



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:59 PM UTC

PDB ID : 4QG0 / pdb_00004qg0
Title : Crystal structure of the tetrameric dGTP/dUTP-bound SAMHD1 (RN206) mutant catalytic core
Authors : Koharudin, L.M.I.; Wu, Y.; DeLucia, M.; Mehrens, J.; Gronenborn, A.M.; Ahn, J.
Deposited on : 2014-05-22
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

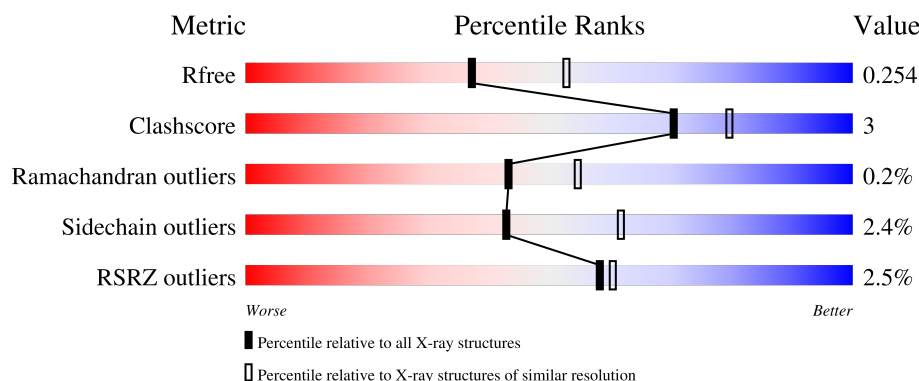
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

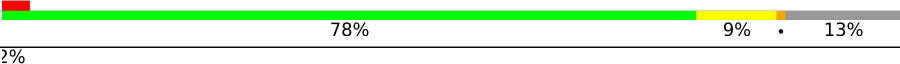
The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	550	 2% 78% 9% 13%
1	B	550	 3% 78% 9% 13%
1	C	550	 2% 75% 12% 13%
1	D	550	 % 80% 7% 13%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 16309 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 Deoxynucleoside triphosphate triphosphohydrolase SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	2	0
			3946	2527	688	711	20			
1	B	480	Total	C	N	O	S	0	2	0
			3938	2521	688	709	20			
1	C	481	Total	C	N	O	S	0	2	0
			3945	2525	686	714	20			
1	D	481	Total	C	N	O	S	0	1	0
			3939	2521	685	713	20			

There are 152 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	77	MET	-	expression tag	UNP Q9Y3Z3
A	78	GLY	-	expression tag	UNP Q9Y3Z3
A	79	SER	-	expression tag	UNP Q9Y3Z3
A	80	SER	-	expression tag	UNP Q9Y3Z3
A	81	HIS	-	expression tag	UNP Q9Y3Z3
A	82	HIS	-	expression tag	UNP Q9Y3Z3
A	83	HIS	-	expression tag	UNP Q9Y3Z3
A	84	HIS	-	expression tag	UNP Q9Y3Z3
A	85	HIS	-	expression tag	UNP Q9Y3Z3
A	86	HIS	-	expression tag	UNP Q9Y3Z3
A	87	SER	-	expression tag	UNP Q9Y3Z3
A	88	SER	-	expression tag	UNP Q9Y3Z3
A	89	GLY	-	expression tag	UNP Q9Y3Z3
A	90	LEU	-	expression tag	UNP Q9Y3Z3
A	91	VAL	-	expression tag	UNP Q9Y3Z3
A	92	PRO	-	expression tag	UNP Q9Y3Z3
A	93	ARG	-	expression tag	UNP Q9Y3Z3
A	94	GLY	-	expression tag	UNP Q9Y3Z3
A	95	SER	-	expression tag	UNP Q9Y3Z3
A	96	HIS	-	expression tag	UNP Q9Y3Z3
A	97	MET	-	expression tag	UNP Q9Y3Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
A	98	ALA	-	expression tag	UNP Q9Y3Z3
A	99	SER	-	expression tag	UNP Q9Y3Z3
A	100	MET	-	expression tag	UNP Q9Y3Z3
A	101	THR	-	expression tag	UNP Q9Y3Z3
A	102	GLY	-	expression tag	UNP Q9Y3Z3
A	103	GLY	-	expression tag	UNP Q9Y3Z3
A	104	GLN	-	expression tag	UNP Q9Y3Z3
A	105	GLN	-	expression tag	UNP Q9Y3Z3
A	106	MET	-	expression tag	UNP Q9Y3Z3
A	107	GLY	-	expression tag	UNP Q9Y3Z3
A	108	ARG	-	expression tag	UNP Q9Y3Z3
A	109	ASP	-	expression tag	UNP Q9Y3Z3
A	110	PRO	-	expression tag	UNP Q9Y3Z3
A	111	ASN	-	expression tag	UNP Q9Y3Z3
A	112	SER	-	expression tag	UNP Q9Y3Z3
A	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
A	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
B	77	MET	-	expression tag	UNP Q9Y3Z3
B	78	GLY	-	expression tag	UNP Q9Y3Z3
B	79	SER	-	expression tag	UNP Q9Y3Z3
B	80	SER	-	expression tag	UNP Q9Y3Z3
B	81	HIS	-	expression tag	UNP Q9Y3Z3
B	82	HIS	-	expression tag	UNP Q9Y3Z3
B	83	HIS	-	expression tag	UNP Q9Y3Z3
B	84	HIS	-	expression tag	UNP Q9Y3Z3
B	85	HIS	-	expression tag	UNP Q9Y3Z3
B	86	HIS	-	expression tag	UNP Q9Y3Z3
B	87	SER	-	expression tag	UNP Q9Y3Z3
B	88	SER	-	expression tag	UNP Q9Y3Z3
B	89	GLY	-	expression tag	UNP Q9Y3Z3
B	90	LEU	-	expression tag	UNP Q9Y3Z3
B	91	VAL	-	expression tag	UNP Q9Y3Z3
B	92	PRO	-	expression tag	UNP Q9Y3Z3
B	93	ARG	-	expression tag	UNP Q9Y3Z3
B	94	GLY	-	expression tag	UNP Q9Y3Z3
B	95	SER	-	expression tag	UNP Q9Y3Z3
B	96	HIS	-	expression tag	UNP Q9Y3Z3
B	97	MET	-	expression tag	UNP Q9Y3Z3
B	98	ALA	-	expression tag	UNP Q9Y3Z3
B	99	SER	-	expression tag	UNP Q9Y3Z3
B	100	MET	-	expression tag	UNP Q9Y3Z3
B	101	THR	-	expression tag	UNP Q9Y3Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	102	GLY	-	expression tag	UNP Q9Y3Z3
B	103	GLY	-	expression tag	UNP Q9Y3Z3
B	104	GLN	-	expression tag	UNP Q9Y3Z3
B	105	GLN	-	expression tag	UNP Q9Y3Z3
B	106	MET	-	expression tag	UNP Q9Y3Z3
B	107	GLY	-	expression tag	UNP Q9Y3Z3
B	108	ARG	-	expression tag	UNP Q9Y3Z3
B	109	ASP	-	expression tag	UNP Q9Y3Z3
B	110	PRO	-	expression tag	UNP Q9Y3Z3
B	111	ASN	-	expression tag	UNP Q9Y3Z3
B	112	SER	-	expression tag	UNP Q9Y3Z3
B	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
B	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
C	77	MET	-	expression tag	UNP Q9Y3Z3
C	78	GLY	-	expression tag	UNP Q9Y3Z3
C	79	SER	-	expression tag	UNP Q9Y3Z3
C	80	SER	-	expression tag	UNP Q9Y3Z3
C	81	HIS	-	expression tag	UNP Q9Y3Z3
C	82	HIS	-	expression tag	UNP Q9Y3Z3
C	83	HIS	-	expression tag	UNP Q9Y3Z3
C	84	HIS	-	expression tag	UNP Q9Y3Z3
C	85	HIS	-	expression tag	UNP Q9Y3Z3
C	86	HIS	-	expression tag	UNP Q9Y3Z3
C	87	SER	-	expression tag	UNP Q9Y3Z3
C	88	SER	-	expression tag	UNP Q9Y3Z3
C	89	GLY	-	expression tag	UNP Q9Y3Z3
C	90	LEU	-	expression tag	UNP Q9Y3Z3
C	91	VAL	-	expression tag	UNP Q9Y3Z3
C	92	PRO	-	expression tag	UNP Q9Y3Z3
C	93	ARG	-	expression tag	UNP Q9Y3Z3
C	94	GLY	-	expression tag	UNP Q9Y3Z3
C	95	SER	-	expression tag	UNP Q9Y3Z3
C	96	HIS	-	expression tag	UNP Q9Y3Z3
C	97	MET	-	expression tag	UNP Q9Y3Z3
C	98	ALA	-	expression tag	UNP Q9Y3Z3
C	99	SER	-	expression tag	UNP Q9Y3Z3
C	100	MET	-	expression tag	UNP Q9Y3Z3
C	101	THR	-	expression tag	UNP Q9Y3Z3
C	102	GLY	-	expression tag	UNP Q9Y3Z3
C	103	GLY	-	expression tag	UNP Q9Y3Z3
C	104	GLN	-	expression tag	UNP Q9Y3Z3
C	105	GLN	-	expression tag	UNP Q9Y3Z3

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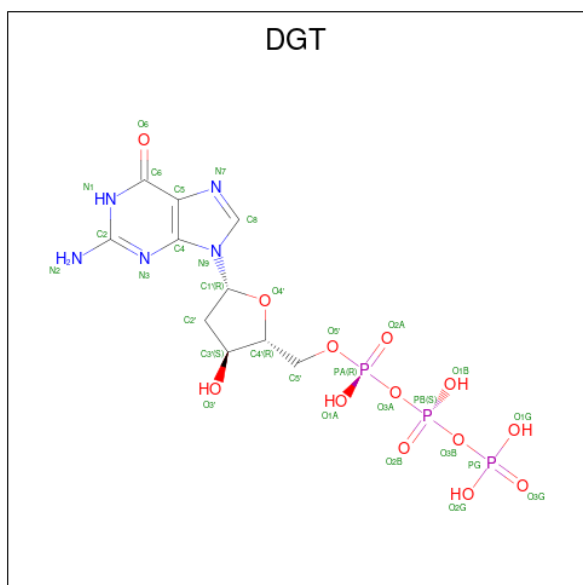
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C	106	MET	-	expression tag	UNP Q9Y3Z3
C	107	GLY	-	expression tag	UNP Q9Y3Z3
C	108	ARG	-	expression tag	UNP Q9Y3Z3
C	109	ASP	-	expression tag	UNP Q9Y3Z3
C	110	PRO	-	expression tag	UNP Q9Y3Z3
C	111	ASN	-	expression tag	UNP Q9Y3Z3
C	112	SER	-	expression tag	UNP Q9Y3Z3
C	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
C	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
D	77	MET	-	expression tag	UNP Q9Y3Z3
D	78	GLY	-	expression tag	UNP Q9Y3Z3
D	79	SER	-	expression tag	UNP Q9Y3Z3
D	80	SER	-	expression tag	UNP Q9Y3Z3
D	81	HIS	-	expression tag	UNP Q9Y3Z3
D	82	HIS	-	expression tag	UNP Q9Y3Z3
D	83	HIS	-	expression tag	UNP Q9Y3Z3
D	84	HIS	-	expression tag	UNP Q9Y3Z3
D	85	HIS	-	expression tag	UNP Q9Y3Z3
D	86	HIS	-	expression tag	UNP Q9Y3Z3
D	87	SER	-	expression tag	UNP Q9Y3Z3
D	88	SER	-	expression tag	UNP Q9Y3Z3
D	89	GLY	-	expression tag	UNP Q9Y3Z3
D	90	LEU	-	expression tag	UNP Q9Y3Z3
D	91	VAL	-	expression tag	UNP Q9Y3Z3
D	92	PRO	-	expression tag	UNP Q9Y3Z3
D	93	ARG	-	expression tag	UNP Q9Y3Z3
D	94	GLY	-	expression tag	UNP Q9Y3Z3
D	95	SER	-	expression tag	UNP Q9Y3Z3
D	96	HIS	-	expression tag	UNP Q9Y3Z3
D	97	MET	-	expression tag	UNP Q9Y3Z3
D	98	ALA	-	expression tag	UNP Q9Y3Z3
D	99	SER	-	expression tag	UNP Q9Y3Z3
D	100	MET	-	expression tag	UNP Q9Y3Z3
D	101	THR	-	expression tag	UNP Q9Y3Z3
D	102	GLY	-	expression tag	UNP Q9Y3Z3
D	103	GLY	-	expression tag	UNP Q9Y3Z3
D	104	GLN	-	expression tag	UNP Q9Y3Z3
D	105	GLN	-	expression tag	UNP Q9Y3Z3
D	106	MET	-	expression tag	UNP Q9Y3Z3
D	107	GLY	-	expression tag	UNP Q9Y3Z3
D	108	ARG	-	expression tag	UNP Q9Y3Z3
D	109	ASP	-	expression tag	UNP Q9Y3Z3

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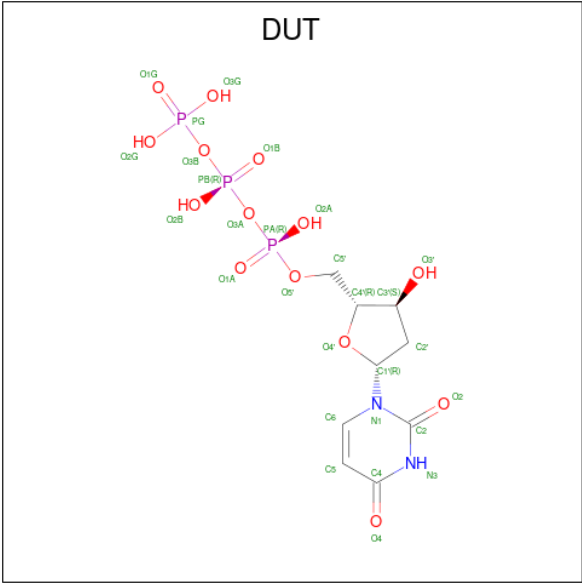
Chain	Residue	Modelled	Actual	Comment	Reference
D	110	PRO	-	expression tag	UNP Q9Y3Z3
D	111	ASN	-	expression tag	UNP Q9Y3Z3
D	112	SER	-	expression tag	UNP Q9Y3Z3
D	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
D	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3

- Molecule 2 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (CCD ID: DGT) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is DEOXYURIDINE-5'-TRIPHOSPHATE (CCD ID: DUT) (formula: $C_9H_{15}N_2O_{14}P_3$).

