



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 02:10 AM UTC

PDB ID : 1N3T / pdb_00001n3t
Title : Biosynthesis of pteridins. Reaction mechanism of GTP cyclohydrolase I
Authors : Rebelo, J.; Auerbach, G.; Bader, G.; Bracher, A.; Nar, H.; Hoesl, C.; Schramek, N.; Kaiser, J.; Bacher, A.; Huber, R.; Fischer, M.
Deposited on : 2002-10-29
Resolution : 3.20 Å(reported)

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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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

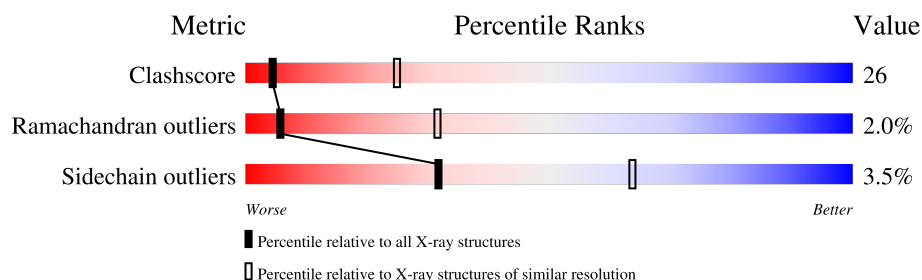
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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))
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	221	
1	B	221	
1	C	221	
1	D	221	
1	E	221	
1	F	221	
1	G	221	
1	H	221	

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Mol	Chain	Length	Quality of chain
1	I	221	 53% 43% .
1	J	221	 53% 42% .
1	K	221	 56% 38% 5%
1	L	221	 55% 40% 5%
1	M	221	 55% 42% .
1	N	221	 50% 48% .
1	O	221	 51% 46% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 26475 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.

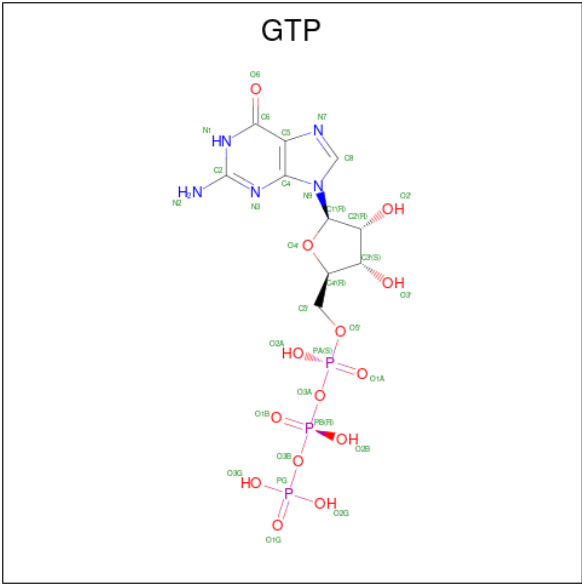
- Molecule 1 is a protein called GTP cyclohydrolase I.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	G	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	H	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	I	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	J	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	K	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	L	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	M	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	N	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	O	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	A	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	B	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	C	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	D	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			
1	E	221	Total	C	N	O	S	0	0	0
			1733	1088	309	328	8			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	SER	CYS	engineered mutation	UNP P0A6T5
B	181	SER	CYS	engineered mutation	UNP P0A6T5
C	181	SER	CYS	engineered mutation	UNP P0A6T5
D	181	SER	CYS	engineered mutation	UNP P0A6T5
E	181	SER	CYS	engineered mutation	UNP P0A6T5
F	181	SER	CYS	engineered mutation	UNP P0A6T5
G	181	SER	CYS	engineered mutation	UNP P0A6T5
H	181	SER	CYS	engineered mutation	UNP P0A6T5
I	181	SER	CYS	engineered mutation	UNP P0A6T5
J	181	SER	CYS	engineered mutation	UNP P0A6T5
K	181	SER	CYS	engineered mutation	UNP P0A6T5
L	181	SER	CYS	engineered mutation	UNP P0A6T5
M	181	SER	CYS	engineered mutation	UNP P0A6T5
N	181	SER	CYS	engineered mutation	UNP P0A6T5
O	181	SER	CYS	engineered mutation	UNP P0A6T5

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	F	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	G	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	H	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	I	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

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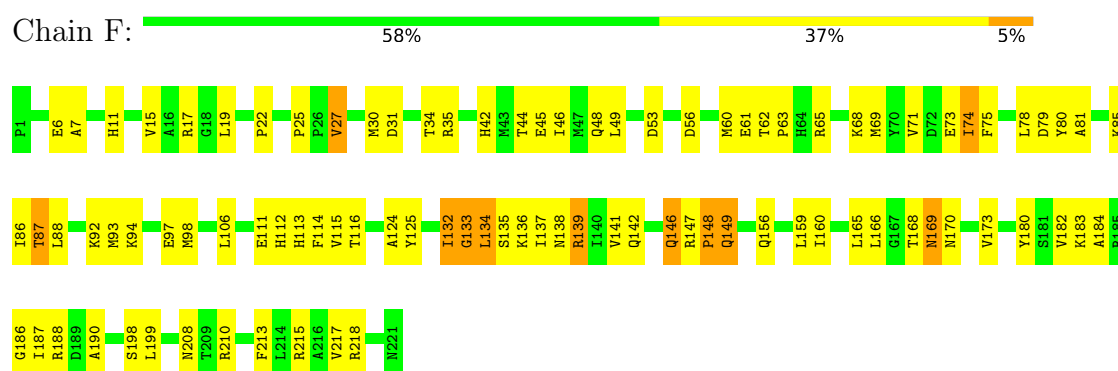
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	J	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	K	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	L	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	M	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	N	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	O	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	A	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	B	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	C	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	D	1	Total 32	C 10	N 5	O 14	P 3	0	0
2	E	1	Total 32	C 10	N 5	O 14	P 3	0	0

3 Residue-property plots

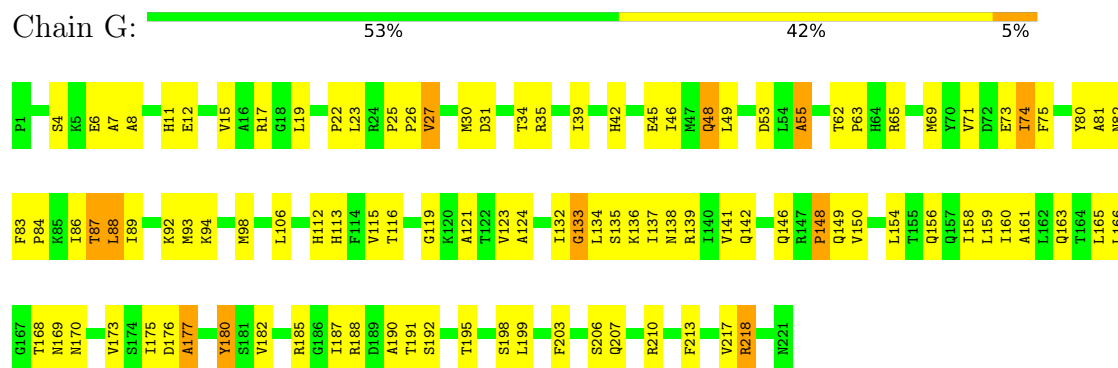
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

