



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:28 AM UTC

PDB ID : 8F73 / pdb_00008f73
Title : Crystal structure of Pseudomonas aeruginosa UDP-glucose phosphorylase in complex with UDP-glucose
Authors : Dirr, L.; Fuehring, J.
Deposited on : 2022-11-17
Resolution : 2.90 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

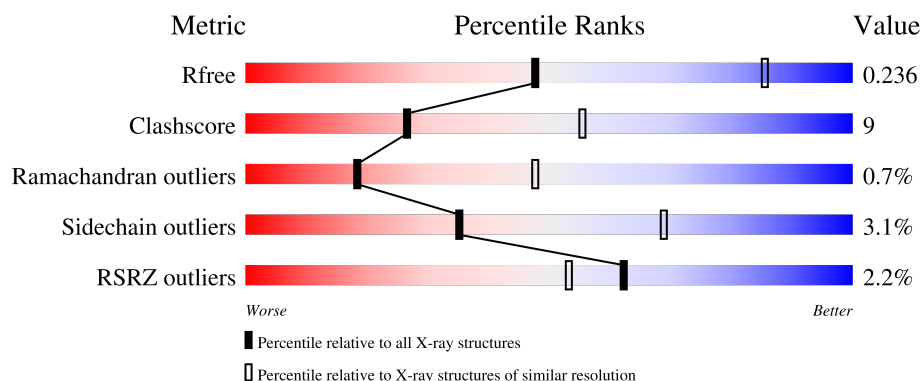
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	285	84% 12% ..
1	B	285	80% 16% ..
1	C	285	74% 21% ..
1	D	285	2% 74% 20% .. 5%
1	E	285	79% 19% ..

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Mol	Chain	Length	Quality of chain
1	F	285	% 80% 14%
1	G	285	6% 61% 31% 5%
1	H	285	4% 70% 25%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 35262 atoms, of which 17592 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 UTP--glucose-1-phosphate uridylyltransferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	275	Total	C	H	N	O	S	62	0	0
			4337	1377	2177	359	410	14			
1	B	280	Total	C	H	N	O	S	39	0	0
			4400	1397	2211	364	414	14			
1	C	274	Total	C	H	N	O	S	40	0	0
			4308	1370	2162	354	408	14			
1	D	272	Total	C	H	N	O	S	49	0	0
			4269	1360	2141	349	405	14			
1	E	281	Total	C	H	N	O	S	72	0	0
			4429	1404	2227	368	416	14			
1	F	275	Total	C	H	N	O	S	61	0	0
			4337	1377	2177	359	410	14			
1	G	271	Total	C	H	N	O	S	42	0	0
			4261	1356	2138	350	403	14			
1	H	274	Total	C	H	N	O	S	89	0	0
			4310	1371	2162	355	408	14			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	ALA	-	expression tag	UNP Q9I291
A	-4	LEU	-	expression tag	UNP Q9I291
A	-3	VAL	-	expression tag	UNP Q9I291
A	-2	PRO	-	expression tag	UNP Q9I291
A	-1	ARG	-	expression tag	UNP Q9I291
A	0	GLY	-	expression tag	UNP Q9I291
A	1	SER	-	expression tag	UNP Q9I291
B	-5	ALA	-	expression tag	UNP Q9I291
B	-4	LEU	-	expression tag	UNP Q9I291
B	-3	VAL	-	expression tag	UNP Q9I291
B	-2	PRO	-	expression tag	UNP Q9I291
B	-1	ARG	-	expression tag	UNP Q9I291
B	0	GLY	-	expression tag	UNP Q9I291

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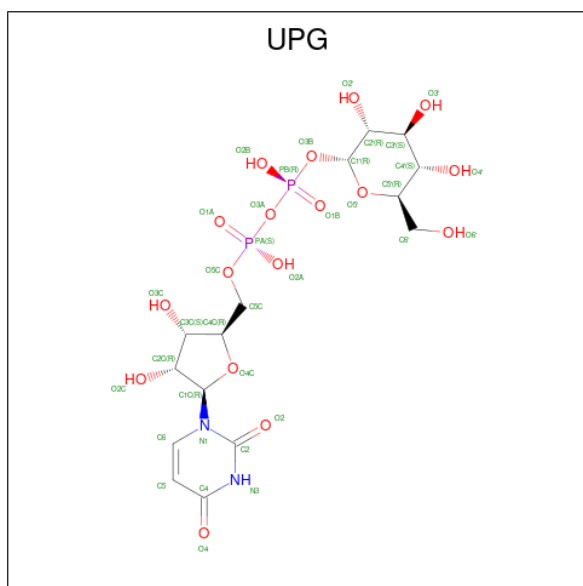
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	SER	-	expression tag	UNP Q9I291
C	-5	ALA	-	expression tag	UNP Q9I291
C	-4	LEU	-	expression tag	UNP Q9I291
C	-3	VAL	-	expression tag	UNP Q9I291
C	-2	PRO	-	expression tag	UNP Q9I291
C	-1	ARG	-	expression tag	UNP Q9I291
C	0	GLY	-	expression tag	UNP Q9I291
C	1	SER	-	expression tag	UNP Q9I291
D	-5	ALA	-	expression tag	UNP Q9I291
D	-4	LEU	-	expression tag	UNP Q9I291
D	-3	VAL	-	expression tag	UNP Q9I291
D	-2	PRO	-	expression tag	UNP Q9I291
D	-1	ARG	-	expression tag	UNP Q9I291
D	0	GLY	-	expression tag	UNP Q9I291
D	1	SER	-	expression tag	UNP Q9I291
E	-5	ALA	-	expression tag	UNP Q9I291
E	-4	LEU	-	expression tag	UNP Q9I291
E	-3	VAL	-	expression tag	UNP Q9I291
E	-2	PRO	-	expression tag	UNP Q9I291
E	-1	ARG	-	expression tag	UNP Q9I291
E	0	GLY	-	expression tag	UNP Q9I291
E	1	SER	-	expression tag	UNP Q9I291
F	-5	ALA	-	expression tag	UNP Q9I291
F	-4	LEU	-	expression tag	UNP Q9I291
F	-3	VAL	-	expression tag	UNP Q9I291
F	-2	PRO	-	expression tag	UNP Q9I291
F	-1	ARG	-	expression tag	UNP Q9I291
F	0	GLY	-	expression tag	UNP Q9I291
F	1	SER	-	expression tag	UNP Q9I291
G	-5	ALA	-	expression tag	UNP Q9I291
G	-4	LEU	-	expression tag	UNP Q9I291
G	-3	VAL	-	expression tag	UNP Q9I291
G	-2	PRO	-	expression tag	UNP Q9I291
G	-1	ARG	-	expression tag	UNP Q9I291
G	0	GLY	-	expression tag	UNP Q9I291
G	1	SER	-	expression tag	UNP Q9I291
H	-5	ALA	-	expression tag	UNP Q9I291
H	-4	LEU	-	expression tag	UNP Q9I291
H	-3	VAL	-	expression tag	UNP Q9I291
H	-2	PRO	-	expression tag	UNP Q9I291
H	-1	ARG	-	expression tag	UNP Q9I291
H	0	GLY	-	expression tag	UNP Q9I291

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Chain	Residue	Modelled	Actual	Comment	Reference
H	1	SER	-	expression tag	UNP Q9I291

- Molecule 2 is URIDINE-5'-DIPHOSPHATE-GLUCOSE (CCD ID: UPG) (formula: $C_{15}H_{24}N_2O_{17}P_2$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			10	2	6	2		
3	C	1	Total	C	H	O	0	0
			10	2	6	2		
3	D	1	Total	C	H	O	0	0
			10	2	6	2		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		
4	E	1	Total	Mg	0	0
			1	1		
4	F	1	Total	Mg	0	0
			1	1		
4	G	1	Total	Mg	0	0
			1	1		
4	H	1	Total	Mg	0	0
			1	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	H	O	0	0
			7	2	3	2		

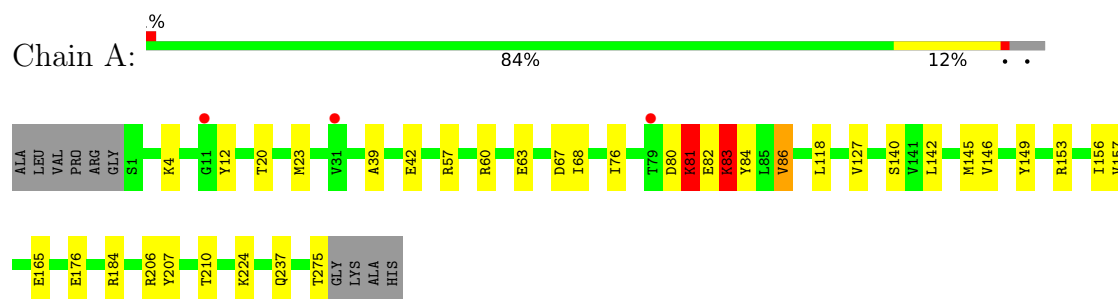
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	18	Total	O	0	0
			18	18		
6	B	25	Total	O	0	0
			25	25		
6	C	13	Total	O	0	0
			13	13		
6	D	8	Total	O	0	0
			8	8		
6	E	14	Total	O	0	0
			14	14		
6	F	10	Total	O	0	0
			10	10		
6	G	10	Total	O	0	0
			10	10		
6	H	4	Total	O	0	0
			4	4		

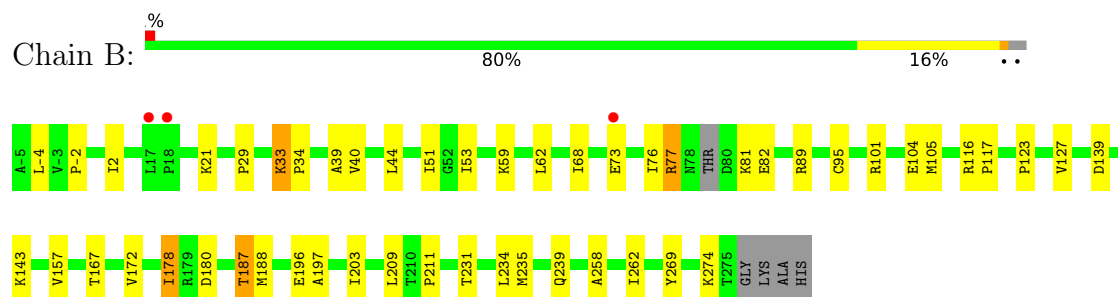
3 Residue-property plots

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.

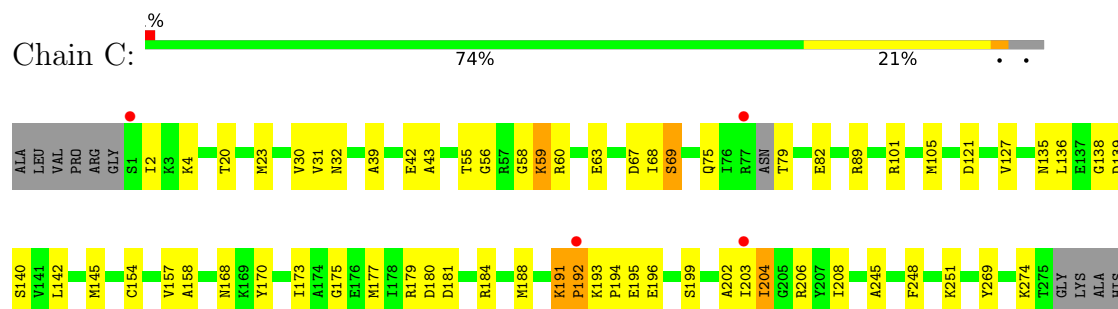
- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase



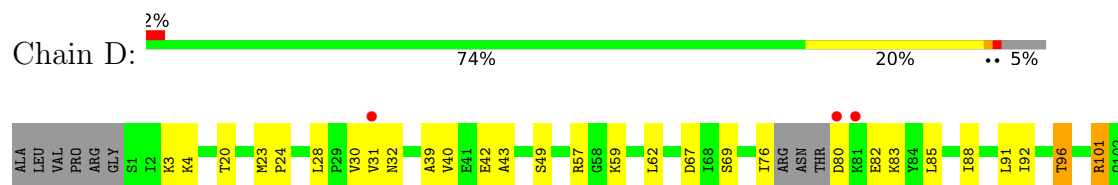
- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase