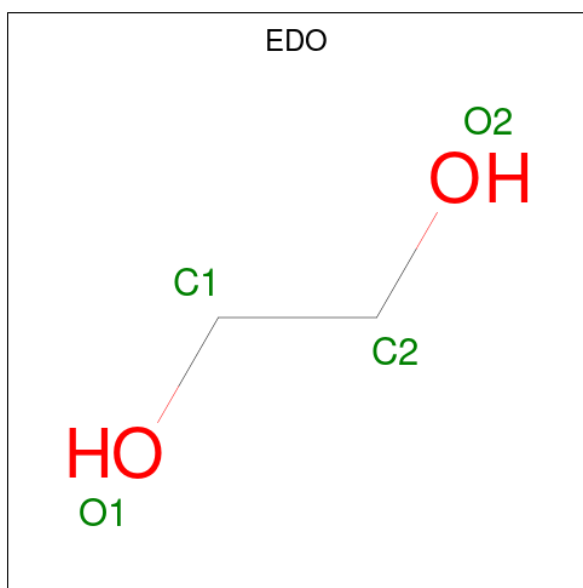


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	A	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	B	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	B	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	C	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	C	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	D	1	Total	C	N	O	P	0	0
			28	9	2	14	3		
3	D	1	Total	C	N	O	P	0	0
			28	9	2	14	3		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	2	Total	Mg	0	0
			2	2		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

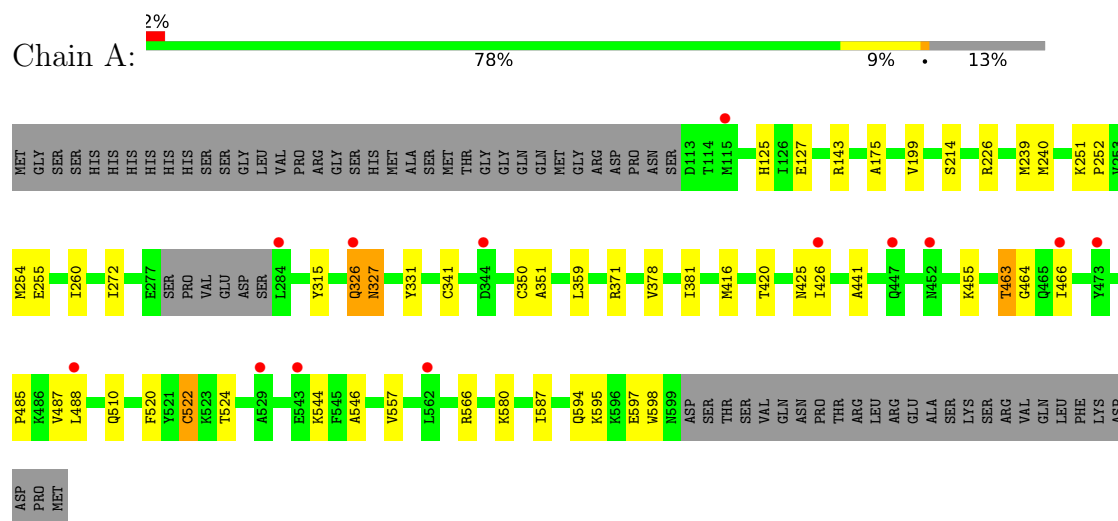
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	43	Total	O	0	0
			43	43		
6	B	45	Total	O	0	0
			45	45		
6	C	39	Total	O	0	0
			39	39		
6	D	55	Total	O	0	0
			55	55		

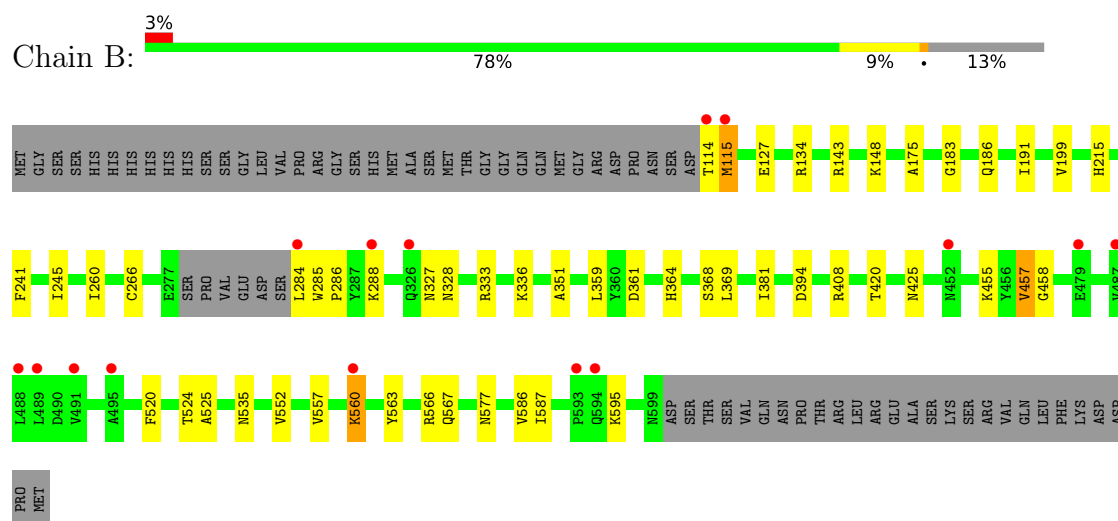
3 Residue-property plots

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.

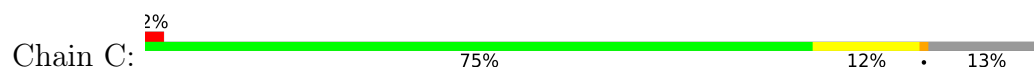
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

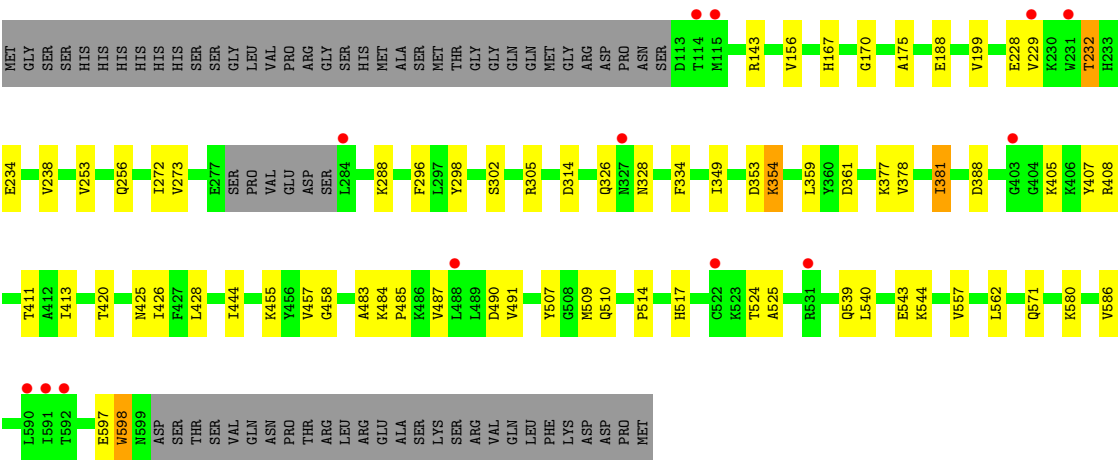


- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

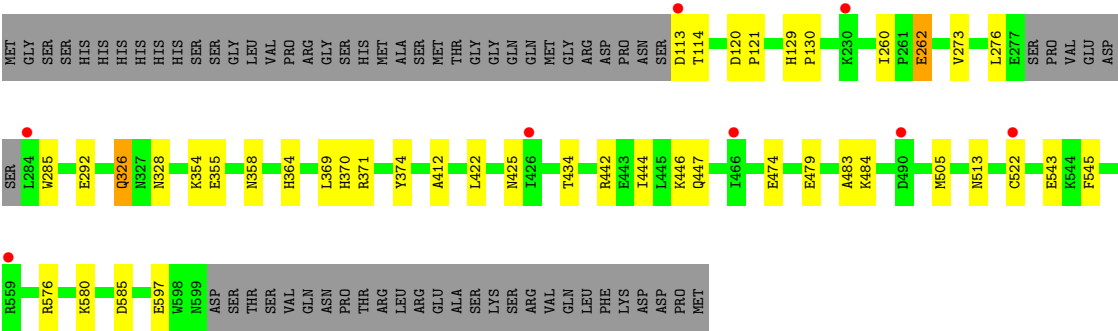
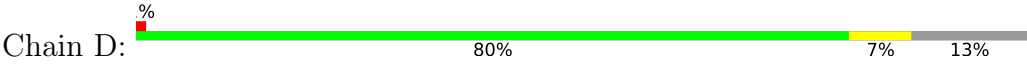


- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1





● Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.86Å 146.60Å 98.69Å 90.00° 114.62° 90.00°	Depositor
Resolution (Å)	42.91 – 2.30 42.91 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (42.91-2.30) 99.8 (42.91-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.29Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.197 , 0.250 0.202 , 0.254	Depositor DCC
R_{free} test set	2011 reflections (1.25%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16309	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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: MG, EDO, DGT, DUT

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	0.45	0/4045	0.84	7/5459 (0.1%)
1	B	0.43	0/4036	0.81	8/5447 (0.1%)
1	C	0.43	0/4043	0.84	7/5457 (0.1%)
1	D	0.49	0/4034	0.84	3/5445 (0.1%)
All	All	0.45	0/16158	0.83	25/21808 (0.1%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	458	GLY	N-CA-C	6.82	120.59	110.42
1	B	458	GLY	N-CA-C	6.44	119.96	110.63
1	A	226	ARG	CA-C-N	6.03	125.91	119.28
1	A	226	ARG	C-N-CA	6.03	125.91	119.28
1	C	484	LYS	CA-C-N	5.61	125.45	119.28
1	C	484	LYS	C-N-CA	5.61	125.45	119.28
1	A	260	ILE	CA-C-N	5.53	124.99	119.24
1	A	260	ILE	C-N-CA	5.53	124.99	119.24
1	A	214	SER	CB-CA-C	-5.47	109.79	117.23
1	B	364	HIS	N-CA-C	-5.42	105.37	111.28
1	C	272	ILE	N-CA-C	5.30	115.51	110.53
1	C	525	ALA	CA-C-N	5.27	125.08	119.28
1	C	525	ALA	C-N-CA	5.27	125.08	119.28
1	A	441	ALA	N-CA-C	-5.27	105.45	111.14
1	B	525	ALA	CA-C-N	5.26	124.97	119.82
1	B	525	ALA	C-N-CA	5.26	124.97	119.82
1	C	510	GLN	CB-CA-C	-5.23	110.12	117.23
1	D	364	HIS	N-CA-C	-5.17	105.64	111.28
1	B	186	GLN	CA-C-N	5.14	125.18	119.32
1	B	186	GLN	C-N-CA	5.14	125.18	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	260	ILE	CA-C-N	5.12	126.24	119.84
1	B	260	ILE	C-N-CA	5.12	126.24	119.84
1	A	272	ILE	N-CA-C	5.08	115.86	110.72
1	D	260	ILE	CA-C-N	5.01	126.10	119.84
1	D	260	ILE	C-N-CA	5.01	126.10	119.84

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	3946	0	3941	31	0
1	B	3938	0	3936	30	0
1	C	3945	0	3935	37	0
1	D	3939	0	3927	26	0
2	A	31	0	12	0	0
2	B	31	0	12	0	0
2	C	31	0	12	1	0
2	D	31	0	12	0	0
3	A	56	0	20	0	0
3	B	56	0	21	2	0
3	C	56	0	21	2	0
3	D	56	0	20	0	0
4	A	3	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	2	0	0	0	0
5	D	4	0	6	1	0
6	A	43	0	0	1	0
6	B	45	0	0	3	0
6	C	39	0	0	3	0
6	D	55	0	0	1	0
All	All	16309	0	15875	109	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 3.