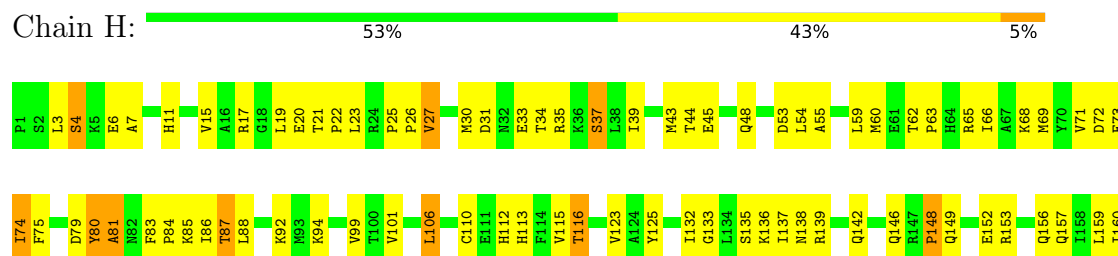
• Molecule 1: GTP cyclohydrolase I

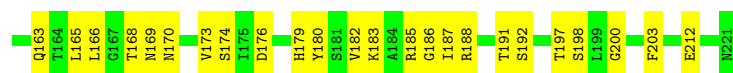


• Molecule 1: GTP cyclohydrolase I

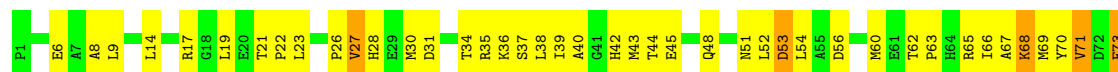


• Molecule 1: GTP cyclohydrolase I





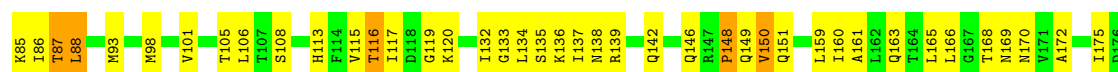
• Molecule 1: GTP cyclohydrolase I



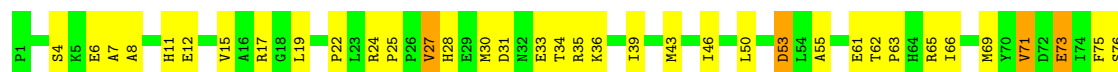
• Molecule 1: GTP cyclohydrolase I

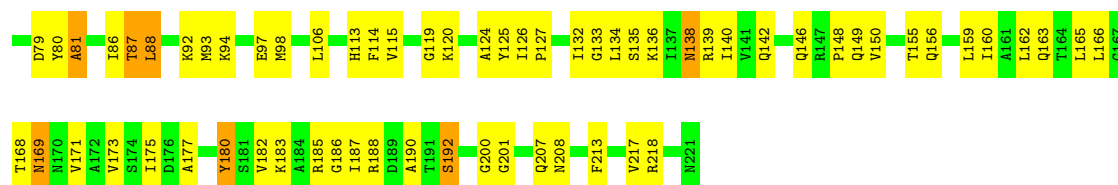


• Molecule 1: GTP cyclohydrolase I

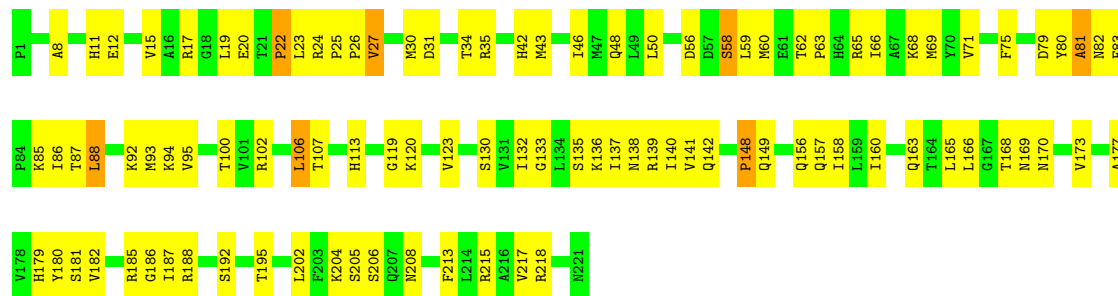


• Molecule 1: GTP cyclohydrolase I

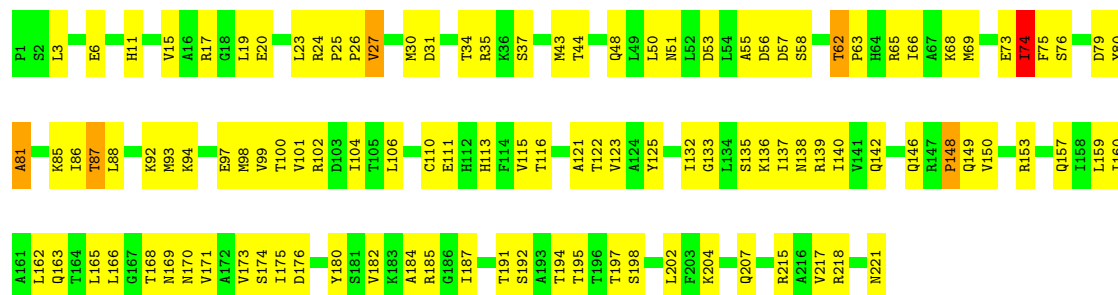




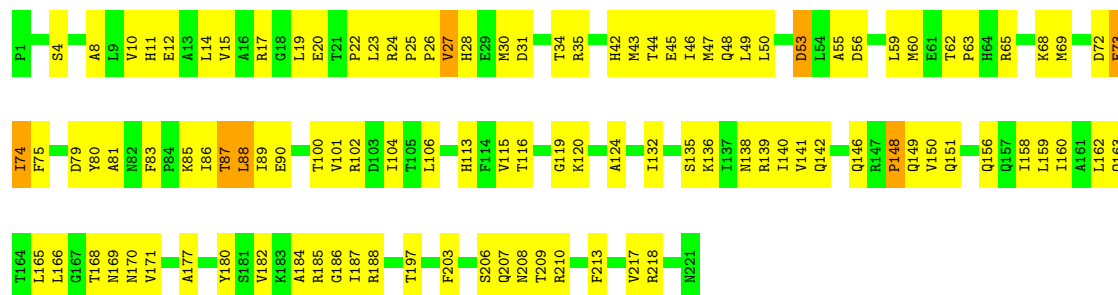
- Molecule 1: GTP cyclohydrolase I



- Molecule 1: GTP cyclohydrolase I

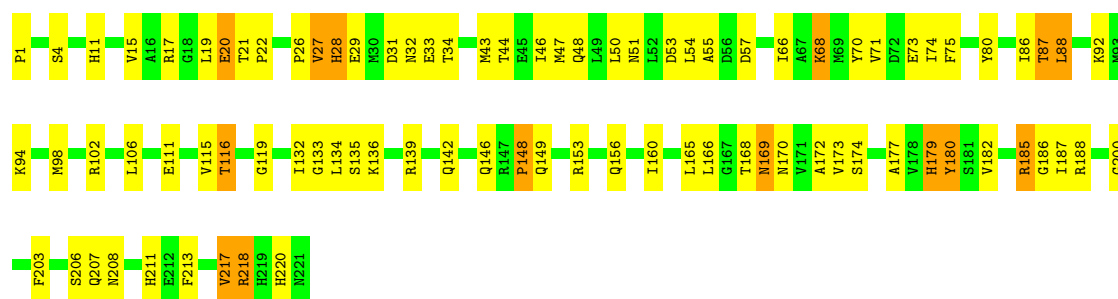


- Molecule 1: GTP cyclohydrolase I

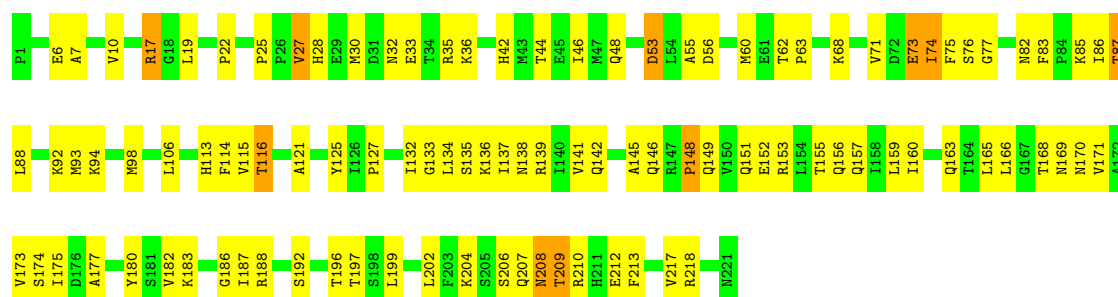


- Molecule 1: GTP cyclohydrolase I

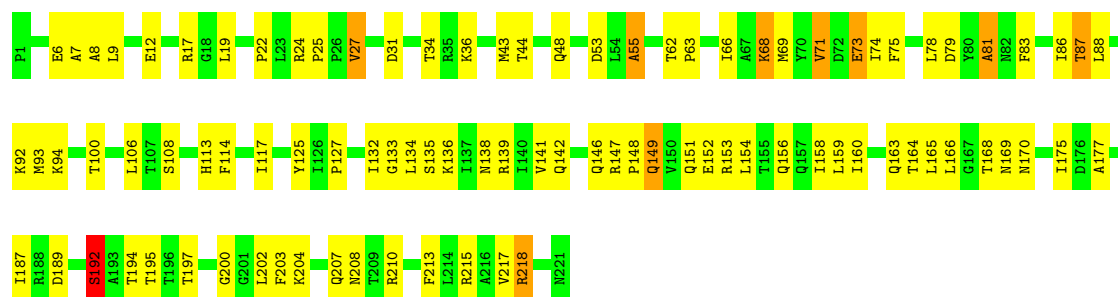




• Molecule 1: GTP cyclohydrolase I



• Molecule 1: GTP cyclohydrolase I

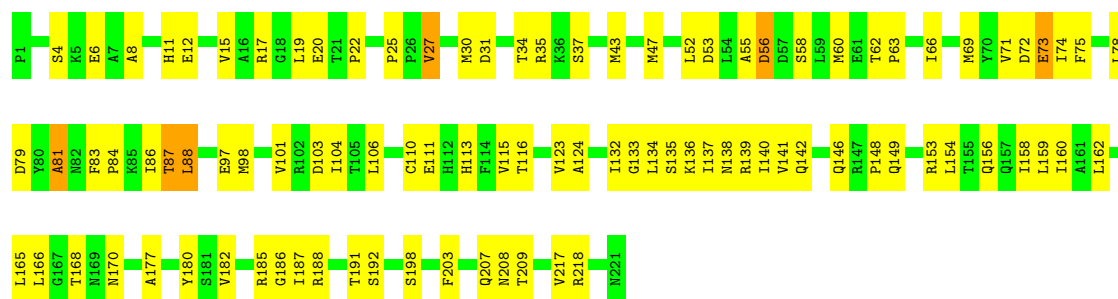


• Molecule 1: GTP cyclohydrolase I



• Molecule 1: GTP cyclohydrolase I

Chain E:



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	223.14Å 317.81Å 132.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.20	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-3.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.185 , 0.228	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	26475	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GTP

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	0.64	0/1761	1.17	16/2384 (0.7%)
1	B	0.60	0/1761	1.03	8/2384 (0.3%)
1	C	0.52	0/1761	1.03	10/2384 (0.4%)
1	D	0.57	0/1761	1.05	12/2384 (0.5%)
1	E	0.59	0/1761	1.05	8/2384 (0.3%)
1	F	0.50	0/1761	1.01	8/2384 (0.3%)
1	G	0.50	0/1761	1.04	13/2384 (0.5%)
1	H	0.50	0/1761	0.99	10/2384 (0.4%)
1	I	0.51	0/1761	1.02	7/2384 (0.3%)
1	J	0.53	0/1761	1.02	9/2384 (0.4%)
1	K	0.53	0/1761	1.01	8/2384 (0.3%)
1	L	0.53	0/1761	1.03	10/2384 (0.4%)
1	M	0.54	0/1761	1.02	9/2384 (0.4%)
1	N	0.53	0/1761	1.02	7/2384 (0.3%)
1	O	0.51	0/1761	0.99	4/2384 (0.2%)
All	All	0.54	0/26415	1.03	139/35760 (0.4%)

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	G	0	1
1	I	0	1
1	L	0	1
All	All	0	4

There are no bond length outliers.

The worst 5 of 139 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	149	GLN	N-CA-C	10.28	124.49	109.69
1	D	149	GLN	N-CA-C	9.84	124.87	109.81
1	G	73	GLU	N-CA-C	9.62	123.31	108.30
1	B	71	VAL	N-CA-C	9.03	119.09	110.42
1	E	149	GLN	N-CA-C	8.72	122.50	109.95

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	180	TYR	Sidechain
1	G	180	TYR	Sidechain
1	I	180	TYR	Sidechain
1	L	180	TYR	Sidechain

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	1733	0	1768	84	0
1	B	1733	0	1768	95	0
1	C	1733	0	1768	93	0
1	D	1733	0	1768	95	0
1	E	1733	0	1768	96	0
1	F	1733	0	1768	114	0
1	G	1733	0	1768	123	0
1	H	1733	0	1768	121	0
1	I	1733	0	1768	119	0
1	J	1733	0	1768	125	0
1	K	1733	0	1768	121	0
1	L	1733	0	1768	119	0
1	M	1733	0	1768	113	0
1	N	1733	0	1768	122	0
1	O	1733	0	1768	120	0
2	A	32	0	11	3	0
2	B	32	0	11	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	32	0	11	2	0
2	D	32	0	11	1	0
2	E	32	0	11	3	0
2	F	32	0	11	2	0
2	G	32	0	11	4	0
2	H	32	0	11	4	0
2	I	32	0	11	3	0
2	J	32	0	11	6	0
2	K	32	0	11	4	0
2	L	32	0	11	4	0
2	M	32	0	11	4	0
2	N	32	0	11	6	0
2	O	32	0	11	3	0
All	All	26475	0	26685	1404	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 26.