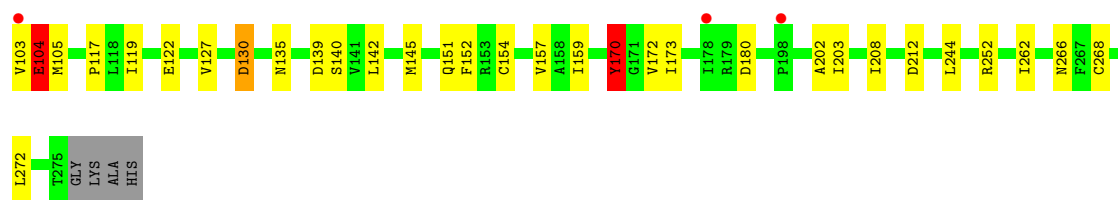


- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase

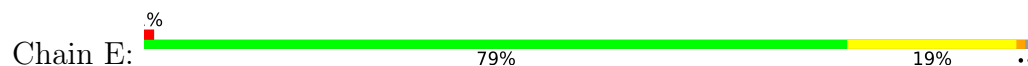


- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase

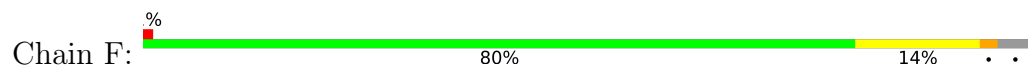




- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase



- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase

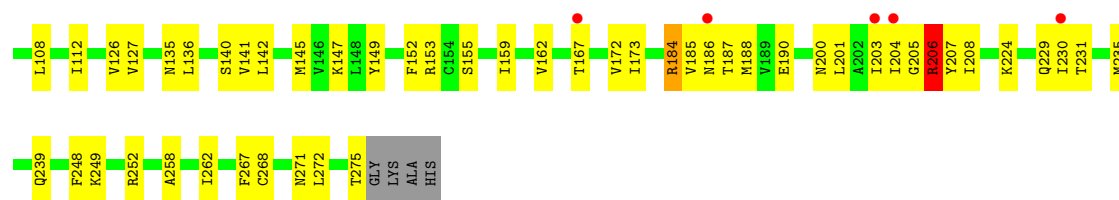


- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase



- Molecule 1: UTP--glucose-1-phosphate uridylyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.72Å 114.85Å 155.76Å 90.00° 97.39° 90.00°	Depositor
Resolution (Å)	46.98 – 2.90 46.98 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.98-2.90) 99.6 (46.98-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.195 , 0.236 0.195 , 0.236	Depositor DCC
R_{free} test set	3230 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	60.0	Xtriage
Anisotropy	0.770	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 49.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	35262	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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: UPG, EDO, MG, ACT

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	0.34	0/2199	0.59	5/2968 (0.2%)
1	B	0.42	3/2228 (0.1%)	0.54	1/3007 (0.0%)
1	C	0.65	3/2184 (0.1%)	0.78	13/2947 (0.4%)
1	D	0.31	0/2166	0.60	7/2923 (0.2%)
1	E	0.31	0/2242	0.52	1/3027 (0.0%)
1	F	0.57	3/2199 (0.1%)	0.82	11/2968 (0.4%)
1	G	0.49	3/2160 (0.1%)	0.74	9/2912 (0.3%)
1	H	0.31	0/2187	0.53	1/2953 (0.0%)
All	All	0.44	12/17565 (0.1%)	0.65	48/23705 (0.2%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	192	PRO	N-CA	20.40	1.69	1.46
1	F	192	PRO	N-CA	17.34	1.68	1.47
1	G	192	PRO	N-CA	15.85	1.67	1.47
1	C	191	LYS	C-N	12.22	1.48	1.33
1	G	193	LYS	C-N	8.65	1.54	1.33
1	F	191	LYS	C-N	8.63	1.44	1.33
1	C	193	LYS	C-N	8.49	1.53	1.33
1	F	193	LYS	C-N	8.21	1.53	1.33
1	G	191	LYS	C-N	6.50	1.49	1.33
1	B	34	PRO	C-O	-5.84	1.17	1.23
1	B	33	LYS	C-O	-5.28	1.18	1.25
1	B	29	PRO	C-O	-5.24	1.17	1.23

All (48) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	191	LYS	CA-C-N	16.41	136.66	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	191	LYS	C-N-CA	16.41	136.66	119.76
1	G	191	LYS	CA-C-N	12.44	135.40	119.84
1	G	191	LYS	C-N-CA	12.44	135.40	119.84
1	C	191	LYS	CA-C-N	12.24	132.31	120.31
1	C	191	LYS	C-N-CA	12.24	132.31	120.31
1	C	195	GLU	CA-C-O	-9.54	109.83	120.32
1	G	193	LYS	CA-C-N	9.06	131.16	119.84
1	G	193	LYS	C-N-CA	9.06	131.16	119.84
1	G	192	PRO	CA-N-CD	-8.47	100.15	112.00
1	A	80	ASP	N-CA-C	-8.18	103.33	112.57
1	C	193	LYS	CA-C-N	8.13	130.00	119.84
1	C	193	LYS	C-N-CA	8.13	130.00	119.84
1	C	195	GLU	O-C-N	8.01	132.72	123.27
1	D	130	ASP	CA-CB-CG	7.94	120.54	112.60
1	F	193	LYS	CA-C-N	7.57	129.31	119.84
1	F	193	LYS	C-N-CA	7.57	129.31	119.84
1	C	69	SER	N-CA-C	7.27	119.42	110.91
1	F	192	PRO	CA-N-CD	-7.08	102.09	112.00
1	D	104	GLU	N-CA-C	-6.96	95.98	110.80
1	C	192	PRO	CA-N-CD	-6.47	102.94	112.00
1	G	196	GLU	N-CA-C	-6.46	102.47	111.52
1	A	86	VAL	CA-C-N	6.43	127.12	119.98
1	A	86	VAL	C-N-CA	6.43	127.12	119.98
1	G	194	PRO	CA-C-N	6.43	133.83	121.54
1	G	194	PRO	C-N-CA	6.43	133.83	121.54
1	G	30	VAL	CA-C-O	-6.32	113.65	121.04
1	A	83	LYS	CB-CA-C	-6.31	109.28	116.54
1	C	193	LYS	N-CA-C	-6.17	100.43	109.50
1	D	170	TYR	CA-C-N	6.13	127.63	120.71
1	D	170	TYR	C-N-CA	6.13	127.63	120.71
1	F	80	ASP	CB-CA-C	6.09	123.55	116.54
1	F	178	ILE	N-CA-C	-6.09	106.87	112.96
1	D	170	TYR	CB-CA-C	-5.85	99.95	110.36
1	C	195	GLU	CA-C-N	5.78	132.58	121.54
1	C	195	GLU	C-N-CA	5.78	132.58	121.54
1	F	30	VAL	N-CA-CB	5.75	117.56	111.00
1	B	81	LYS	N-CA-C	-5.52	103.31	111.81
1	H	206	ARG	N-CA-C	-5.50	99.32	108.34
1	F	195	GLU	CA-C-N	5.41	131.87	121.54
1	F	195	GLU	C-N-CA	5.41	131.87	121.54
1	E	204	ILE	CB-CA-C	5.37	117.50	111.59
1	F	193	LYS	N-CA-C	-5.35	101.63	109.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	194	PRO	CA-C-N	5.26	130.41	123.05
1	C	194	PRO	C-N-CA	5.26	130.41	123.05
1	D	69	SER	N-CA-C	5.26	117.06	110.91
1	A	81	LYS	N-CA-C	-5.05	107.09	113.20
1	D	104	GLU	CB-CA-C	5.03	120.43	110.42

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	2160	2177	2179	19	1
1	B	2189	2211	2208	35	3
1	C	2146	2162	2161	42	1
1	D	2128	2141	2141	37	1
1	E	2202	2227	2227	41	2
1	F	2160	2177	2179	31	5
1	G	2123	2138	2136	72	2
1	H	2148	2162	2160	52	3
2	A	36	22	22	1	0
2	B	36	22	22	2	0
2	C	36	22	22	3	0
2	D	36	22	22	1	0
2	E	36	22	22	1	0
2	F	36	22	22	2	0
2	G	36	22	22	1	0
2	H	36	22	22	1	0
3	A	4	6	6	0	0
3	C	4	6	6	1	0
3	D	4	6	6	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	B	4	3	3	0	0
6	A	18	0	0	2	0
6	B	25	0	0	0	0
6	C	13	0	0	2	0
6	D	8	0	0	4	0
6	E	14	0	0	6	0
6	F	10	0	0	0	0
6	G	10	0	0	2	0
6	H	4	0	0	0	0
All	All	17670	17592	17588	324	9

