2.2
1	B	394	LEU	2.2
1	B	344	PHE	2.2
1	B	46	SER	2.2
1	B	352	GLU	2.2
1	B	404	VAL	2.1
1	A	298	PHE	2.1
1	B	317	SER	2.1
1	B	340	PRO	2.1
1	A	494	GLU	2.1
1	A	343	GLY	2.1
1	B	3	GLU	2.1
1	B	77	SER	2.1
1	B	338	LEU	2.0
1	A	382	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

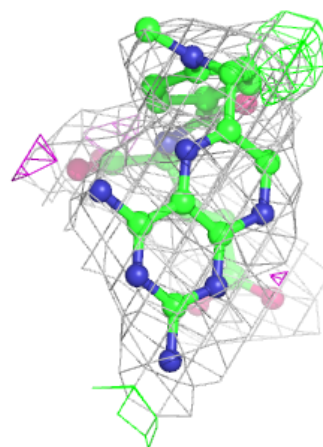
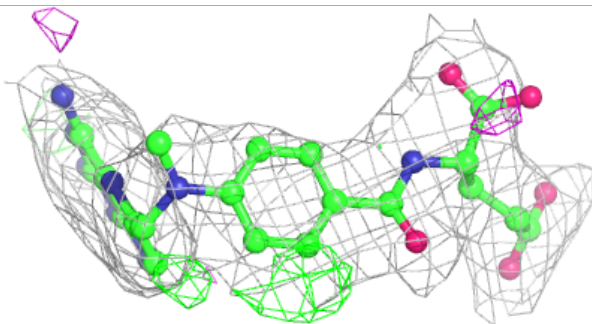
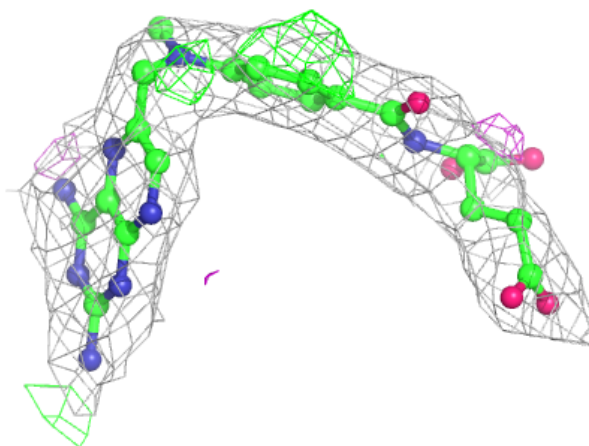
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MTX	B	602	33/33	0.89	0.25	51,61,83,85	0
3	MTX	A	602	33/33	0.89	0.29	46,58,80,84	0
2	NDP	B	601	48/48	0.93	0.18	53,70,95,107	0
2	NDP	A	601	48/48	0.96	0.18	49,62,78,83	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 MTX B 602:**

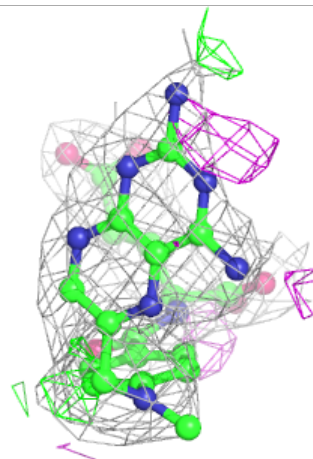
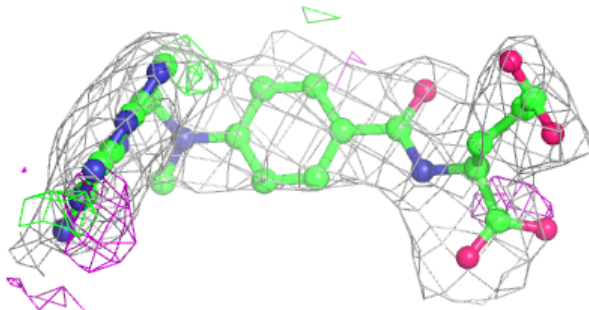
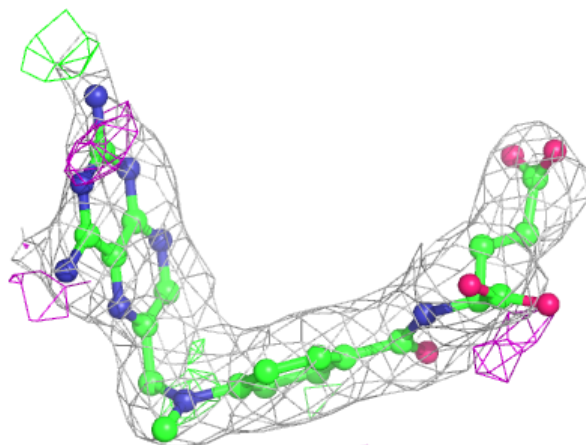
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