The worst 5 of 1404 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:H:168:THR:HG22	1:H:170:ASN:H	1.14	1.06
1:A:86:ILE:HD13	1:A:165:LEU:HD13	1.37	1.05
1:O:168:THR:HG22	1:O:170:ASN:H	1.23	1.03
1:D:187:ILE:HD11	1:E:139:ARG:HG2	1.41	1.03
1:F:168:THR:HG22	1:F:170:ASN:H	1.21	1.02

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	219/221 (99%)	193 (88%)	22 (10%)	4 (2%)	6	34
1	B	219/221 (99%)	187 (85%)	25 (11%)	7 (3%)	3	21
1	C	219/221 (99%)	192 (88%)	25 (11%)	2 (1%)	14	47
1	D	219/221 (99%)	196 (90%)	20 (9%)	3 (1%)	9	39
1	E	219/221 (99%)	196 (90%)	18 (8%)	5 (2%)	5	29
1	F	219/221 (99%)	193 (88%)	20 (9%)	6 (3%)	4	25
1	G	219/221 (99%)	196 (90%)	19 (9%)	4 (2%)	6	34
1	H	219/221 (99%)	194 (89%)	22 (10%)	3 (1%)	9	39
1	I	219/221 (99%)	195 (89%)	19 (9%)	5 (2%)	5	29
1	J	219/221 (99%)	192 (88%)	23 (10%)	4 (2%)	6	34
1	K	219/221 (99%)	191 (87%)	23 (10%)	5 (2%)	5	29
1	L	219/221 (99%)	196 (90%)	18 (8%)	5 (2%)	5	29
1	M	219/221 (99%)	187 (85%)	29 (13%)	3 (1%)	9	39
1	N	219/221 (99%)	187 (85%)	27 (12%)	5 (2%)	5	29
1	O	219/221 (99%)	187 (85%)	26 (12%)	6 (3%)	4	25
All	All	3285/3315 (99%)	2882 (88%)	336 (10%)	67 (2%)	6	31

5 of 67 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	53	ASP
1	F	53	ASP
1	F	80	TYR
1	F	116	THR
1	G	27	VAL

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	194/194 (100%)	184 (95%)	10 (5%)	21	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	194/194 (100%)	189 (97%)	5 (3%)	40	69
1	C	194/194 (100%)	188 (97%)	6 (3%)	35	66
1	D	194/194 (100%)	188 (97%)	6 (3%)	35	66
1	E	194/194 (100%)	189 (97%)	5 (3%)	40	69
1	F	194/194 (100%)	188 (97%)	6 (3%)	35	66
1	G	194/194 (100%)	187 (96%)	7 (4%)	31	63
1	H	194/194 (100%)	185 (95%)	9 (5%)	24	58
1	I	194/194 (100%)	189 (97%)	5 (3%)	40	69
1	J	194/194 (100%)	188 (97%)	6 (3%)	35	66
1	K	194/194 (100%)	185 (95%)	9 (5%)	24	58
1	L	194/194 (100%)	187 (96%)	7 (4%)	31	63
1	M	194/194 (100%)	187 (96%)	7 (4%)	31	63
1	N	194/194 (100%)	187 (96%)	7 (4%)	31	63
1	O	194/194 (100%)	187 (96%)	7 (4%)	31	63
All	All	2910/2910 (100%)	2808 (96%)	102 (4%)	32	64

5 of 102 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	N	37	SER
1	A	20	GLU
1	E	73	GLU
1	N	68	LYS
1	O	68	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 64 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	163	GLN
1	D	208	ASN
1	K	146	GLN
1	K	138	ASN
1	E	48	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

15 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GTP	M	417	-	33,34,34	3.03	13 (39%)	50,54,54	2.94	18 (36%)
2	GTP	O	419	-	33,34,34	2.99	12 (36%)	50,54,54	2.95	18 (36%)
2	GTP	N	418	-	33,34,34	3.13	14 (42%)	50,54,54	2.95	16 (32%)
2	GTP	K	420	-	33,34,34	3.11	16 (48%)	50,54,54	3.03	18 (36%)
2	GTP	I	413	-	33,34,34	3.17	14 (42%)	50,54,54	2.93	17 (34%)
2	GTP	E	424	-	33,34,34	3.04	14 (42%)	50,54,54	2.75	17 (34%)
2	GTP	A	425	-	33,34,34	3.22	14 (42%)	50,54,54	2.80	18 (36%)
2	GTP	J	414	-	33,34,34	3.02	14 (42%)	50,54,54	2.93	18 (36%)
2	GTP	D	423	-	33,34,34	3.11	13 (39%)	50,54,54	2.94	17 (34%)
2	GTP	B	421	-	33,34,34	3.33	15 (45%)	50,54,54	2.93	18 (36%)
2	GTP	C	422	-	33,34,34	3.29	13 (39%)	50,54,54	2.99	19 (38%)
2	GTP	H	412	-	33,34,34	3.03	15 (45%)	50,54,54	2.94	17 (34%)
2	GTP	G	411	-	33,34,34	3.08	14 (42%)	50,54,54	2.96	17 (34%)
2	GTP	L	416	-	33,34,34	3.18	13 (39%)	50,54,54	2.96	18 (36%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GTP	F	415	-	33,34,34	3.23	13 (39%)	50,54,54	2.99	18 (36%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GTP	M	417	-	-	4/22/38/38	0/3/3/3
2	GTP	O	419	-	-	7/22/38/38	0/3/3/3
2	GTP	N	418	-	-	4/22/38/38	0/3/3/3
2	GTP	K	420	-	-	4/22/38/38	0/3/3/3
2	GTP	I	413	-	-	4/22/38/38	0/3/3/3
2	GTP	E	424	-	-	6/22/38/38	0/3/3/3
2	GTP	A	425	-	-	4/22/38/38	0/3/3/3
2	GTP	J	414	-	-	4/22/38/38	0/3/3/3
2	GTP	D	423	-	-	4/22/38/38	0/3/3/3
2	GTP	B	421	-	-	4/22/38/38	0/3/3/3
2	GTP	C	422	-	-	5/22/38/38	0/3/3/3
2	GTP	H	412	-	-	4/22/38/38	0/3/3/3
2	GTP	G	411	-	-	4/22/38/38	0/3/3/3
2	GTP	L	416	-	-	4/22/38/38	0/3/3/3
2	GTP	F	415	-	-	4/22/38/38	0/3/3/3

The worst 5 of 207 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	422	GTP	PB-O3B	9.65	1.69	1.59
2	F	415	GTP	PB-O3B	9.02	1.69	1.59
2	B	421	GTP	PB-O3B	8.92	1.69	1.59
2	L	416	GTP	PB-O3B	8.83	1.69	1.59
2	C	422	GTP	C6-N1	8.79	1.55	1.38

The worst 5 of 264 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	417	GTP	N9-C8-N7	-8.54	97.55	113.40
2	G	411	GTP	N9-C8-N7	-8.50	97.64	113.40

Continued on next page...

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	419	GTP	N9-C8-N7	-8.44	97.75	113.40
2	F	415	GTP	N9-C8-N7	-8.39	97.84	113.40
2	N	418	GTP	N9-C8-N7	-8.37	97.87	113.40

There are no chirality outliers.

5 of 66 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	412	GTP	O4'-C4'-C5'-O5'
2	I	413	GTP	O4'-C4'-C5'-O5'
2	K	420	GTP	O4'-C4'-C5'-O5'
2	L	416	GTP	O4'-C4'-C5'-O5'
2	M	417	GTP	O4'-C4'-C5'-O5'

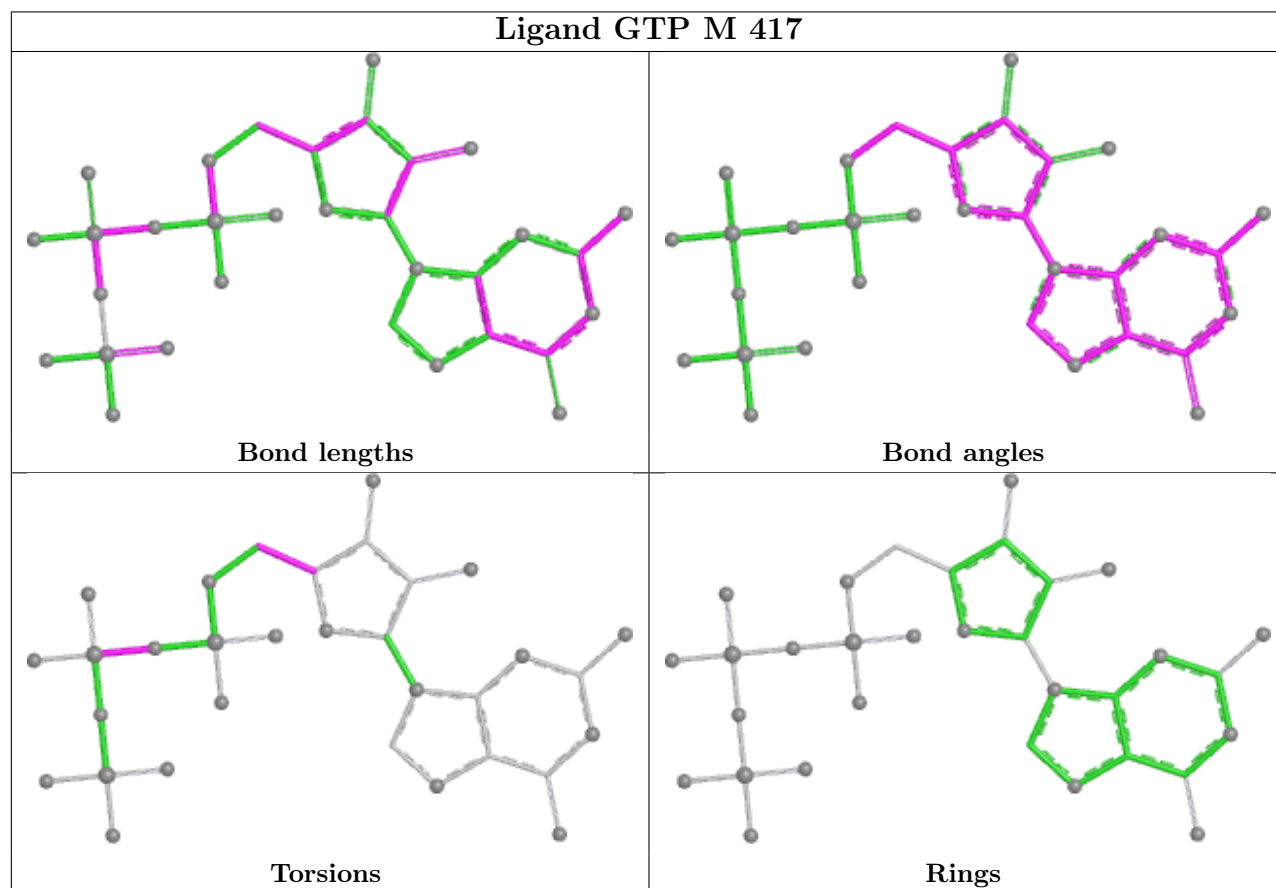
There are no ring outliers.

