wwPDB X-ray Structure Validation Summary Report (i)

Mar 3, 2024 – 06:19 PM EST

PDB ID	:	6DWD
Title	:	SAMHD1 Bound to Clofarabine-TP in the Catalytic Pocket and Allosteric Pocket
Authors	:	Knecht, K.M.; Buzovetsky, O.; Schneider, C.; Thomas, D.; Srikanth, V.; Kaderali, L.; Tofoleanu, F.; Reiss, K.; Ferreiros, N.; Geisslinger, G.; Batista, V.S.; Ji, X.; Cinatl, J.; Keppler, O.T.; Xiong, Y.
Deposited on	:	2018-06-26
Resolution	:	1.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (1)) were used in the production of this report:

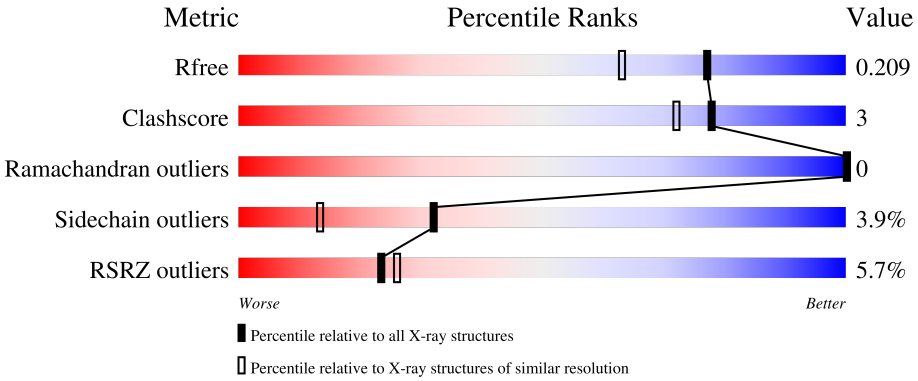
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 <=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	6% 75% 8% • 15%
1	B	550	6% 78% 8% • 13%
1	C	550	5% 80% 6% • 13%
1	D	550	3% 79% 8% • 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GLY	A	716	-	X	-	-
7	SIN	D	714	-	X	-	-

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 17162 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	D	481	Total	C	N	O	S	0	1	0
			3939	2520	686	712	21			
1	C	481	Total	C	N	O	S	0	2	0
			3941	2522	686	712	21			
1	B	481	Total	C	N	O	S	0	0	0
			3933	2517	685	711	20			
1	A	466	Total	C	N	O	S	0	0	0
			3810	2436	665	689	20			

There are 152 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	77	MET	-	initiating methionine	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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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
C	77	MET	-	initiating methionine	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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
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
B	77	MET	-	initiating methionine	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
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

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Chain	Residue	Modelled	Actual	Comment	Reference
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
A	77	MET	-	initiating methionine	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
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

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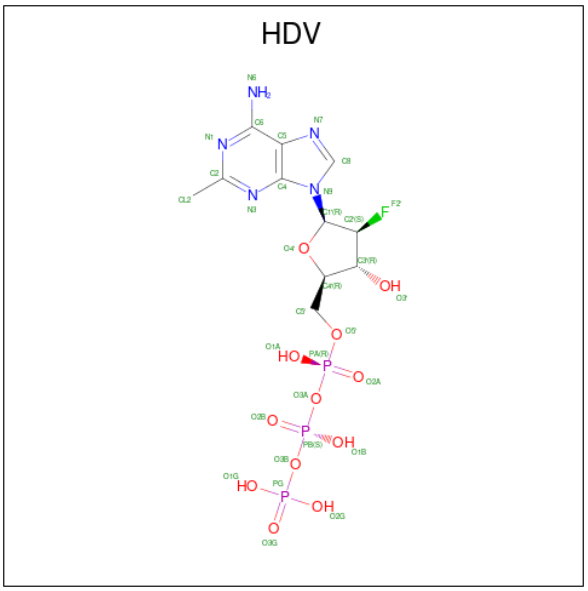
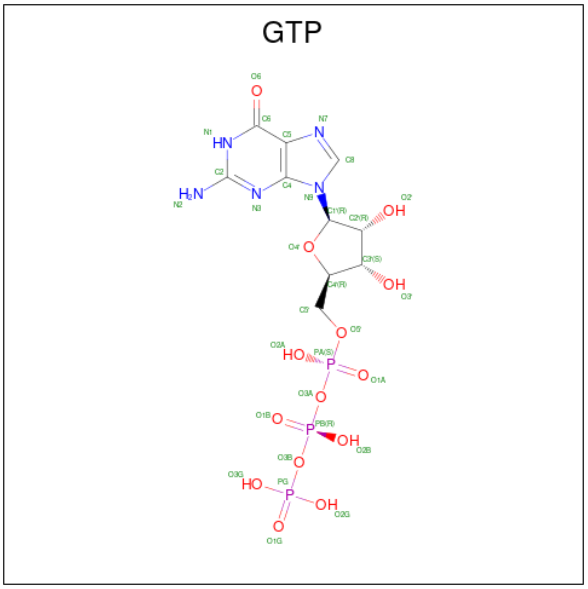
Chain	Residue	Modelled	Actual	Comment	Reference
A	110	PRO	-	expression tag	UNP Q9Y3Z3
A	111	ASN	-	expression tag	UNP Q9Y3Z3
A	112	SER	-	expression tag	UNP Q9Y3Z3
A	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
A	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3

- Molecule 2 is 9-{2-deoxy-2-fluoro-5-O-[(R)-hydroxy{[(S)-hydroxy(phosphonooxy)phosphoryl]oxy}phosphoryl]-beta-D-arabinofuranosyl}-2-me thyl-9H-purin-6-amine (three-letter code: HDV) (formula: $C_{11}H_{17}FN_5O_{12}P_3$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	D	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	D	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	C	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	C	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	B	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	B	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	A	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		
2	A	1	Total	C	F	N	O	P	0	0
			32	11	1	5	12	3		

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: $C_{10}H_{16}N_5O_{14}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
3	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
3	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
3	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

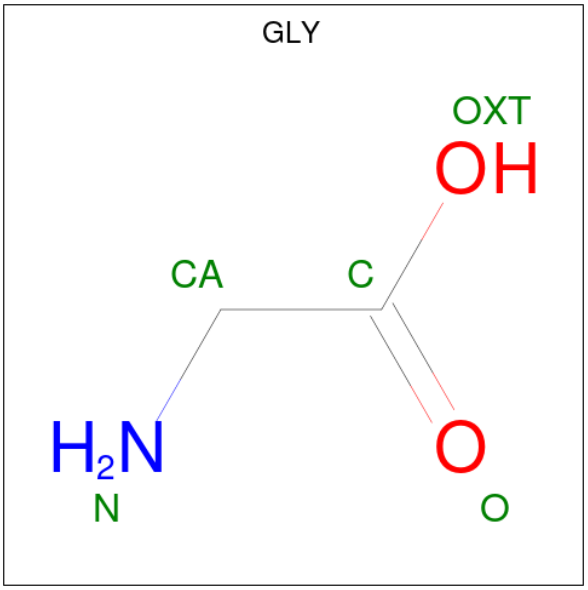
- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

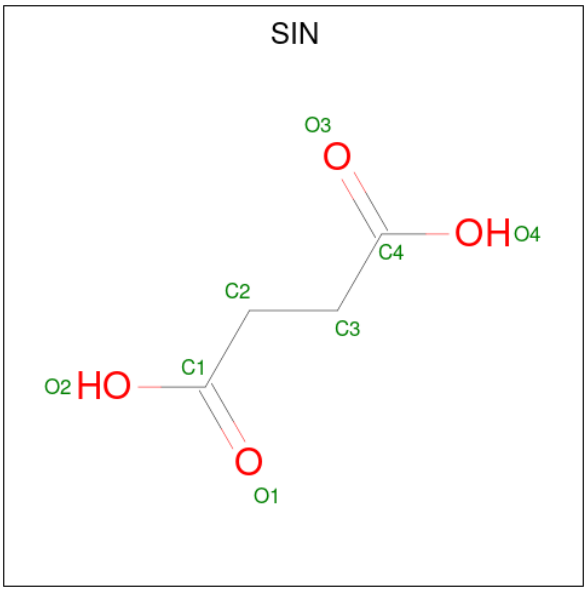
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	D	6	Total Na 6 6	0	0
5	C	4	Total Na 4 4	0	0
5	B	4	Total Na 4 4	0	0
5	A	9	Total Na 9 9	0	0

- Molecule 6 is GLYCINE (three-letter code: GLY) (formula: $C_2H_5NO_2$).



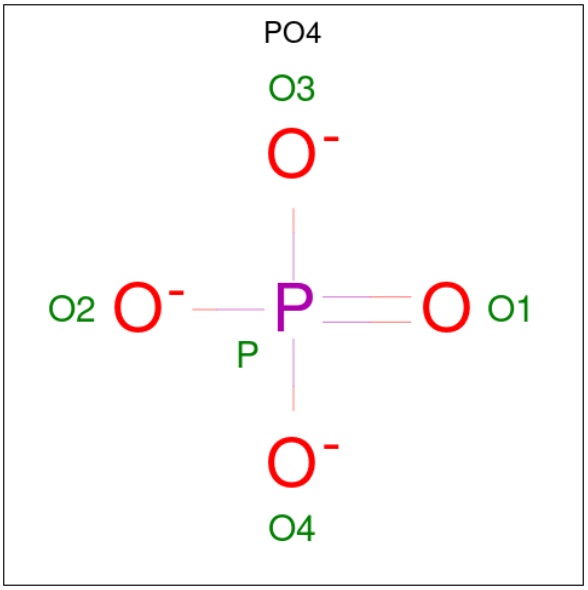
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total C N O 5 2 1 2	0	0
6	C	1	Total C N O 5 2 1 2	0	0
6	C	1	Total C N O 5 2 1 2	0	0
6	B	1	Total C N O 5 2 1 2	0	0
6	A	1	Total C N O 5 2 1 2	0	0
6	A	1	Total C N O 5 2 1 2	0	0
6	A	1	Total C N O 5 2 1 2	0	0

- Molecule 7 is SUCCINIC ACID (three-letter code: SIN) (formula: $C_4H_6O_4$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	D	1	Total C O 8 4 4	0	0
7	B	1	Total C O 8 4 4	0	0

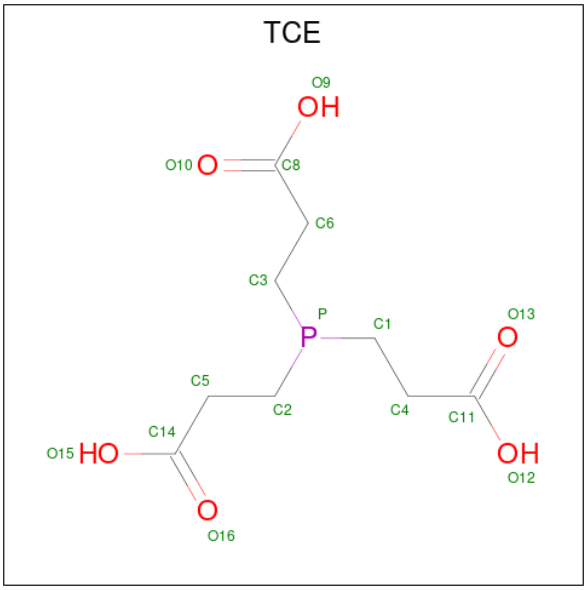
- Molecule 8 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	C	1	Total O P 5 4 1	0	0

- Molecule 9 is 3,3',3''-phosphanetriyltripropanoic acid (three-letter code: TCE) (formula:

$C_9H_{15}O_6P$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total C O P 16 9 6 1	0	0

- Molecule 10 is water.