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (324) 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:192:PRO:CA	1:C:192:PRO:N	1.69	1.43
1:F:192:PRO:N	1:F:192:PRO:CA	1.68	1.42
1:G:192:PRO:N	1:G:192:PRO:CA	1.67	1.34
2:E:301:UPG:O2A	6:E:401:HOH:O	1.59	1.20
1:E:130:ASP:OD2	6:E:401:HOH:O	1.62	1.16
1:B:73:GLU:HG2	1:B:89:ARG:HH22	1.24	0.99
1:G:176:GLU:OE1	1:G:186:ASN:ND2	2.04	0.91
1:D:130:ASP:OD1	6:D:401:HOH:O	1.89	0.89
2:C:301:UPG:H6	2:C:301:UPG:H5C2	1.55	0.88
1:D:130:ASP:CG	6:D:401:HOH:O	2.15	0.88
2:G:301:UPG:O2A	6:G:401:HOH:O	1.91	0.88
1:D:130:ASP:OD2	6:D:401:HOH:O	1.93	0.85
1:D:3:LYS:NZ	1:D:122:GLU:OE2	2.10	0.84
2:B:301:UPG:H6	2:B:301:UPG:H5C2	1.59	0.83
1:G:173:ILE:HG22	1:G:188:MET:HG3	1.60	0.83
1:H:204:ILE:HD13	1:H:248:PHE:CZ	2.11	0.83
1:G:171:GLY:O	6:G:402:HOH:O	1.96	0.82
1:C:68:ILE:HD11	1:G:103:VAL:HG13	1.59	0.82
1:E:76:ILE:HB	1:E:79:THR:HG21	1.60	0.81
2:D:301:UPG:O1B	6:D:401:HOH:O	1.99	0.80
1:G:159:ILE:HG21	1:G:173:ILE:HD11	1.61	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ILE:HD11	1:H:103:VAL:HG22	1.61	0.79
1:F:23:MET:HE3	1:F:28:LEU:HD23	1.62	0.79
2:A:301:UPG:O1B	6:A:401:HOH:O	2.02	0.78
1:C:127:VAL:HG13	1:C:206:ARG:NH1	1.99	0.76
1:H:190:GLU:OE2	2:H:301:UPG:O2'	2.03	0.76
2:F:301:UPG:H6	2:F:301:UPG:H5C2	1.69	0.74
1:E:76:ILE:HB	1:E:79:THR:CG2	2.17	0.74
1:B:73:GLU:O	1:B:77:ARG:HG3	1.88	0.74
1:B:73:GLU:HG2	1:B:89:ARG:NH2	2.02	0.72
1:F:173:ILE:HG22	1:F:188:MET:HG2	1.71	0.72
1:G:221:GLU:N	1:G:221:GLU:OE1	2.23	0.72
1:G:82:GLU:OE1	1:G:82:GLU:N	2.24	0.71
1:H:43:ALA:HA	1:H:142:LEU:HD23	1.73	0.71
1:E:188:MET:N	1:E:235:MET:SD	2.64	0.70
1:H:204:ILE:HD13	1:H:248:PHE:CE1	2.26	0.70
1:H:135:ASN:ND2	1:H:249:LYS:O	2.24	0.69
1:B:269:TYR:CE2	1:B:274:LYS:HE2	2.26	0.69
1:E:130:ASP:OD2	6:E:402:HOH:O	2.09	0.68
1:G:136:LEU:HD23	1:G:136:LEU:H	1.59	0.68
1:A:42:GLU:HG3	1:A:140:SER:HB2	1.75	0.67
1:B:104:GLU:N	1:B:104:GLU:OE1	2.29	0.66
1:D:135:ASN:ND2	1:D:139:ASP:O	2.29	0.66
1:A:81:LYS:O	1:A:81:LYS:HG2	1.96	0.66
1:B:40:VAL:HG13	1:B:51:ILE:HD13	1.78	0.64
1:C:157:VAL:HG22	1:C:203:ILE:HG23	1.80	0.64
1:G:144:GLN:O	1:G:148:LEU:HG	1.98	0.64
1:A:63:GLU:O	1:A:67:ASP:OD1	2.15	0.64
1:D:24:PRO:HG2	1:D:62:LEU:HD21	1.80	0.64
1:F:26:GLU:OE2	2:F:301:UPG:O2C	2.15	0.64
1:F:3:LYS:HE2	1:F:50:GLU:CD	2.22	0.64
1:F:15:ARG:NH1	1:F:256:GLY:O	2.31	0.63
1:G:192:PRO:N	1:G:192:PRO:C	2.54	0.63
1:B:269:TYR:CZ	1:B:274:LYS:HE2	2.34	0.62
1:E:130:ASP:CG	6:E:402:HOH:O	2.41	0.62
1:G:173:ILE:HD12	1:G:245:ALA:HB2	1.82	0.61
1:C:59:LYS:NZ	1:G:67:ASP:OD2	2.32	0.61
1:C:135:ASN:ND2	1:C:139:ASP:O	2.33	0.61
1:E:130:ASP:OD1	6:E:402:HOH:O	2.16	0.61
1:G:76:ILE:HD11	1:G:80:ASP:HB2	1.82	0.61
1:H:39:ALA:HB1	1:H:127:VAL:HG11	1.83	0.61
1:G:176:GLU:OE1	1:G:176:GLU:N	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:GLU:OE1	1:C:89:ARG:NH2	2.34	0.61
1:H:45:GLU:OE1	1:H:140:SER:OG	2.12	0.60
1:G:173:ILE:CD1	1:G:245:ALA:HB2	2.31	0.60
1:G:72:LEU:O	1:G:76:ILE:HG22	2.01	0.60
1:H:204:ILE:CD1	1:H:248:PHE:CE1	2.84	0.60
1:A:39:ALA:HB1	1:A:127:VAL:HG11	1.84	0.60
1:H:159:ILE:HG22	1:H:203:ILE:HA	1.84	0.59
1:F:192:PRO:HG2	1:F:197:ALA:HB2	1.84	0.59
1:H:36:ILE:HD13	1:H:62:LEU:HD11	1.83	0.59
1:F:192:PRO:N	1:F:192:PRO:C	2.56	0.59
1:H:24:PRO:HG3	1:H:62:LEU:HD23	1.85	0.59
1:C:251:LYS:NZ	6:C:402:HOH:O	2.33	0.59
1:C:67:ASP:OD2	1:G:59:LYS:NZ	2.32	0.59
1:A:145:MET:HE3	1:A:156:ILE:HG22	1.85	0.58
1:H:136:LEU:H	1:H:136:LEU:HD23	1.68	0.58
1:C:179:ARG:O	1:C:180:ASP:C	2.47	0.57
1:C:63:GLU:OE2	1:C:101:ARG:NH1	2.36	0.57
1:G:20:THR:HG23	1:G:57:ARG:HH22	1.69	0.57
1:D:39:ALA:HB1	1:D:127:VAL:HG11	1.86	0.57
1:D:20:THR:O	1:D:20:THR:HG22	2.03	0.57
1:D:180:ASP:OD1	1:D:180:ASP:N	2.35	0.57
1:G:42:GLU:OE1	1:G:206:ARG:NH2	2.35	0.57
1:A:153:ARG:NH2	1:C:121:ASP:OD1	2.37	0.57
1:F:3:LYS:HE2	1:F:50:GLU:OE2	2.05	0.57
1:E:76:ILE:HD12	1:E:76:ILE:O	2.05	0.56
1:F:39:ALA:HB1	1:F:127:VAL:HG11	1.86	0.56
1:H:76:ILE:HG23	1:H:76:ILE:O	2.04	0.56
1:B:187:THR:HA	1:B:235:MET:HE1	1.86	0.56
1:C:157:VAL:HG22	1:C:203:ILE:CG2	2.36	0.56
1:B:167:THR:HG23	1:B:172:VAL:HG21	1.87	0.56
1:E:179:ARG:HG2	1:E:180:ASP:OD1	2.06	0.55
1:F:173:ILE:HG22	1:F:188:MET:CG	2.35	0.55
1:B:39:ALA:HB1	1:B:127:VAL:HG11	1.89	0.55
1:H:267:PHE:O	1:H:271:ASN:ND2	2.36	0.55
1:B:68:ILE:HG23	1:B:68:ILE:O	2.06	0.55
1:G:229:GLN:OE1	1:G:229:GLN:N	2.39	0.55
1:C:55:THR:HB	1:C:59:LYS:HB2	1.89	0.54
1:C:180:ASP:OD1	1:C:181:ASP:N	2.39	0.54
1:C:39:ALA:HB1	1:C:127:VAL:HG11	1.89	0.54
1:A:68:ILE:H	1:A:68:ILE:HD12	1.72	0.54
1:G:145:MET:HA	1:G:148:LEU:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:269:TYR:CZ	1:G:274:LYS:HE3	2.42	0.54
1:C:173:ILE:HG22	1:C:188:MET:HG2	1.90	0.54
1:B:105:MET:HG3	2:B:301:UPG:O4	2.08	0.53
1:C:79:THR:HA	1:C:82:GLU:HG3	1.90	0.53
1:D:23:MET:HE3	1:D:28:LEU:HD23	1.89	0.53
1:G:39:ALA:HB1	1:G:127:VAL:HG11	1.91	0.53
1:A:157:VAL:HG23	1:A:207:TYR:CD2	2.43	0.53
1:D:43:ALA:HA	1:D:142:LEU:HD23	1.90	0.53
1:C:192:PRO:N	1:C:192:PRO:C	2.61	0.53
1:C:42:GLU:HG3	1:C:140:SER:HB2	1.91	0.53
1:A:12:TYR:O	1:A:57:ARG:NH2	2.42	0.52
1:F:42:GLU:HG3	1:F:140:SER:HB2	1.90	0.52
1:E:20:THR:HG23	1:E:20:THR:O	2.09	0.52
1:E:159:ILE:HD13	1:E:245:ALA:HB1	1.91	0.52
1:H:167:THR:HG23	1:H:172:VAL:HG21	1.92	0.52
1:G:108:LEU:HD11	1:G:128:LEU:HD11	1.92	0.52
1:C:136:LEU:HD12	1:C:136:LEU:H	1.74	0.52
1:D:4:LYS:NZ	1:D:119:ILE:O	2.40	0.52
1:D:42:GLU:HG3	1:D:140:SER:HB2	1.91	0.52
1:H:162:VAL:O	1:H:200:ASN:OD1	2.28	0.52
1:F:178:ILE:N	1:F:178:ILE:HD12	2.25	0.52
1:G:174:ALA:HB3	1:G:187:THR:OG1	2.10	0.52
1:E:219:GLN:N	1:E:219:GLN:OE1	2.43	0.51
1:G:175:GLY:HA2	1:G:186:ASN:ND2	2.25	0.51
1:A:210:THR:OG1	1:A:237:GLN:NE2	2.44	0.51
1:G:159:ILE:CG2	1:G:173:ILE:HD11	2.37	0.51
1:C:173:ILE:O	1:C:199:SER:OG	2.20	0.51
1:E:252:ARG:NH1	1:E:254:ASP:HB2	2.26	0.51
1:G:212:ASP:OD1	1:G:237:GLN:NE2	2.44	0.51
1:E:103:VAL:HG13	1:E:104:GLU:HG2	1.93	0.51
1:B:188:MET:HE3	1:B:231:THR:HG23	1.94	0.50
1:D:23:MET:CE	1:D:28:LEU:HD23	2.42	0.50
1:A:20:THR:O	1:A:57:ARG:NH1	2.45	0.50
1:G:145:MET:HE3	1:G:156:ILE:HG22	1.93	0.50
1:G:199:SER:O	1:G:199:SER:OG	2.23	0.50
1:H:275:THR:O	1:H:275:THR:OG1	2.30	0.50
3:D:302:EDO:H12	1:E:101:ARG:HB2	1.94	0.50
1:D:96:THR:HG21	1:E:117:PRO:HG3	1.94	0.49
1:F:43:ALA:HA	1:F:142:LEU:HD23	1.93	0.49
1:G:173:ILE:CD1	1:G:245:ALA:CB	2.89	0.49
1:D:173:ILE:HD12	1:D:173:ILE:C	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:185:VAL:HG11	1:H:188:MET:HE2	1.94	0.49
1:B:167:THR:CG2	1:B:172:VAL:HG21	2.43	0.49
1:G:42:GLU:HG2	1:G:140:SER:HB2	1.93	0.49
1:B:180:ASP:OD1	1:B:180:ASP:N	2.45	0.49
1:F:2:ILE:O	1:F:2:ILE:HG23	2.11	0.49
1:C:43:ALA:HA	1:C:142:LEU:HD23	1.93	0.49
1:H:200:ASN:OD1	1:H:200:ASN:O	2.31	0.49
1:E:-2:PRO:HD2	1:E:211:PRO:HG2	1.95	0.49
1:G:173:ILE:HG23	1:G:203:ILE:HD11	1.95	0.49
1:G:179:ARG:HE	1:H:184:ARG:HD3	1.78	0.49
1:H:24:PRO:HG3	1:H:62:LEU:CD2	2.42	0.49
1:E:159:ILE:HD12	1:E:201:LEU:HD22	1.94	0.48
1:G:130:ASP:OD1	1:G:130:ASP:N	2.46	0.48
1:G:43:ALA:HA	1:G:142:LEU:HD23	1.94	0.48
1:G:204:ILE:HG22	1:G:205:GLY:N	2.28	0.48
1:D:170:TYR:CD1	1:D:170:TYR:N	2.82	0.48
1:D:24:PRO:HG2	1:D:62:LEU:CD2	2.43	0.48
1:G:20:THR:HG23	1:G:57:ARG:NH2	2.27	0.48
1:C:68:ILE:CD1	1:G:103:VAL:HG13	2.39	0.48
1:G:173:ILE:CG2	1:G:203:ILE:HD11	2.43	0.48
1:H:147:LYS:HA	1:H:147:LYS:HE2	1.95	0.48
1:H:204:ILE:HG21	1:H:252:ARG:HD2	1.96	0.48
1:G:24:PRO:HG2	1:G:62:LEU:HG	1.96	0.48
1:B:-4:LEU:HD22	1:E:178:ILE:HG21	1.95	0.47
1:B:40:VAL:HG13	1:B:51:ILE:CD1	2.43	0.47
1:B:2:ILE:HG12	1:B:123:PRO:HB2	1.96	0.47
1:H:145:MET:SD	1:H:206:ARG:HB3	2.54	0.47
1:D:117:PRO:HG2	1:E:96:THR:HG21	1.96	0.47
1:E:180:ASP:OD1	1:E:180:ASP:N	2.48	0.47
1:G:222:PRO:HA	1:G:227:GLU:O	2.14	0.47
1:C:20:THR:HG22	1:C:20:THR:O	2.13	0.47
1:F:23:MET:CE	1:F:28:LEU:HD23	2.37	0.47
1:G:40:VAL:HG12	1:G:91:LEU:HD22	1.97	0.47
1:G:221:GLU:H	1:G:221:GLU:CD	2.21	0.47
1:A:176:GLU:HG3	1:A:184:ARG:HB3	1.96	0.47
1:E:42:GLU:HG3	1:E:140:SER:HB2	1.97	0.47
1:G:197:ALA:O	1:G:199:SER:N	2.48	0.47
1:C:269:TYR:O	1:C:274:LYS:HG3	2.15	0.47
1:E:179:ARG:HE	1:E:182:ILE:HD12	1.80	0.47
1:B:178:ILE:HD13	1:B:178:ILE:N	2.31	0.46
1:G:106:LYS:O	1:G:228:ILE:N	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:204:ILE:CD1	1:H:248:PHE:CZ	2.90	0.46
1:C:168:ASN:O	1:C:191:LYS:HG2	2.15	0.46
1:F:212:ASP:OD1	1:F:237:GLN:NE2	2.49	0.46
1:E:68:ILE:CD1	1:E:92:ILE:HG21	2.45	0.46
1:D:59:LYS:HD2	1:D:101:ARG:NH2	2.29	0.46
1:H:4:LYS:NZ	1:H:50:GLU:OE1	2.48	0.46
1:B:-2:PRO:HD2	1:B:211:PRO:HG2	1.98	0.46
1:D:145:MET:HE2	1:D:208:ILE:HG12	1.98	0.46
1:G:159:ILE:CG2	1:G:245:ALA:HB1	2.46	0.46
1:B:139:ASP:HB3	1:B:143:LYS:HD2	1.97	0.46
1:F:3:LYS:CE	1:F:50:GLU:OE2	2.63	0.46
1:G:172:VAL:HG12	1:G:173:ILE:N	2.30	0.46
1:F:167:THR:HG21	1:F:194:PRO:HA	1.98	0.46
1:F:57:ARG:HH11	1:F:57:ARG:HG3	1.81	0.46
1:D:88:ILE:O	1:D:92:ILE:HD12	2.16	0.45
1:H:230:ILE:HG23	1:H:231:THR:N	2.29	0.45
1:B:258:ALA:O	1:B:262:ILE:HG12	2.16	0.45
1:C:30:VAL:O	1:C:32:ASN:N	2.48	0.45
1:C:138:GLY:O	6:C:401:HOH:O	2.20	0.45
1:D:262:ILE:O	1:D:266:ASN:ND2	2.42	0.45
1:D:62:LEU:H	1:D:62:LEU:HD22	1.82	0.45
1:D:76:ILE:O	1:D:80:ASP:N	2.49	0.45
1:C:20:THR:HG23	1:C:23:MET:O	2.17	0.45
1:D:101:ARG:CD	1:D:101:ARG:N	2.80	0.45
1:E:116:ARG:NH1	6:E:403:HOH:O	2.49	0.45
1:H:112:ILE:HD12	1:H:230:ILE:HD12	1.98	0.45
1:H:206:ARG:HD3	1:H:206:ARG:HA	1.80	0.45
1:E:201:LEU:HD21	1:E:247:LYS:HE2	1.98	0.45
1:F:2:ILE:HD13	1:F:48:LEU:CD2	2.47	0.45
1:F:72:LEU:HD21	1:F:85:LEU:HD13	1.98	0.45
1:G:141:VAL:O	1:G:145:MET:HG3	2.17	0.45
1:G:13:GLY:O	1:G:20:THR:HG21	2.17	0.44
1:G:188:MET:HG2	1:G:231:THR:HG22	1.98	0.44
1:C:175:GLY:HA3	1:C:184:ARG:O	2.17	0.44
1:B:235:MET:CE	1:B:239:GLN:OE1	2.65	0.44
1:C:4:LYS:NZ	3:C:302:EDO:H21	2.32	0.44
1:D:154:CYS:SG	1:D:244:LEU:HG	2.58	0.44
1:F:2:ILE:HD13	1:F:48:LEU:HD22	1.99	0.44
1:F:178:ILE:HD11	1:F:184:ARG:HB2	1.99	0.44
1:G:113:LEU:O	1:G:113:LEU:HD23	2.17	0.44
1:H:186:ASN:OD1	1:H:187:THR:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:80:ASP:C	1:E:82:GLU:H	2.25	0.44
1:E:212:ASP:OD1	1:E:212:ASP:N	2.51	0.44
1:G:172:VAL:HG11	1:G:198:PRO:O	2.17	0.44
1:G:179:ARG:HH21	1:H:184:ARG:HD2	1.82	0.44
1:D:252:ARG:HD3	1:D:252:ARG:C	2.43	0.44
1:F:85:LEU:O	1:F:89:ARG:HG2	2.18	0.44
1:C:170:TYR:CG	1:C:202:ALA:HB1	2.53	0.44
1:C:170:TYR:CB	1:C:202:ALA:HB1	2.48	0.44
1:A:224:LYS:HE3	1:B:196:GLU:HA	2.00	0.44
1:D:103:VAL:O	1:D:104:GLU:HB3	2.18	0.44
1:D:159:ILE:HA	1:D:202:ALA:O	2.17	0.44
1:E:200:ASN:OD1	1:E:200:ASN:O	2.35	0.44
1:G:182:ILE:HG23	1:G:244:LEU:HD13	1.99	0.44
1:H:112:ILE:HD13	1:H:126:VAL:HG21	2.00	0.44
1:E:116:ARG:N	1:E:117:PRO:HD2	2.33	0.43
1:H:230:ILE:CG2	1:H:231:THR:N	2.81	0.43
1:C:58:GLY:C	1:C:60:ARG:H	2.27	0.43
1:C:145:MET:CE	1:C:158:ALA:HB2	2.48	0.43
1:F:157:VAL:HG12	1:F:203:ILE:HG23	2.01	0.43
1:D:40:VAL:HG12	1:D:91:LEU:HD22	2.00	0.43
1:D:157:VAL:CG2	1:D:203:ILE:HG23	2.49	0.43
1:E:39:ALA:HB1	1:E:127:VAL:HG11	2.00	0.43
1:G:209:LEU:HD22	1:G:234:LEU:HD11	2.01	0.43
1:B:167:THR:HG21	1:B:197:ALA:CB	2.48	0.43
1:G:224:LYS:O	1:G:229:GLN:NE2	2.52	0.43
1:H:200:ASN:C	1:H:201:LEU:HD22	2.44	0.43
1:A:142:LEU:O	1:A:146:VAL:HG23	2.19	0.43
1:G:109:GLY:N	1:G:228:ILE:O	2.52	0.43
1:G:251:LYS:HD2	1:G:252:ARG:N	2.34	0.43
1:B:209:LEU:CD2	1:B:234:LEU:HD21	2.49	0.43
1:B:269:TYR:CE2	1:B:274:LYS:CE	3.00	0.43
1:D:104:GLU:O	1:D:105:MET:SD	2.76	0.43
1:E:24:PRO:HD3	1:E:58:GLY:O	2.19	0.43
1:B:235:MET:HE3	1:B:239:GLN:OE1	2.19	0.42
1:G:185:VAL:HG12	1:G:238:ALA:HB1	2.01	0.42
1:A:83:LYS:HB3	1:A:84:TYR:H	1.52	0.42
1:C:69:SER:O	1:C:69:SER:OG	2.35	0.42
1:C:204:ILE:HD12	1:C:248:PHE:CE1	2.54	0.42
1:E:80:ASP:O	1:E:82:GLU:N	2.52	0.42
1:E:85:LEU:O	1:E:89:ARG:HG3	2.18	0.42
1:H:173:ILE:HD12	1:H:173:ILE:C	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:239:GLN:OE1	1:H:239:GLN:N	2.52	0.42
1:A:206:ARG:NH2	6:A:402:HOH:O	2.52	0.42
1:B:157:VAL:HG13	1:B:203:ILE:HG23	2.01	0.42
1:C:105:MET:HG2	2:C:301:UPG:O4	2.19	0.42
1:G:136:LEU:H	1:G:136:LEU:CD2	2.29	0.42
1:H:155:SER:HA	1:H:208:ILE:O	2.19	0.42
1:C:157:VAL:HG12	1:C:245:ALA:HA	2.02	0.42
1:E:188:MET:H	1:E:235:MET:CE	2.31	0.42
1:G:259:GLU:O	1:G:262:ILE:HG22	2.19	0.42
1:H:108:LEU:O	1:H:112:ILE:HG13	2.19	0.42
1:A:4:LYS:HE2	1:A:118:LEU:O	2.19	0.42
1:B:53:ILE:HG21	1:B:62:LEU:HD13	2.02	0.42
1:C:56:GLY:O	1:C:59:LYS:HB3	2.18	0.42
1:E:108:LEU:HD12	1:E:108:LEU:O	2.20	0.42
1:B:44:LEU:HD13	1:B:95:CYS:SG	2.60	0.42
1:F:135:ASN:ND2	1:F:139:ASP:O	2.51	0.42
1:G:168:ASN:OD1	1:G:168:ASN:N	2.46	0.42
1:H:205:GLY:HA2	1:H:207:TYR:CE1	2.54	0.42
1:D:82:GLU:HG2	1:D:85:LEU:HD12	2.02	0.42
1:B:187:THR:HA	1:B:235:MET:CE	2.48	0.42
1:H:149:TYR:O	1:H:153:ARG:N	2.50	0.42
1:H:258:ALA:O	1:H:262:ILE:HG12	2.20	0.42
1:E:134:LEU:HD12	1:E:253:PHE:HE2	1.85	0.41
1:F:192:PRO:CG	1:F:197:ALA:HB2	2.48	0.41
1:H:141:VAL:O	1:H:145:MET:HG3	2.20	0.41
1:G:20:THR:O	1:G:21:LYS:C	2.63	0.41
1:G:112:ILE:CD1	1:G:126:VAL:HG11	2.50	0.41
1:H:42:GLU:OE1	1:H:206:ARG:NH1	2.53	0.41
1:E:76:ILE:O	1:E:76:ILE:CG1	2.69	0.41
1:G:267:PHE:CD1	1:G:267:PHE:C	2.98	0.41
1:H:12:TYR:CE1	1:H:105:MET:HG3	2.55	0.41
1:C:105:MET:CG	2:C:301:UPG:O4	2.68	0.41
1:E:242:CYS:SG	1:E:244:LEU:HD13	2.60	0.41
1:G:179:ARG:HG2	1:G:180:ASP:H	1.86	0.41
1:F:167:THR:CG2	1:F:194:PRO:HA	2.50	0.41
1:H:152:PHE:N	1:H:152:PHE:CD1	2.89	0.41
1:E:91:LEU:O	1:E:95:CYS:HB2	2.21	0.41
1:G:12:TYR:CD2	1:G:105:MET:SD	3.14	0.41
1:G:85:LEU:O	1:G:89:ARG:HG3	2.20	0.41
1:G:172:VAL:CG1	1:G:173:ILE:N	2.84	0.41
1:H:85:LEU:O	1:H:89:ARG:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:224:LYS:O	1:H:229:GLN:NE2	2.51	0.41
1:B:101:ARG:N	1:B:101:ARG:HD2	2.36	0.41
1:H:90:ARG:HG2	1:H:94:GLU:HG3	2.02	0.41
1:D:268:CYS:O	1:D:272:LEU:HB2	2.21	0.40
1:D:151:GLN:HB3	1:D:152:PHE:CD1	2.56	0.40
1:H:159:ILE:HG22	1:H:203:ILE:CA	2.51	0.40
1:B:68:ILE:O	1:B:68:ILE:CG2	2.70	0.40
1:B:116:ARG:N	1:B:117:PRO:CD	2.85	0.40
1:E:170:TYR:CB	1:E:202:ALA:HB1	2.51	0.40
1:F:30:VAL:O	1:F:30:VAL:CG2	2.69	0.40
1:A:149:TYR:CE1	1:A:153:ARG:HA	2.57	0.40
1:D:30:VAL:O	1:D:32:ASN:N	2.51	0.40
1:F:20:THR:HA	1:F:23:MET:O	2.22	0.40
1:G:182:ILE:C	1:G:183:PHE:CD1	2.99	0.40
1:H:187:THR:HA	1:H:235:MET:SD	2.61	0.40
1:H:268:CYS:O	1:H:272:LEU:HB2	2.22	0.40