All (109) 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:328:ASN:HA	1:D:326:GLN:HG3	1.42	1.01
3:B:702:DUT:O2G	6:B:830:HOH:O	1.80	0.98
1:A:331:TYR:OH	6:A:820:HOH:O	2.00	0.78
3:C:703:DUT:O2A	6:C:807:HOH:O	2.01	0.76
1:A:425:ASN:OD1	1:D:425:ASN:ND2	2.20	0.72
1:C:298:TYR:O	6:C:835:HOH:O	2.08	0.71
1:C:354:LYS:HE2	3:C:701:DUT:O2A	1.90	0.71
1:C:302:SER:O	6:C:835:HOH:O	2.09	0.70
1:B:328:ASN:HA	1:D:326:GLN:CG	2.20	0.69
1:C:234:GLU:HB3	1:C:273:VAL:HG12	1.76	0.68
1:B:455:LYS:HG2	1:B:557:VAL:HG12	1.75	0.68
1:C:232:THR:OG1	1:C:234:GLU:OE1	2.14	0.65
1:D:355:GLU:OE1	1:D:358:ASN:ND2	2.29	0.65
1:B:327:ASN:O	1:D:326:GLN:HG3	1.97	0.64
1:A:326:GLN:OE1	1:A:327:ASN:N	2.22	0.63
1:C:543:GLU:HG3	1:C:544:LYS:HG3	1.81	0.63
1:A:175:ALA:HB1	1:A:199:VAL:HG12	1.81	0.61
1:D:328:ASN:ND2	6:D:805:HOH:O	2.29	0.61
1:C:580:LYS:HD3	1:C:598:TRP:HB3	1.83	0.60
1:A:252:PRO:HA	1:A:255:GLU:HG2	1.86	0.58
1:A:127:GLU:HG3	1:B:336:LYS:HE3	1.87	0.56
1:B:566:ARG:HD3	1:B:587:ILE:HB	1.87	0.56
1:A:239:MET:HE2	1:A:416:MET:HE3	1.88	0.56
1:A:566:ARG:HD3	1:A:587:ILE:HB	1.88	0.55
1:A:326:GLN:HG3	1:C:326[A]:GLN:HG2	1.88	0.55
1:B:148:LYS:NZ	6:B:827:HOH:O	2.37	0.53
1:A:544:LYS:HE2	1:C:539:GLN:HB3	1.91	0.52
1:A:326:GLN:HG3	1:C:326[A]:GLN:CG	2.39	0.52
1:B:175:ALA:HB1	1:B:199:VAL:HG12	1.91	0.51
1:C:353:ASP:OD1	1:C:354:LYS:N	2.43	0.51
1:B:425:ASN:HB2	1:C:428:LEU:HD13	1.93	0.51
1:A:341:CYS:HB2	1:A:350:CYS:SG	2.51	0.51
1:A:371:ARG:NH2	1:C:361:ASP:OD1	2.32	0.49
1:B:143:ARG:HD2	1:B:420:THR:HA	1.92	0.49
1:C:455:LYS:HG2	1:C:557:VAL:HG12	1.92	0.49
1:C:408:ARG:N	1:C:411:THR:OG1	2.39	0.49
1:B:115:MET:HE2	1:B:127:GLU:HB3	1.95	0.49
1:D:597:GLU:OE1	1:D:597:GLU:N	2.44	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:586:VAL:HG21	1:D:522:CYS:SG	2.54	0.48
1:A:522:CYS:SG	1:A:524:THR:OG1	2.71	0.47
1:B:351:ALA:O	1:B:520:PHE:HA	2.14	0.47
1:B:359:LEU:HD23	1:B:359:LEU:HA	1.74	0.47
1:A:240:MET:HE2	1:A:416:MET:SD	2.55	0.47
1:A:597:GLU:OE1	1:A:597:GLU:N	2.28	0.47
1:B:361:ASP:OD2	1:D:371:ARG:NH2	2.46	0.46
1:C:507:TYR:HB3	1:C:514:PRO:HG3	1.97	0.46
1:B:560:LYS:HD2	1:B:560:LYS:HA	1.48	0.46
1:B:381:ILE:HD12	1:B:381:ILE:HA	1.82	0.46
1:A:510:GLN:O	1:A:546:ALA:HB2	2.16	0.45
1:A:455:LYS:HG2	1:A:557:VAL:HG12	1.97	0.45
1:A:143:ARG:HD2	1:A:420:THR:HA	1.98	0.45
1:B:183:GLY:HA2	1:B:191:ILE:HD12	1.98	0.45
1:B:215:HIS:HE2	3:B:702:DUT:PA	2.40	0.45
1:C:485:PRO:HA	1:C:571:GLN:HE21	1.82	0.45
1:B:285:TRP:HA	1:B:286:PRO:HD3	1.80	0.45
1:B:586:VAL:HG11	1:D:522:CYS:SG	2.57	0.44
1:C:405:LYS:HD2	1:C:407:TYR:CZ	2.52	0.44
1:D:374:TYR:O	5:D:706:EDO:O2	2.36	0.44
1:A:351:ALA:O	1:A:520:PHE:HA	2.17	0.44
1:A:580:LYS:HD2	1:A:598:TRP:HB3	2.00	0.44
1:B:457:VAL:HG22	1:B:552:VAL:O	2.17	0.44
1:D:483:ALA:O	1:D:484:LYS:HD2	2.18	0.44
1:C:167:HIS:ND1	1:C:314:ASP:OD2	2.48	0.43
1:D:479:GLU:OE1	1:D:576:ARG:NH1	2.48	0.43
1:C:509:MET:HE1	1:C:517:HIS:ND1	2.33	0.43
1:A:251:LYS:O	1:A:254:MET:HB2	2.19	0.43
1:A:595:LYS:HZ3	1:A:597:GLU:CD	2.27	0.43
1:B:425:ASN:OD1	1:C:425:ASN:ND2	2.50	0.43
1:D:444:ILE:O	1:D:447:GLN:HB2	2.19	0.43
1:A:463:THR:HG22	1:A:464:GLY:H	1.82	0.43
1:D:262:GLU:H	1:D:262:GLU:CD	2.27	0.43
1:C:305:ARG:NH1	1:C:517:HIS:HD2	2.17	0.42
1:D:120:ASP:OD1	1:D:121:PRO:HD2	2.19	0.42
1:A:125:HIS:CE1	1:B:333:ARG:HB2	2.54	0.42
1:A:326:GLN:O	1:C:328:ASN:HB2	2.19	0.42
6:B:805:HOH:O	1:D:543:GLU:HG2	2.20	0.42
1:C:175:ALA:HB1	1:C:199:VAL:HG12	2.01	0.42
1:C:296:PHE:HB2	1:C:349:ILE:HG13	2.01	0.42
1:B:241:PHE:O	1:B:245:ILE:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:228:GLU:OE1	1:C:228:GLU:N	2.42	0.42
1:C:156:VAL:O	2:C:702:DGT:H8	2.20	0.42
1:B:266:CYS:SG	1:B:288:LYS:HE3	2.59	0.42
1:B:369:LEU:HD23	1:B:369:LEU:HA	1.92	0.42
1:D:369:LEU:HD23	1:D:369:LEU:HA	1.90	0.42
1:C:377:LYS:HG3	1:C:378:VAL:HG23	2.01	0.42
1:D:513:ASN:HB2	1:D:545:PHE:CE2	2.55	0.42
1:D:580:LYS:NZ	1:D:585:ASP:OD2	2.41	0.41
1:A:381:ILE:HD12	1:A:381:ILE:HA	1.86	0.41
1:A:522:CYS:SG	1:C:586:VAL:HG11	2.60	0.41
1:B:563:TYR:O	1:B:567:GLN:HG2	2.20	0.41
1:C:359:LEU:HD23	1:C:359:LEU:HA	1.81	0.41
1:C:483:ALA:O	1:C:485:PRO:HD3	2.20	0.41
1:D:113:ASP:HB3	1:D:114:THR:H	1.74	0.41
1:A:485:PRO:O	1:A:487:VAL:N	2.50	0.41
1:D:370:HIS:HB3	1:D:505:MET:HE1	2.02	0.41
1:C:143:ARG:HD2	1:C:420:THR:HA	2.02	0.41
1:A:378:VAL:O	1:A:381:ILE:HG22	2.20	0.41
1:C:253:VAL:O	1:C:256:GLN:HB3	2.21	0.41
1:D:442:ARG:HG2	1:D:446:LYS:HE2	2.02	0.41
1:A:463:THR:O	1:A:466:ILE:HB	2.21	0.41
1:B:134:ARG:HA	1:B:134:ARG:HD3	1.91	0.41
1:B:394:ASP:O	1:B:408:ARG:HD2	2.21	0.41
1:C:170:GLY:HA3	1:C:314:ASP:OD1	2.20	0.41
1:C:388:ASP:OD1	1:C:444:ILE:HD13	2.20	0.41
1:C:381:ILE:HD12	1:C:381:ILE:HA	1.75	0.41
1:D:129:HIS:CG	1:D:130:PRO:HD2	2.56	0.41
1:C:334:PHE:CE2	1:C:359:LEU:HD21	2.57	0.40
1:D:285:TRP:CD2	1:D:292:GLU:HG2	2.57	0.40
1:D:412:ALA:HB3	1:D:422:LEU:HD22	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	479/550 (87%)	467 (98%)	12 (2%)	0	100	100
1	B	478/550 (87%)	465 (97%)	12 (2%)	1 (0%)	43	55
1	C	479/550 (87%)	460 (96%)	18 (4%)	1 (0%)	43	55
1	D	478/550 (87%)	467 (98%)	10 (2%)	1 (0%)	43	55
All	All	1914/2200 (87%)	1859 (97%)	52 (3%)	3 (0%)	43	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	595	LYS
1	C	598	TRP
1	D	262	GLU

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	429/488 (88%)	420 (98%)	9 (2%)	47	66
1	B	428/488 (88%)	418 (98%)	10 (2%)	44	63
1	C	429/488 (88%)	411 (96%)	18 (4%)	26	40
1	D	428/488 (88%)	422 (99%)	6 (1%)	59	76
All	All	1714/1952 (88%)	1671 (98%)	43 (2%)	43	60

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	315	TYR
1	A	326	GLN
1	A	327	ASN
1	A	359	LEU
1	A	426	ILE

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Mol	Chain	Res	Type
1	A	463	THR
1	A	488	LEU
1	A	522	CYS
1	A	594	GLN
1	B	114	THR
1	B	115	MET
1	B	284	LEU
1	B	368	SER
1	B	457	VAL
1	B	524	THR
1	B	535[A]	ASN
1	B	535[B]	ASN
1	B	560	LYS
1	B	577	ASN
1	C	188	GLU
1	C	229	VAL
1	C	232	THR
1	C	238	VAL
1	C	288	LYS
1	C	354	LYS
1	C	381	ILE
1	C	413	ILE
1	C	426	ILE
1	C	457	VAL
1	C	487	VAL
1	C	490	ASP
1	C	491	VAL
1	C	524	THR
1	C	540	LEU
1	C	562	LEU
1	C	597[A]	GLU
1	C	597[B]	GLU
1	D	273	VAL
1	D	276	LEU
1	D	326	GLN
1	D	354	LYS
1	D	434	THR
1	D	474	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	322	HIS
1	A	539	GLN
1	A	567	GLN
1	A	571	GLN
1	A	594	GLN
1	B	140	GLN
1	B	248	ASN
1	B	326	GLN
1	B	567	GLN
1	B	571	GLN
1	C	149	GLN
1	C	210	HIS
1	C	370	HIS
1	C	375	GLN
1	C	380	ASN
1	C	517	HIS
1	D	149	GLN
1	D	200	GLN
1	D	210	HIS
1	D	235	GLN
1	D	243	HIS
1	D	364	HIS
1	D	571	GLN
1	D	577	ASN
1	D	594	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 7 are monoatomic - leaving 13 for Mogul analysis.

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	DUT	C	701	4	28,29,29	2.28	6 (21%)	41,45,45	1.62	6 (14%)
3	DUT	B	702	4	28,29,29	2.31	7 (25%)	41,45,45	1.59	5 (12%)
3	DUT	A	703	4	28,29,29	2.30	9 (32%)	41,45,45	1.66	4 (9%)
3	DUT	D	703	4	28,29,29	2.30	7 (25%)	41,45,45	1.55	5 (12%)
3	DUT	A	702	4	28,29,29	2.29	7 (25%)	41,45,45	1.58	6 (14%)
2	DGT	B	701	4	32,33,33	2.38	7 (21%)	48,52,52	1.92	12 (25%)
3	DUT	D	701	4	28,29,29	2.23	7 (25%)	41,45,45	1.63	4 (9%)
3	DUT	B	703	4	28,29,29	2.33	7 (25%)	41,45,45	1.67	5 (12%)
2	DGT	C	702	4	32,33,33	2.57	9 (28%)	48,52,52	1.97	11 (22%)
3	DUT	C	703	-	28,29,29	2.35	7 (25%)	41,45,45	1.62	7 (17%)
5	EDO	D	706	-	3,3,3	0.53	0	2,2,2	0.28	0
2	DGT	D	702	4	32,33,33	2.49	8 (25%)	48,52,52	1.97	15 (31%)
2	DGT	A	701	4	32,33,33	2.55	10 (31%)	48,52,52	1.92	14 (29%)

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	DUT	C	701	4	-	5/22/34/34	0/2/2/2
3	DUT	B	702	4	-	6/22/34/34	0/2/2/2
3	DUT	A	703	4	-	3/22/34/34	0/2/2/2
3	DUT	D	703	4	-	3/22/34/34	0/2/2/2
3	DUT	A	702	4	-	5/22/34/34	0/2/2/2
2	DGT	B	701	4	-	9/22/34/34	0/3/3/3
3	DUT	D	701	4	-	2/22/34/34	0/2/2/2
3	DUT	B	703	4	-	3/22/34/34	0/2/2/2
2	DGT	C	702	4	-	9/22/34/34	0/3/3/3
3	DUT	C	703	-	-	4/22/34/34	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	D	706	-	-	1/1/1/1	-
2	DGT	D	702	4	-	9/22/34/34	0/3/3/3
2	DGT	A	701	4	-	9/22/34/34	0/3/3/3

All (91) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	702	DGT	O6-C6	9.33	1.41	1.23
2	A	701	DGT	O6-C6	8.97	1.40	1.23
2	D	702	DGT	O6-C6	8.73	1.40	1.23
2	B	701	DGT	O6-C6	8.69	1.40	1.23
3	B	703	DUT	O4-C4	8.27	1.40	1.24
3	A	703	DUT	O4-C4	8.10	1.40	1.24
3	C	701	DUT	O4-C4	8.07	1.40	1.24
3	B	702	DUT	O4-C4	7.90	1.39	1.24
3	C	703	DUT	O4-C4	7.87	1.39	1.24
3	D	703	DUT	O4-C4	7.86	1.39	1.24
3	D	701	DUT	O4-C4	7.75	1.39	1.24
3	A	702	DUT	O4-C4	7.67	1.39	1.24
2	A	701	DGT	C2-N2	6.25	1.48	1.34
2	C	702	DGT	C2-N2	6.21	1.48	1.34
2	D	702	DGT	C2-N2	6.16	1.48	1.34
2	B	701	DGT	C2-N2	5.80	1.47	1.34
3	C	703	DUT	PB-O3A	-5.43	1.53	1.59
3	A	702	DUT	PB-O3A	-5.15	1.53	1.59
3	B	702	DUT	PB-O3A	-5.06	1.54	1.59
3	D	703	DUT	PB-O3A	-4.98	1.54	1.59
3	C	701	DUT	PB-O3A	-4.94	1.54	1.59
2	A	701	DGT	PB-O3B	-4.71	1.54	1.59
3	B	703	DUT	PB-O3A	-4.62	1.54	1.59
2	C	702	DGT	PB-O3B	-4.33	1.54	1.59
3	A	703	DUT	PB-O3A	-4.05	1.55	1.59
2	B	701	DGT	PB-O3B	-3.79	1.55	1.59
2	D	702	DGT	PA-O3A	-3.75	1.55	1.59
3	D	701	DUT	PB-O3A	-3.67	1.55	1.59
2	D	702	DGT	PB-O3B	-3.66	1.55	1.59
3	D	701	DUT	O3'-C3'	-3.65	1.35	1.43
2	C	702	DGT	O4'-C1'	3.55	1.50	1.42
2	A	701	DGT	O4'-C1'	3.55	1.50	1.42
2	D	702	DGT	O4'-C1'	3.47	1.50	1.42
3	A	702	DUT	O3'-C3'	-3.43	1.36	1.43
3	A	703	DUT	O3'-C3'	-3.35	1.36	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	DGT	O4'-C1'	3.35	1.49	1.42
3	C	703	DUT	O3'-C3'	-3.35	1.36	1.43
3	B	702	DUT	O3'-C3'	-3.32	1.36	1.43
3	D	703	DUT	O3'-C3'	-3.13	1.36	1.43
3	C	701	DUT	C6-C5	3.02	1.42	1.35
3	B	703	DUT	C6-N1	-3.01	1.30	1.38
2	C	702	DGT	O3'-C3'	-2.94	1.37	1.43
3	D	703	DUT	C2-N1	2.89	1.43	1.38
3	C	701	DUT	O3'-C3'	-2.88	1.37	1.43
3	C	701	DUT	C2-N1	2.85	1.42	1.38
3	A	703	DUT	C2-N1	2.81	1.42	1.38
2	A	701	DGT	O3'-C3'	-2.80	1.37	1.43
3	D	703	DUT	C6-C5	2.80	1.41	1.35
3	B	703	DUT	O3'-C3'	-2.77	1.37	1.43
3	A	703	DUT	C6-N1	-2.76	1.31	1.38
3	B	702	DUT	C2-N1	2.72	1.42	1.38
3	A	703	DUT	C6-C5	2.71	1.41	1.35
3	D	703	DUT	C6-N1	-2.70	1.31	1.38
3	B	702	DUT	C6-C5	2.70	1.41	1.35
3	C	703	DUT	C6-N1	-2.67	1.31	1.38
3	A	702	DUT	C6-N1	-2.64	1.31	1.38
3	D	701	DUT	C6-C5	2.63	1.41	1.35
3	D	701	DUT	C6-N1	-2.63	1.31	1.38
3	C	703	DUT	C6-C5	2.62	1.41	1.35
2	A	701	DGT	PA-O3A	-2.60	1.56	1.59
3	B	703	DUT	C4-N3	-2.57	1.34	1.38
3	B	702	DUT	C6-N1	-2.55	1.32	1.38
3	A	702	DUT	C6-C5	2.52	1.40	1.35
3	C	701	DUT	C6-N1	-2.52	1.32	1.38
3	B	703	DUT	C6-C5	2.48	1.40	1.35
2	D	702	DGT	O3'-C3'	-2.46	1.38	1.43
3	C	703	DUT	C2'-C3'	-2.43	1.46	1.52
2	B	701	DGT	O3'-C3'	-2.36	1.38	1.43
2	C	702	DGT	O4'-C4'	-2.35	1.39	1.45
2	D	702	DGT	PB-O3A	-2.32	1.57	1.59
3	D	701	DUT	C2-N1	2.29	1.42	1.38
2	C	702	DGT	PB-O3A	-2.29	1.57	1.59
3	A	702	DUT	C2'-C3'	-2.27	1.47	1.52
3	D	701	DUT	C4-N3	-2.26	1.34	1.38
3	A	702	DUT	C2-N1	2.21	1.41	1.38
2	A	701	DGT	PB-O3A	-2.20	1.57	1.59
3	A	703	DUT	C4-N3	-2.19	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	703	DUT	C2-N1	2.18	1.41	1.38
3	D	703	DUT	C2'-C3'	-2.18	1.47	1.52
2	B	701	DGT	C6-N1	-2.13	1.34	1.38
2	A	701	DGT	C6-N1	-2.12	1.34	1.38
2	C	702	DGT	C2'-C3'	-2.09	1.47	1.52
3	A	703	DUT	C2'-C3'	-2.08	1.47	1.52
3	B	702	DUT	C2'-C3'	-2.07	1.47	1.52
2	D	702	DGT	O4'-C4'	-2.07	1.40	1.45
2	A	701	DGT	O4'-C4'	-2.07	1.40	1.45
2	C	702	DGT	C6-N1	-2.06	1.35	1.38
3	B	703	DUT	C2-N1	2.04	1.41	1.38
2	B	701	DGT	O4'-C4'	-2.02	1.40	1.45
2	A	701	DGT	C5-N7	-2.00	1.35	1.39
3	A	703	DUT	C5'-C4'	-2.00	1.45	1.51