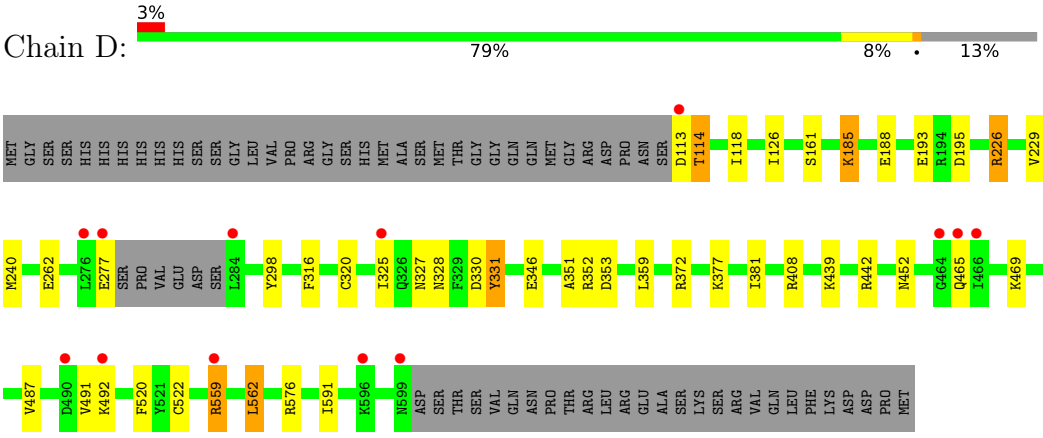
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	D	287	Total O 287 287	0	0
10	C	252	Total O 252 252	0	0
10	B	227	Total O 227 227	0	0
10	A	284	Total O 284 284	0	0

3 Residue-property plots (i)

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

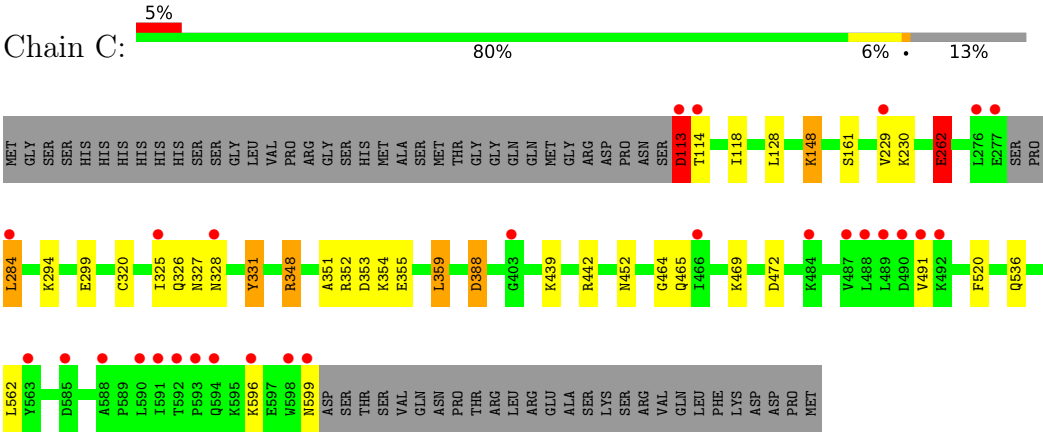
• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain D:



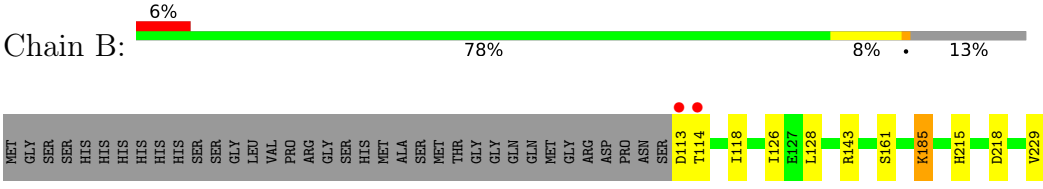
• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

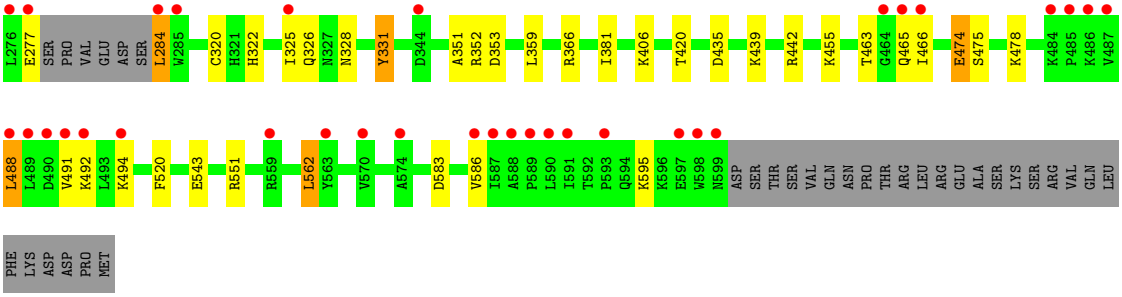
Chain C:



• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

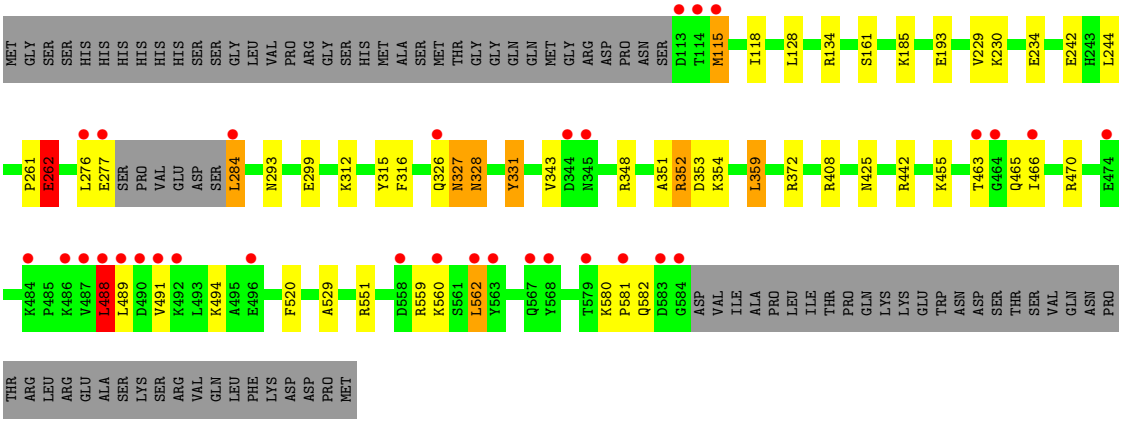
Chain B:





• Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	80.48Å 142.23Å 98.79Å 90.00° 114.10° 90.00°	Depositor
Resolution (Å)	50.00 – 1.70 48.19 – 1.70	Depositor EDS
% Data completeness (in resolution range)	94.0 (50.00-1.70) 94.0 (48.19-1.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.60 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.176 , 0.202 0.185 , 0.209	Depositor DCC
R_{free} test set	10589 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	19.4	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 41.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	17162	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.25% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: HDV, GTP, MG, NA, TCE, PO4, SIN

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	0.99	8/3898 (0.2%)	1.05	18/5258 (0.3%)
1	B	0.89	2/4025 (0.0%)	0.98	9/5433 (0.2%)
1	C	0.95	4/4039 (0.1%)	1.02	12/5452 (0.2%)
1	D	0.95	5/4031 (0.1%)	1.00	15/5441 (0.3%)
All	All	0.94	19/15993 (0.1%)	1.01	54/21584 (0.3%)

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	C	0	1

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	161	SER	CB-OG	12.94	1.59	1.42
1	C	388	ASP	CB-CG	-10.55	1.29	1.51
1	B	161	SER	CB-OG	8.31	1.53	1.42
1	C	113	ASP	CB-CG	7.16	1.66	1.51
1	D	226	ARG	CG-CD	-7.16	1.34	1.51

The worst 5 of 54 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	328	ASN	N-CA-CB	-8.86	94.66	110.60
1	C	388	ASP	CB-CG-OD1	-8.75	110.43	118.30
1	D	442	ARG	NE-CZ-NH1	8.44	124.52	120.30
1	A	328	ASN	CB-CA-C	-8.35	93.70	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	353	ASP	CB-CG-OD1	8.20	125.68	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	464	GLY	Peptide

5.2 Too-close contacts (i)

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	3810	0	3793	30	0
1	B	3933	0	3921	23	0
1	C	3941	0	3932	27	0
1	D	3939	0	3925	23	0
2	A	64	0	0	0	0
2	B	64	0	0	3	0
2	C	64	0	0	1	0
2	D	64	0	0	1	0
3	B	32	0	12	0	0
3	C	32	0	12	0	0
3	D	64	0	24	0	0
4	A	3	0	0	0	0
4	B	1	0	0	0	0
4	C	4	0	0	0	0
4	D	2	0	0	0	0
5	A	9	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	6	0	0	0	0
6	A	15	0	6	0	0
6	B	5	0	2	1	0
6	C	10	0	4	0	0
6	D	5	0	2	0	0
7	B	8	0	4	0	0
7	D	8	0	4	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	C	5	0	0	0	0
9	A	16	0	12	2	0
10	A	284	0	0	5	0
10	B	227	0	0	6	0
10	C	252	0	0	7	0
10	D	287	0	0	10	0
All	All	17162	0	15653	92	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.

The worst 5 of 92 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:C:328[B]:ASN:HD22	1:A:328:ASN:ND2	1.36	1.20
1:C:452:ASN:HB3	10:C:1023:HOH:O	1.49	1.12
2:D:703:HDV:CL2	10:D:830:HOH:O	2.02	1.07
2:B:701:HDV:CL2	10:B:842:HOH:O	2.03	1.03
1:C:465:GLN:HE22	1:C:544:LYS:HE2	1.20	1.01

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	462/550 (84%)	452 (98%)	10 (2%)	0	100	100
1	B	477/550 (87%)	469 (98%)	8 (2%)	0	100	100
1	C	479/550 (87%)	473 (99%)	6 (1%)	0	100	100
1	D	478/550 (87%)	470 (98%)	8 (2%)	0	100	100
All	All	1896/2200 (86%)	1864 (98%)	32 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	413/488 (85%)	395 (96%)	18 (4%)	28	11
1	B	427/488 (88%)	408 (96%)	19 (4%)	28	11
1	C	429/488 (88%)	414 (96%)	15 (4%)	36	17
1	D	428/488 (88%)	414 (97%)	14 (3%)	38	19
All	All	1697/1952 (87%)	1631 (96%)	66 (4%)	32	13

5 of 66 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	359	LEU
1	A	465	GLN
1	A	562	LEU
1	C	469	LYS
1	C	439	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	328	ASN
1	A	577	ASN
1	A	364	HIS
1	C	535	ASN
1	A	326	GLN

5.3.3 RNA

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)

Of 56 ligands modelled in this entry, 33 are monoatomic - leaving 23 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	$\#\|Z\| > 2$
2	HDV	B	701	4	27,34,34	2.12	10 (37%)	34,54,54	1.85	7 (20%)
6	GLY	D	713	-	4,4,4	1.13	1 (25%)	3,4,4	1.46	0
2	HDV	A	702	4	27,34,34	1.82	9 (33%)	34,54,54	1.92	7 (20%)
6	GLY	C	713	-	4,4,4	0.99	0	3,4,4	1.25	0
6	GLY	A	715	-	4,4,4	1.10	0	3,4,4	0.63	0
3	GTP	D	704	4	26,34,34	1.05	1 (3%)	32,54,54	1.14	2 (6%)
7	SIN	B	710	-	7,7,7	1.37	0	8,8,8	1.37	1 (12%)
2	HDV	C	701	4	27,34,34	2.21	6 (22%)	34,54,54	2.36	11 (32%)
3	GTP	C	702	4	26,34,34	1.38	4 (15%)	32,54,54	1.45	5 (15%)
7	SIN	D	714	5	7,7,7	1.20	0	8,8,8	2.80	3 (37%)
6	GLY	A	716	-	4,4,4	1.24	1 (25%)	3,4,4	2.29	1 (33%)
3	GTP	D	702	4	26,34,34	1.30	5 (19%)	32,54,54	1.45	5 (15%)
2	HDV	C	704	4	27,34,34	2.01	7 (25%)	34,54,54	1.55	8 (23%)
2	HDV	B	702	4	27,34,34	1.88	6 (22%)	34,54,54	2.33	8 (23%)
6	GLY	C	712	-	4,4,4	1.19	0	3,4,4	1.73	1 (33%)
2	HDV	A	704	4	27,34,34	1.94	9 (33%)	34,54,54	2.43	17 (50%)
6	GLY	B	709	-	4,4,4	0.68	0	3,4,4	1.94	1 (33%)
2	HDV	D	703	4	27,34,34	2.50	8 (29%)	34,54,54	1.92	8 (23%)
8	PO4	C	714	-	4,4,4	0.90	0	6,6,6	1.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
9	TCE	A	718	-	12,15,15	1.37	1 (8%)	12,18,18	3.37	5 (41%)
2	HDV	D	701	4	27,34,34	1.97	7 (25%)	34,54,54	2.74	14 (41%)
6	GLY	A	717	-	4,4,4	0.92	0	3,4,4	1.44	0
3	GTP	B	703	4	26,34,34	1.22	3 (11%)	32,54,54	1.43	4 (12%)

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	HDV	B	701	4	-	5/18/38/38	0/3/3/3
6	GLY	D	713	-	-	2/2/2/2	-
2	HDV	A	702	4	-	1/18/38/38	0/3/3/3
6	GLY	C	713	-	-	0/2/2/2	-
6	GLY	A	715	-	-	0/2/2/2	-
3	GTP	D	704	4	-	2/18/38/38	0/3/3/3
7	SIN	B	710	-	-	3/5/5/5	-
2	HDV	C	701	4	-	4/18/38/38	0/3/3/3
3	GTP	C	702	4	-	5/18/38/38	0/3/3/3
7	SIN	D	714	5	-	5/5/5/5	-
6	GLY	A	716	-	-	2/2/2/2	-
3	GTP	D	702	4	-	2/18/38/38	0/3/3/3
2	HDV	C	704	4	-	3/18/38/38	0/3/3/3
2	HDV	B	702	4	-	4/18/38/38	0/3/3/3
6	GLY	C	712	-	-	2/2/2/2	-
2	HDV	A	704	4	-	4/18/38/38	0/3/3/3
6	GLY	B	709	-	-	0/2/2/2	-
2	HDV	D	703	4	-	1/18/38/38	0/3/3/3
9	TCE	A	718	-	-	9/15/15/15	-
2	HDV	D	701	4	-	3/18/38/38	0/3/3/3
6	GLY	A	717	-	-	2/2/2/2	-
3	GTP	B	703	4	-	5/18/38/38	0/3/3/3

The worst 5 of 78 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	703	HDV	O4'-C1'	9.41	1.54	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	704	HDV	O4'-C1'	6.58	1.50	1.41
2	B	701	HDV	O4'-C1'	6.28	1.49	1.41
2	C	701	HDV	C2'-C3'	-5.75	1.44	1.52
2	D	701	HDV	O4'-C1'	5.33	1.48	1.41

The worst 5 of 108 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	702	HDV	C2-N3-C4	7.71	121.79	115.52
2	D	701	HDV	CL2-C2-N1	6.52	127.34	117.15
9	A	718	TCE	C1-P-C3	6.44	121.47	100.95
2	A	702	HDV	C5-C6-N1	-6.36	116.84	121.01
7	D	714	SIN	C2-C3-C4	6.26	127.06	113.60

There are no chirality outliers.