15 monomers are involved in 51 short contacts:

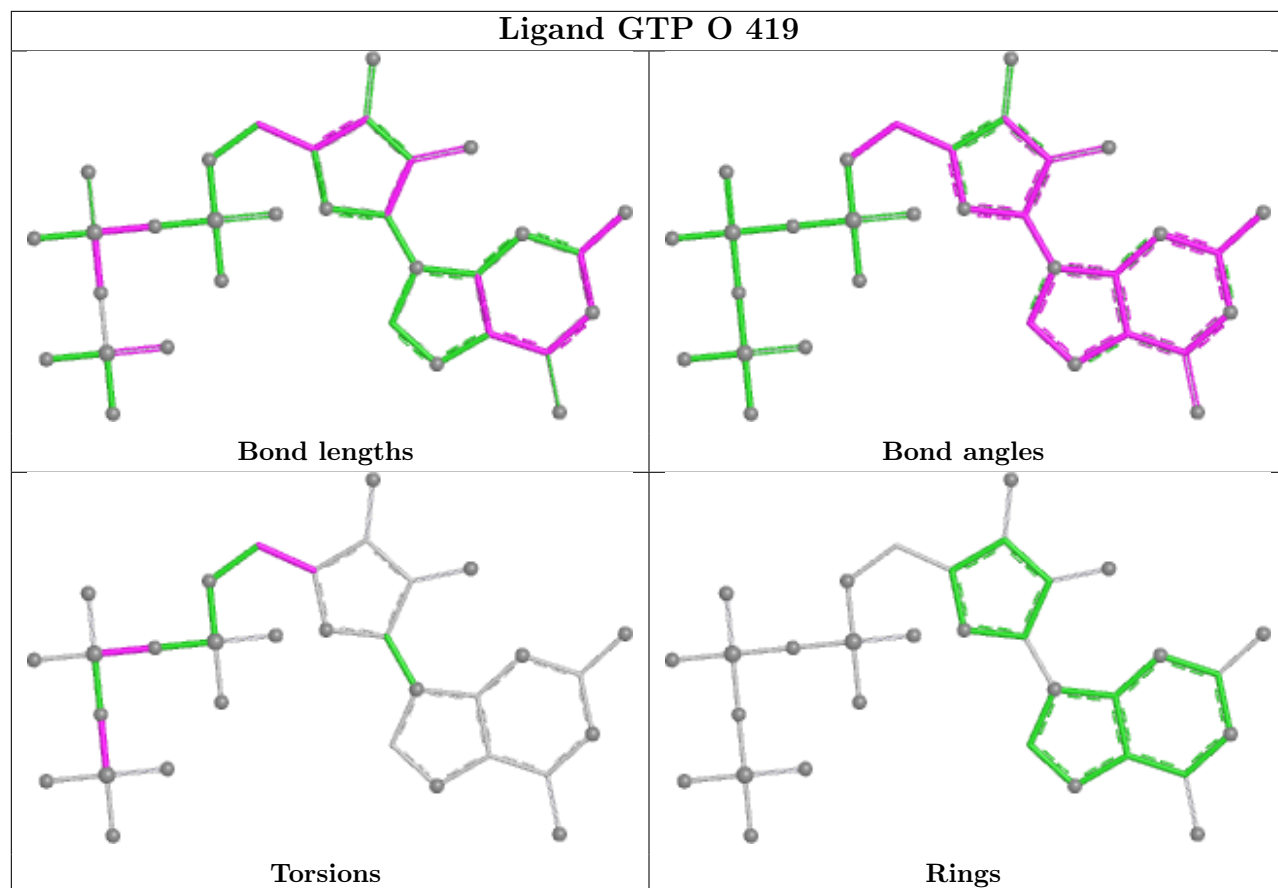
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	M	417	GTP	4	0
2	O	419	GTP	3	0
2	N	418	GTP	6	0
2	K	420	GTP	4	0
2	I	413	GTP	3	0
2	E	424	GTP	3	0
2	A	425	GTP	3	0
2	J	414	GTP	6	0
2	D	423	GTP	1	0
2	B	421	GTP	2	0
2	C	422	GTP	2	0
2	H	412	GTP	4	0
2	G	411	GTP	4	0
2	L	416	GTP	4	0
2	F	415	GTP	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

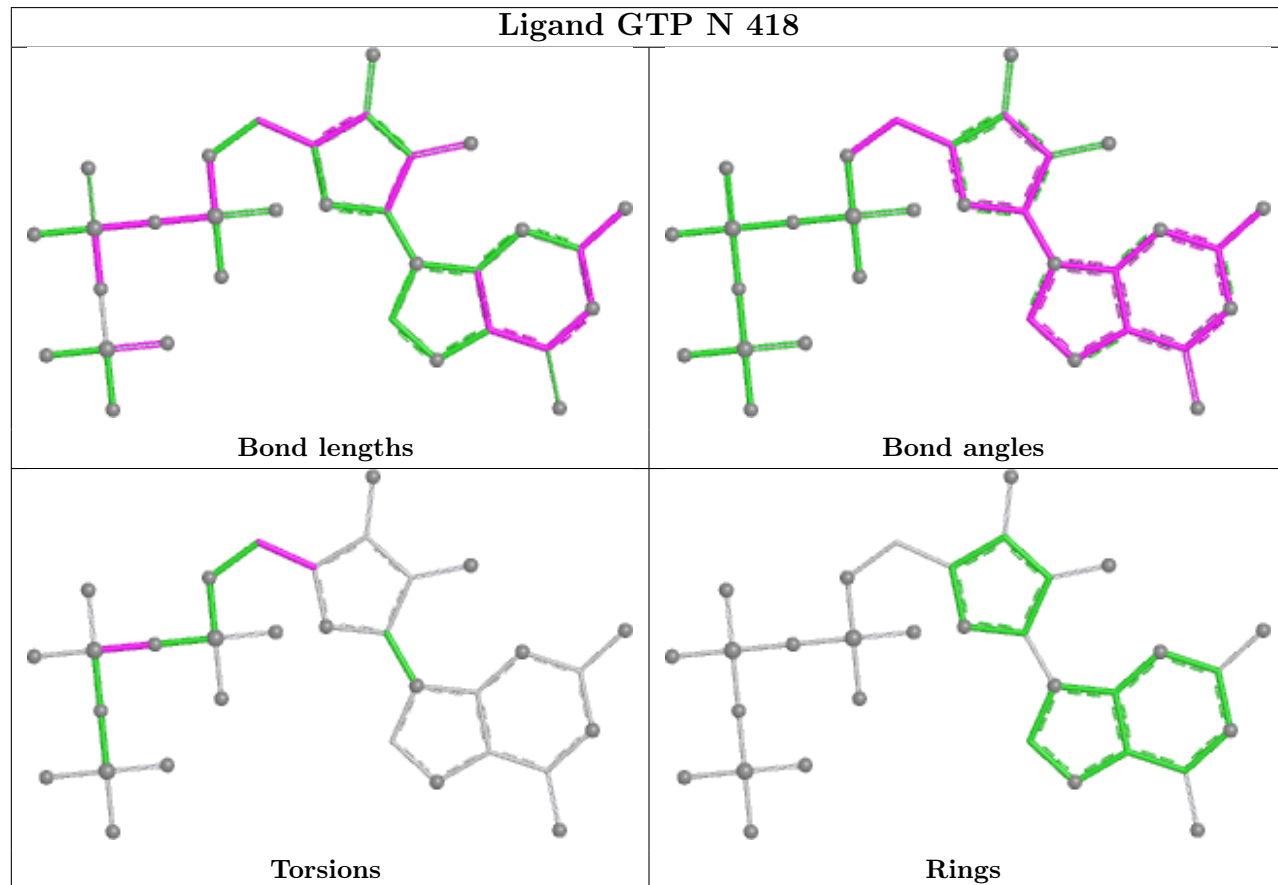
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