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	702	DGT	C5-C4-N3	-6.42	118.17	128.39
2	B	701	DGT	C5-C4-N3	-6.05	118.77	128.39
3	A	703	DUT	C4-N3-C2	-6.03	119.13	126.61
2	A	701	DGT	C5-C4-N3	-5.99	118.86	128.39
2	D	702	DGT	C2-N3-C4	5.96	122.56	112.30
3	D	701	DUT	C4-N3-C2	-5.87	119.32	126.61
2	D	702	DGT	C5-C4-N3	-5.81	119.14	128.39
2	C	702	DGT	C2-N3-C4	5.79	122.28	112.30
3	C	703	DUT	C4-N3-C2	-5.65	119.60	126.61
3	B	702	DUT	C4-N3-C2	-5.62	119.64	126.61
3	B	703	DUT	C4-N3-C2	-5.61	119.65	126.61
2	B	701	DGT	C2-N3-C4	5.60	121.94	112.30
2	A	701	DGT	C2-N3-C4	5.59	121.92	112.30
3	A	702	DUT	C4-N3-C2	-5.42	119.88	126.61
3	C	701	DUT	C4-N3-C2	-5.36	119.96	126.61
3	D	703	DUT	C4-N3-C2	-5.32	120.01	126.61
3	A	703	DUT	N3-C2-N1	4.76	121.09	114.89
3	B	703	DUT	N3-C2-N1	4.70	121.01	114.89
3	C	701	DUT	N3-C2-N1	4.64	120.93	114.89
3	B	702	DUT	N3-C2-N1	4.50	120.74	114.89
2	C	702	DGT	N9-C4-N3	4.42	134.79	125.95
3	D	701	DUT	N3-C2-N1	4.39	120.61	114.89
3	D	703	DUT	N3-C2-N1	4.31	120.50	114.89
3	C	703	DUT	N3-C2-N1	4.11	120.25	114.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	701	DUT	C5-C4-N3	4.01	120.42	114.80
3	A	702	DUT	N3-C2-N1	3.93	120.00	114.89
2	B	701	DGT	N9-C4-N3	3.85	133.65	125.95
3	A	702	DUT	C5-C4-N3	3.84	120.18	114.80
2	A	701	DGT	N9-C4-N3	3.80	133.54	125.95
3	A	703	DUT	C5-C4-N3	3.73	120.02	114.80
3	C	703	DUT	C5-C4-N3	3.72	120.01	114.80
3	D	703	DUT	C5-C4-N3	3.64	119.90	114.80
2	D	702	DGT	N9-C4-N3	3.52	132.98	125.95
3	B	702	DUT	C5-C4-N3	3.50	119.70	114.80
3	B	703	DUT	C5-C4-N3	3.42	119.58	114.80
2	D	702	DGT	O1G-PG-O3B	3.39	116.01	104.64
2	C	702	DGT	O1G-PG-O3B	3.33	115.81	104.64
3	C	701	DUT	C5-C4-N3	3.29	119.41	114.80
3	C	703	DUT	O4-C4-C5	-3.04	119.92	125.16
2	B	701	DGT	O1G-PG-O3B	2.95	114.52	104.64
2	B	701	DGT	C6-C5-N7	2.90	135.56	130.29
2	D	702	DGT	C6-C5-N7	2.90	135.56	130.29
3	A	702	DUT	O4-C4-C5	-2.88	120.19	125.16
3	C	701	DUT	O2A-PA-O3A	2.88	115.05	107.27
2	C	702	DGT	C8-N7-C5	2.87	109.38	104.26
2	D	702	DGT	O4'-C1'-N9	2.87	112.96	107.86
3	B	702	DUT	O4-C4-C5	-2.81	120.31	125.16
2	A	701	DGT	O4'-C1'-N9	2.74	112.73	107.86
2	A	701	DGT	O1G-PG-O3B	2.74	113.81	104.64
2	A	701	DGT	C6-C5-N7	2.71	135.23	130.29
2	B	701	DGT	C4-C5-N7	-2.67	106.44	110.67
2	B	701	DGT	C8-N7-C5	2.66	109.00	104.26
2	A	701	DGT	C8-N7-C5	2.65	108.98	104.26
2	B	701	DGT	C5-C6-N1	2.64	119.97	113.25
2	D	702	DGT	N2-C2-N1	2.63	122.31	116.76
2	A	701	DGT	C4-C5-N7	-2.63	106.51	110.67
2	D	702	DGT	O6-C6-C5	-2.61	119.66	126.53
2	C	702	DGT	N9-C8-N7	-2.59	108.60	113.40
2	C	702	DGT	O4'-C1'-N9	2.59	112.46	107.86
2	C	702	DGT	C4-C5-N7	-2.58	106.58	110.67
2	D	702	DGT	C5-C6-N1	2.53	119.70	113.25
2	D	702	DGT	C4-C5-N7	-2.52	106.68	110.67
3	A	702	DUT	O5'-C5'-C4'	2.51	117.53	108.99
3	B	703	DUT	O2B-PB-O3A	2.48	113.98	107.27
2	B	701	DGT	N9-C8-N7	-2.47	108.81	113.40
2	C	702	DGT	C6-C5-N7	2.47	134.79	130.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	DGT	C2-N1-C6	-2.47	120.63	125.11
2	A	701	DGT	O5'-C5'-C4'	2.44	117.31	108.99
3	C	703	DUT	O2-C2-N1	-2.43	119.64	122.80
3	B	703	DUT	O5'-C5'-C4'	2.40	117.17	108.99
2	B	701	DGT	O5'-C5'-C4'	2.40	117.17	108.99
2	D	702	DGT	O5'-C5'-C4'	2.39	117.12	108.99
2	A	701	DGT	N9-C8-N7	-2.39	108.98	113.40
3	C	703	DUT	O5'-C5'-C4'	2.37	117.07	108.99
2	A	701	DGT	C5-C6-N1	2.36	119.25	113.25
3	A	702	DUT	O2-C2-N1	-2.34	119.75	122.80
2	D	702	DGT	C8-N7-C5	2.33	108.41	104.26
2	C	702	DGT	C5-C6-N1	2.32	119.17	113.25
2	C	702	DGT	O5'-C5'-C4'	2.31	116.86	108.99
2	A	701	DGT	O1B-PB-O3A	2.30	113.50	107.27
3	C	703	DUT	O2B-PB-O3B	2.30	113.50	107.27
2	D	702	DGT	N9-C8-N7	-2.23	109.27	113.40
3	D	701	DUT	O2-C2-N1	-2.23	119.90	122.80
3	D	703	DUT	O4'-C1'-N1	2.22	111.81	107.86
3	A	703	DUT	O5'-C5'-C4'	2.22	116.56	108.99
2	B	701	DGT	O6-C6-C5	-2.20	120.72	126.53
3	C	701	DUT	O2B-PB-O3B	2.19	113.20	107.27
2	A	701	DGT	C2'-C3'-C4'	2.19	107.24	102.80
3	C	701	DUT	O2-C2-N3	-2.14	117.54	121.49
2	D	702	DGT	C2-N1-C6	-2.12	121.26	125.11
3	B	702	DUT	O5'-C5'-C4'	2.07	116.04	108.99
2	A	701	DGT	O6-C6-C5	-2.05	121.12	126.53
2	D	702	DGT	N1-C2-N3	-2.05	119.58	123.32
3	D	703	DUT	O4-C4-C5	-2.04	121.64	125.16

There are no chirality outliers.

All (68) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	DGT	PB-O3B-PG-O1G
2	B	701	DGT	PB-O3B-PG-O1G
2	C	702	DGT	PB-O3B-PG-O1G
2	D	702	DGT	PB-O3B-PG-O1G
3	A	702	DUT	C5'-O5'-PA-O1A
3	B	702	DUT	PB-O3A-PA-O5'
2	C	702	DGT	PB-O3B-PG-O3G
3	B	703	DUT	PB-O3A-PA-O1A
3	C	701	DUT	PG-O3B-PB-O1B

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Mol	Chain	Res	Type	Atoms
3	D	701	DUT	PA-O3A-PB-O1B
3	A	702	DUT	C4'-C5'-O5'-PA
3	D	703	DUT	C4'-C5'-O5'-PA
3	B	702	DUT	C4'-C5'-O5'-PA
3	C	703	DUT	C4'-C5'-O5'-PA
3	A	702	DUT	PB-O3A-PA-O5'
3	C	703	DUT	PB-O3A-PA-O5'
3	D	703	DUT	PB-O3A-PA-O5'
3	B	702	DUT	PG-O3B-PB-O2B
2	B	701	DGT	C2'-C1'-N9-C8
2	C	702	DGT	C2'-C1'-N9-C8
2	D	702	DGT	C2'-C1'-N9-C8
3	A	702	DUT	C5'-O5'-PA-O2A
3	A	702	DUT	C5'-O5'-PA-O3A
2	A	701	DGT	PG-O3B-PB-O1B
2	B	701	DGT	PG-O3B-PB-O1B
2	B	701	DGT	PB-O3A-PA-O1A
2	C	702	DGT	PG-O3B-PB-O1B
2	D	702	DGT	PG-O3B-PB-O1B
2	D	702	DGT	PB-O3A-PA-O1A
3	B	702	DUT	PA-O3A-PB-O2B
3	B	703	DUT	PA-O3A-PB-O1B
3	C	701	DUT	PB-O3A-PA-O2A
2	A	701	DGT	C2'-C1'-N9-C4
2	B	701	DGT	C2'-C1'-N9-C4
2	C	702	DGT	C2'-C1'-N9-C4
2	D	702	DGT	C2'-C1'-N9-C4
2	B	701	DGT	PB-O3B-PG-O3G
2	A	701	DGT	PB-O3A-PA-O1A
2	C	702	DGT	PB-O3A-PA-O1A
3	A	703	DUT	PA-O3A-PB-O1B
3	C	703	DUT	PG-O3B-PB-O1B
2	A	701	DGT	C4'-C5'-O5'-PA
2	A	701	DGT	PB-O3B-PG-O3G
2	D	702	DGT	PB-O3B-PG-O3G
3	C	703	DUT	PB-O3B-PG-O3G
2	A	701	DGT	C2'-C1'-N9-C8
2	B	701	DGT	C4'-C5'-O5'-PA
2	D	702	DGT	C4'-C5'-O5'-PA
2	A	701	DGT	PG-O3B-PB-O2B
2	B	701	DGT	PG-O3B-PB-O2B
2	B	701	DGT	PB-O3A-PA-O2A

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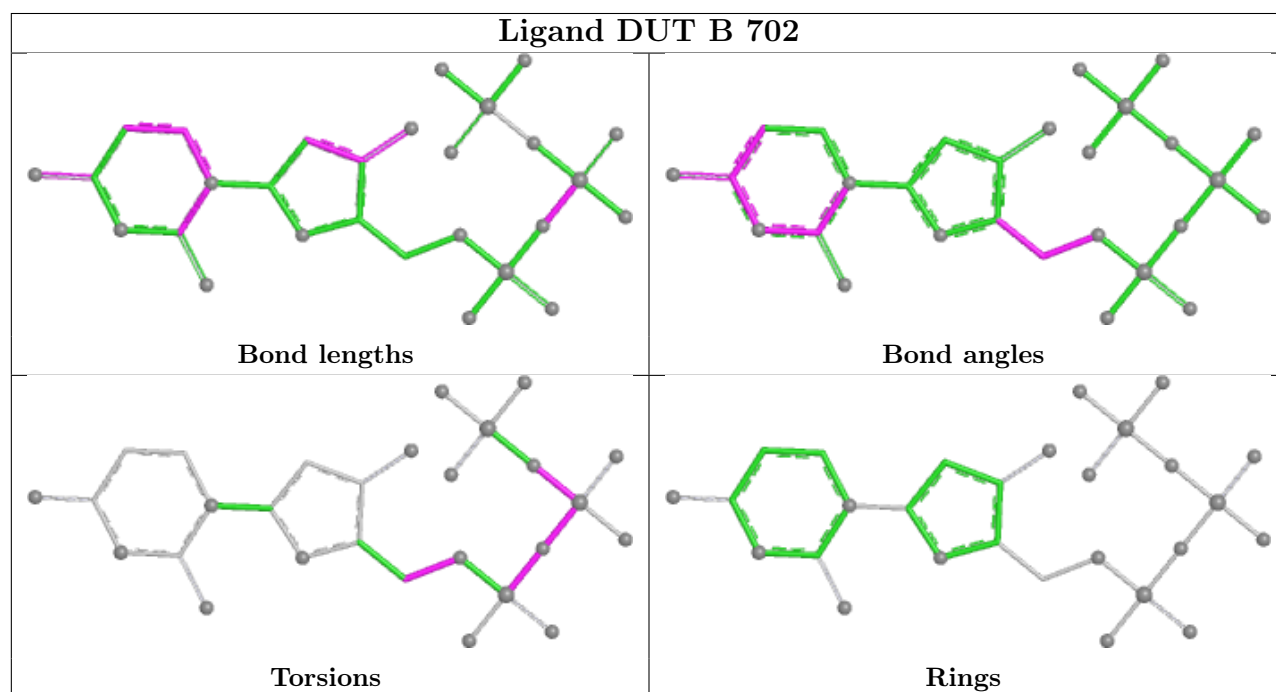
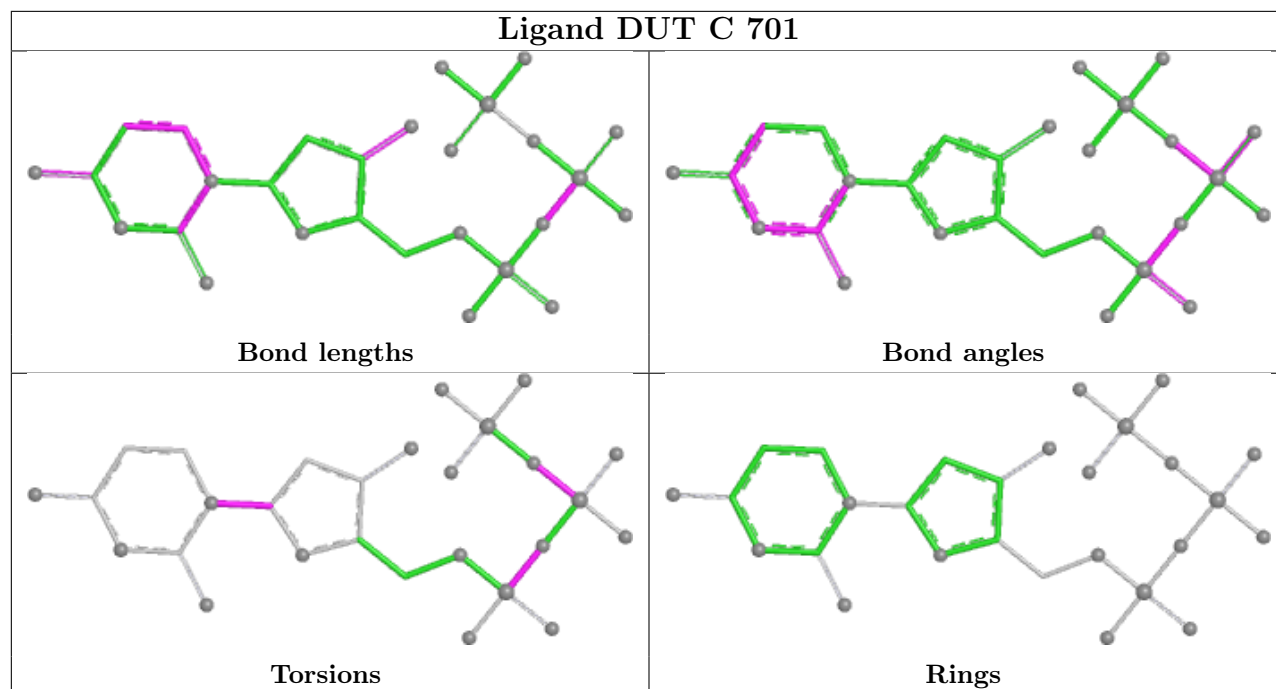
Mol	Chain	Res	Type	Atoms
2	C	702	DGT	PG-O3B-PB-O2B
2	C	702	DGT	PB-O3A-PA-O2A
2	D	702	DGT	PG-O3B-PB-O2B
3	A	703	DUT	PB-O3A-PA-O2A
3	B	702	DUT	PA-O3A-PB-O1B
3	C	701	DUT	PB-O3A-PA-O1A
3	D	703	DUT	PA-O3A-PB-O1B
5	D	706	EDO	O1-C1-C2-O2
3	C	701	DUT	O4'-C1'-N1-C2
2	C	702	DGT	C4'-C5'-O5'-PA
3	C	701	DUT	O4'-C1'-N1-C6
2	A	701	DGT	PB-O3A-PA-O2A
2	D	702	DGT	PB-O3A-PA-O2A
3	A	703	DUT	PG-O3B-PB-O2B
3	B	702	DUT	PG-O3B-PB-O1B
3	B	703	DUT	PB-O3A-PA-O2A
3	D	701	DUT	PB-O3A-PA-O2A