5 of 64 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	701	HDV	C3'-C4'-C5'-O5'
2	C	701	HDV	C3'-C4'-C5'-O5'
2	B	701	HDV	PB-O3B-PG-O1G
2	B	702	HDV	C3'-C4'-C5'-O5'
2	A	704	HDV	C3'-C4'-C5'-O5'

There are no ring outliers.

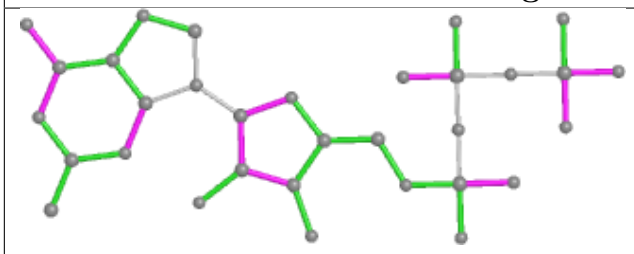
7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	HDV	1	0
7	D	714	SIN	2	0
2	C	704	HDV	1	0
2	B	702	HDV	2	0
6	B	709	GLY	1	0
2	D	703	HDV	1	0
9	A	718	TCE	2	0

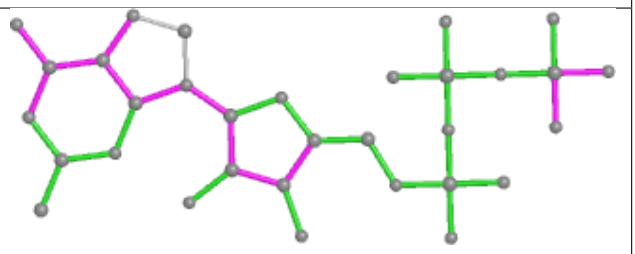
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

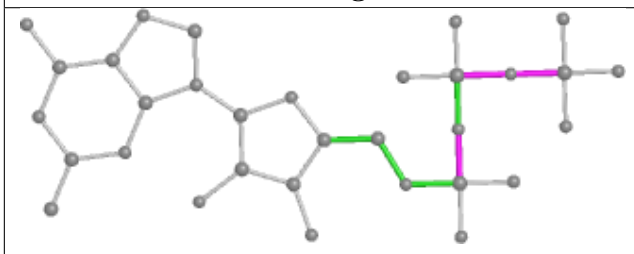
Ligand HDV B 701



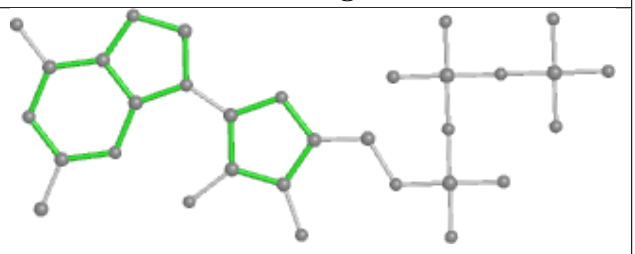
Bond lengths



Bond angles

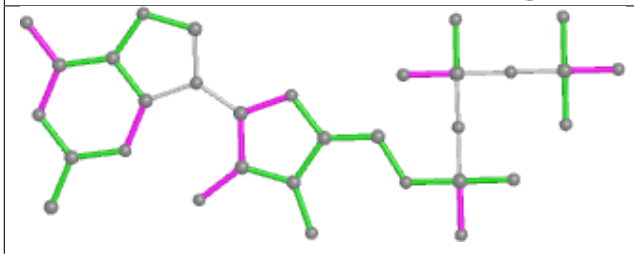


Torsions



Rings

Ligand HDV A 702



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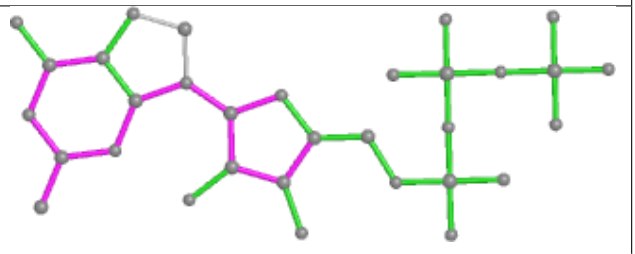

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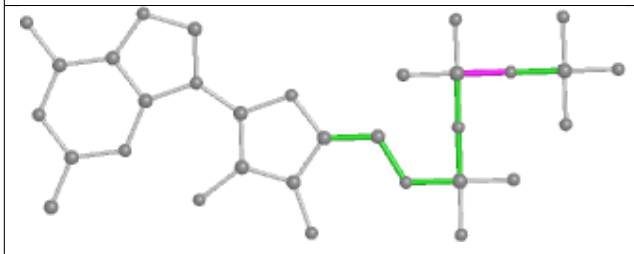

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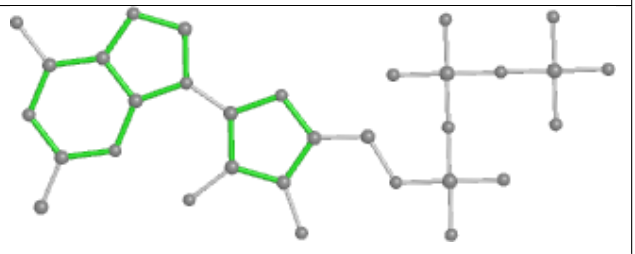
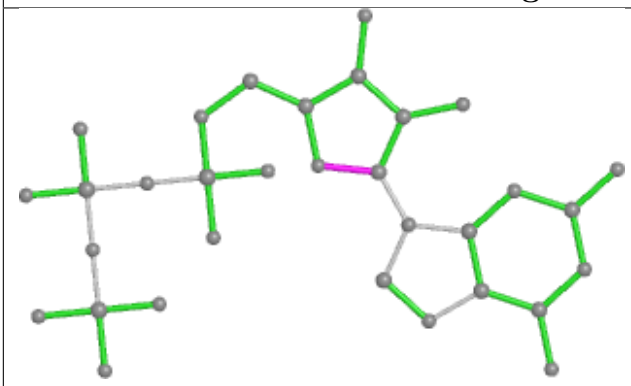
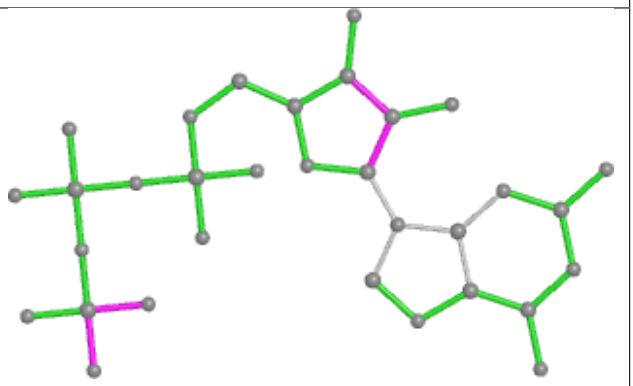


Rings

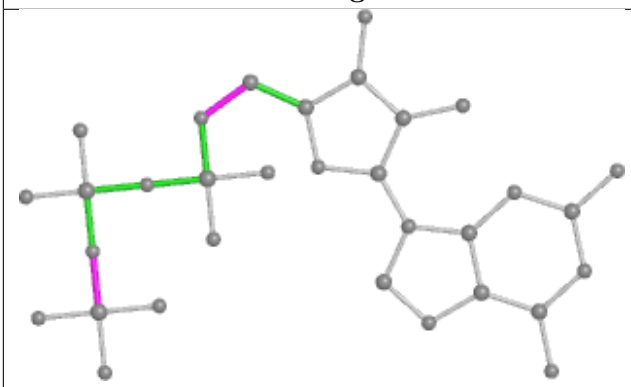
Ligand GTP D 704



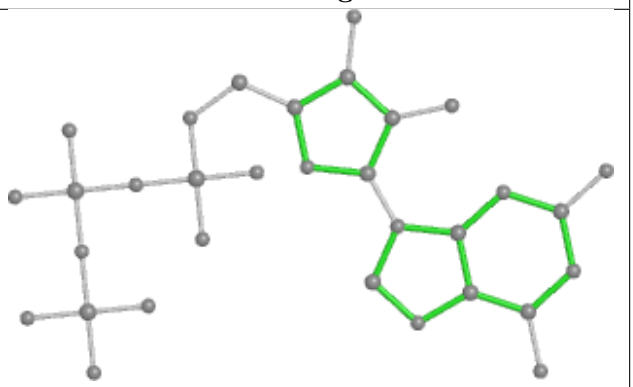
Bond lengths



Bond angles

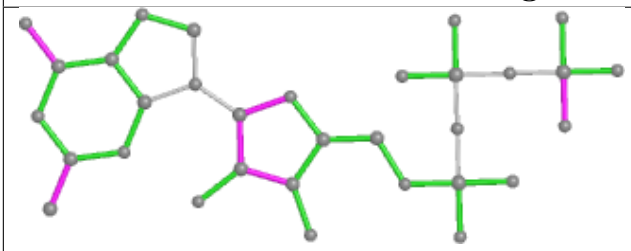


Torsions

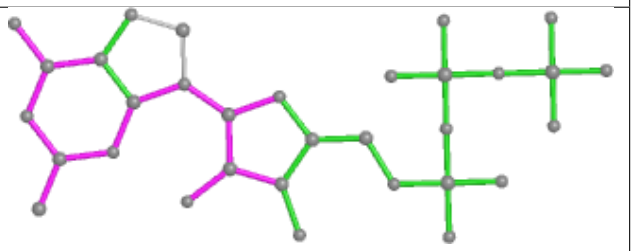


Rings

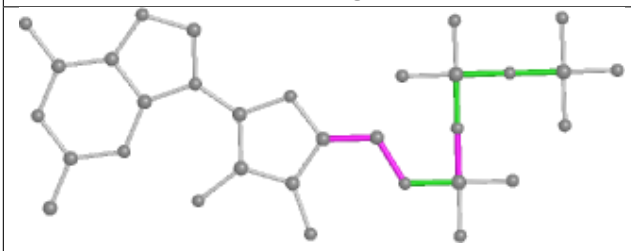
Ligand HDV C 701



Bond lengths



Bond angles



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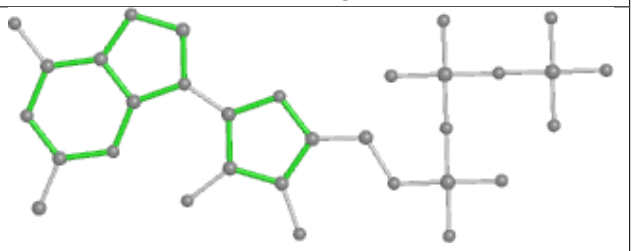
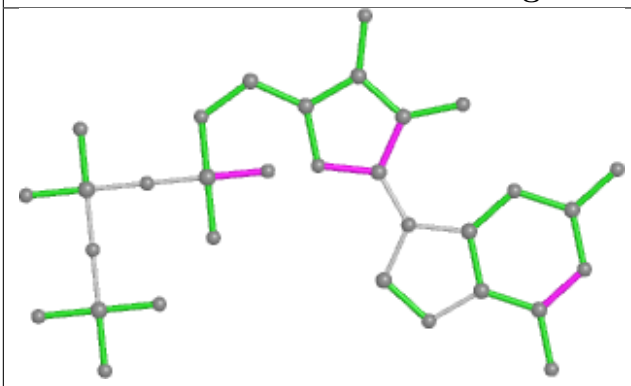
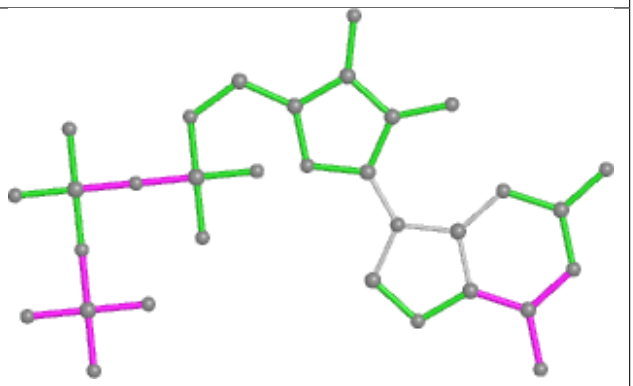