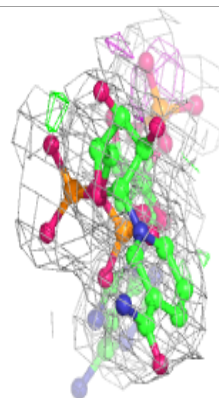
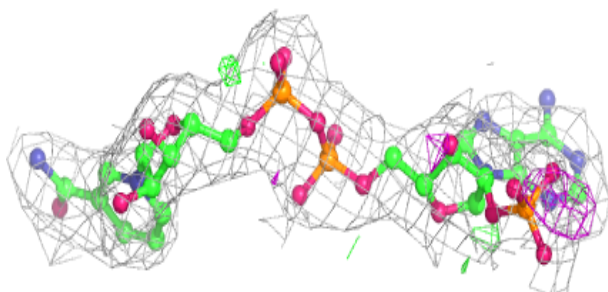
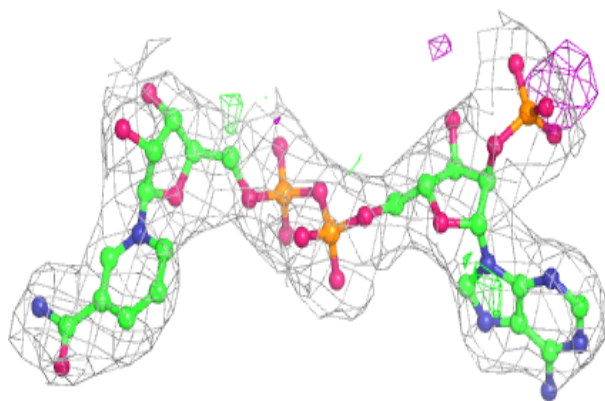
**Electron density around MTX A 602:**

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

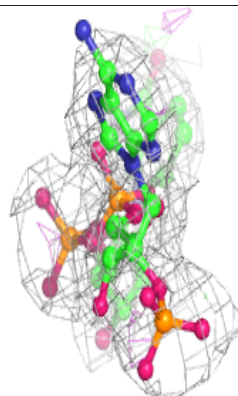
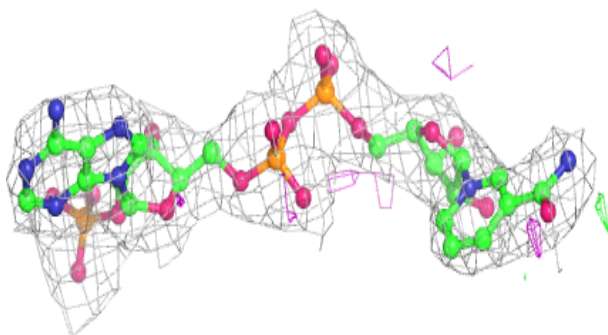
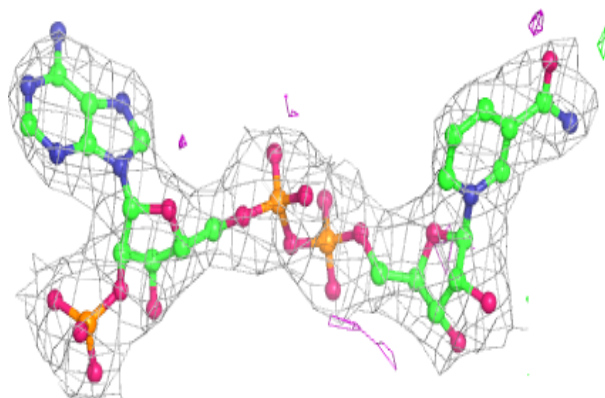


**Electron density around NDP B 601:**

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

**Electron density around NDP 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)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.