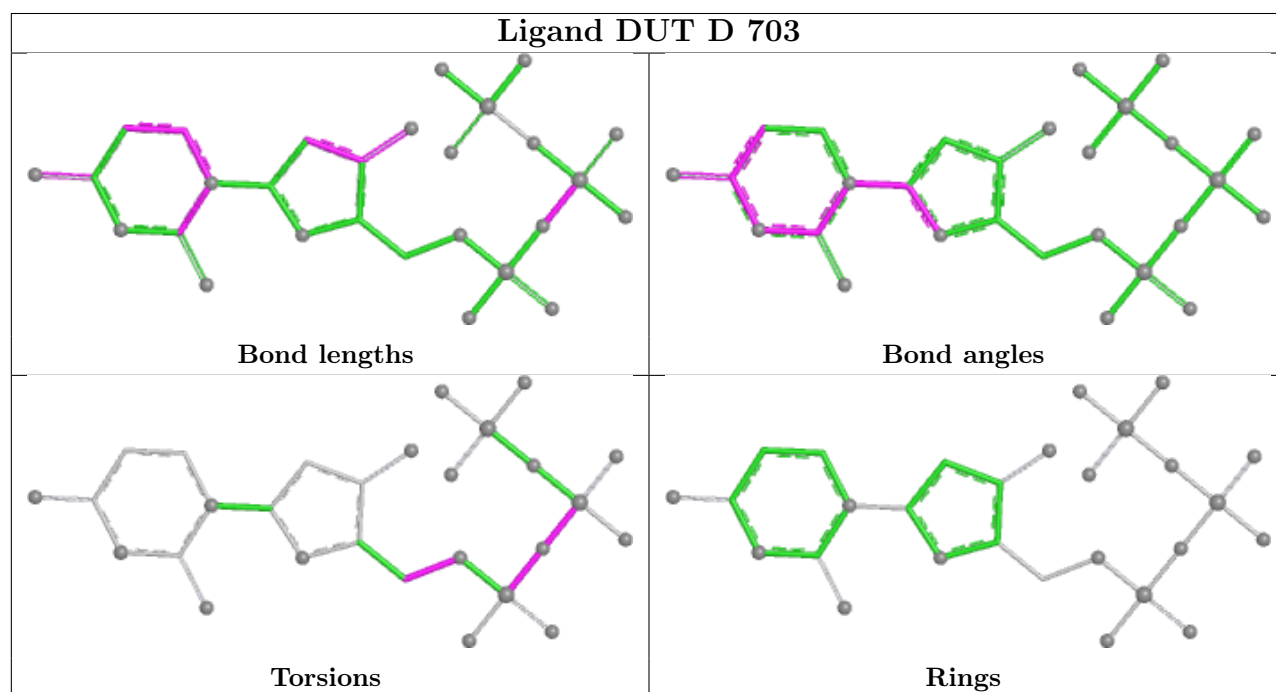
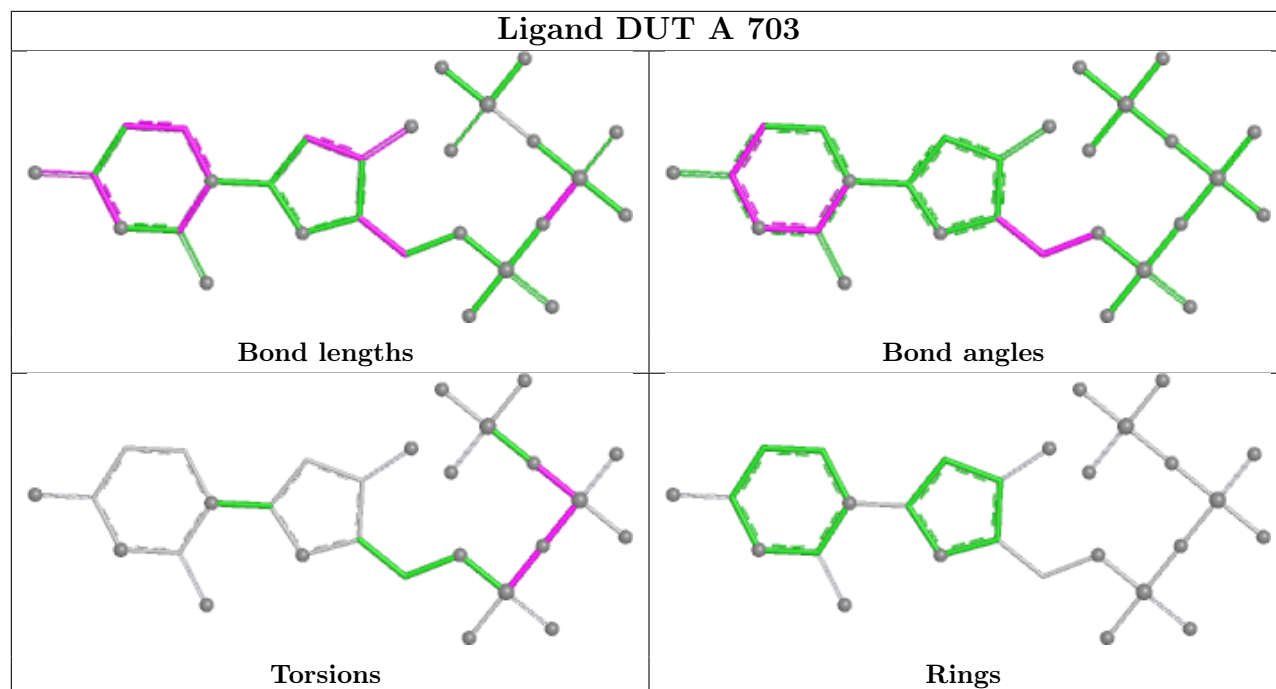
There are no ring outliers.

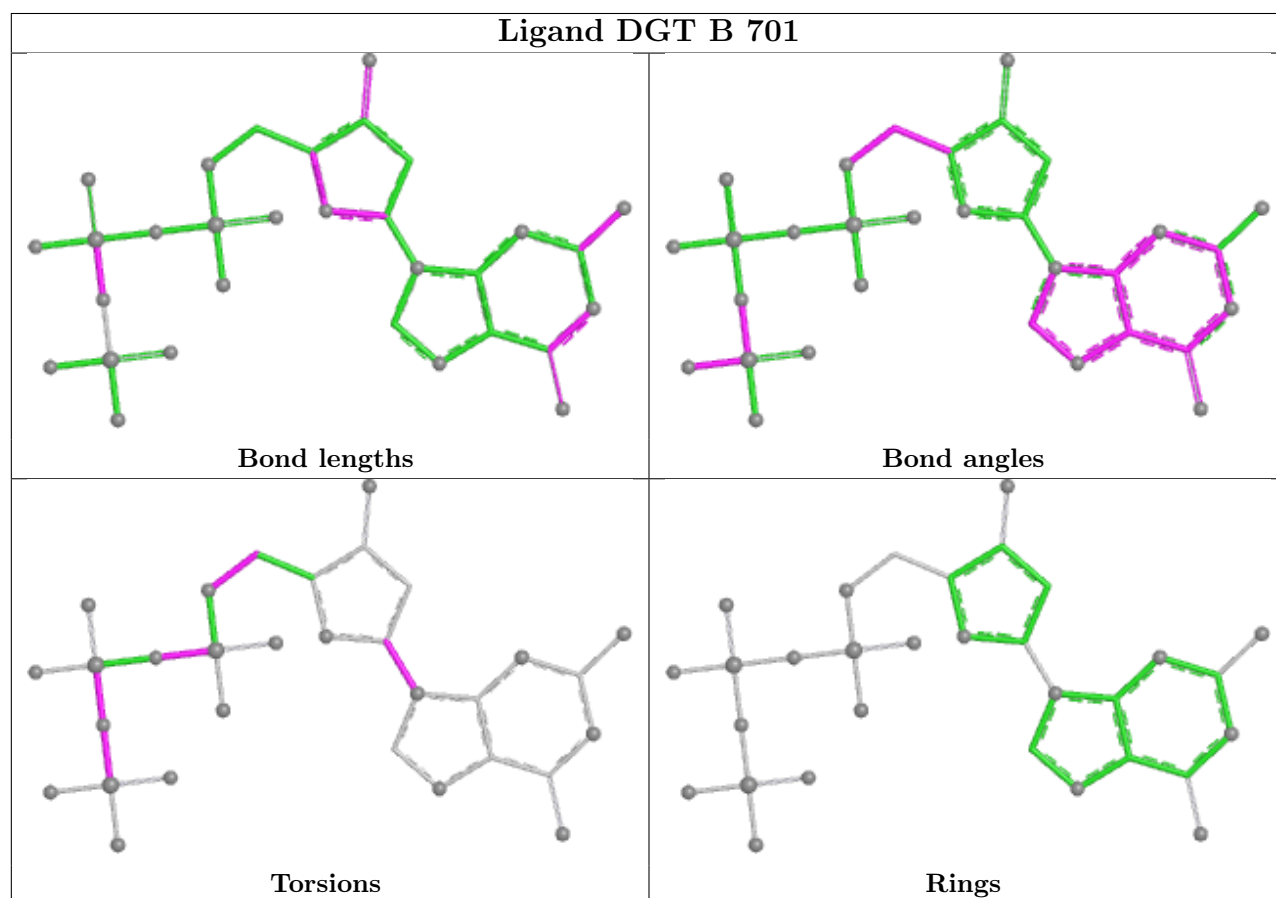
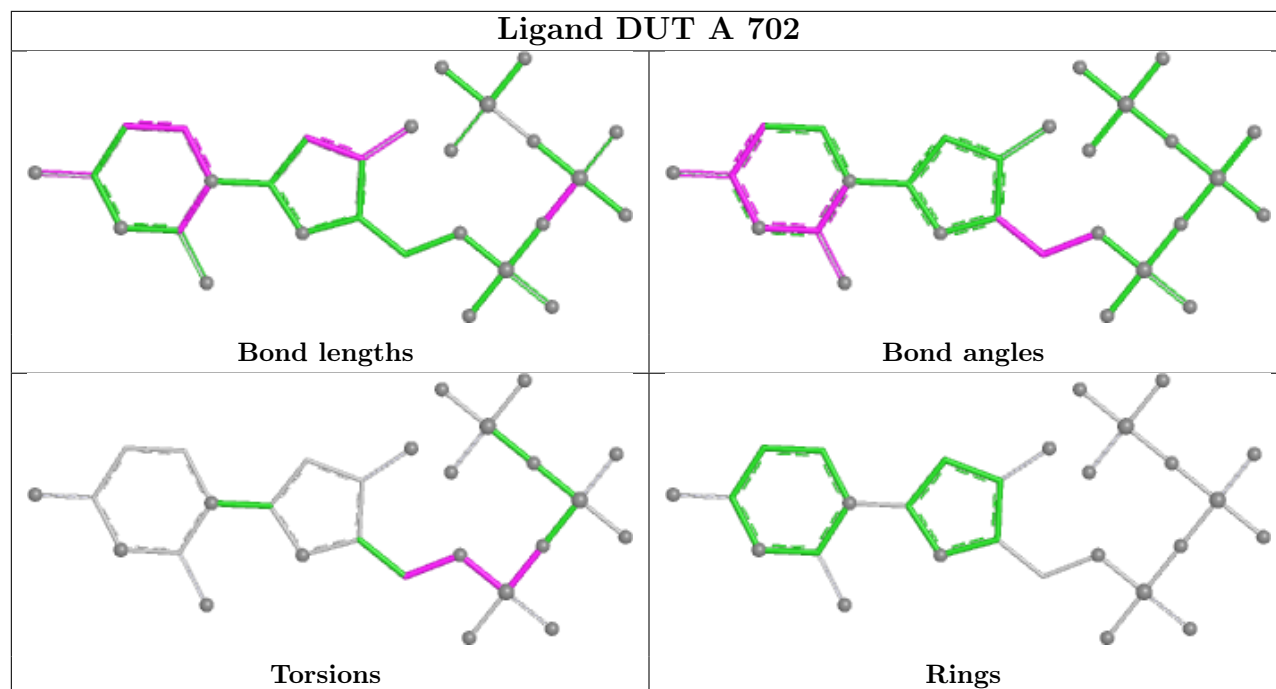
5 monomers are involved in 6 short contacts:

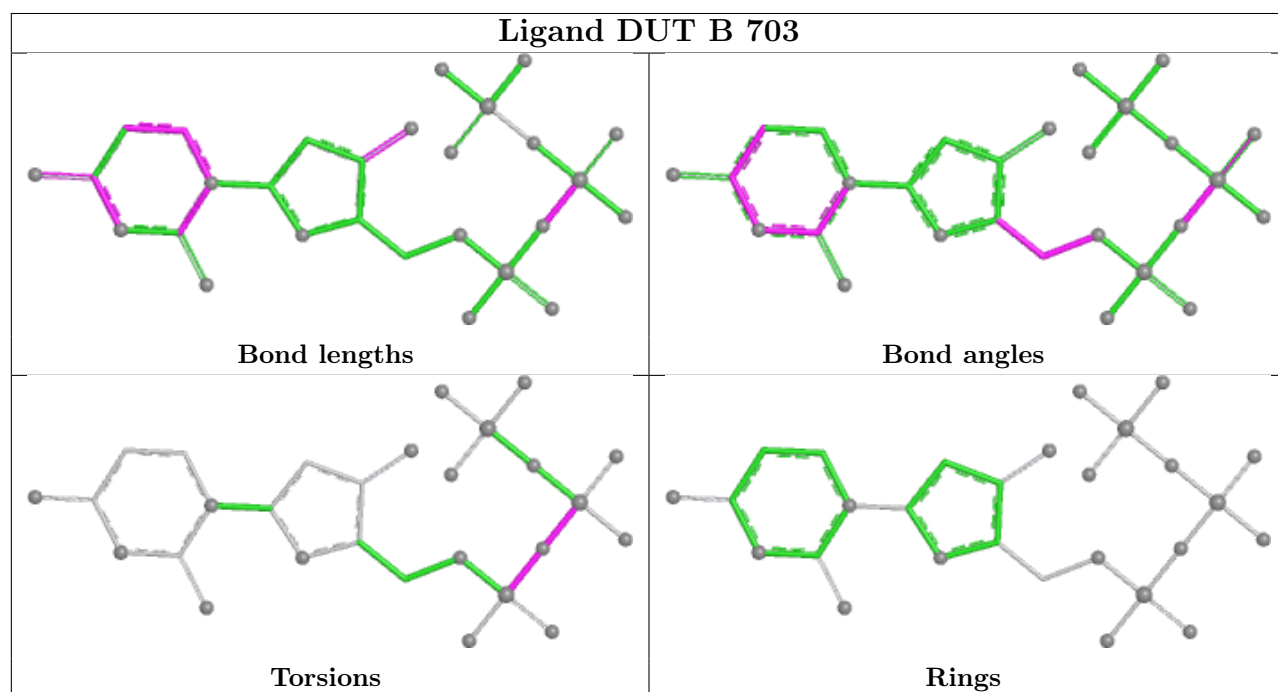
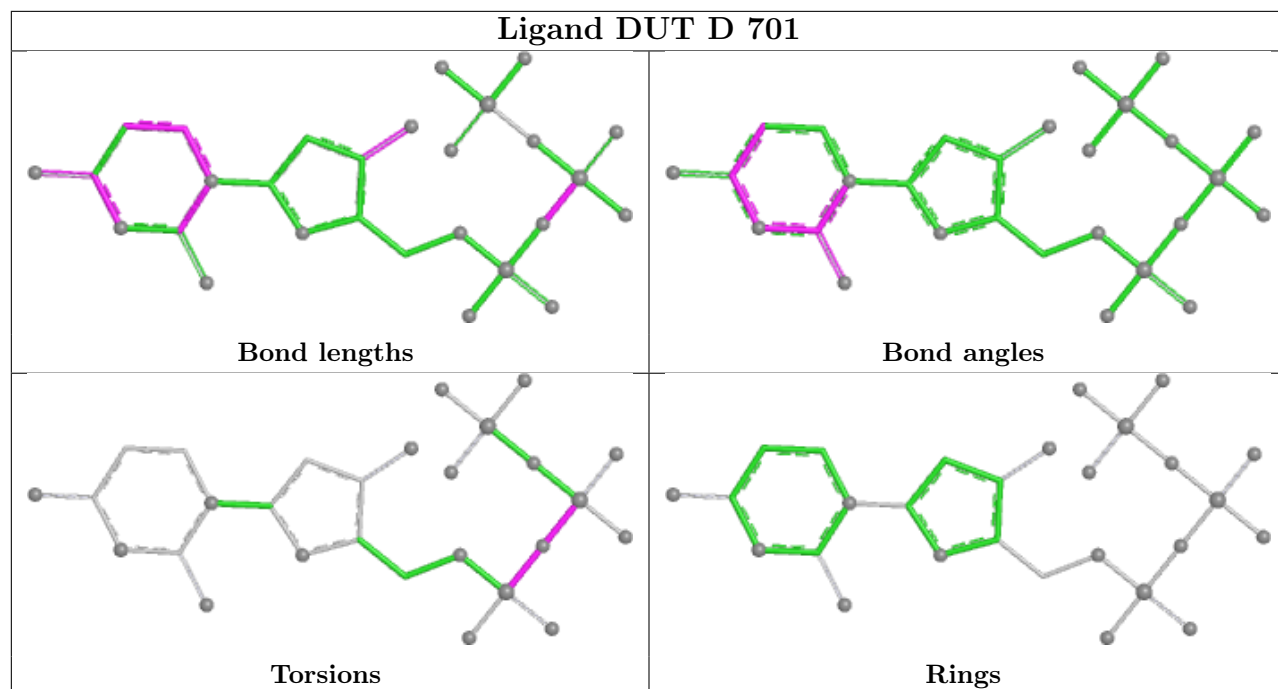
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	701	DUT	1	0
3	B	702	DUT	2	0
2	C	702	DGT	1	0
3	C	703	DUT	1	0
5	D	706	EDO	1	0

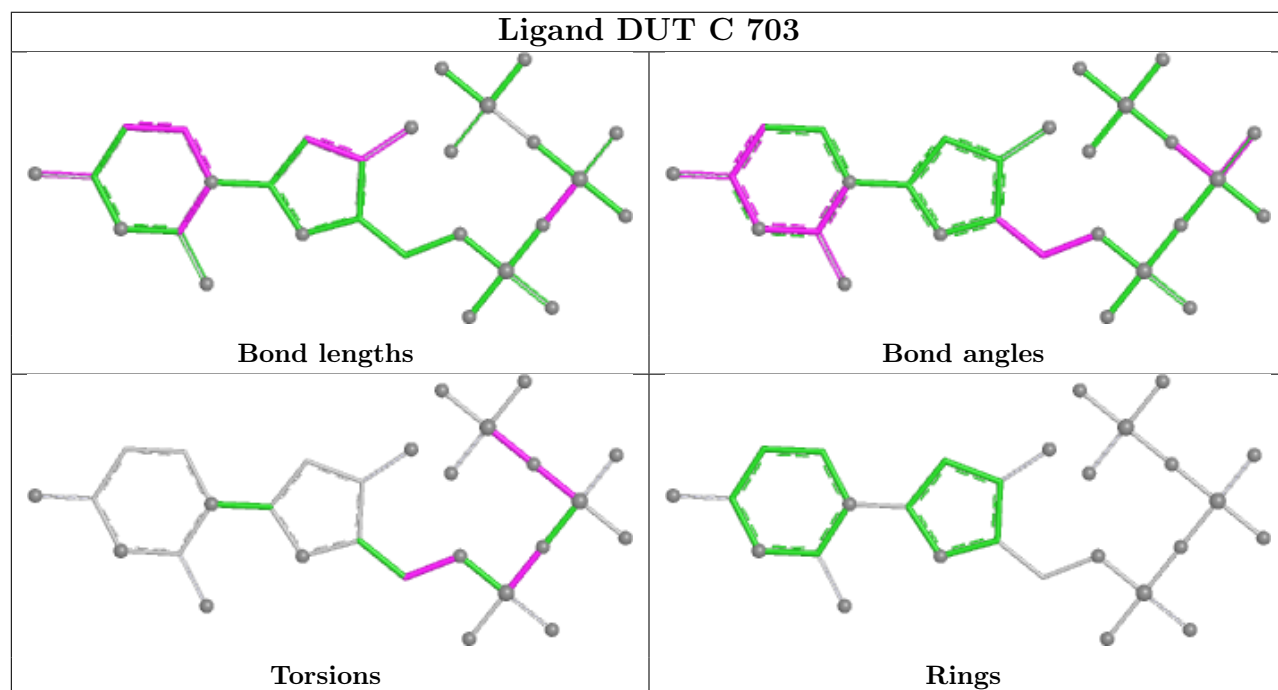
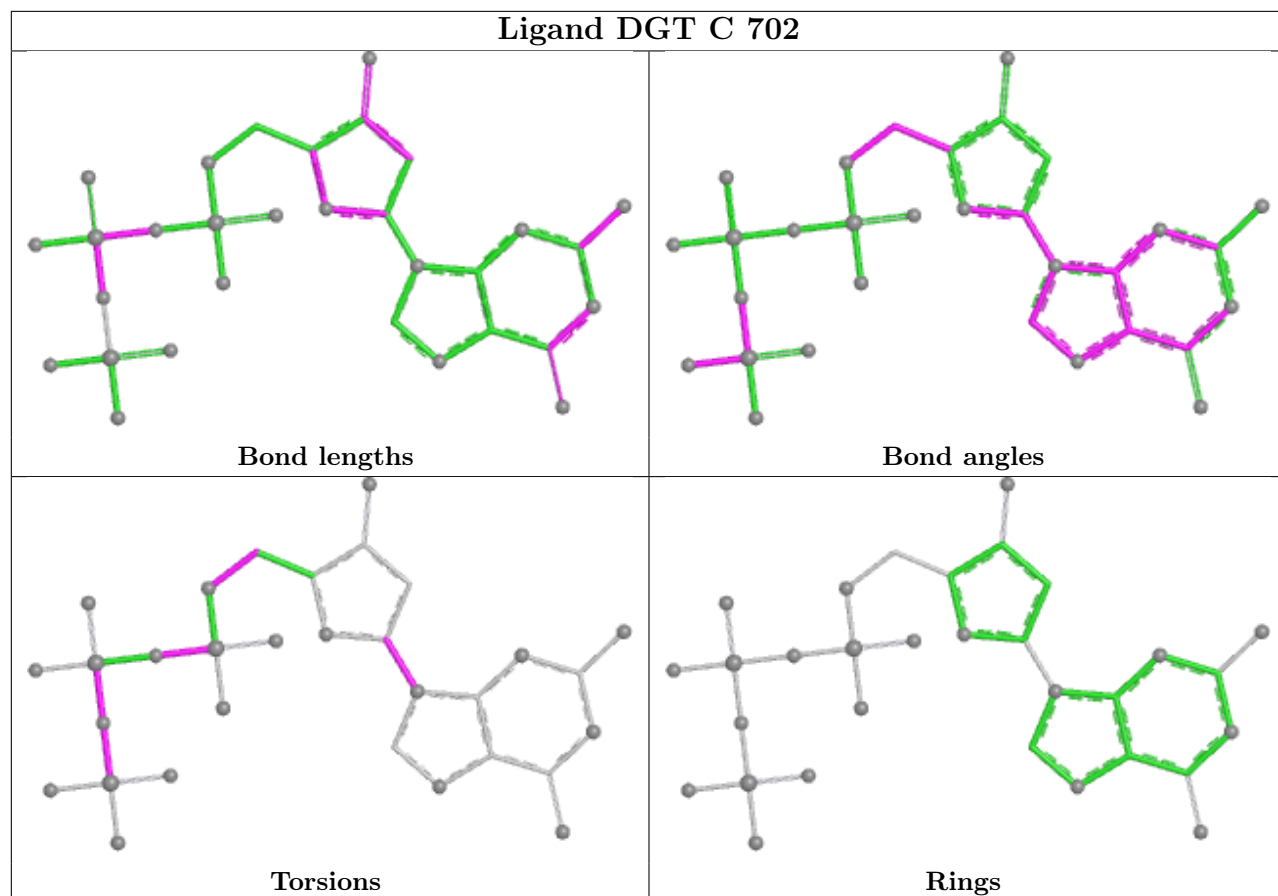
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

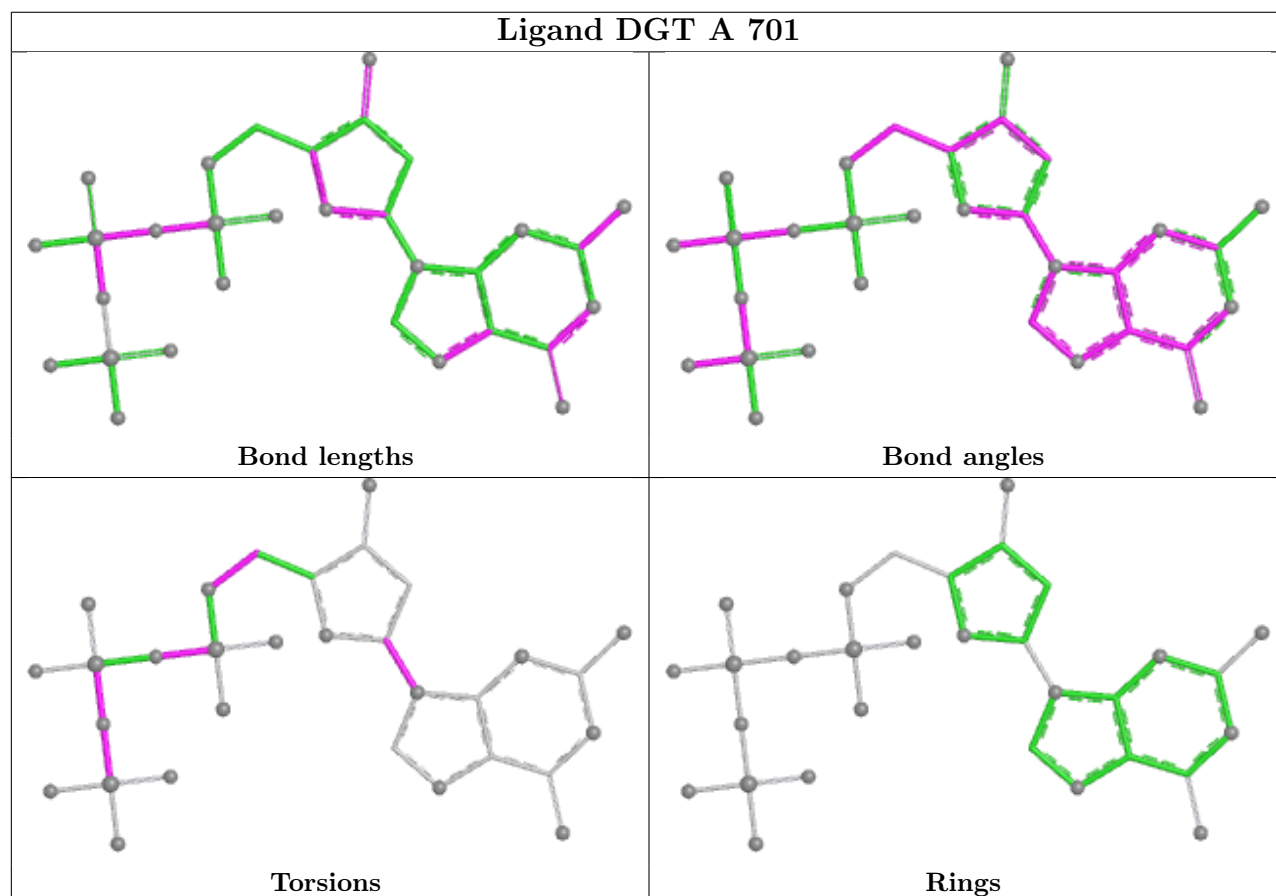
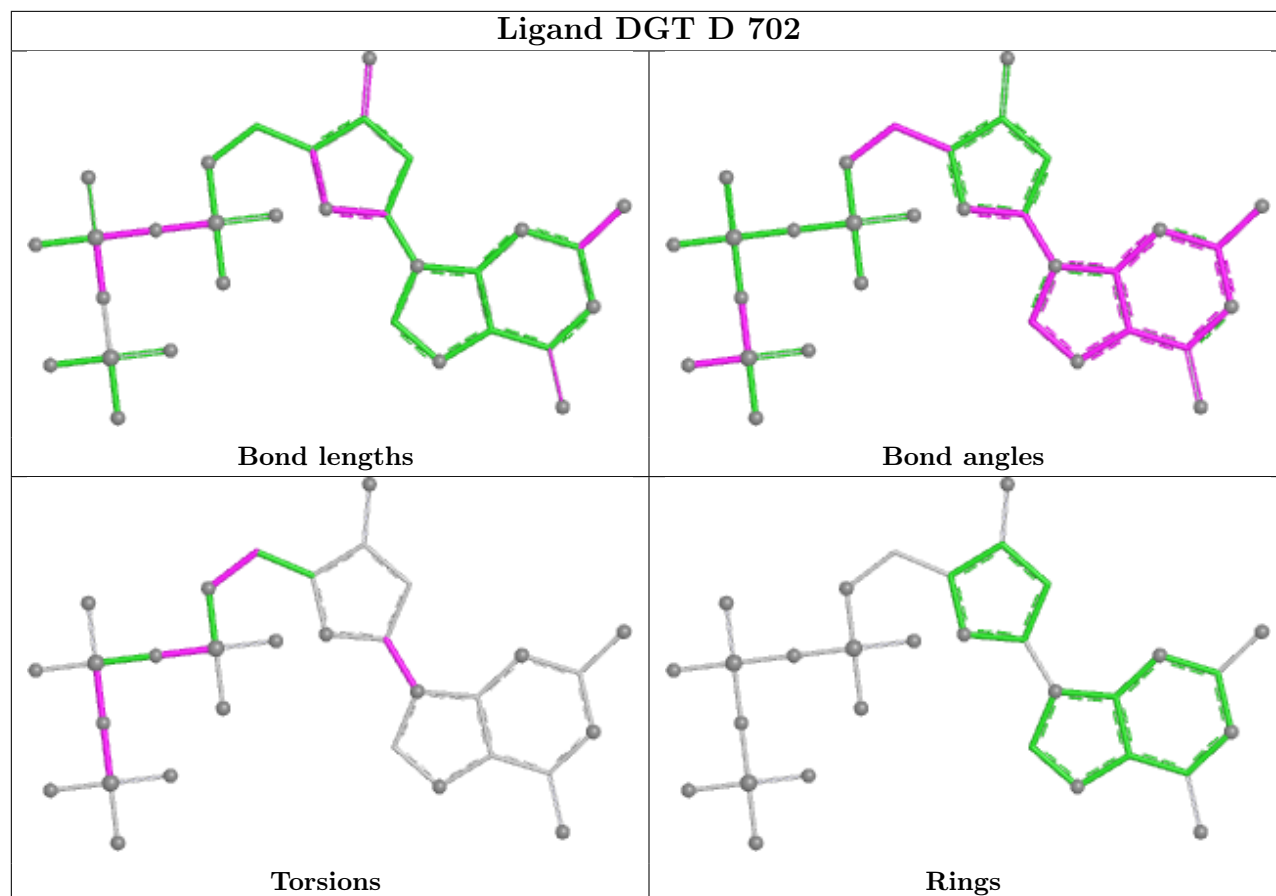