All (9) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:LYS:NZ	1:F:67:ASP:OD2[1_655]	1.94	0.26
1:F:32:ASN:HD22	1:H:17:LEU:O[2_555]	1.35	0.25
1:F:32:ASN:ND2	1:H:17:LEU:O[2_555]	1.97	0.23
1:E:101:ARG:NH2	1:G:71:GLU:OE2[2_656]	2.00	0.20
1:A:67:ASP:O	1:B:21:LYS:HZ1[2_645]	1.53	0.07
1:E:101:ARG:HH21	1:G:71:GLU:OE2[2_656]	1.53	0.07
1:C:75:GLN:OE1	1:D:57:ARG:NH1[2_646]	2.16	0.04
1:F:21:LYS:HZ1	1:H:67:ASP:O[2_555]	1.58	0.02
1:B:59:LYS:HZ3	1:F:67:ASP:OD2[1_655]	1.59	0.01

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	273/285 (96%)	259 (95%)	14 (5%)	0	100	100
1	B	276/285 (97%)	264 (96%)	10 (4%)	2 (1%)	18	48
1	C	270/285 (95%)	254 (94%)	13 (5%)	3 (1%)	11	36
1	D	268/285 (94%)	253 (94%)	14 (5%)	1 (0%)	30	59
1	E	279/285 (98%)	266 (95%)	11 (4%)	2 (1%)	18	48
1	F	273/285 (96%)	255 (93%)	16 (6%)	2 (1%)	18	48
1	G	265/285 (93%)	245 (92%)	17 (6%)	3 (1%)	11	36
1	H	272/285 (95%)	264 (97%)	6 (2%)	2 (1%)	18	48
All	All	2176/2280 (95%)	2060 (95%)	101 (5%)	15 (1%)	18	48