Rings

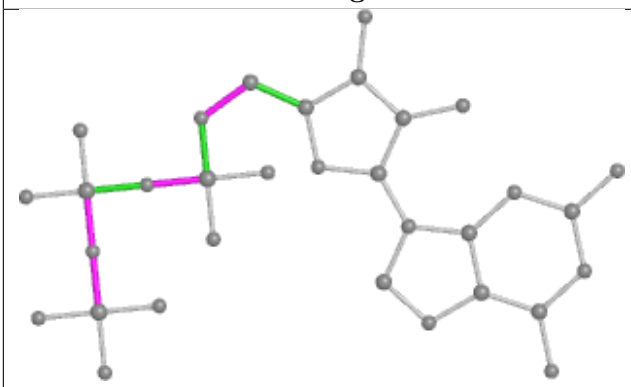
Ligand GTP C 702



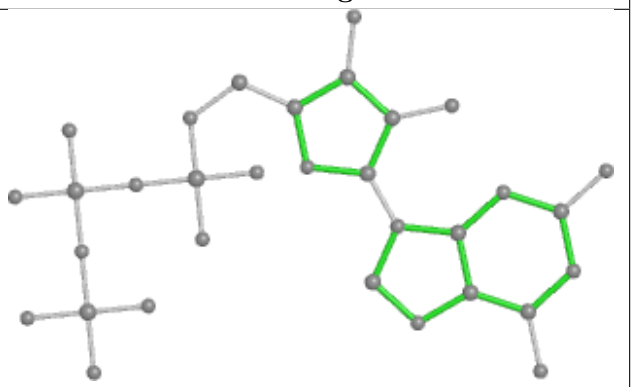
Bond lengths



Bond angles

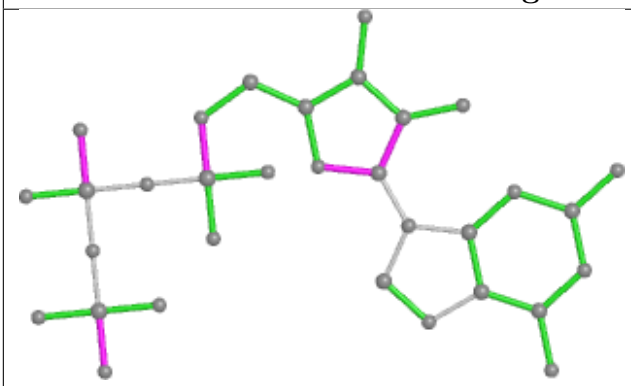


Torsions

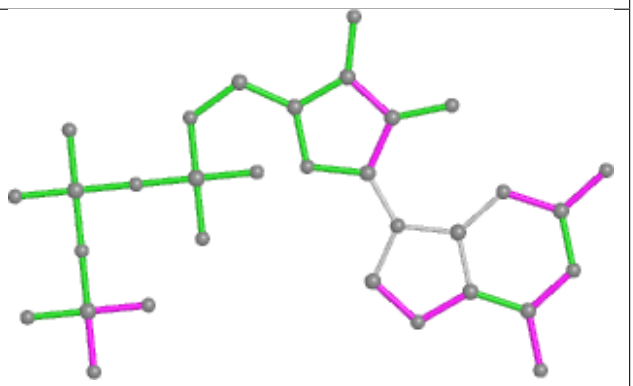


Rings

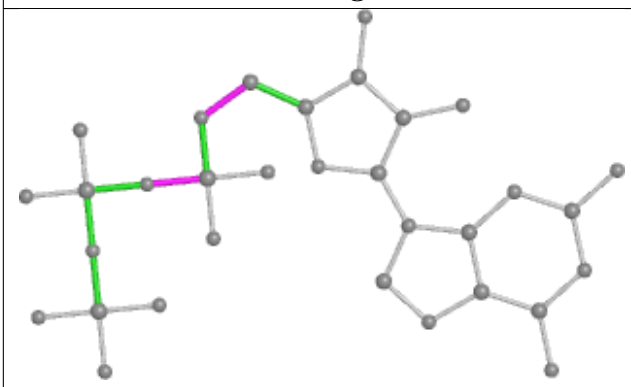
Ligand GTP D 702



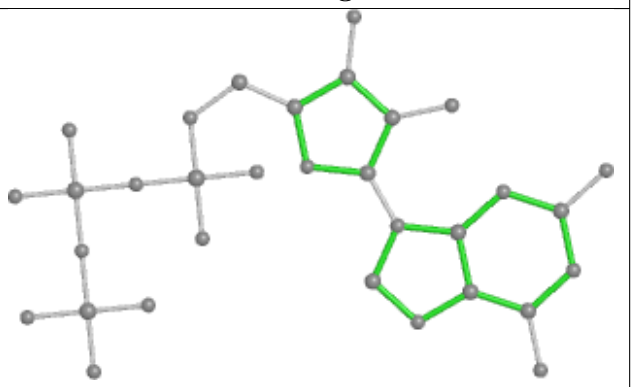
Bond lengths



Bond angles

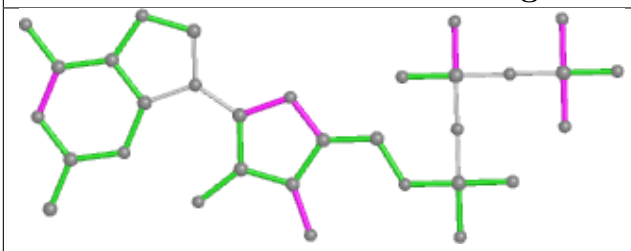


Torsions

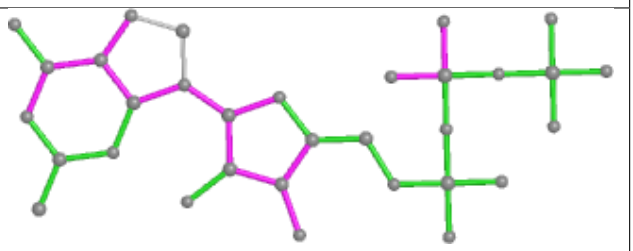


Rings

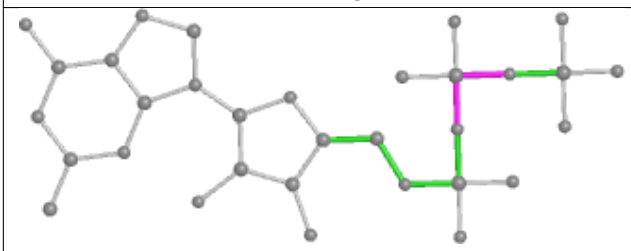
Ligand HDV C 704



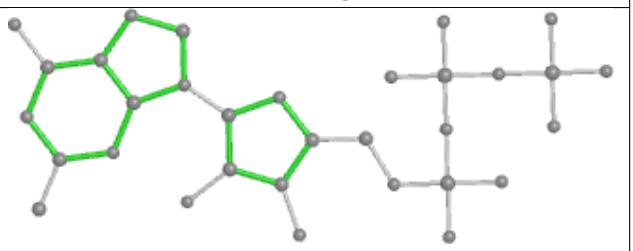
Bond lengths



Bond angles

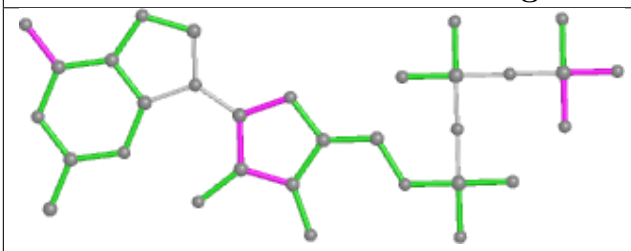


Torsions

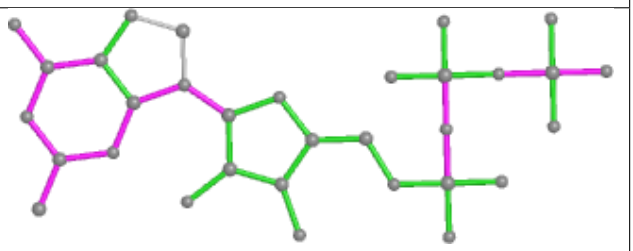


Rings

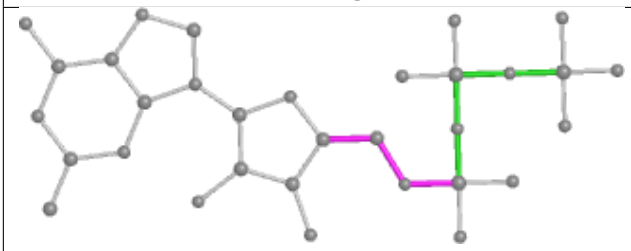
Ligand HDV B 702



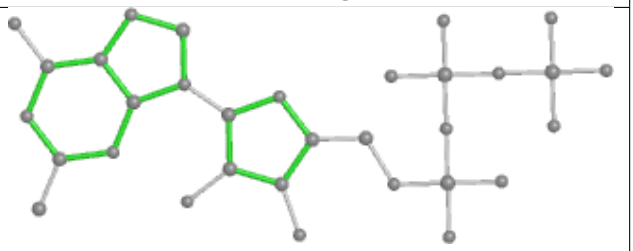
Bond lengths



Bond angles

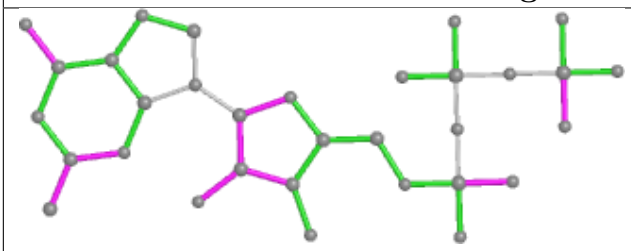


Torsions

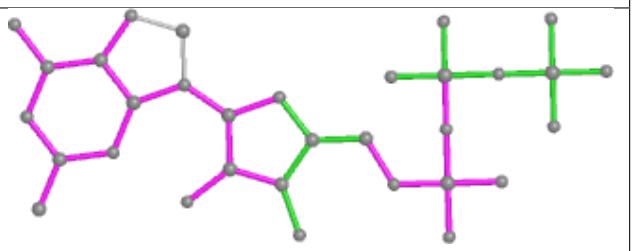


Rings

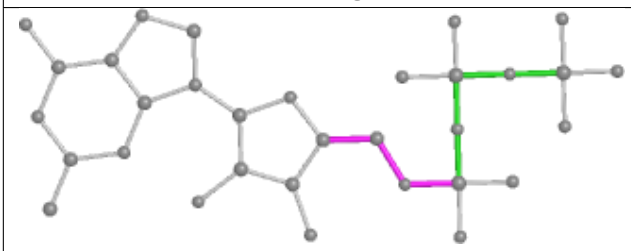
Ligand HDV A 704



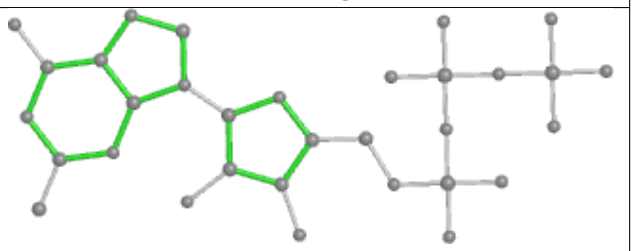
Bond lengths



Bond angles

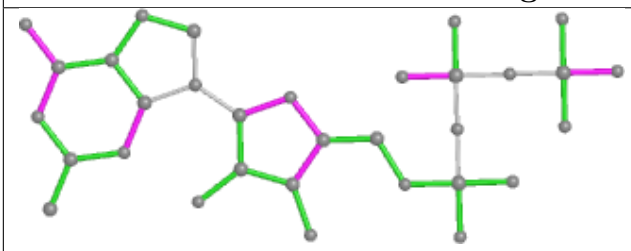


Torsions

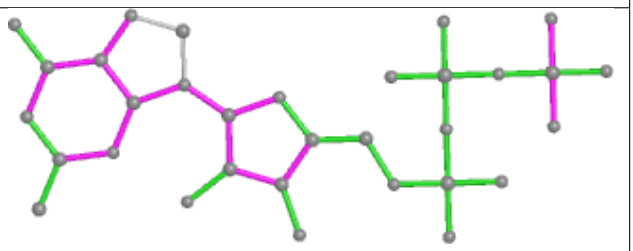


Rings

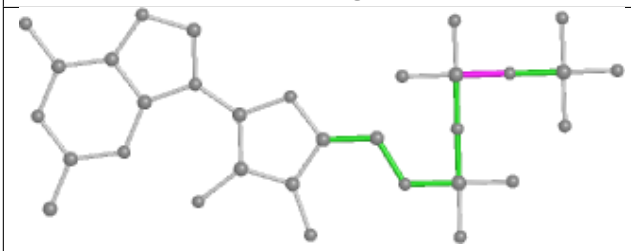
Ligand HDV D 703



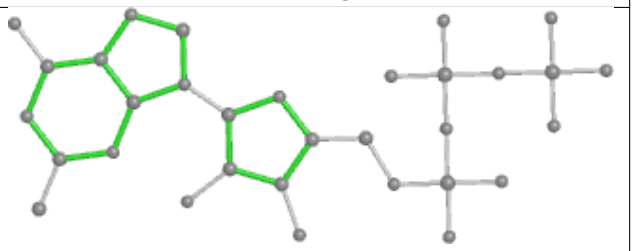
Bond lengths



Bond angles

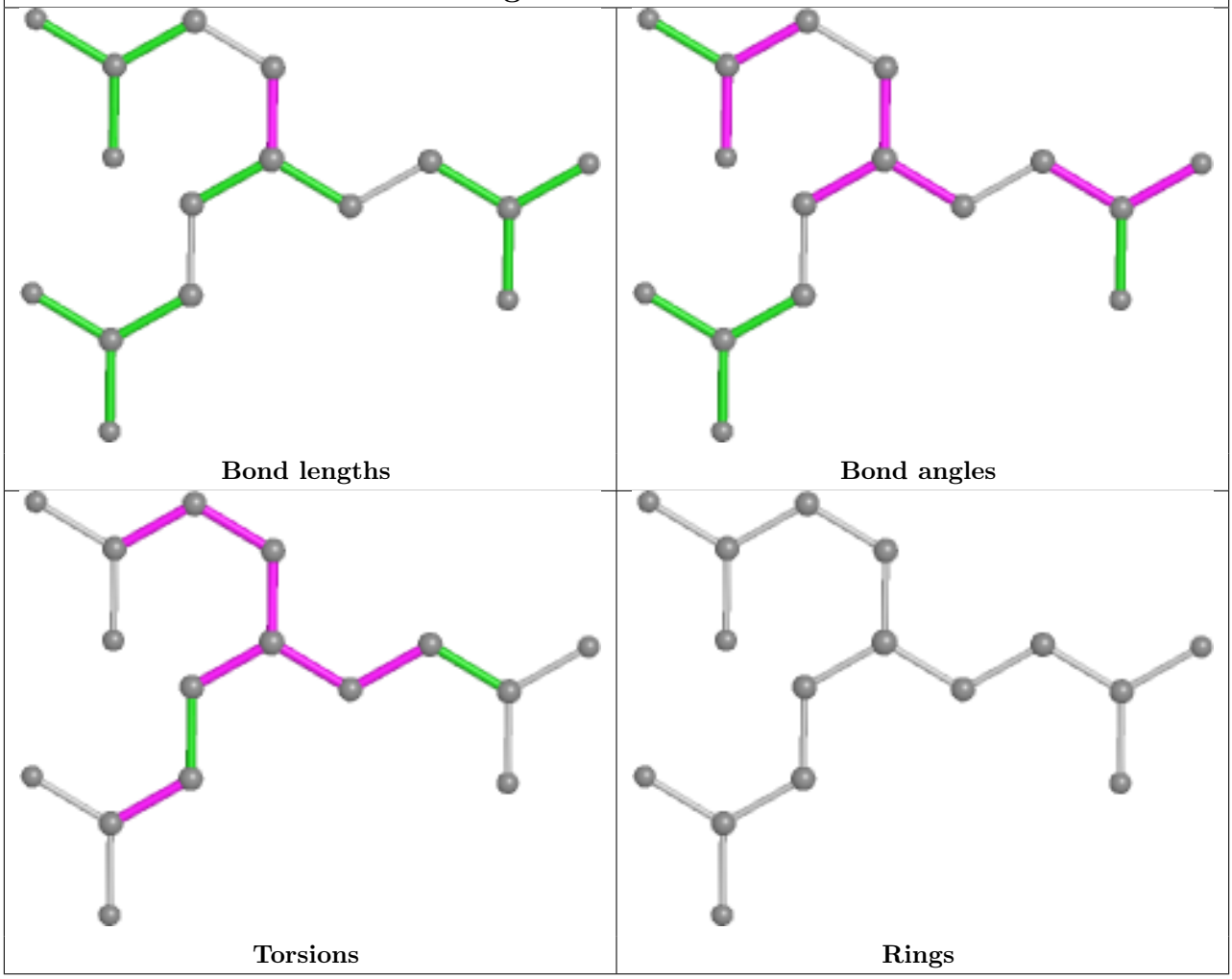


Torsions

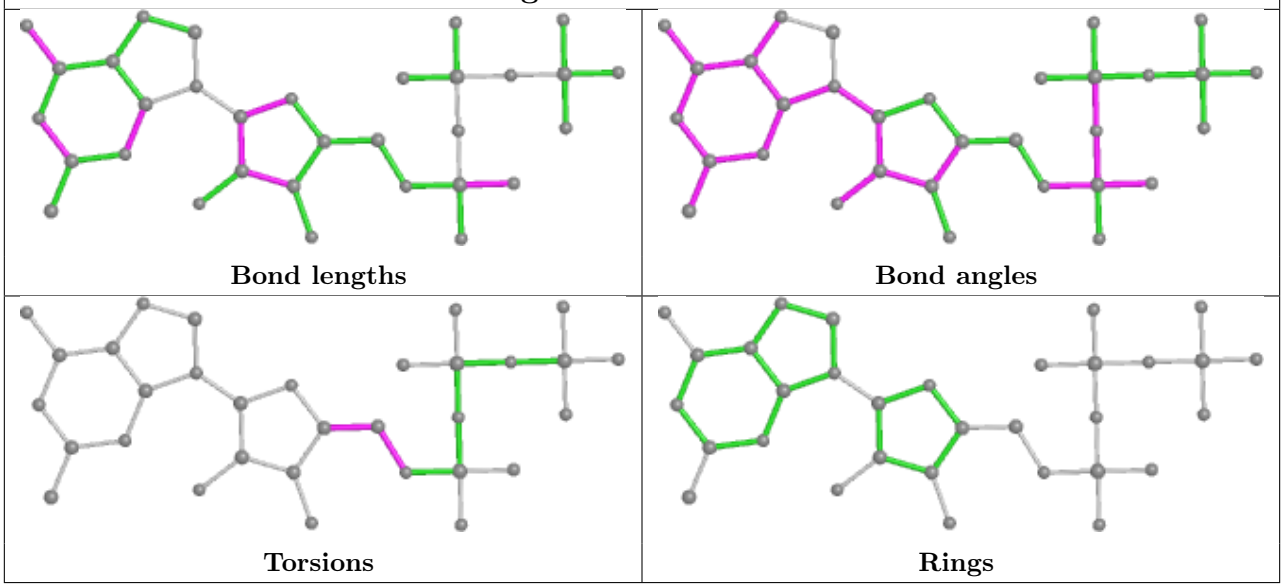


Rings

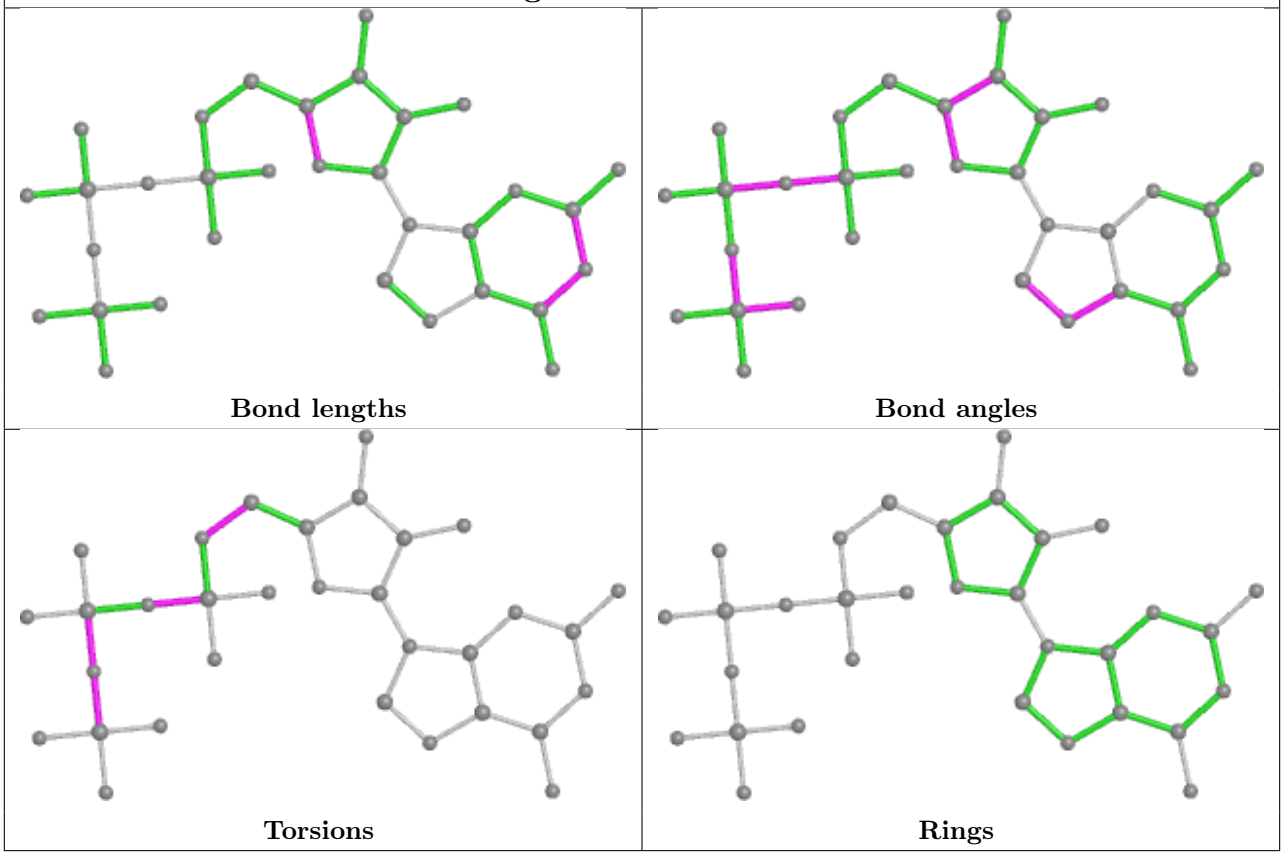
Ligand TCE A 718



Ligand HDV D 701



Ligand GTP B 703



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	466/550 (84%)	0.17	32 (6%)	16	19	10, 23, 62, 133	0
1	B	481/550 (87%)	0.22	35 (7%)	15	17	12, 27, 61, 98	0
1	C	481/550 (87%)	0.11	28 (5%)	23	25	12, 25, 62, 91	0
1	D	481/550 (87%)	-0.07	14 (2%)	51	56	9, 22, 52, 110	0
All	All	1909/2200 (86%)	0.11	109 (5%)	23	26	9, 24, 61, 133	0

The worst 5 of 109 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	LEU	16.8
1	C	487	VAL	9.8
1	C	488	LEU	9.5