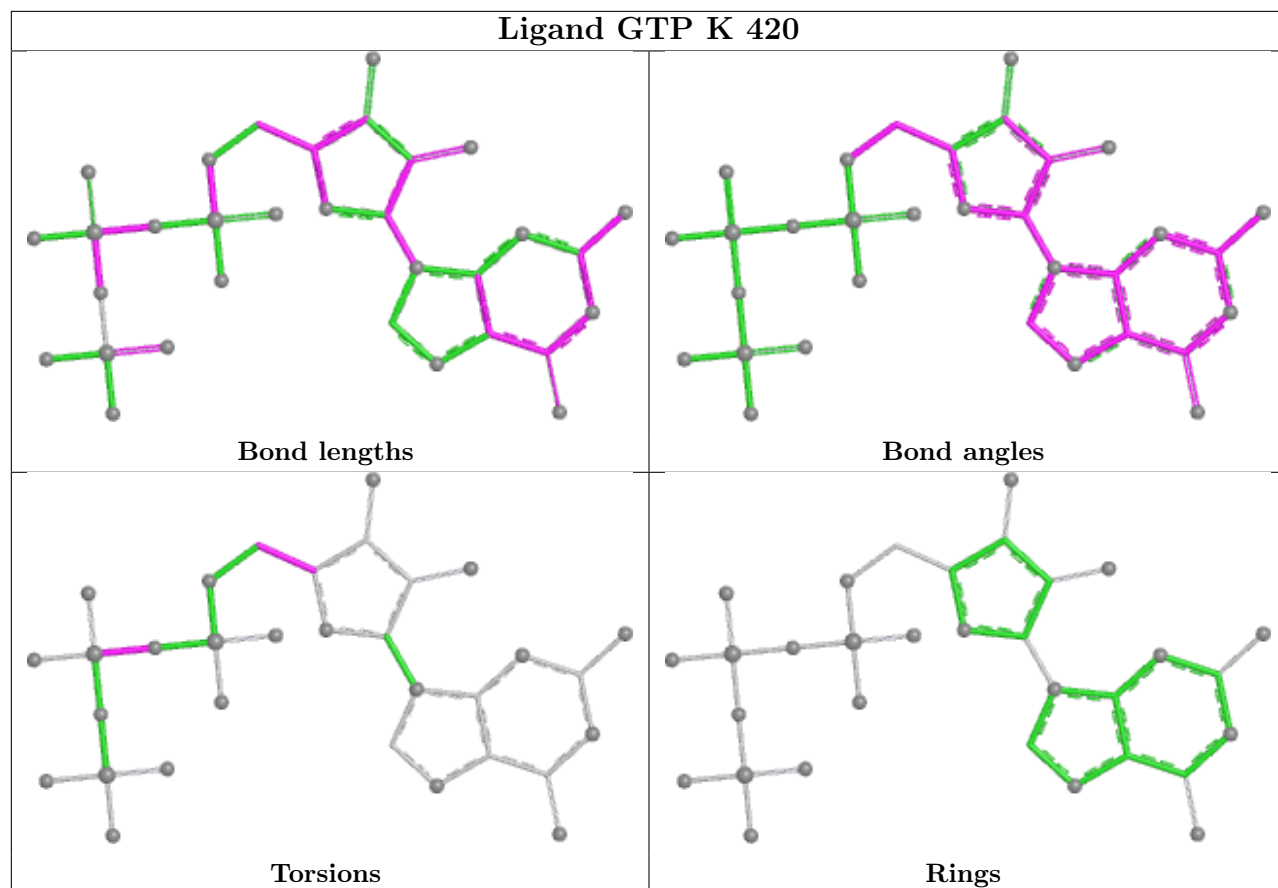
Ligand GTP O 419



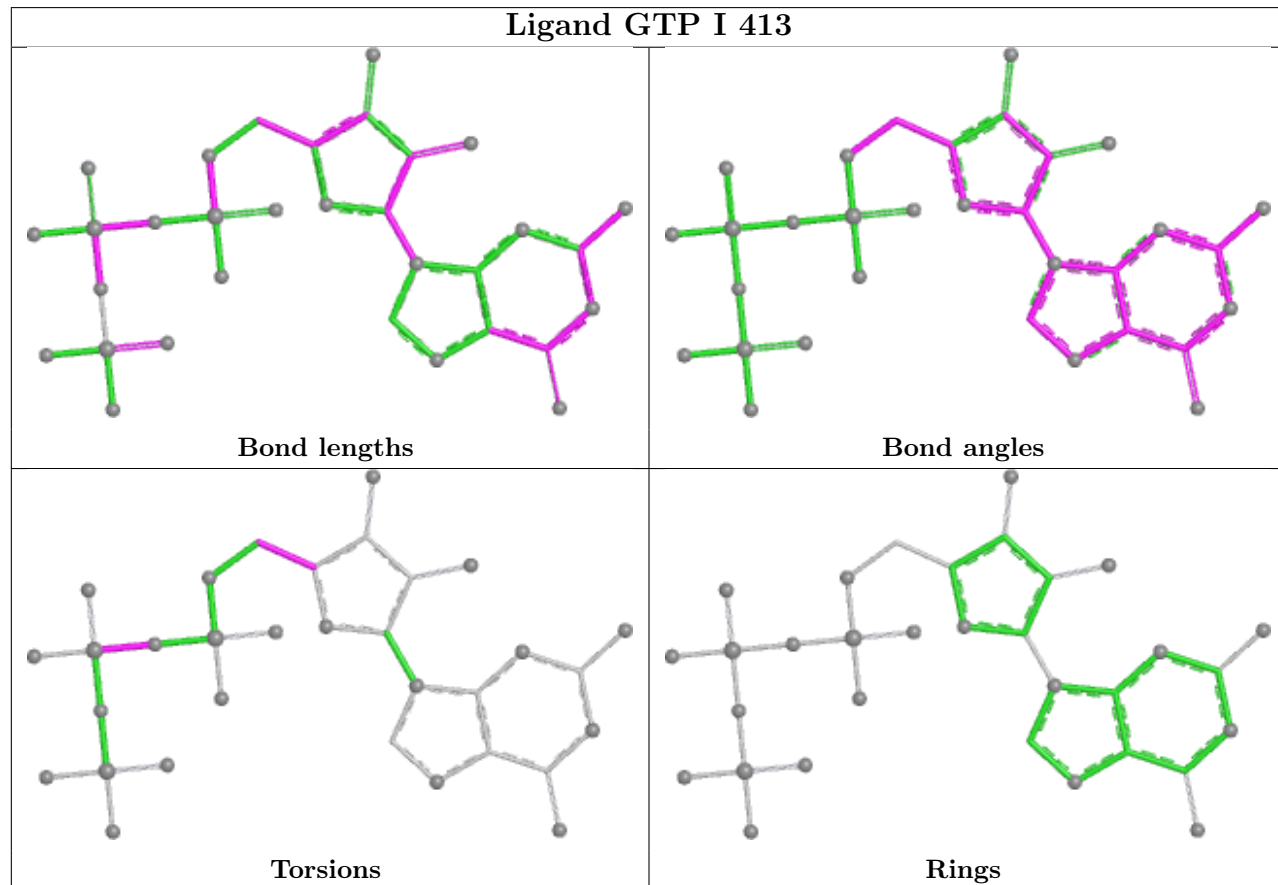
Ligand GTP N 418



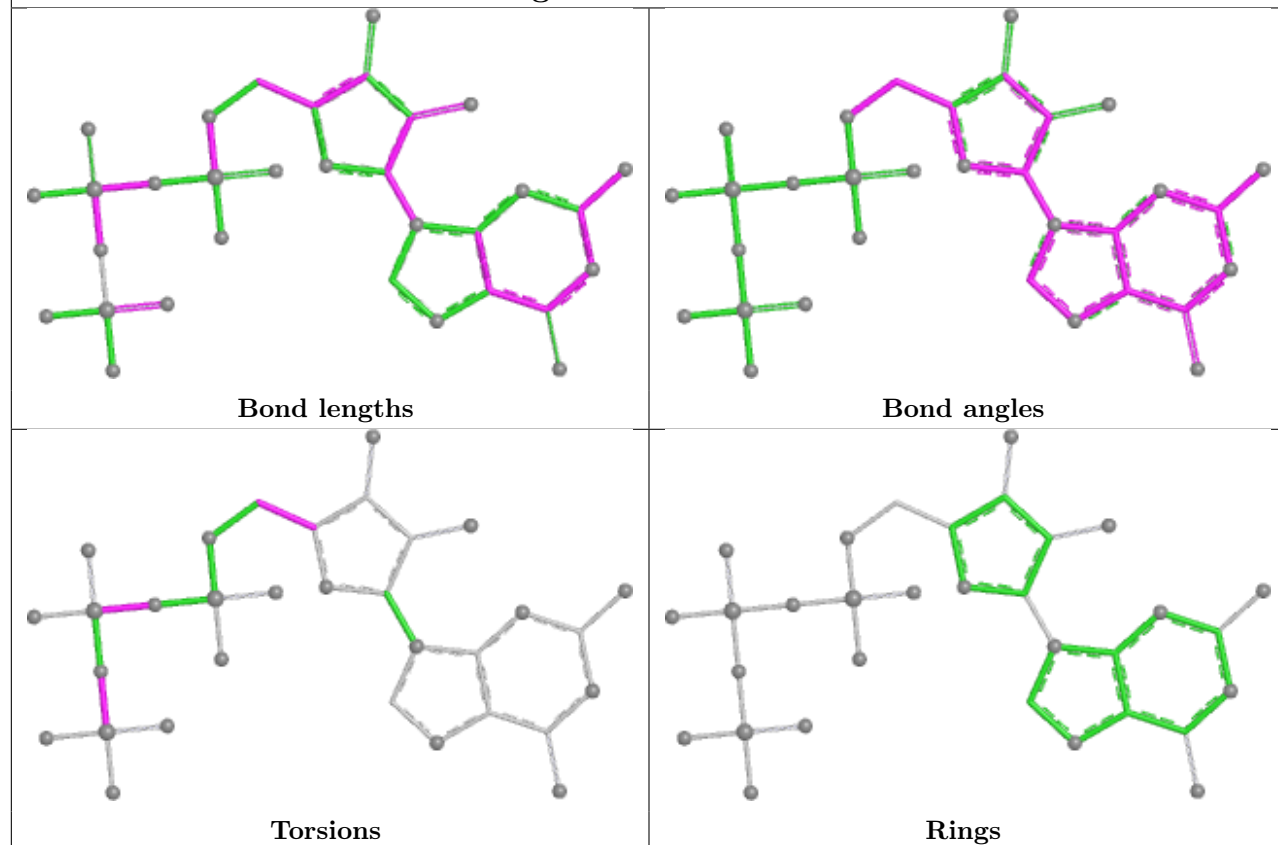
Ligand GTP K 420



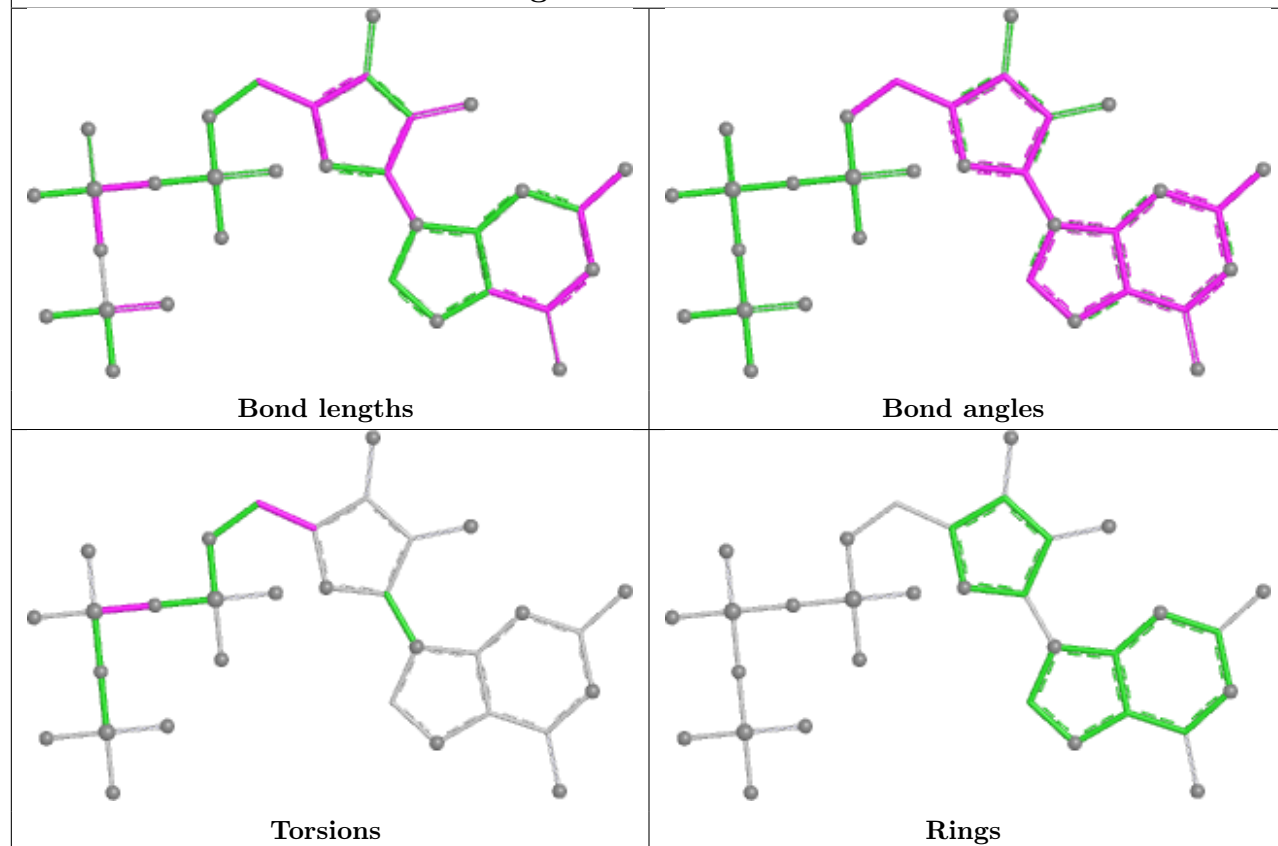