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	195	GLU
1	H	79	THR
1	C	31	VAL
1	C	196	GLU
1	D	104	GLU
1	E	81	LYS
1	B	77	ARG
1	F	79	THR
1	C	59	LYS
1	F	31	VAL
1	G	21	LYS
1	B	76	ILE
1	H	31	VAL
1	E	31	VAL
1	G	198	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	233/239 (98%)	224 (96%)	9 (4%)	28	63
1	B	235/239 (98%)	231 (98%)	4 (2%)	53	82
1	C	231/239 (97%)	226 (98%)	5 (2%)	45	77
1	D	229/239 (96%)	219 (96%)	10 (4%)	25	59
1	E	237/239 (99%)	230 (97%)	7 (3%)	36	70
1	F	233/239 (98%)	227 (97%)	6 (3%)	40	73
1	G	228/239 (95%)	218 (96%)	10 (4%)	25	59
1	H	231/239 (97%)	224 (97%)	7 (3%)	36	70
All	All	1857/1912 (97%)	1799 (97%)	58 (3%)	35	69

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	MET
1	A	60	ARG
1	A	76	ILE
1	A	81	LYS
1	A	82	GLU
1	A	83	LYS
1	A	86	VAL
1	A	165	GLU
1	A	275	THR
1	B	33	LYS
1	B	82	GLU
1	B	178	ILE
1	B	187	THR
1	C	2	ILE
1	C	154	CYS
1	C	177	MET
1	C	204	ILE
1	C	208	ILE
1	D	31	VAL
1	D	49	SER
1	D	67	ASP
1	D	83	LYS
1	D	96	THR
1	D	101	ARG
1	D	104	GLU
1	D	170	TYR
1	D	172	VAL

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Mol	Chain	Res	Type
1	D	212	ASP
1	E	20	THR
1	E	25	LYS
1	E	60	ARG
1	E	73	GLU
1	E	167	THR
1	E	208	ILE
1	E	219	GLN
1	F	2	ILE
1	F	57	ARG
1	F	59	LYS
1	F	195	GLU
1	F	224	LYS
1	F	275	THR
1	G	68	ILE
1	G	94	GLU
1	G	112	ILE
1	G	114	THR
1	G	161	GLU
1	G	168	ASN
1	G	188	MET
1	G	192	PRO
1	G	195	GLU
1	G	213	ILE
1	H	14	THR
1	H	67	ASP
1	H	68	ILE
1	H	93	ASP
1	H	96	THR
1	H	184	ARG
1	H	206	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	266	ASN
1	A	271	ASN
1	C	200	ASN
1	C	266	ASN
1	D	160	GLN
1	E	271	ASN
1	H	266	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 20 ligands modelled in this entry, 8 are monoatomic - leaving 12 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	UPG	E	301	4	37,38,38	0.46	0	55,58,58	0.39	0
2	UPG	H	301	4	37,38,38	0.84	0	55,58,58	0.76	2 (3%)
2	UPG	A	301	4	37,38,38	0.42	0	55,58,58	0.46	0
2	UPG	D	301	4	37,38,38	1.05	3 (8%)	55,58,58	0.72	0
5	ACT	B	303	-	3,3,3	1.15	0	3,3,3	1.40	0
3	EDO	D	302	-	3,3,3	0.39	0	2,2,2	0.45	0
2	UPG	C	301	4	37,38,38	1.10	2 (5%)	55,58,58	0.86	4 (7%)
2	UPG	B	301	4	37,38,38	1.35	3 (8%)	55,58,58	0.74	1 (1%)
2	UPG	G	301	4	37,38,38	0.46	0	55,58,58	0.48	0
3	EDO	A	302	-	3,3,3	0.44	0	2,2,2	0.38	0
2	UPG	F	301	4	37,38,38	0.93	0	55,58,58	0.78	0
3	EDO	C	302	-	3,3,3	0.44	0	2,2,2	0.30	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	UPG	E	301	4	-	7/23/59/59	0/3/3/3
2	UPG	H	301	4	-	3/23/59/59	0/3/3/3
2	UPG	A	301	4	-	3/23/59/59	0/3/3/3
2	UPG	D	301	4	-	3/23/59/59	0/3/3/3
3	EDO	D	302	-	-	1/1/1/1	-
2	UPG	C	301	4	-	9/23/59/59	0/3/3/3
2	UPG	B	301	4	-	3/23/59/59	0/3/3/3
2	UPG	G	301	4	-	4/23/59/59	0/3/3/3
3	EDO	A	302	-	-	1/1/1/1	-
2	UPG	F	301	4	-	5/23/59/59	0/3/3/3
3	EDO	C	302	-	-	0/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	301	UPG	PA-O3A	-3.90	1.55	1.59
2	B	301	UPG	PB-O3A	-2.84	1.56	1.59
2	B	301	UPG	C3'-C2'	-2.84	1.45	1.52
2	C	301	UPG	C4'-C5'	-2.65	1.47	1.53
2	D	301	UPG	C4'-C5'	-2.26	1.48	1.53
2	D	301	UPG	C3'-C2'	-2.24	1.46	1.52
2	D	301	UPG	PA-O3A	-2.10	1.57	1.59
2	C	301	UPG	C3'-C2'	-2.03	1.47	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	301	UPG	C1'-O5'-C5'	2.73	119.06	113.72
2	C	301	UPG	O2A-PA-O1A	2.41	123.64	112.44
2	H	301	UPG	O2'-C2'-C3'	-2.23	105.12	110.38
2	B	301	UPG	O2A-PA-O1A	2.20	122.69	112.44
2	C	301	UPG	O5'-C1'-O3B	2.17	114.20	111.36
2	H	301	UPG	O3B-PB-O1B	-2.12	103.03	109.81
2	C	301	UPG	O3B-PB-O1B	-2.10	103.08	109.81

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	301	UPG	C5C-O5C-PA-O2A

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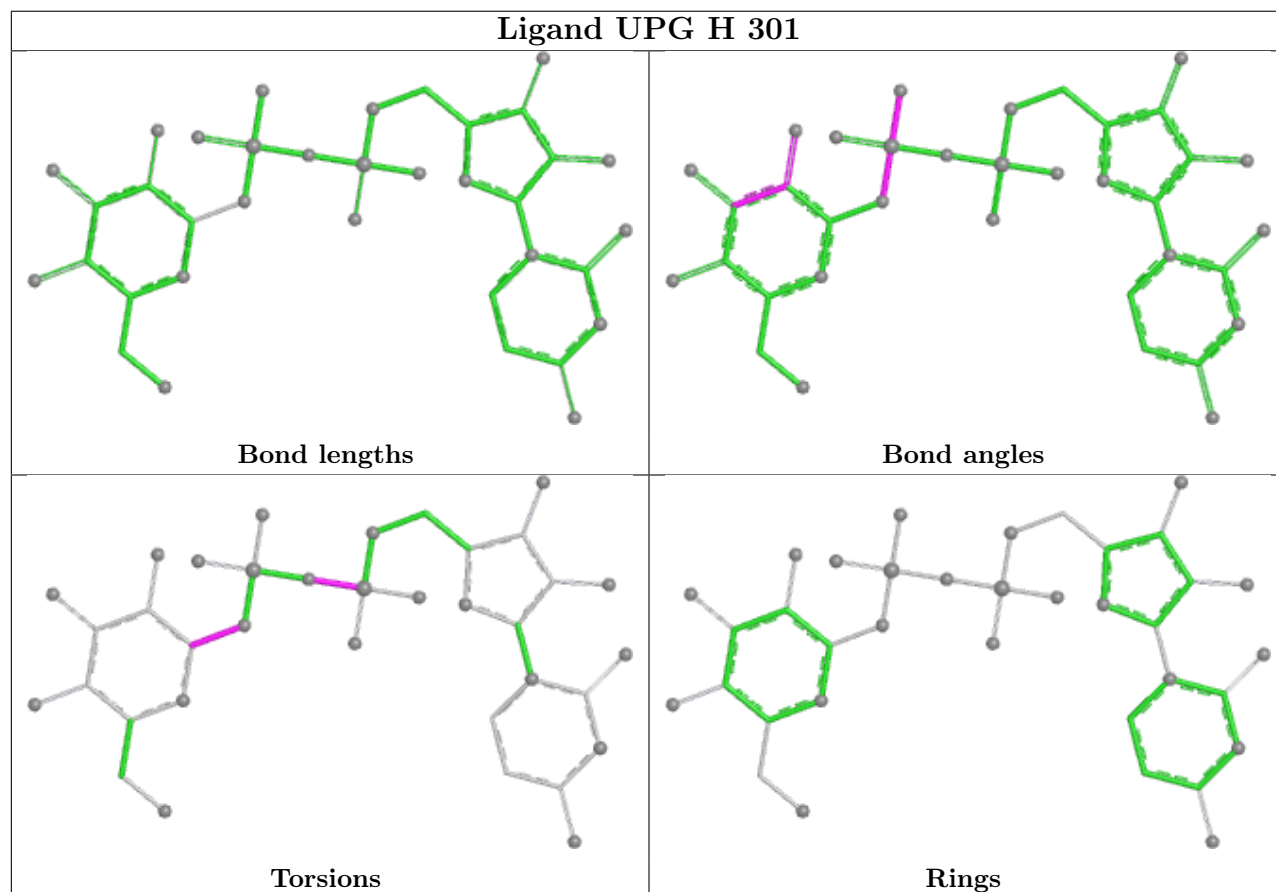
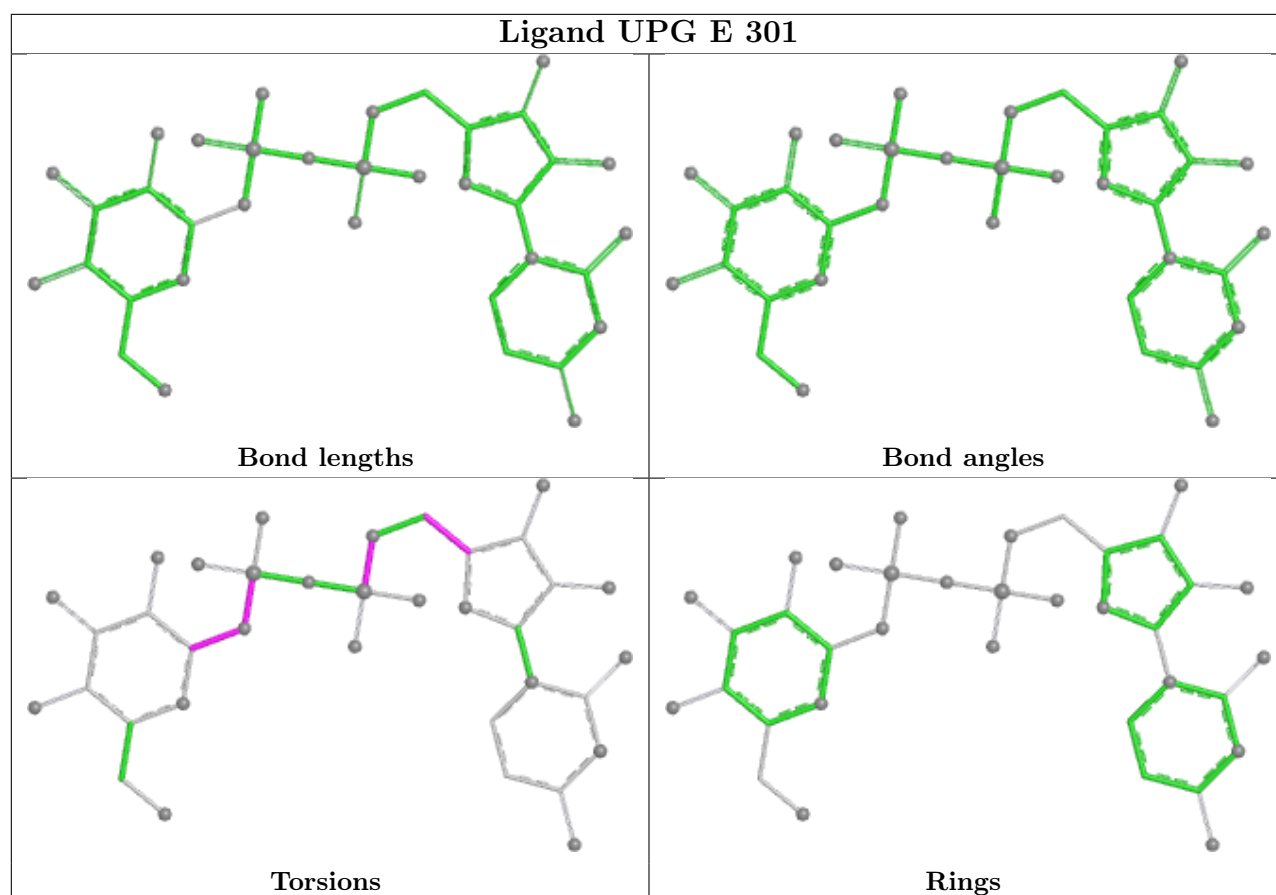
Mol	Chain	Res	Type	Atoms
2	C	301	UPG	C5C-O5C-PA-O3A
2	C	301	UPG	O5'-C1'-O3B-PB
2	E	301	UPG	C5C-O5C-PA-O2A
2	E	301	UPG	C1'-O3B-PB-O3A
2	F	301	UPG	C5C-O5C-PA-O2A
2	D	301	UPG	C2'-C1'-O3B-PB
2	E	301	UPG	C2'-C1'-O3B-PB
2	F	301	UPG	C2'-C1'-O3B-PB
2	H	301	UPG	C2'-C1'-O3B-PB
2	A	301	UPG	O5'-C5'-C6'-O6'
3	D	302	EDO	O1-C1-C2-O2
2	B	301	UPG	C2'-C1'-O3B-PB
2	C	301	UPG	C2'-C1'-O3B-PB
2	G	301	UPG	C2'-C1'-O3B-PB
2	E	301	UPG	O4C-C4C-C5C-O5C
2	G	301	UPG	O4C-C4C-C5C-O5C
2	D	301	UPG	O5'-C5'-C6'-O6'
3	A	302	EDO	O1-C1-C2-O2
2	E	301	UPG	C3C-C4C-C5C-O5C
2	A	301	UPG	C2'-C1'-O3B-PB
2	C	301	UPG	C5C-O5C-PA-O1A
2	E	301	UPG	C5C-O5C-PA-O1A
2	E	301	UPG	C5C-O5C-PA-O3A
2	F	301	UPG	C5C-O5C-PA-O3A
2	B	301	UPG	PB-O3A-PA-O2A
2	D	301	UPG	PB-O3A-PA-O2A
2	G	301	UPG	PB-O3A-PA-O2A
2	C	301	UPG	O5'-C5'-C6'-O6'
2	A	301	UPG	C4'-C5'-C6'-O6'
2	G	301	UPG	PB-O3A-PA-O1A
2	C	301	UPG	O4C-C4C-C5C-O5C
2	F	301	UPG	C4'-C5'-C6'-O6'
2	B	301	UPG	PB-O3A-PA-O1A
2	C	301	UPG	PA-O3A-PB-O1B
2	H	301	UPG	PB-O3A-PA-O1A
2	H	301	UPG	PB-O3A-PA-O2A
2	C	301	UPG	PA-O3A-PB-O2B
2	F	301	UPG	O4C-C4C-C5C-O5C