1	A	489	LEU	8.9
1	A	490	ASP	7.9

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

There are no monosaccharides in this entry.

6.4 Ligands

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^{th} 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
6	GLY	C	713	5/5	0.45	0.22	50,53,61,66	0
6	GLY	A	715	5/5	0.66	0.24	37,40,52,62	0
6	GLY	D	713	5/5	0.69	0.17	47,48,52,52	0
9	TCE	A	718	16/16	0.70	0.26	50,64,95,97	0
6	GLY	C	712	5/5	0.74	0.18	40,42,48,58	0
6	GLY	A	716	5/5	0.75	0.19	47,51,62,64	0
7	SIN	B	710	8/8	0.76	0.18	43,48,54,56	0
5	NA	A	714	1/1	0.76	0.12	55,55,55,55	0
5	NA	C	710	1/1	0.83	0.17	48,48,48,48	0
5	NA	A	711	1/1	0.85	0.15	54,54,54,54	0
5	NA	D	712	1/1	0.85	0.15	49,49,49,49	0
5	NA	B	706	1/1	0.85	0.15	46,46,46,46	0
4	MG	C	707	1/1	0.86	0.12	49,49,49,49	0
5	NA	A	709	1/1	0.87	0.14	49,49,49,49	0
5	NA	D	709	1/1	0.87	0.29	48,48,48,48	0
7	SIN	D	714	8/8	0.88	0.18	44,48,55,59	0
5	NA	D	708	1/1	0.88	0.13	48,48,48,48	0
6	GLY	B	709	5/5	0.88	0.14	29,31,38,39	0
5	NA	A	712	1/1	0.89	0.10	48,48,48,48	0
5	NA	A	713	1/1	0.89	0.18	47,47,47,47	0
5	NA	A	707	1/1	0.90	0.09	39,39,39,39	0
5	NA	A	708	1/1	0.90	0.13	40,40,40,40	0
5	NA	D	711	1/1	0.91	0.10	47,47,47,47	0
5	NA	B	708	1/1	0.91	0.09	53,53,53,53	0
5	NA	A	710	1/1	0.92	0.27	49,49,49,49	0
6	GLY	A	717	5/5	0.93	0.19	33,33,41,45	0
5	NA	C	711	1/1	0.93	0.19	47,47,47,47	0
5	NA	D	707	1/1	0.93	0.23	52,52,52,52	0
8	PO4	C	714	5/5	0.93	0.16	61,64,68,73	0
5	NA	B	707	1/1	0.93	0.12	42,42,42,42	0
4	MG	D	706	1/1	0.94	0.11	45,45,45,45	0
5	NA	A	706	1/1	0.94	0.16	34,34,34,34	0
5	NA	C	709	1/1	0.94	0.06	34,34,34,34	0
5	NA	B	705	1/1	0.95	0.23	44,44,44,44	0
5	NA	C	708	1/1	0.95	0.13	38,38,38,38	0
5	NA	D	710	1/1	0.95	0.17	50,50,50,50	0
2	HDV	B	702	32/32	0.97	0.07	15,25,32,36	0
4	MG	A	705	1/1	0.98	0.04	18,18,18,18	0
2	HDV	A	704	32/32	0.98	0.06	11,20,27,30	0
2	HDV	C	701	32/32	0.98	0.07	14,22,29,32	0
2	HDV	D	701	32/32	0.98	0.07	11,18,26,29	0
4	MG	A	701	1/1	0.98	0.05	15,15,15,15	0
2	HDV	D	703	32/32	0.99	0.08	11,15,20,26	0