Ligand GTP I 413



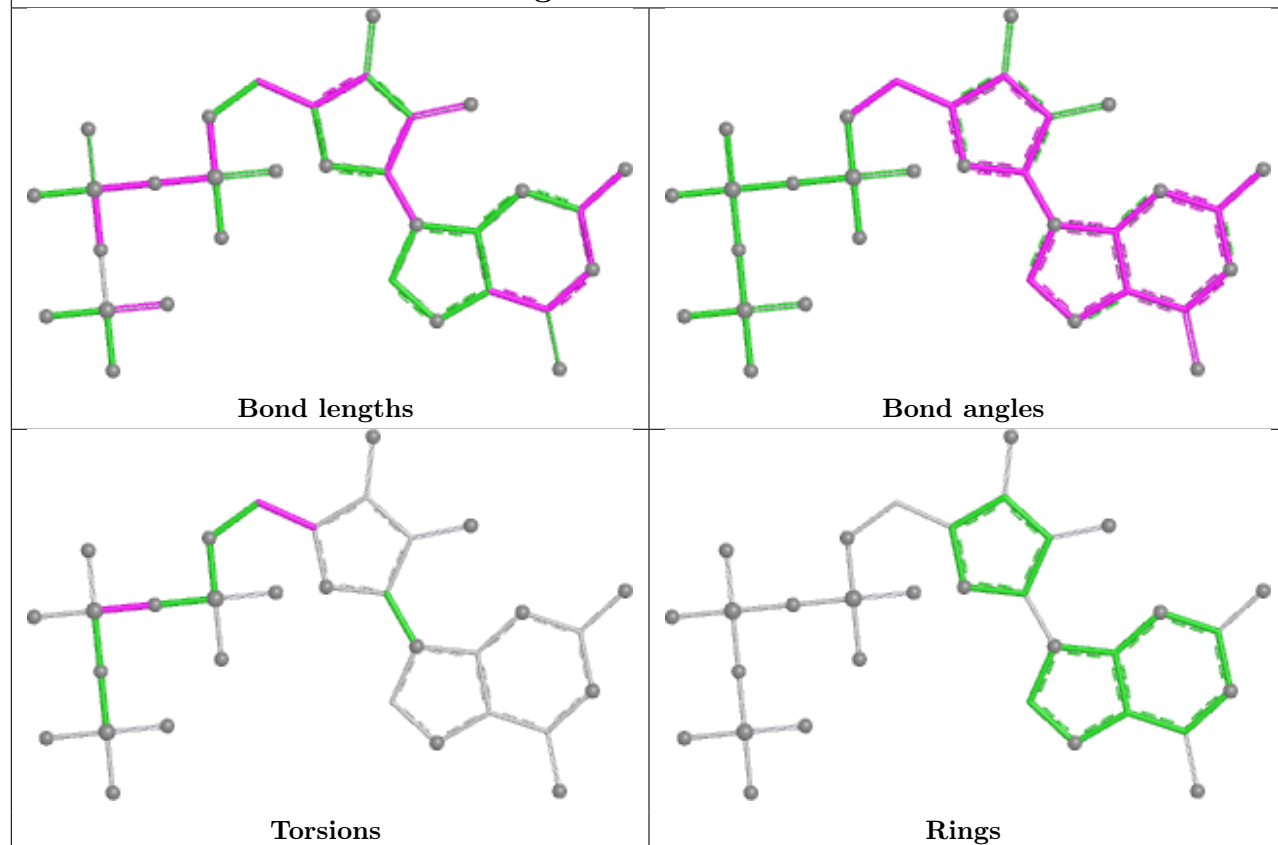
Ligand GTP E 424



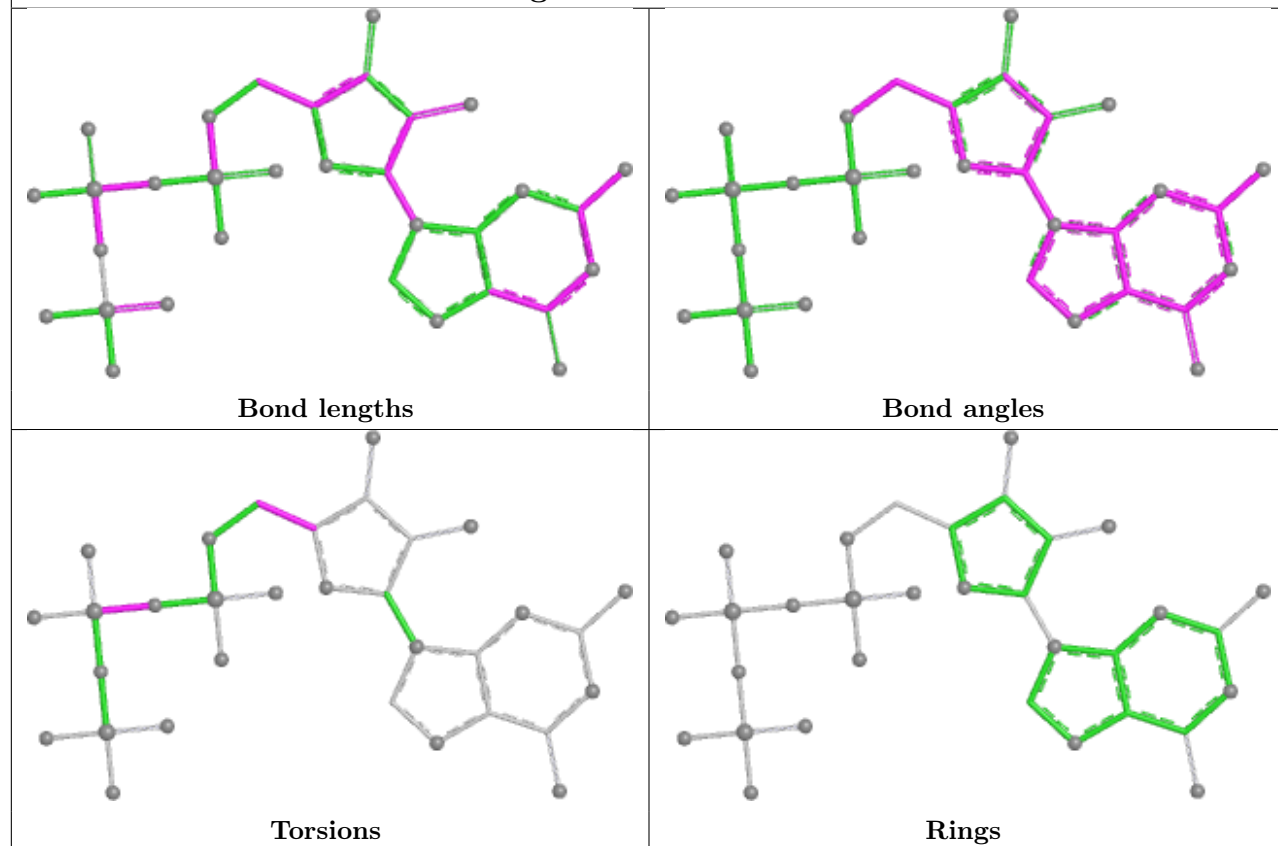
Ligand GTP A 425



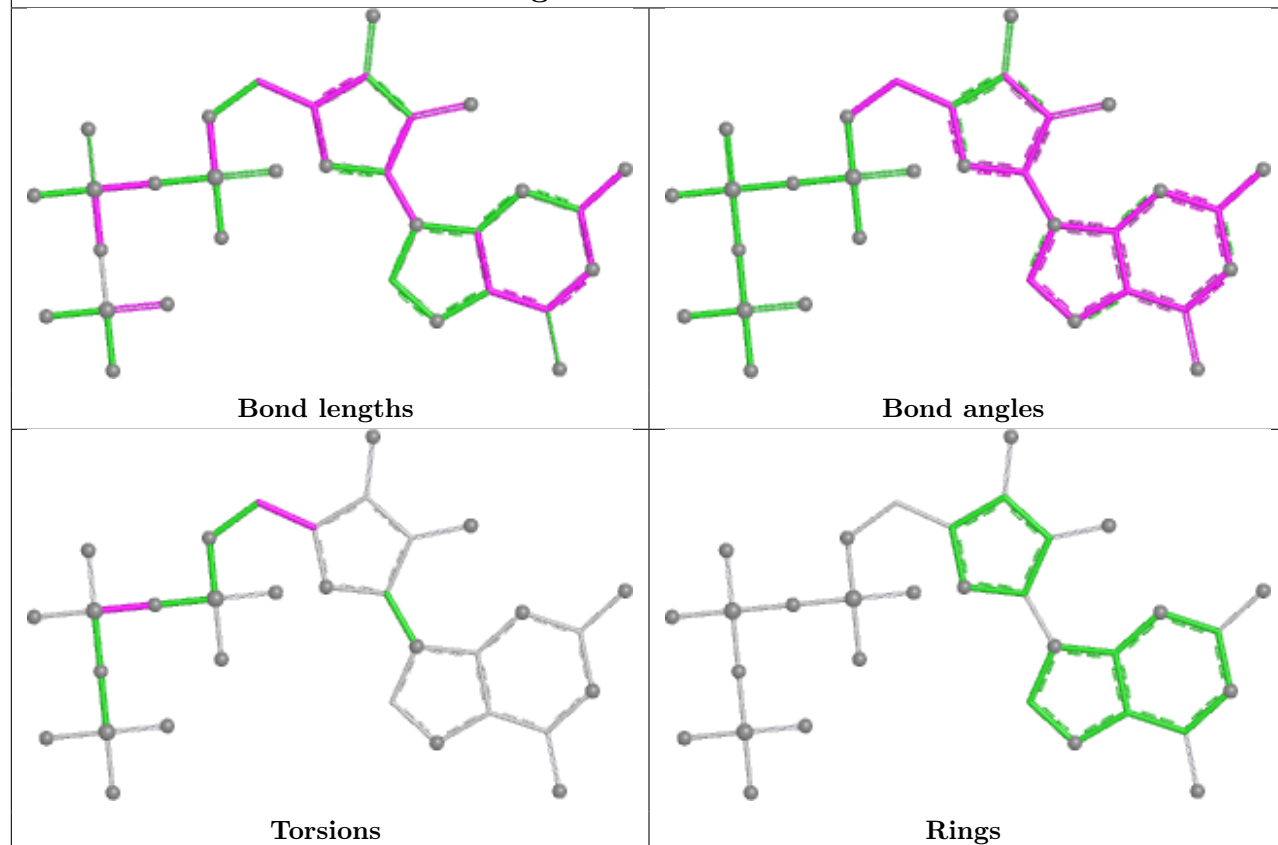
Ligand GTP J 414