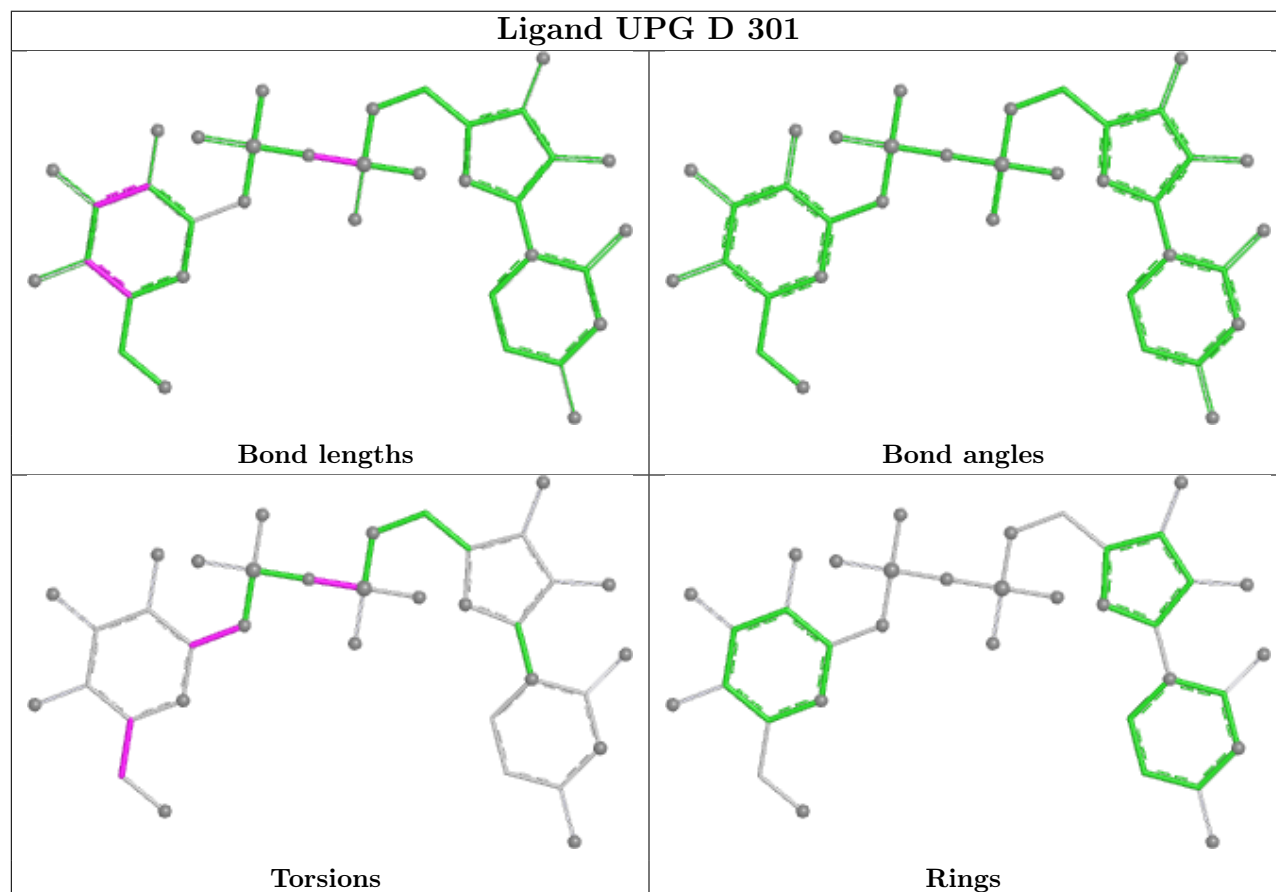
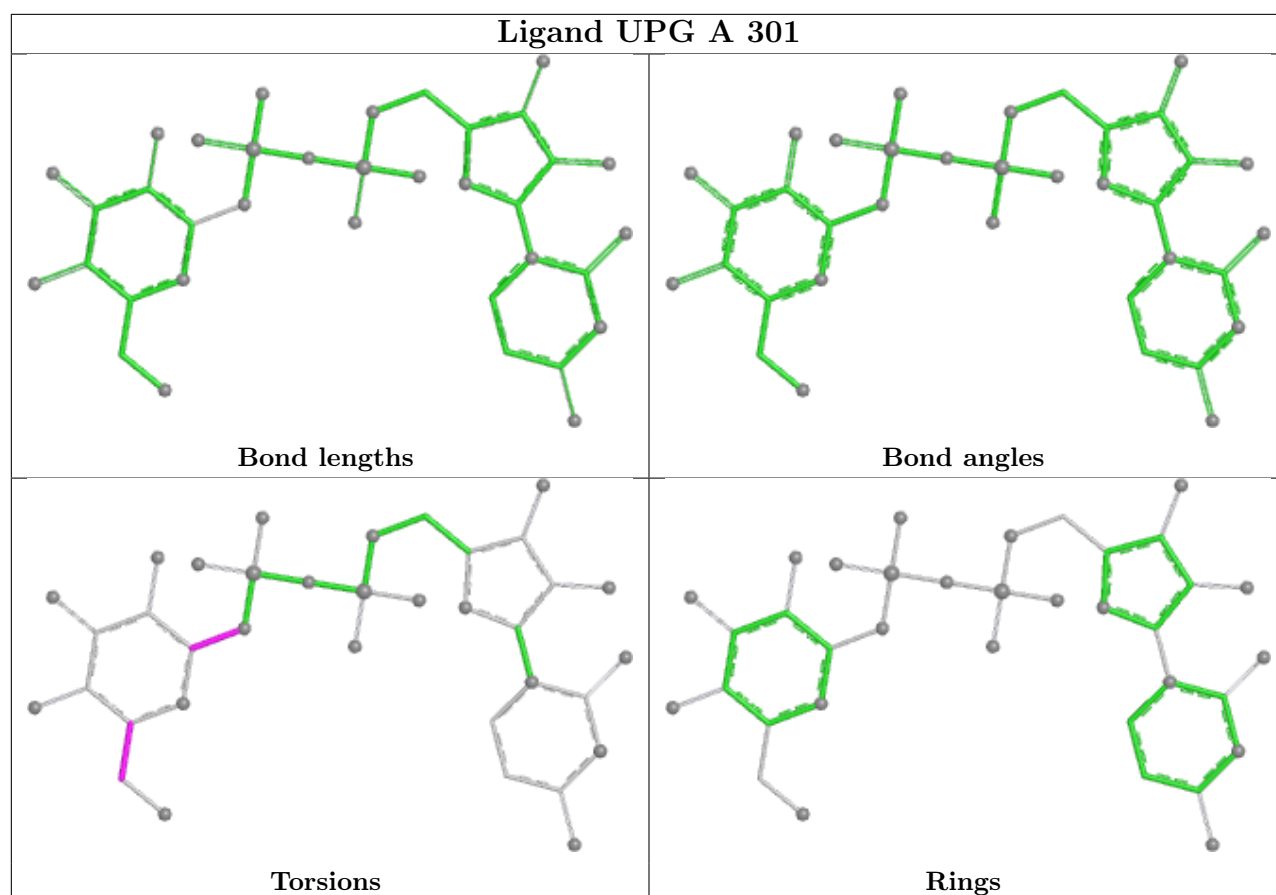
There are no ring outliers.