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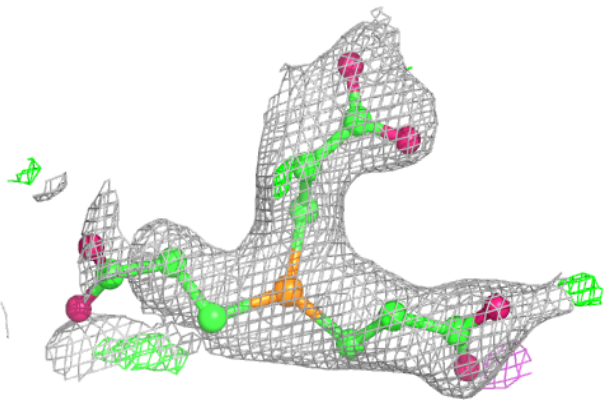
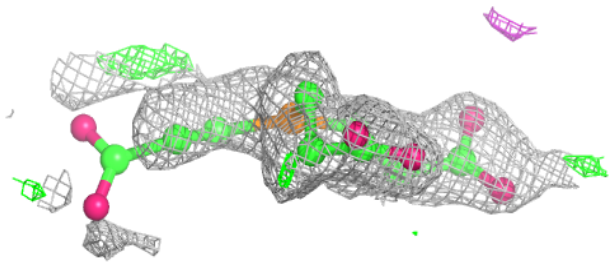
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	MG	B	704	1/1	0.99	0.06	24,24,24,24	0
2	HDV	A	702	32/32	0.99	0.06	13,16,19,25	0
4	MG	A	703	1/1	0.99	0.05	18,18,18,18	0
2	HDV	C	704	32/32	0.99	0.06	9,12,16,21	0
3	GTP	D	702	32/32	0.99	0.07	10,12,19,22	0
3	GTP	D	704	32/32	0.99	0.06	9,11,14,16	0
3	GTP	C	702	32/32	0.99	0.05	13,16,23,25	0
3	GTP	B	703	32/32	0.99	0.05	12,13,23,24	0
4	MG	D	705	1/1	0.99	0.03	17,17,17,17	0
2	HDV	B	701	32/32	0.99	0.07	10,14,18,26	0
4	MG	C	703	1/1	0.99	0.06	16,16,16,16	0
4	MG	C	706	1/1	0.99	0.02	22,22,22,22	0
4	MG	C	705	1/1	1.00	0.04	12,12,12,12	0

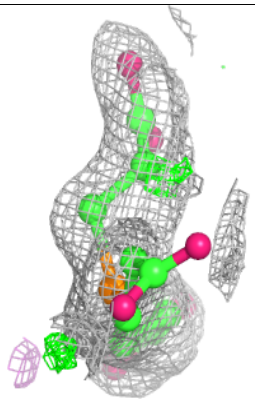
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around TCE A 718:

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

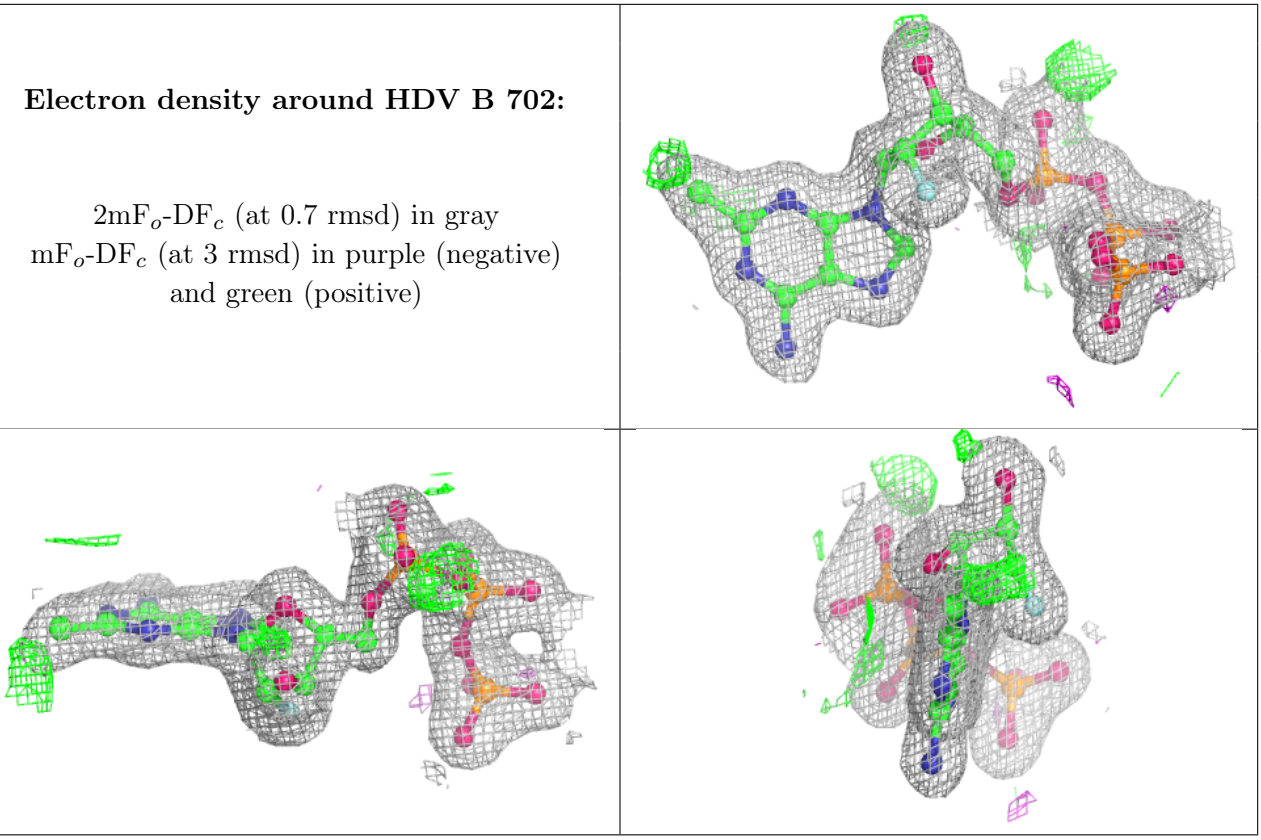






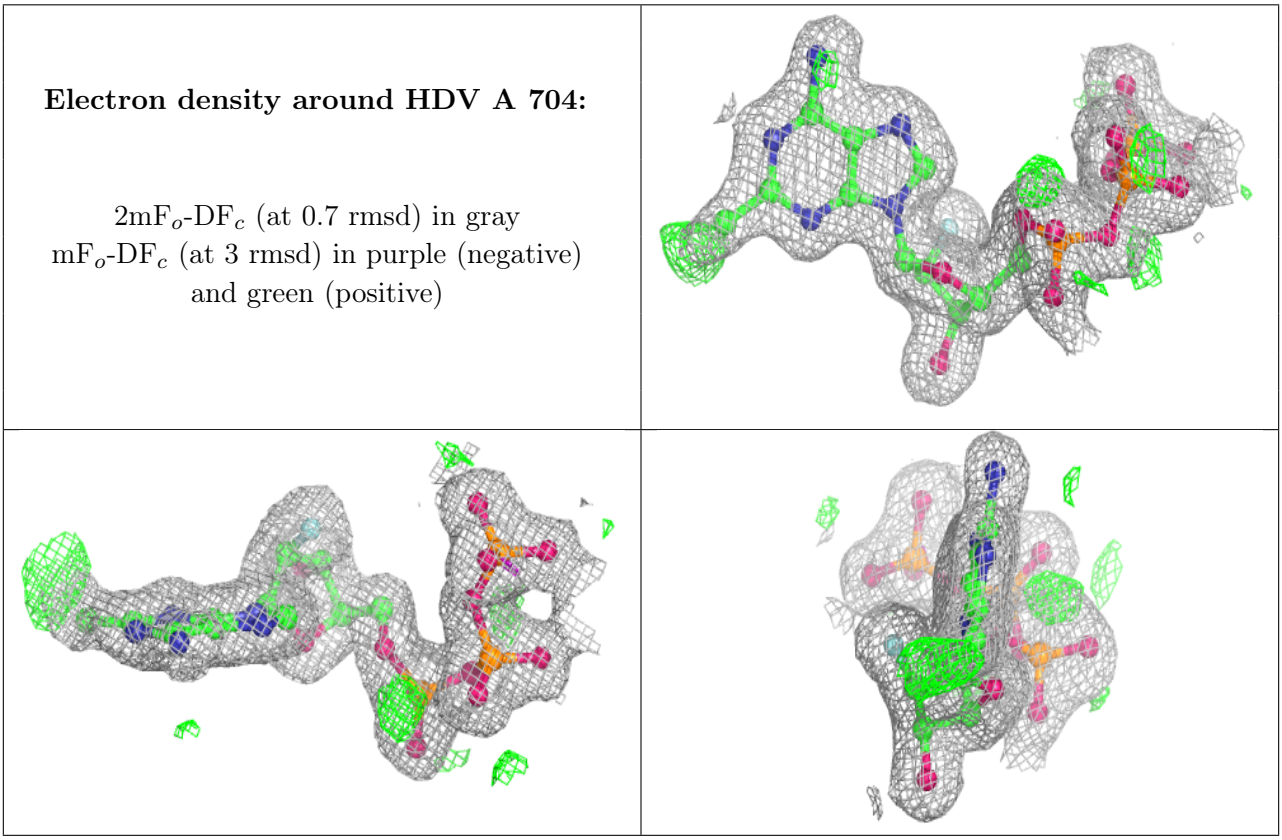
Electron density around HDV B 702:

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



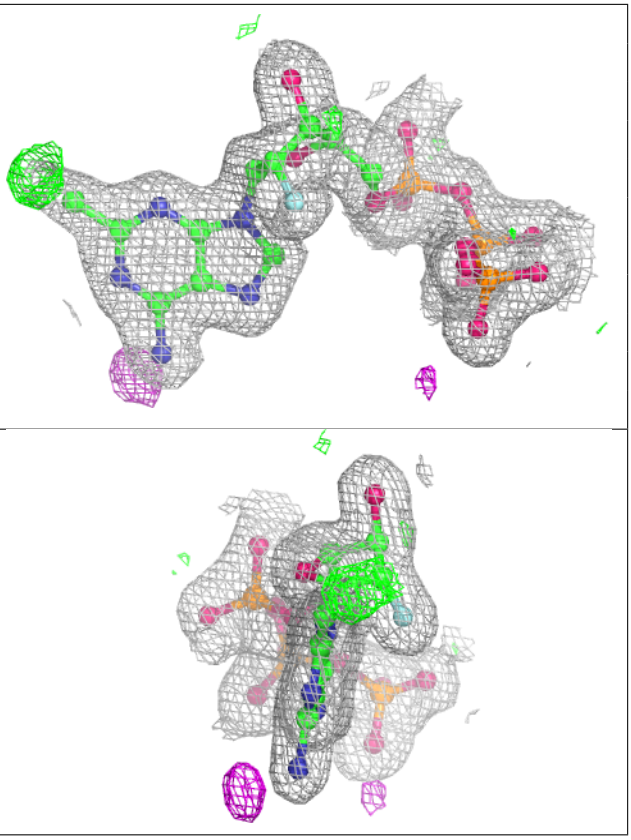
Electron density around HDV A 704:

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



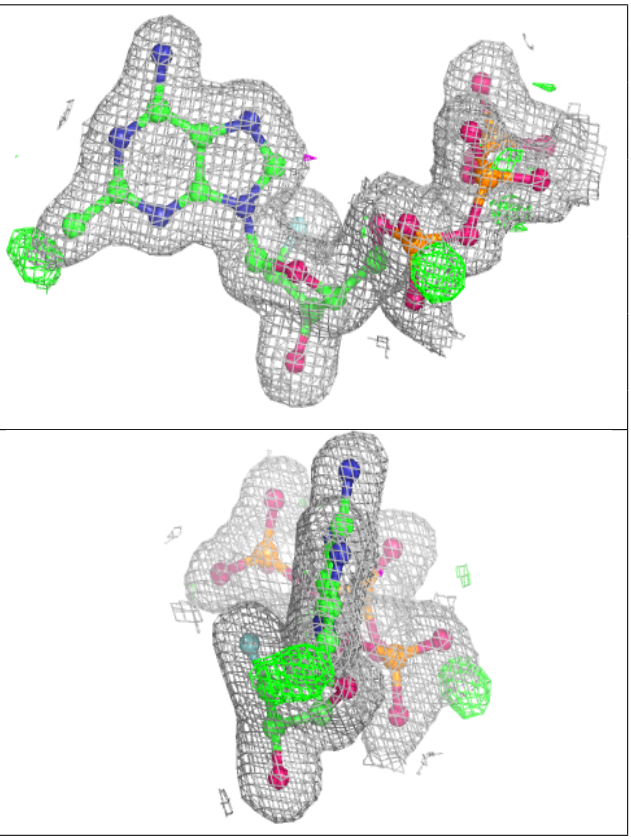
Electron density around HDV C 701:

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



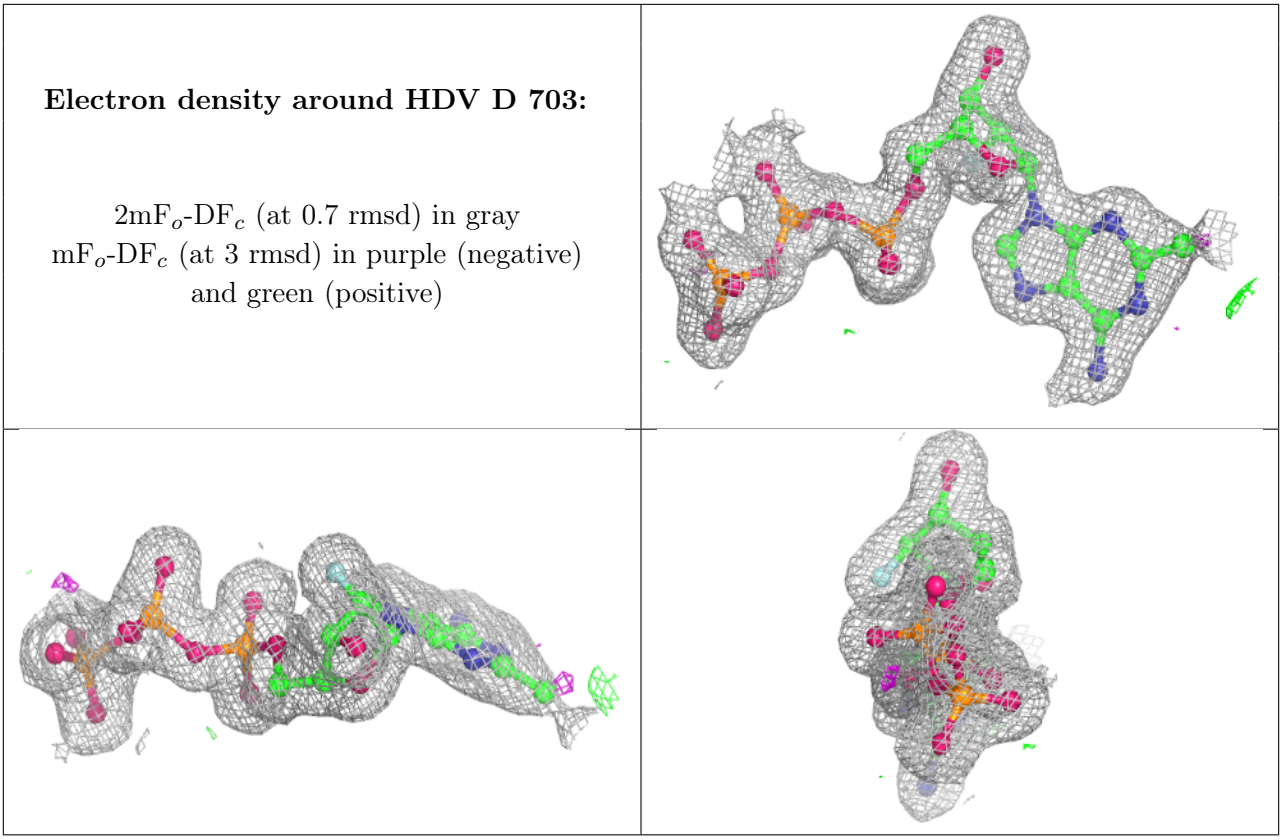
Electron density around HDV D 701:

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



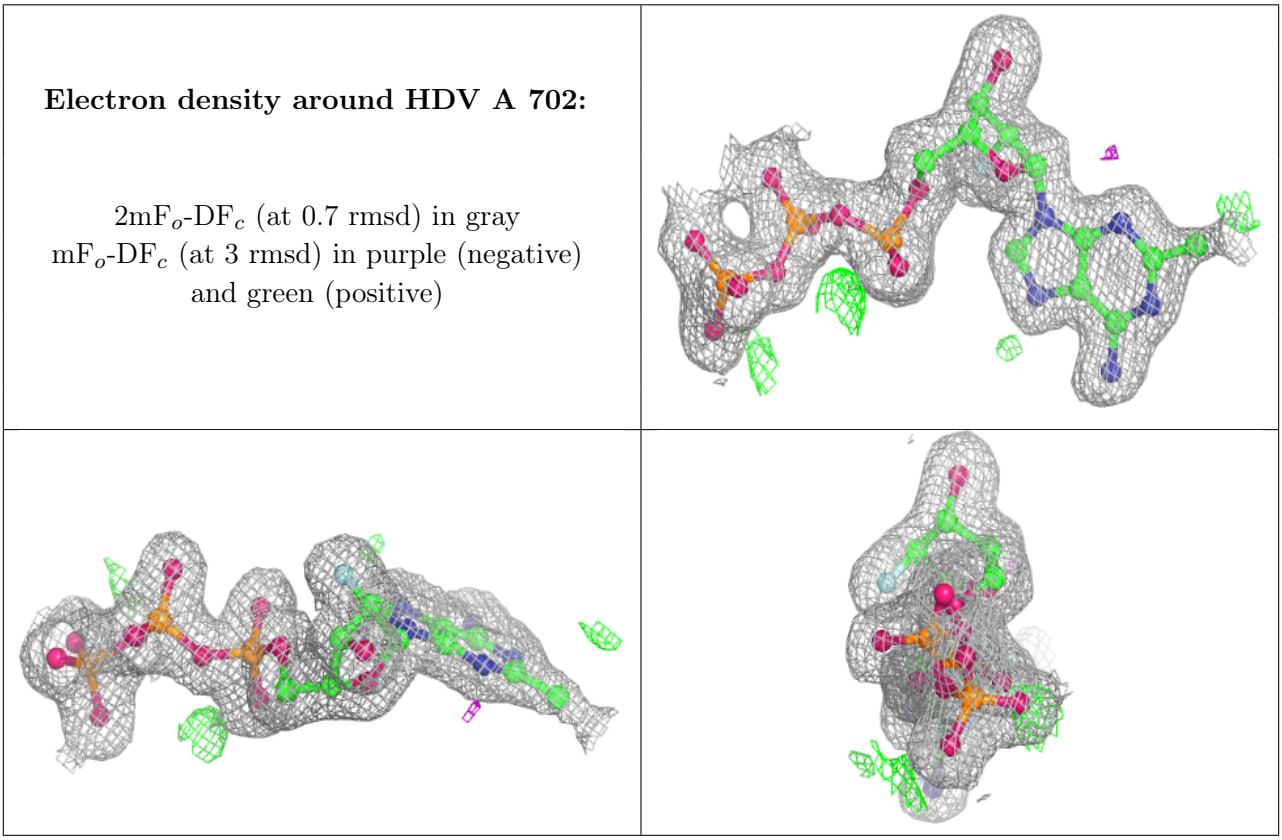
Electron density around HDV D 703:

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



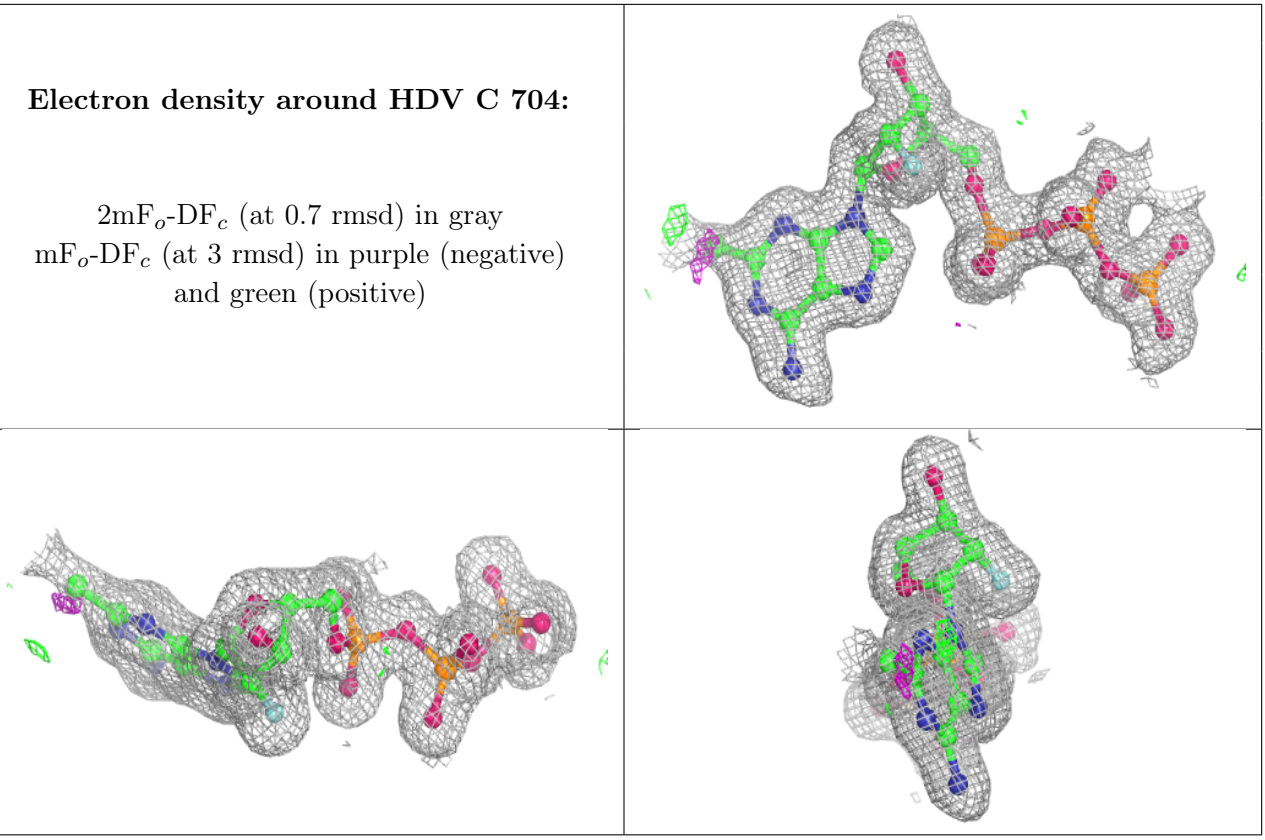
Electron density around HDV A 702:

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



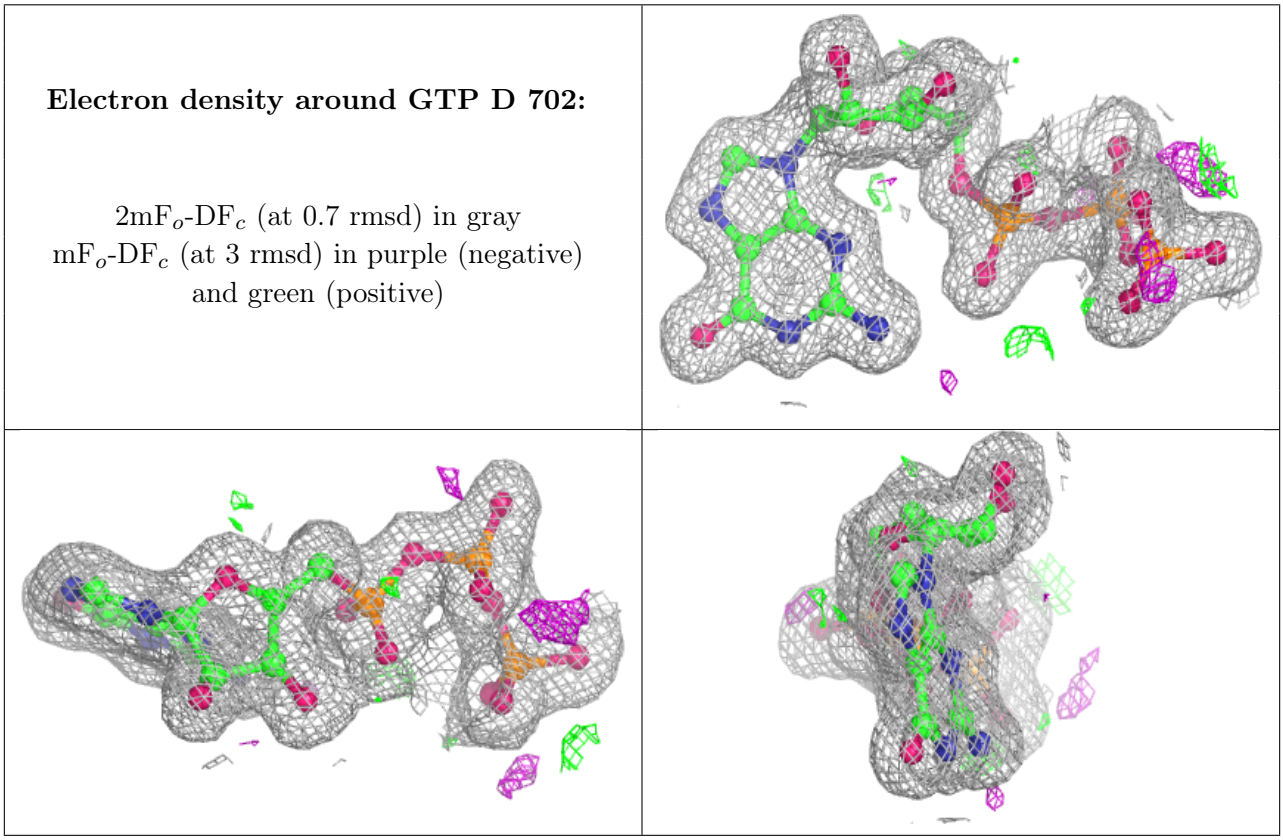
Electron density around HDV C 704:

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



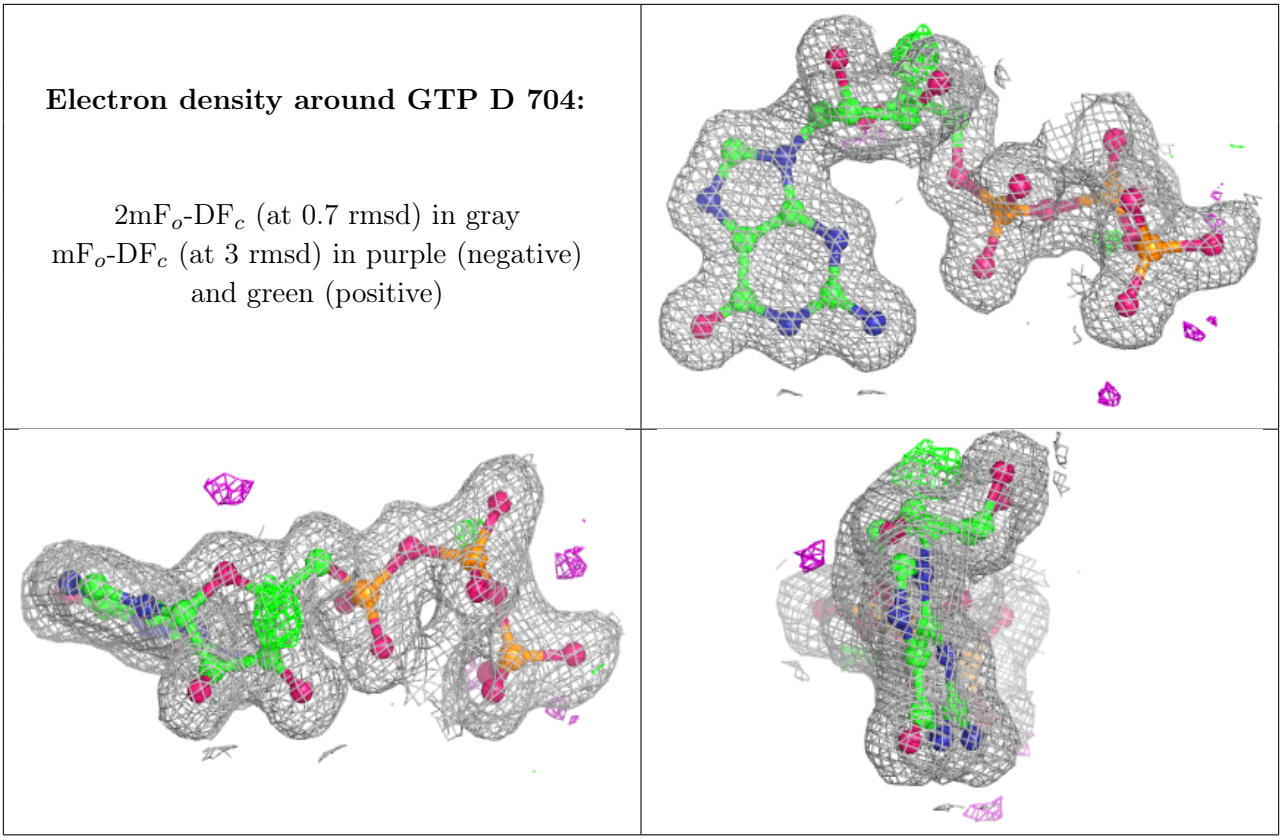
Electron density around GTP D 702:

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



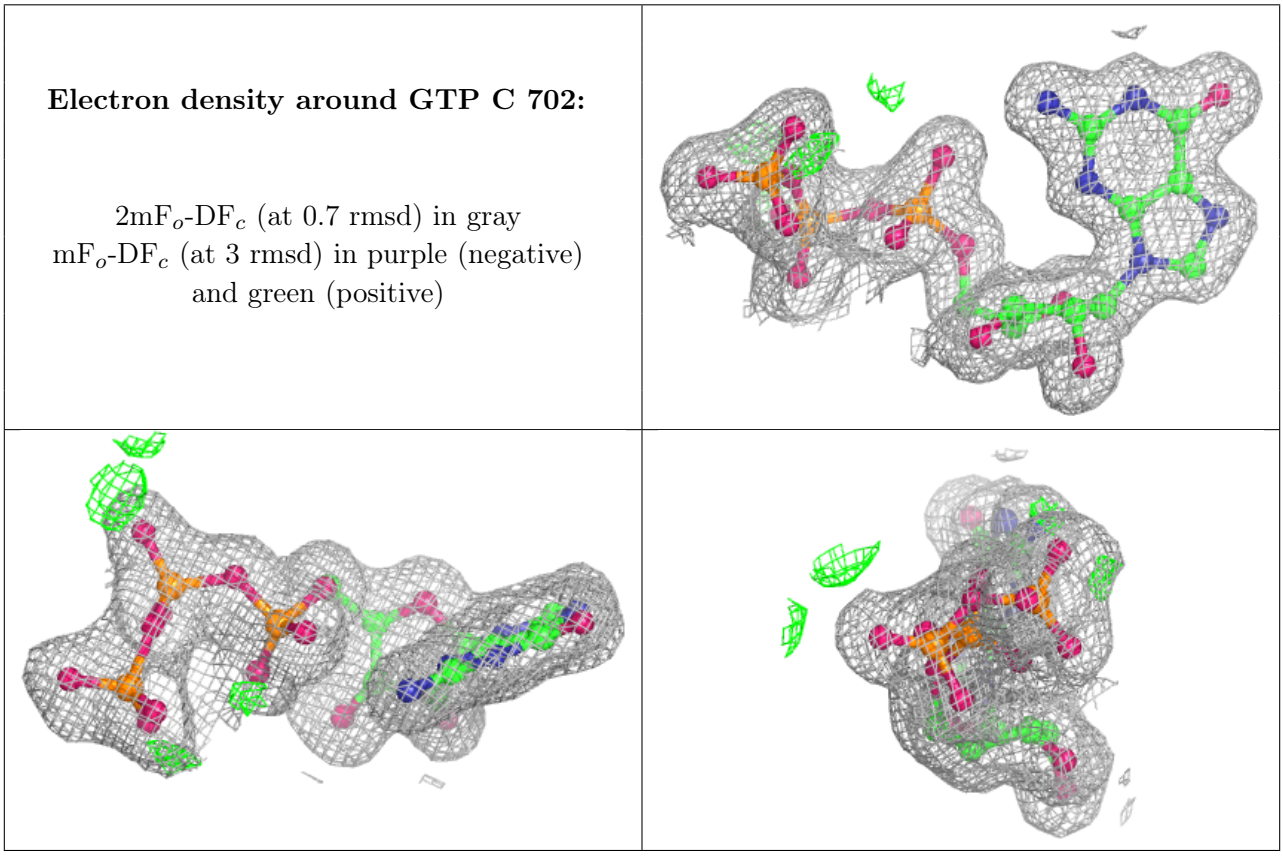
Electron density around GTP D 704:

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



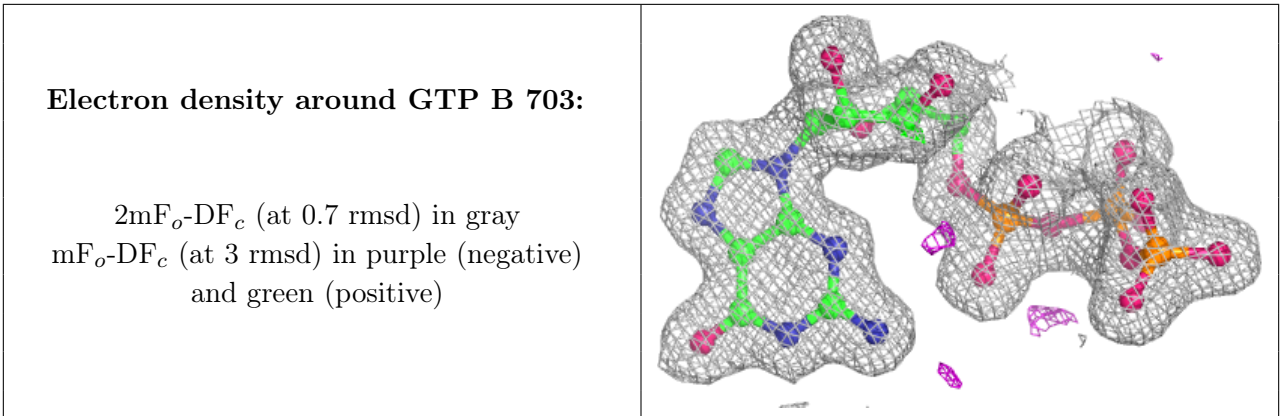
Electron density around GTP C 702:

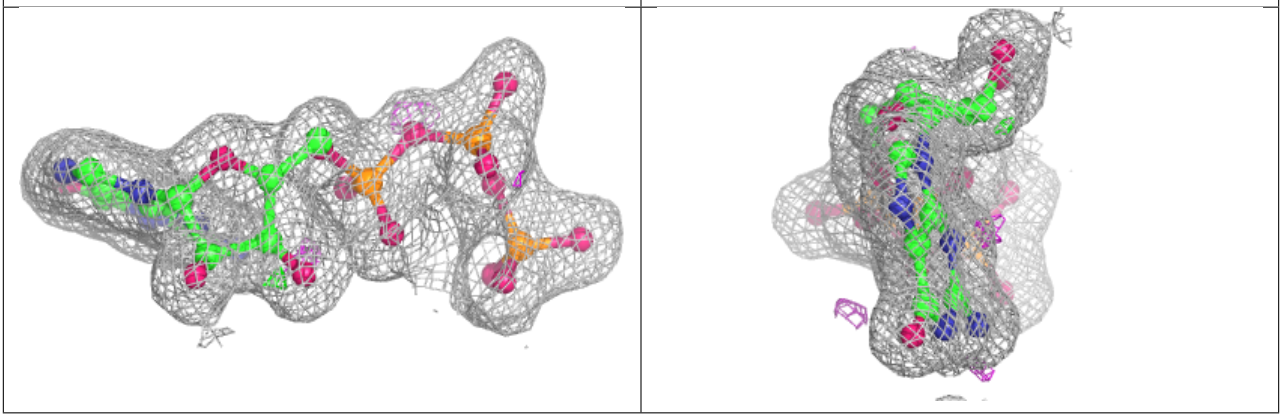
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



Electron density around GTP B 703:

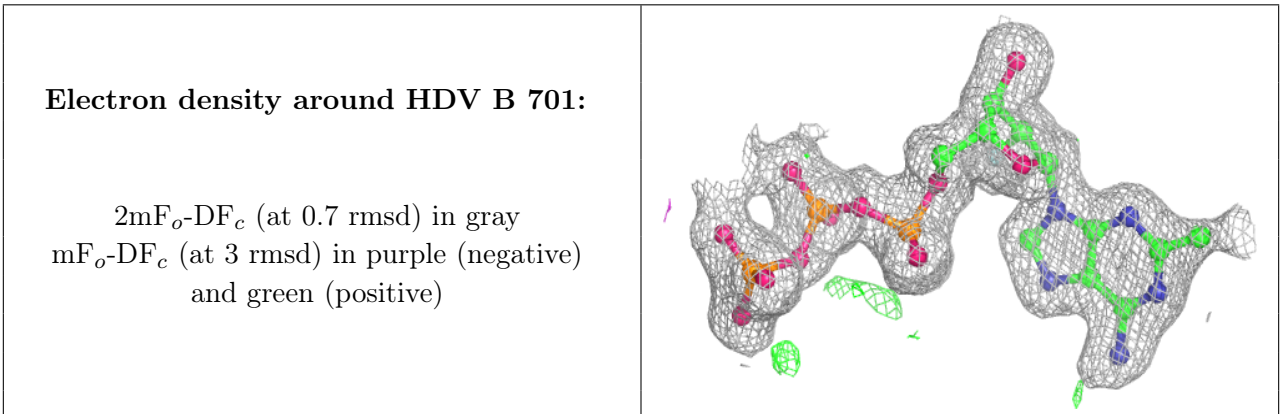
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

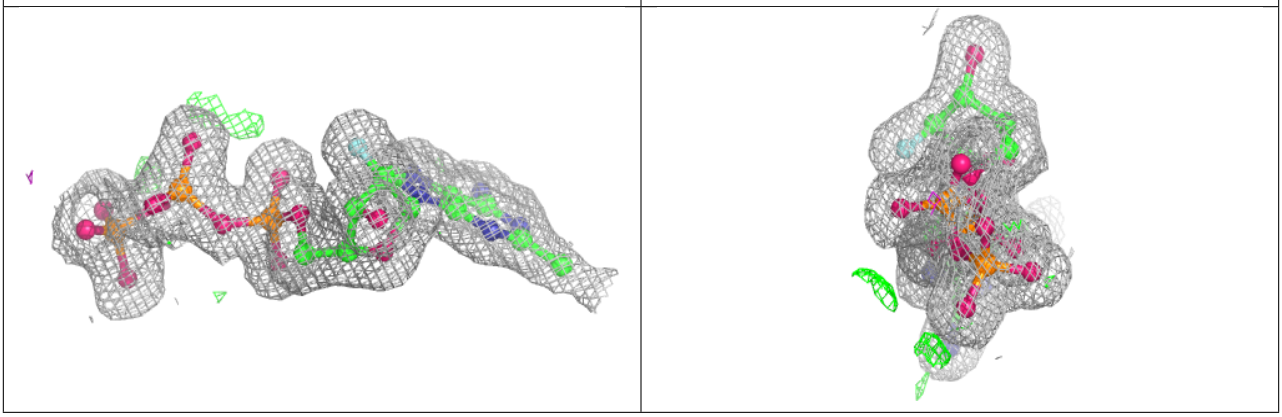




Electron density around HDV B 701:

$2mF_o$-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.