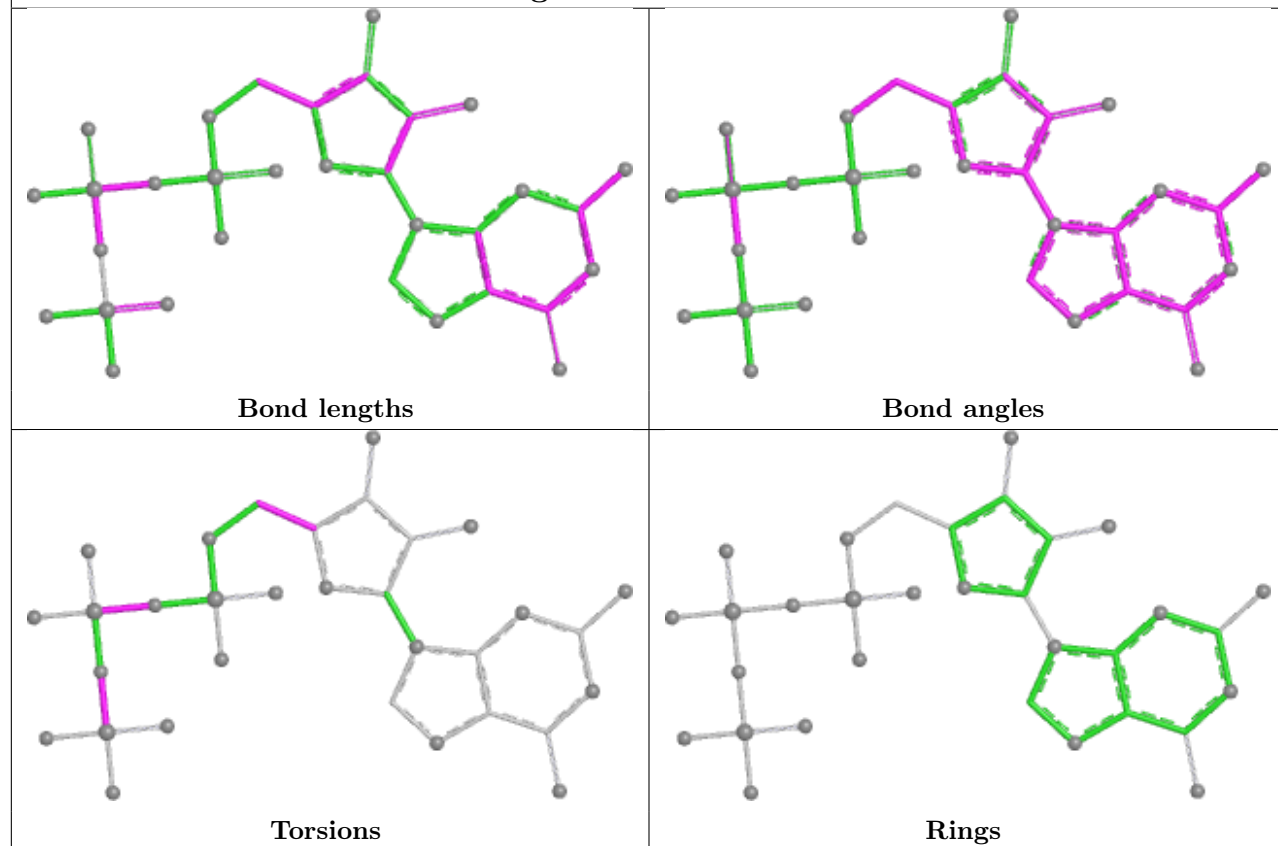
Ligand GTP D 423



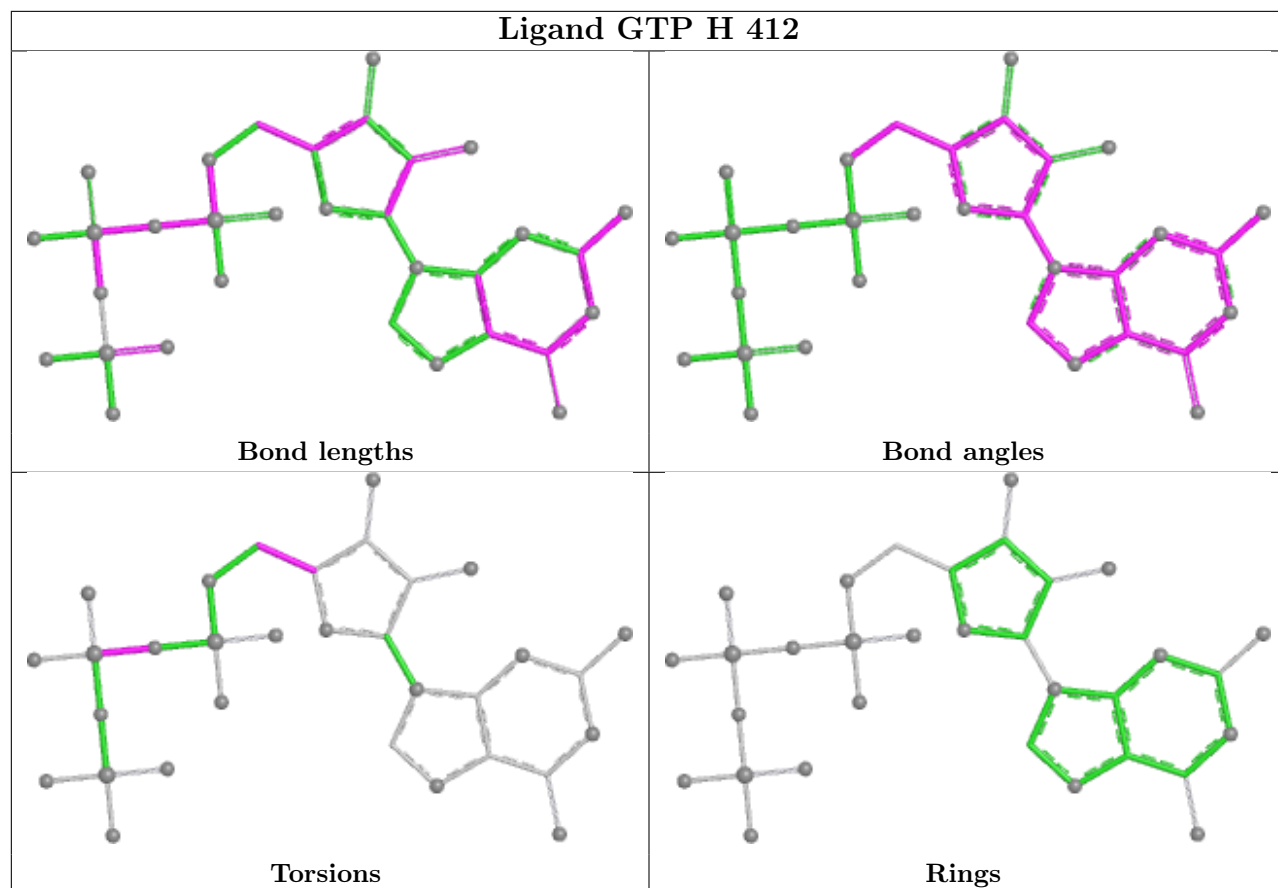
Ligand GTP B 421



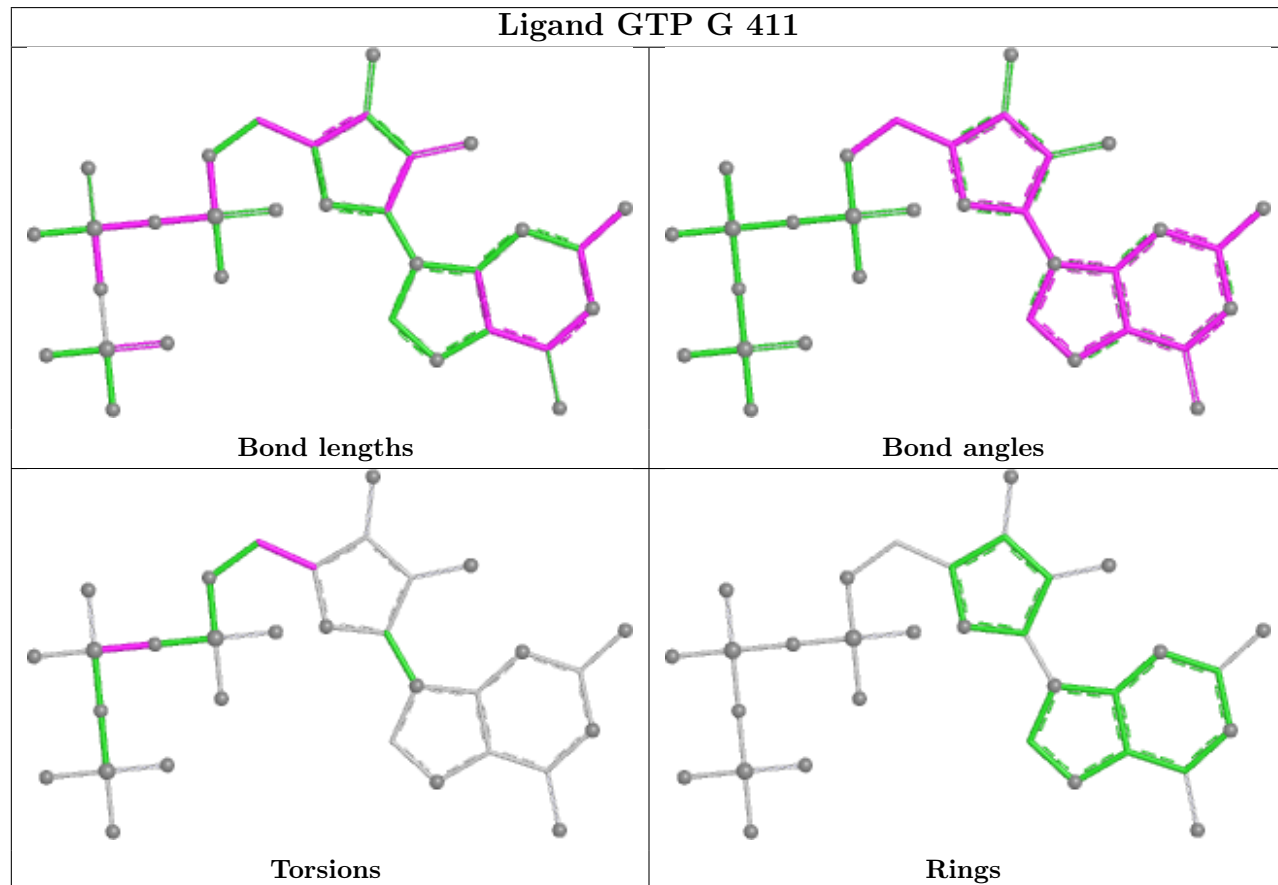
Ligand GTP C 422



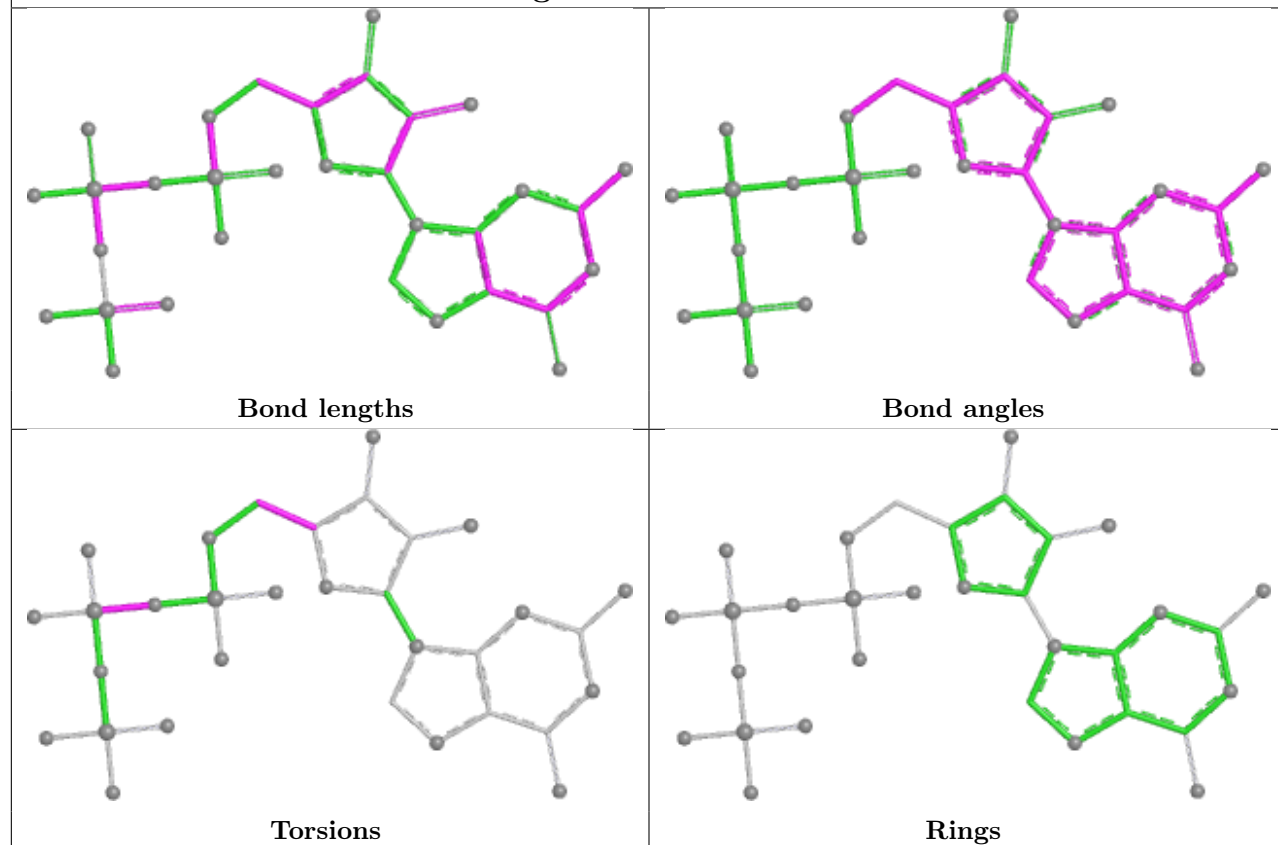