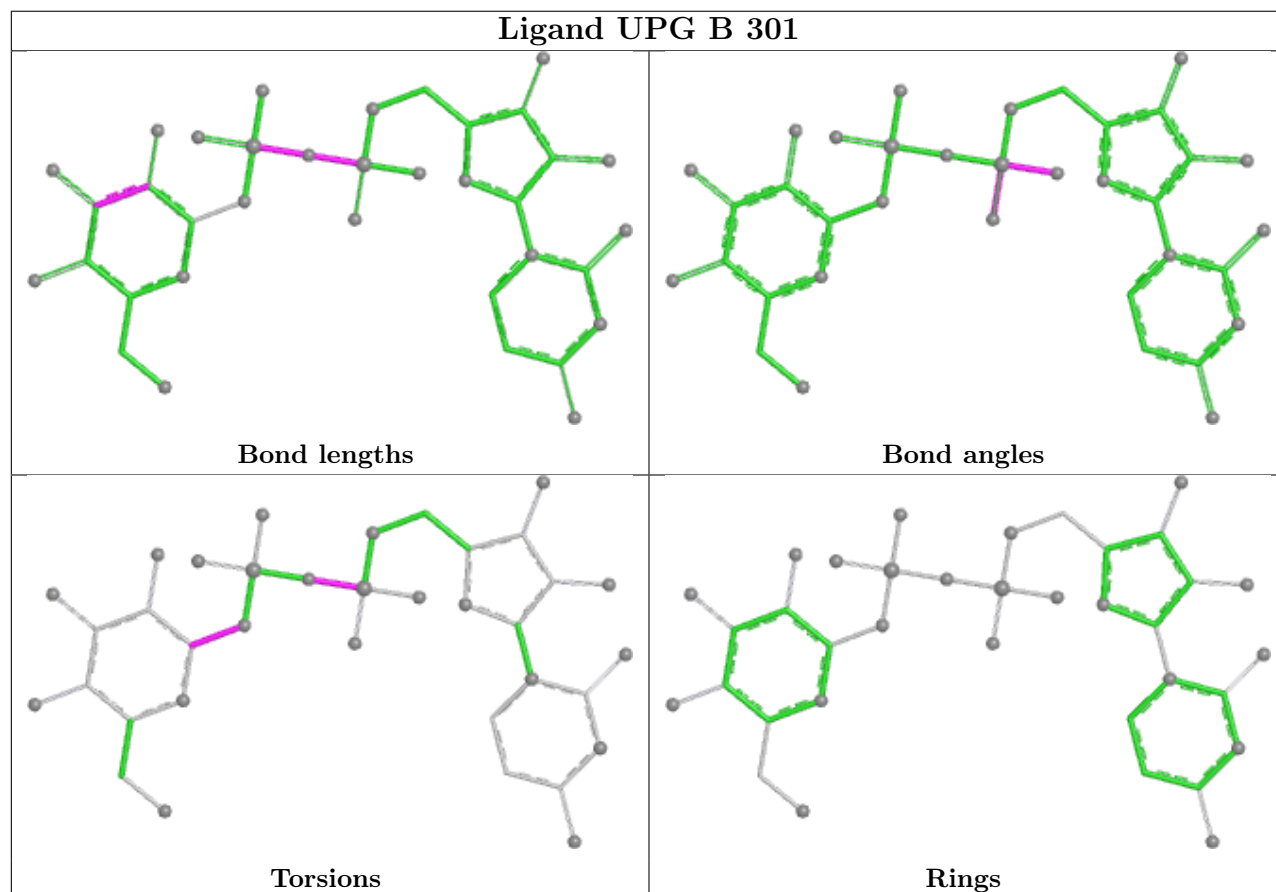
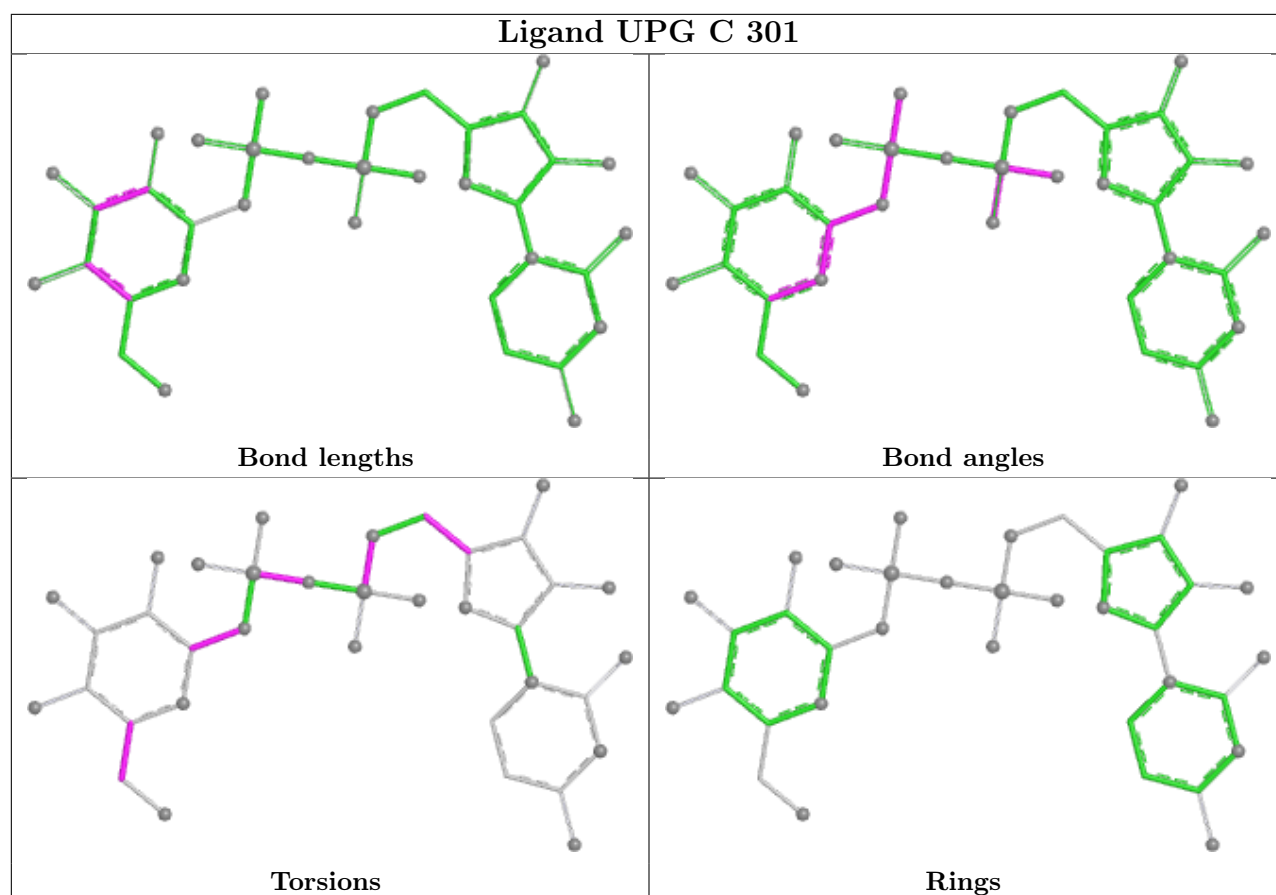
10 monomers are involved in 14 short contacts:

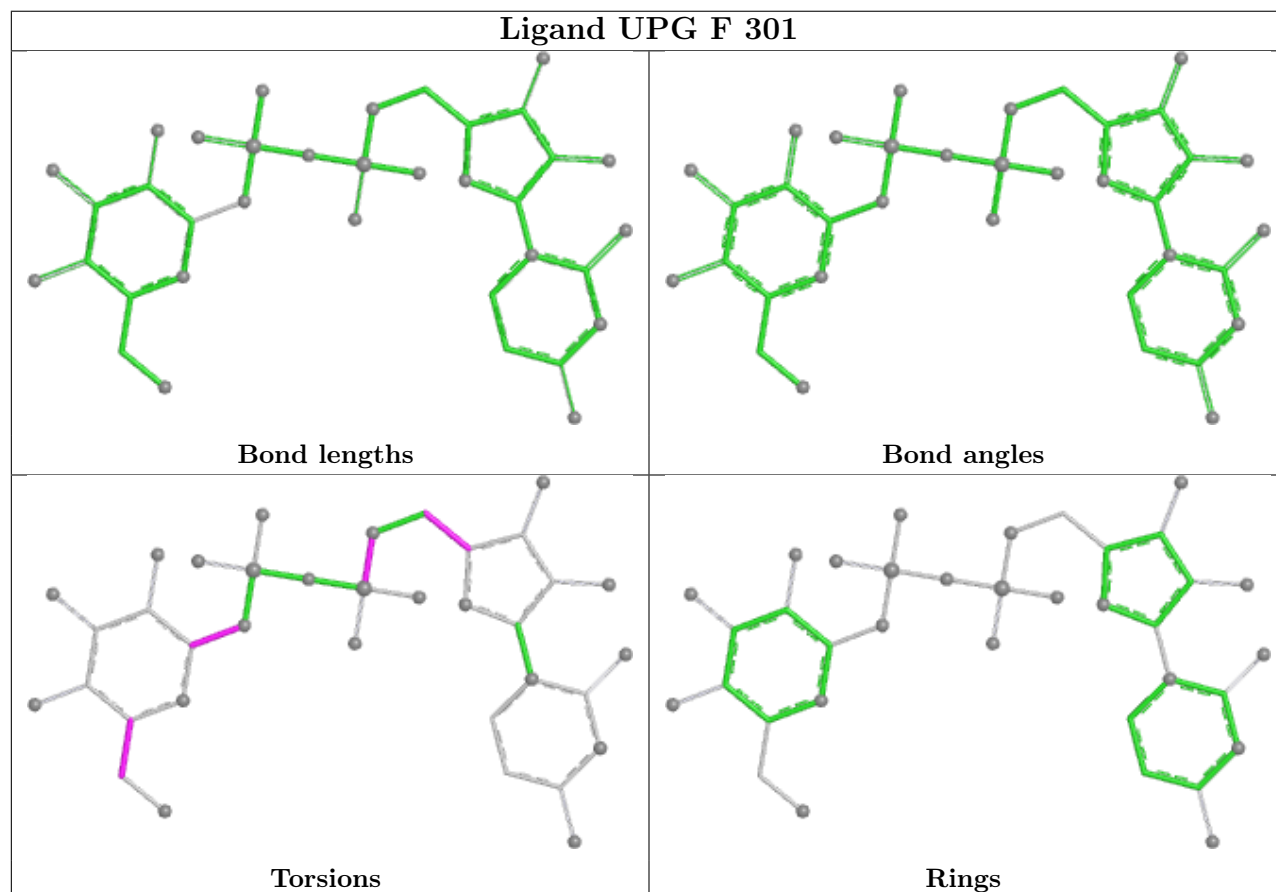
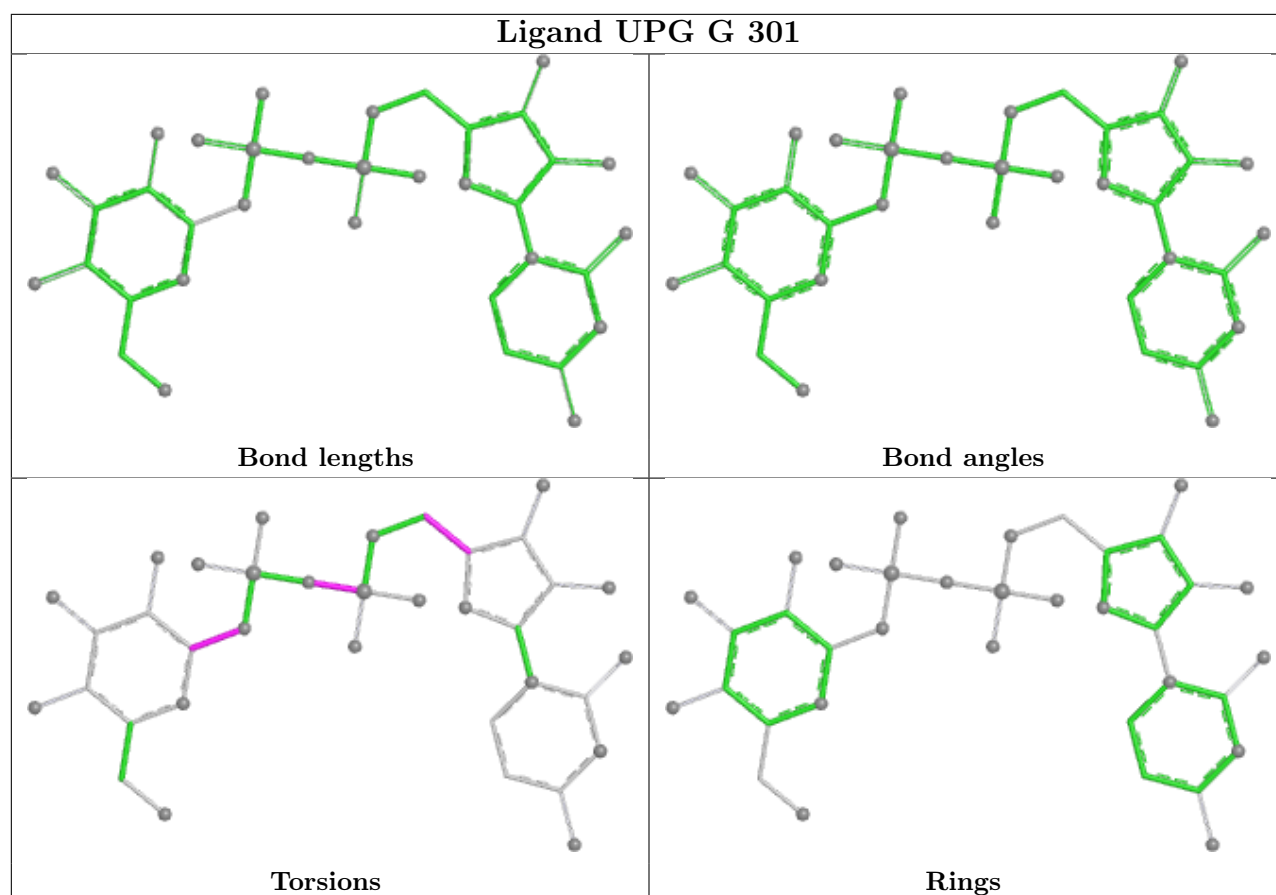
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	301	UPG	1	0
2	H	301	UPG	1	0
2	A	301	UPG	1	0
2	D	301	UPG	1	0
3	D	302	EDO	1	0
2	C	301	UPG	3	0
2	B	301	UPG	2	0
2	G	301	UPG	1	0
2	F	301	UPG	2	0
3	C	302	EDO	1	0