5.7 Other polymers [i](#)

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	481/550 (87%)	0.25	13 (2%) 56 58	20, 36, 56, 81	334 (69%)
1	B	480/550 (87%)	0.14	15 (3%) 51 53	21, 36, 54, 85	286 (59%)
1	C	481/550 (87%)	0.32	13 (2%) 56 58	20, 37, 56, 94	341 (70%)
1	D	481/550 (87%)	-0.02	8 (1%) 69 70	18, 31, 46, 64	315 (65%)
All	All	1923/2200 (87%)	0.17	49 (2%) 58 60	18, 35, 54, 94	1276 (66%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	284	LEU	3.7
1	B	284	LEU	3.6
1	B	593	PRO	3.3
1	A	466	ILE	3.2
1	C	403	GLY	3.2
1	A	284	LEU	3.2
1	C	114	THR	3.0
1	B	560	LYS	2.8
1	A	447	GLN	2.7
1	B	594	GLN	2.7
1	A	115	MET	2.6
1	D	284	LEU	2.6
1	A	529	ALA	2.6
1	A	488	LEU	2.6
1	B	115	MET	2.5
1	C	229	VAL	2.5
1	C	231	TRP	2.5
1	D	522	CYS	2.4
1	A	543	GLU	2.4
1	B	489	LEU	2.4
1	C	488	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	590	LEU	2.4
1	B	488	LEU	2.4
1	A	326	GLN	2.4
1	B	452	ASN	2.3
1	B	114	THR	2.3
1	D	559	ARG	2.3
1	A	426	ILE	2.3
1	C	115	MET	2.2
1	C	531	ARG	2.2
1	C	592	THR	2.2
1	B	487	VAL	2.2
1	B	495	ALA	2.2
1	B	288	LYS	2.2
1	A	473	TYR	2.2
1	D	490	ASP	2.2
1	D	466	ILE	2.1
1	B	479	GLU	2.1
1	C	522	CYS	2.1
1	C	591	ILE	2.1
1	D	426	ILE	2.1
1	B	326	GLN	2.0
1	A	344	ASP	2.0
1	A	452	ASN	2.0
1	A	562	LEU	2.0
1	D	113	ASP	2.0
1	B	491	VAL	2.0
1	C	327	ASN	2.0
1	D	230	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

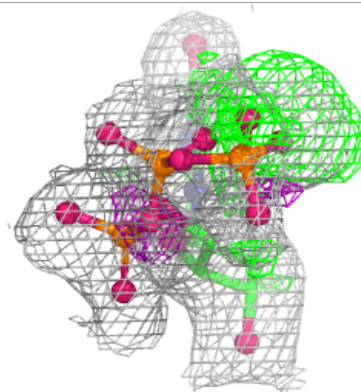
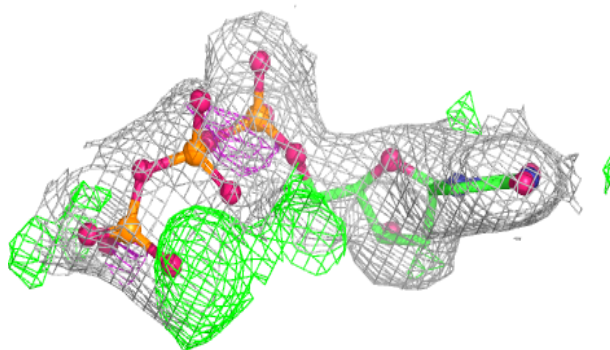
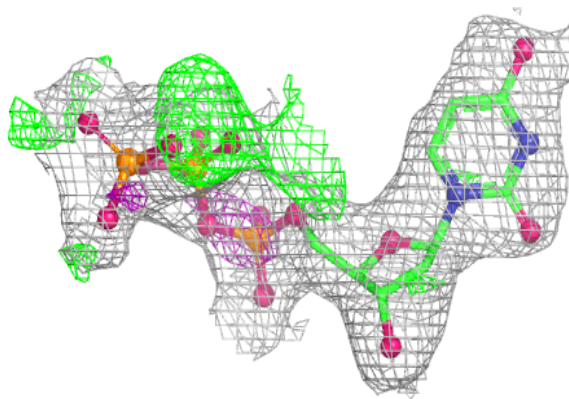
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MG	D	705	1/1	0.73	0.19	58,58,58,58	0
4	MG	B	704	1/1	0.77	0.18	61,61,61,61	0
3	DUT	C	703	28/28	0.78	0.13	25,38,57,78	28
3	DUT	D	703	28/28	0.85	0.11	18,29,54,59	28
5	EDO	D	706	4/4	0.85	0.14	30,30,35,44	4
4	MG	C	704	1/1	0.86	0.08	42,42,42,42	0
3	DUT	B	702	28/28	0.90	0.11	23,32,52,57	28
4	MG	A	705	1/1	0.90	0.04	26,26,26,26	0
3	DUT	A	702	28/28	0.90	0.11	24,33,53,59	28
4	MG	A	706	1/1	0.92	0.10	62,62,62,62	0
2	DGT	B	701	31/31	0.96	0.06	25,30,39,40	31
3	DUT	C	701	28/28	0.96	0.07	22,31,39,43	0
2	DGT	A	701	31/31	0.97	0.06	26,31,36,44	31
3	DUT	A	703	28/28	0.97	0.06	26,33,43,44	0
3	DUT	D	701	28/28	0.97	0.06	22,29,38,41	28
2	DGT	C	702	31/31	0.97	0.06	30,34,42,45	0
3	DUT	B	703	28/28	0.97	0.06	22,27,33,36	28
2	DGT	D	702	31/31	0.98	0.05	20,27,31,35	0
4	MG	A	704	1/1	0.99	0.02	32,32,32,32	1
4	MG	D	704	1/1	0.99	0.02	31,31,31,31	0

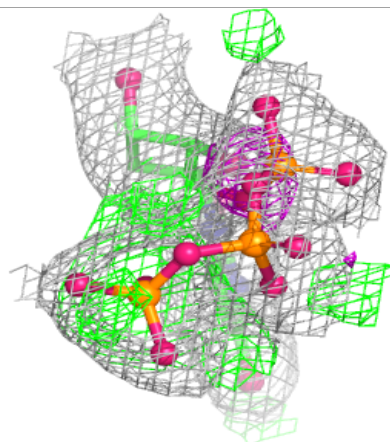
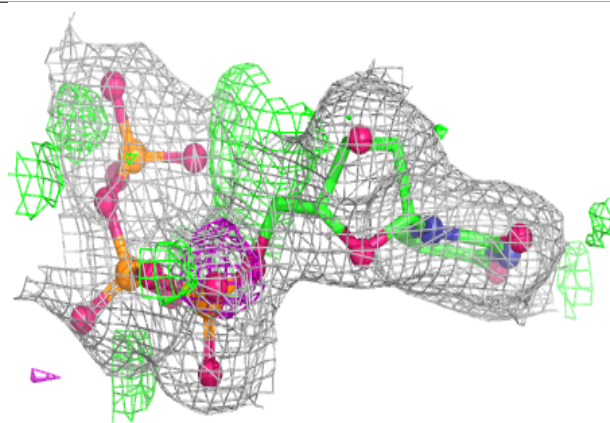
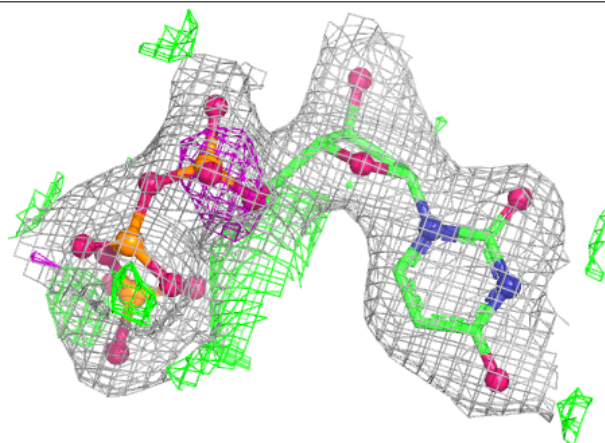
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around DUT C 703:

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

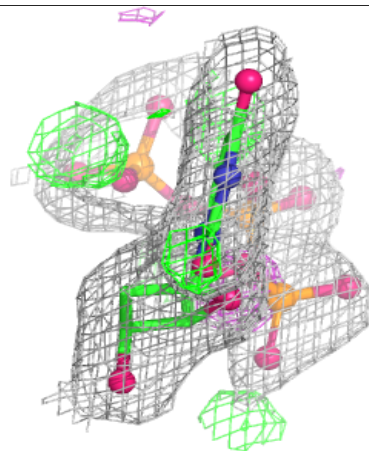
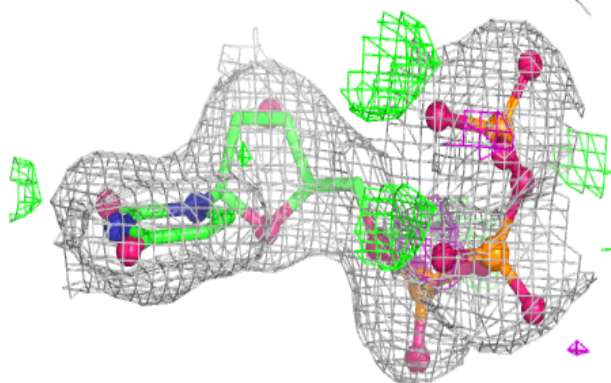
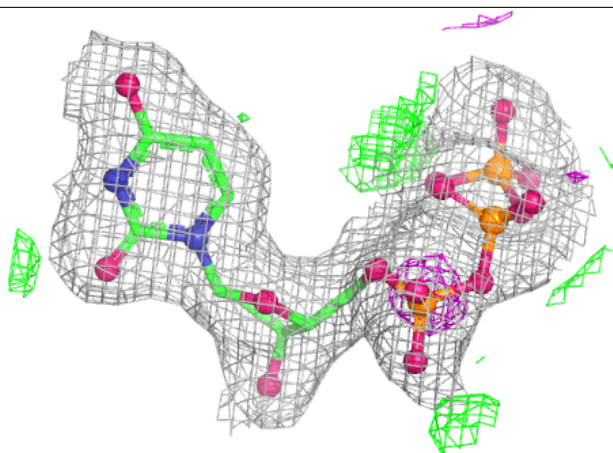
**Electron density around DUT D 703:**

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

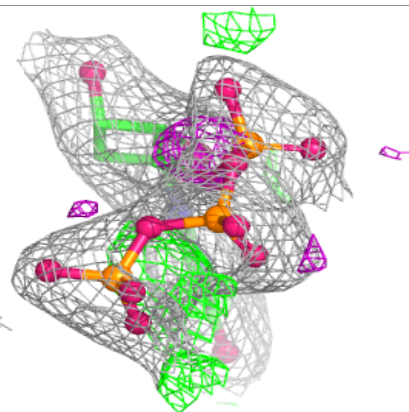
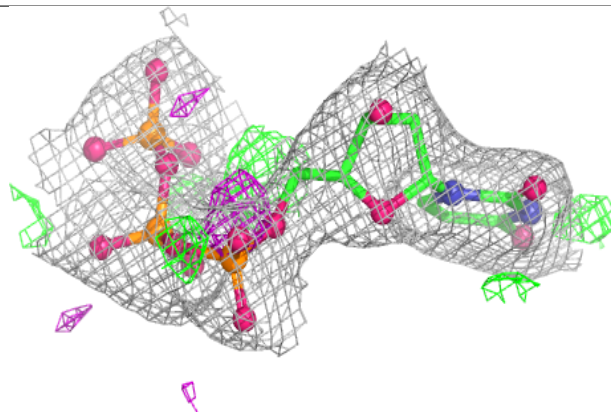
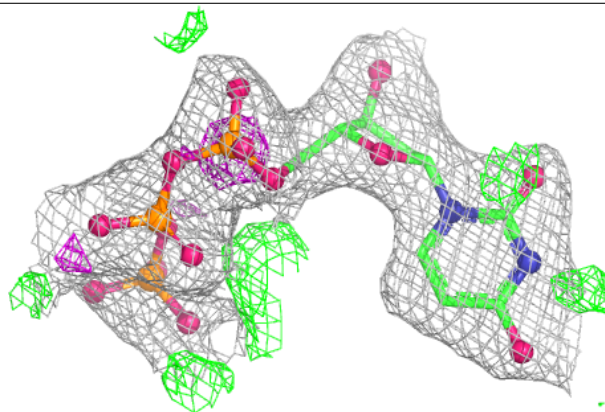


Electron density around DUT B 702:

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

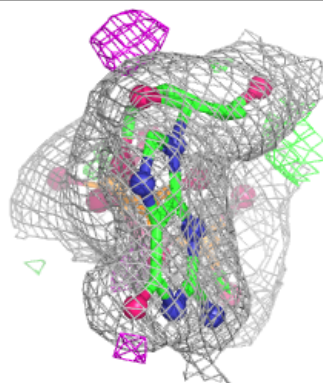
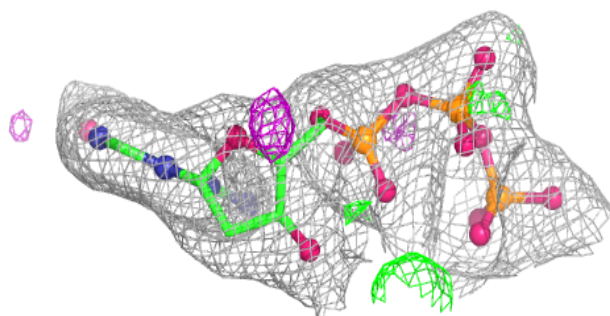
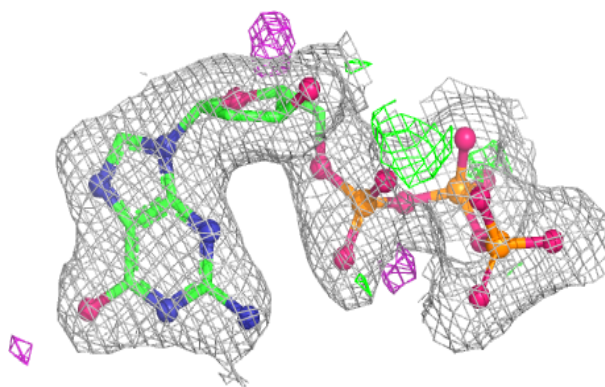
**Electron density around DUT A 702:**

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and green (positive)

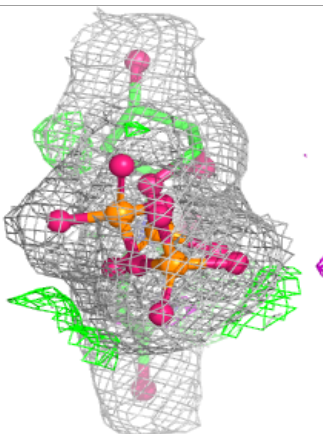
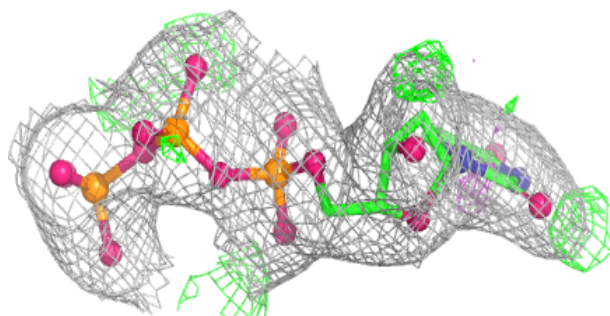
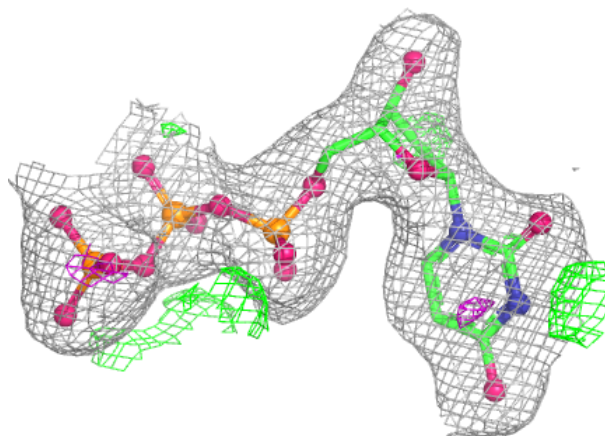


Electron density around DGT B 701:

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

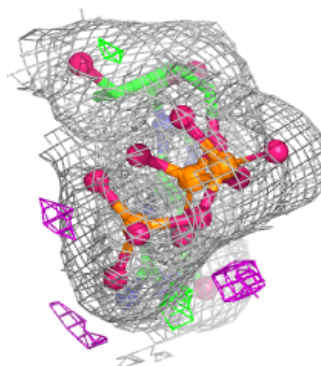
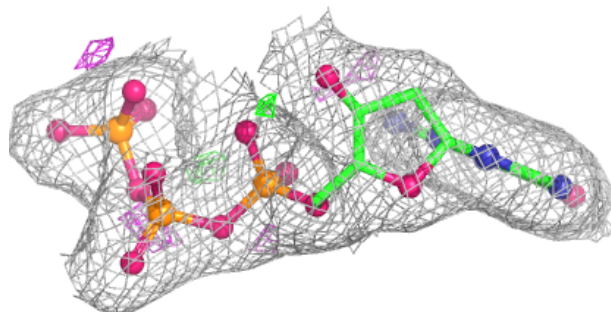
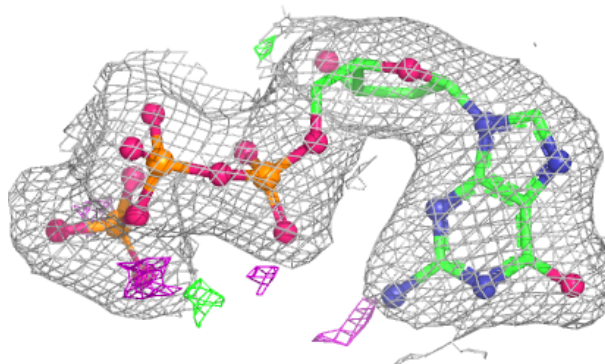
**Electron density around DUT C 701:**

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

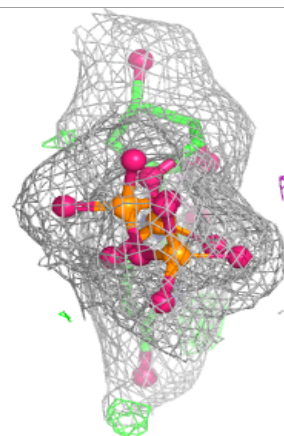
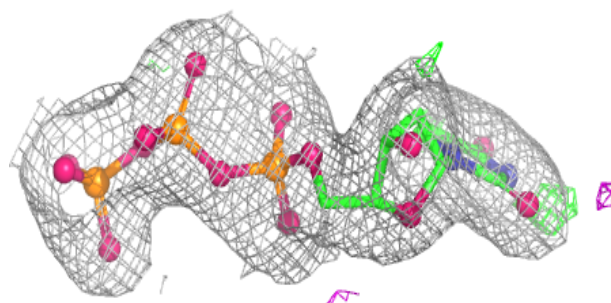
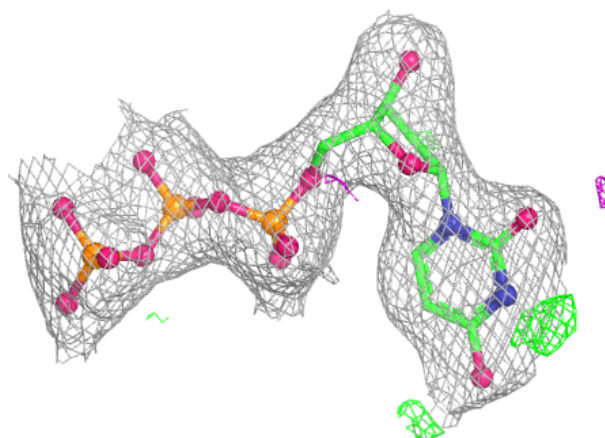


Electron density around DGT A 701:

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

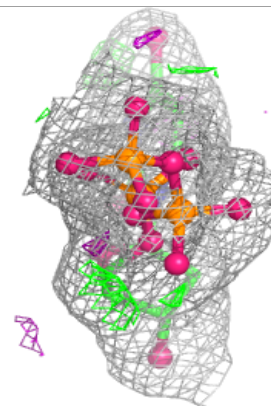
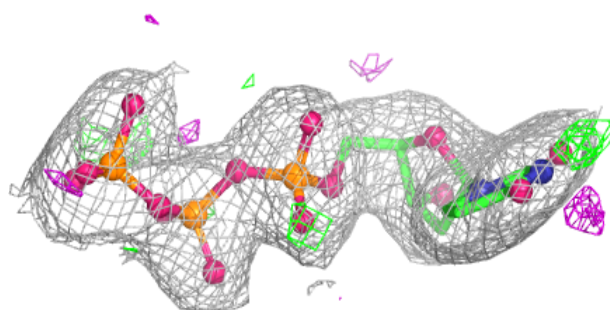
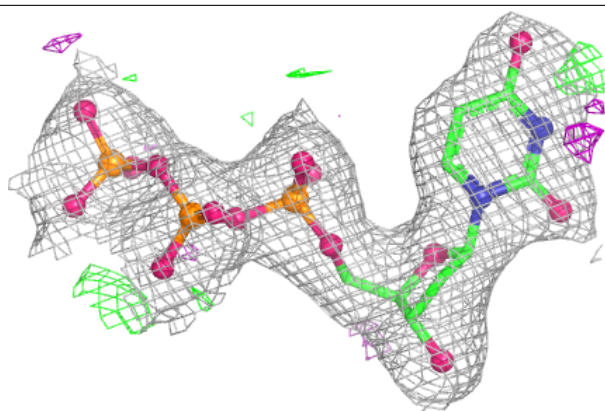
**Electron density around DUT A 703:**

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

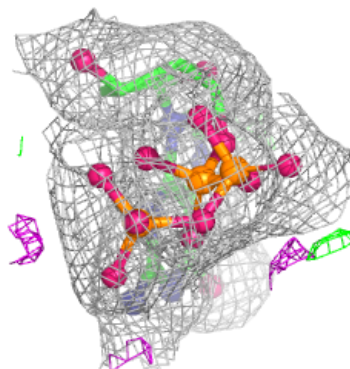
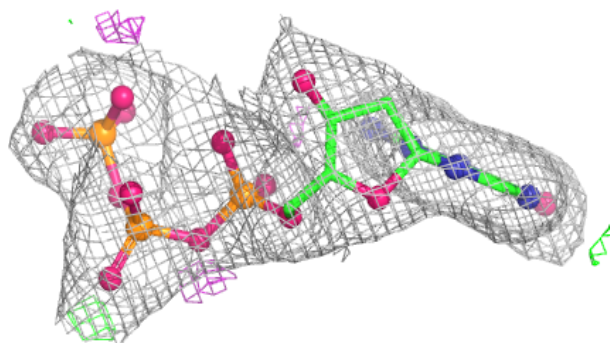
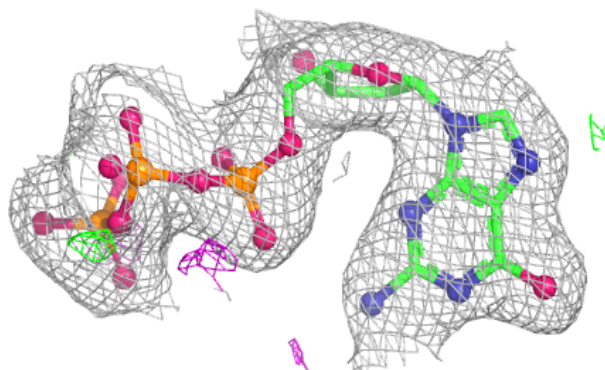


Electron density around DUT D 701:

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

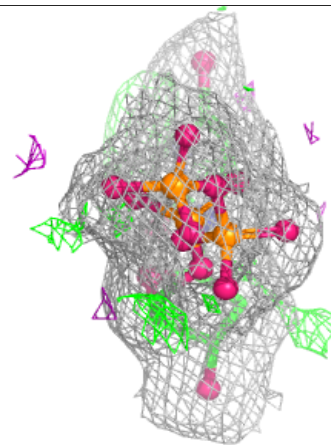
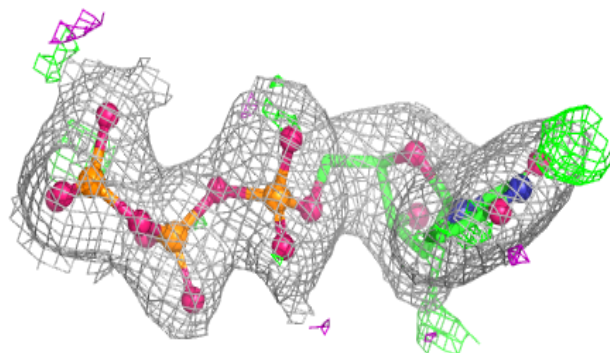
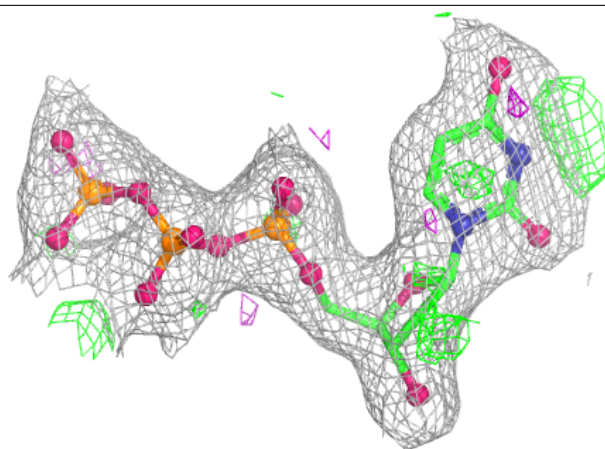
**Electron density around DGT C 702:**

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

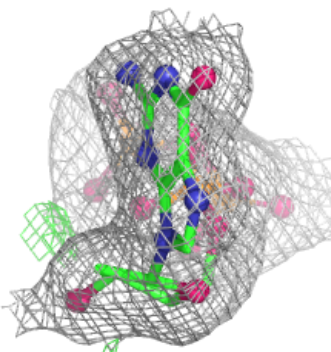
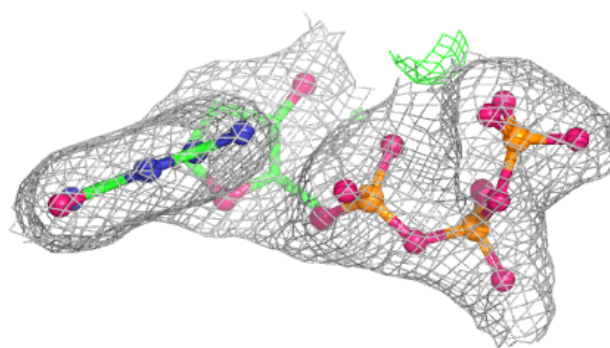
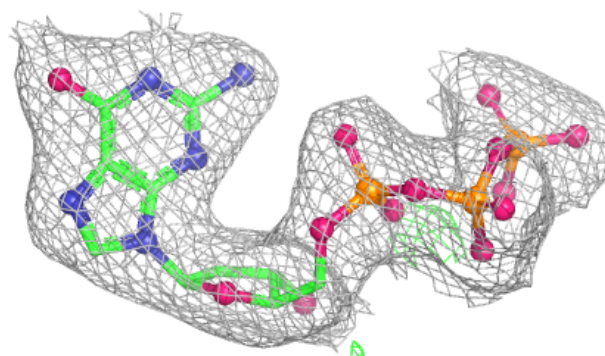


Electron density around DUT B 703:

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

**Electron density around DGT D 702:**

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



6.5 Other polymers [i](#)

There are no such residues in this entry.