Ligand GTP H 412



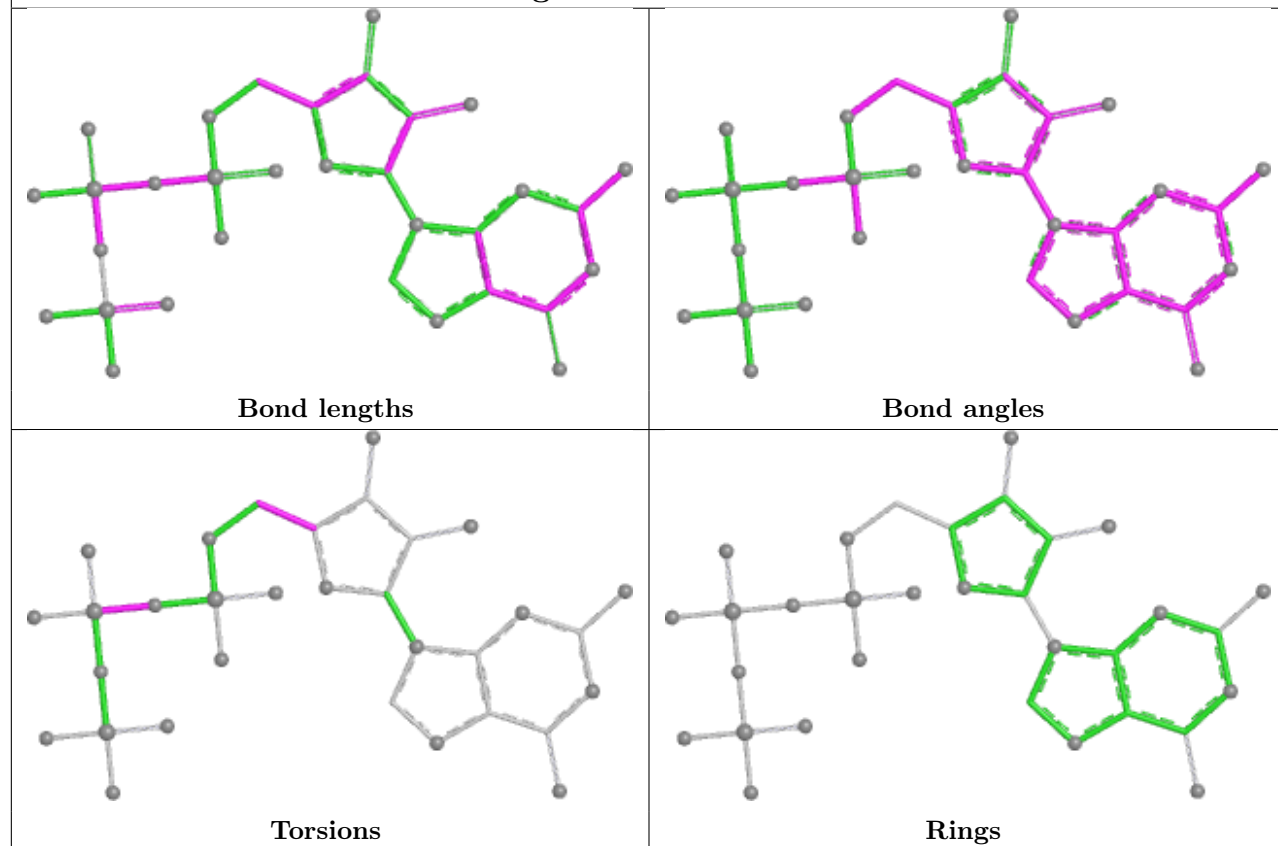
Ligand GTP G 411



Ligand GTP L 416



Ligand GTP F 415



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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.