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	275/285 (96%)	-0.10	3 (1%) 78 71	26, 57, 85, 117	5 (1%)
1	B	280/285 (98%)	-0.19	3 (1%) 78 71	35, 56, 84, 142	5 (1%)
1	C	274/285 (96%)	0.02	4 (1%) 72 64	40, 66, 99, 140	4 (1%)
1	D	272/285 (95%)	0.05	6 (2%) 62 53	36, 67, 95, 119	6 (2%)
1	E	281/285 (98%)	0.02	2 (0%) 84 79	29, 67, 100, 132	7 (2%)
1	F	275/285 (96%)	0.04	2 (0%) 84 79	40, 69, 100, 142	6 (2%)
1	G	271/285 (95%)	0.50	17 (6%) 26 20	33, 77, 113, 141	4 (1%)
1	H	274/285 (96%)	0.33	11 (4%) 42 34	34, 76, 108, 150	8 (2%)
All	All	2202/2280 (96%)	0.08	48 (2%) 62 53	26, 67, 103, 150	45 (2%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	77	ARG	4.3
1	B	17	LEU	4.0
1	G	202	ALA	3.8
1	G	199	SER	3.6
1	B	18	PRO	3.6
1	G	84	TYR	3.5
1	H	84	TYR	3.2
1	G	173	ILE	3.2
1	G	223	GLY	2.8
1	G	148	LEU	2.8
1	D	80	ASP	2.7
1	G	134	LEU	2.6
1	G	80	ASP	2.6
1	H	81	LYS	2.6
1	H	204	ILE	2.6
1	H	230	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	77	ARG	2.5
1	F	240	ASP	2.5
1	G	136	LEU	2.5
1	C	192	PRO	2.5
1	G	12	TYR	2.5
1	G	172	VAL	2.4
1	G	191	LYS	2.4
1	H	167	THR	2.3
1	D	178	ILE	2.3
1	H	59	LYS	2.3
1	G	175	GLY	2.3
1	H	186	ASN	2.3
1	D	103	VAL	2.3
1	C	1	SER	2.2
1	D	198	PRO	2.2
1	A	79	THR	2.2
1	A	11	GLY	2.2
1	C	203	ILE	2.2
1	G	244	LEU	2.1
1	E	77	ARG	2.1
1	F	82	GLU	2.1
1	A	31	VAL	2.1
1	G	243	VAL	2.1
1	D	81	LYS	2.1
1	G	76	ILE	2.1
1	H	203	ILE	2.1
1	B	73	GLU	2.1
1	H	105	MET	2.1
1	H	77	ARG	2.1
1	D	31	VAL	2.0
1	E	118	LEU	2.0
1	H	68	ILE	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 ⓘ

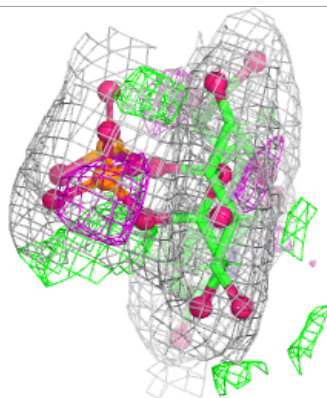
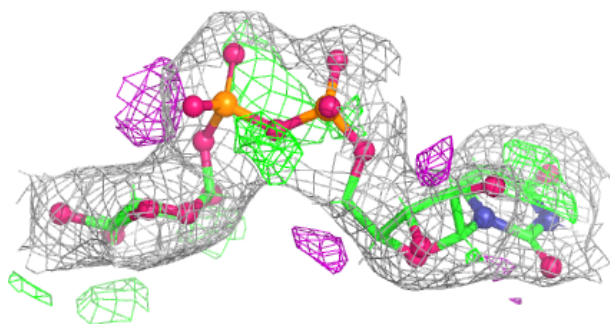
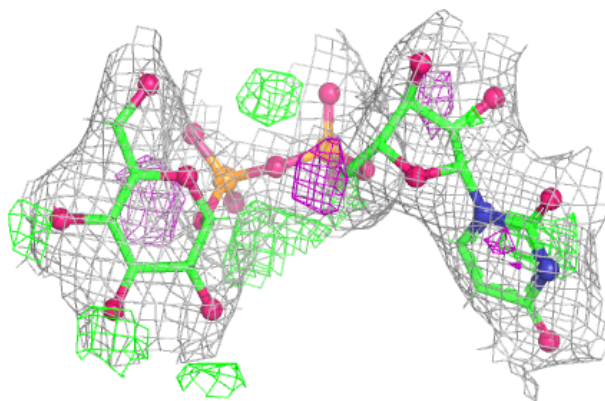
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	ACT	B	303	4/4	0.77	0.25	67,79,99,99	0
3	EDO	A	302	4/4	0.78	0.21	63,78,93,93	0
4	MG	H	302	1/1	0.88	0.08	66,66,66,66	0
3	EDO	C	302	4/4	0.89	0.17	55,68,84,85	0
2	UPG	E	301	36/36	0.90	0.12	42,60,81,89	0
2	UPG	G	301	36/36	0.92	0.10	44,67,86,100	0
4	MG	D	303	1/1	0.94	0.06	68,68,68,68	0
4	MG	G	302	1/1	0.94	0.12	57,57,57,57	0
2	UPG	H	301	36/36	0.95	0.07	44,68,92,100	0
2	UPG	F	301	36/36	0.95	0.07	49,62,84,89	0
2	UPG	B	301	36/36	0.95	0.09	36,57,73,130	0
3	EDO	D	302	4/4	0.95	0.20	74,97,129,141	0
2	UPG	C	301	36/36	0.96	0.08	43,56,72,76	0
4	MG	E	302	1/1	0.97	0.11	52,52,52,52	0
4	MG	F	302	1/1	0.97	0.09	61,61,61,61	0
2	UPG	A	301	36/36	0.97	0.07	35,49,63,67	0
4	MG	C	303	1/1	0.97	0.08	45,45,45,45	0
2	UPG	D	301	36/36	0.97	0.06	37,58,73,86	0
4	MG	A	303	1/1	0.98	0.04	52,52,52,52	0
4	MG	B	302	1/1	0.99	0.03	51,51,51,51	0

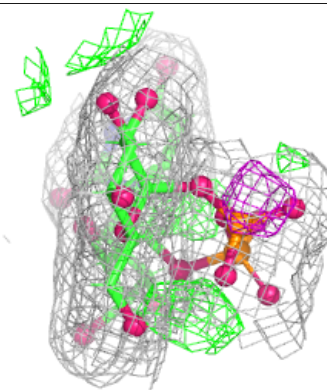
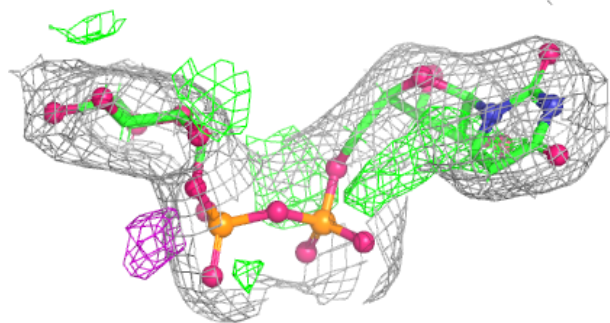
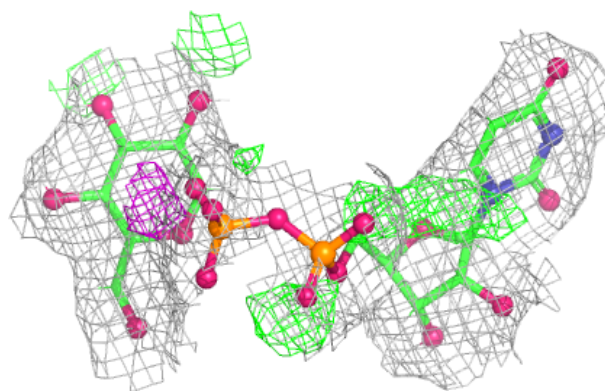
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around UPG E 301:

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

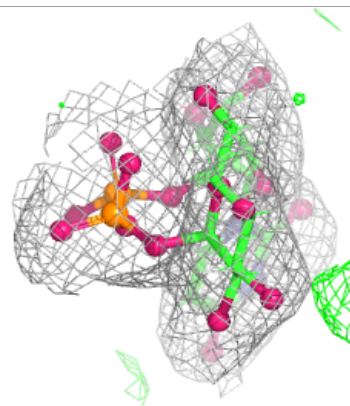
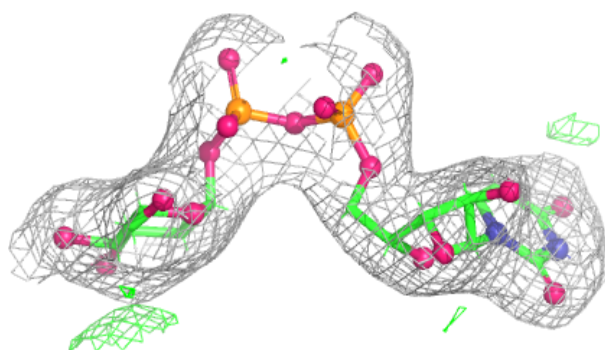
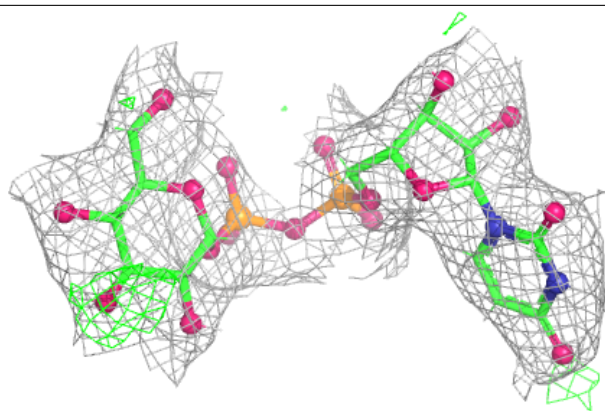
**Electron density around UPG G 301:**

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

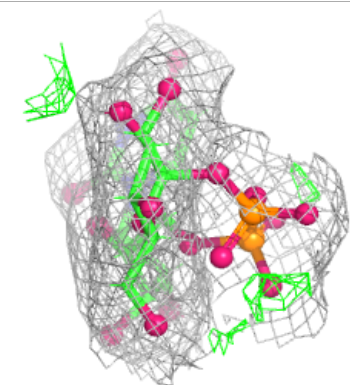
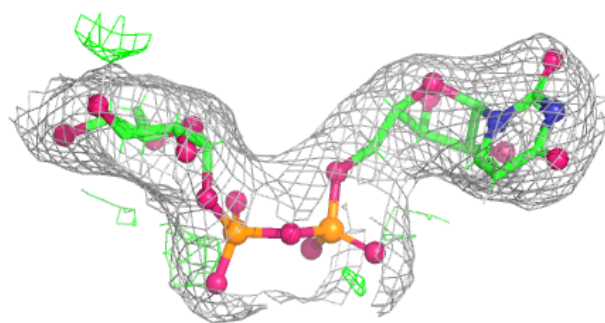
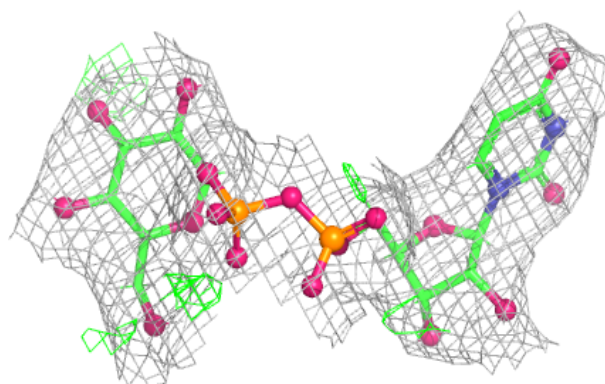


Electron density around UPG H 301:

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

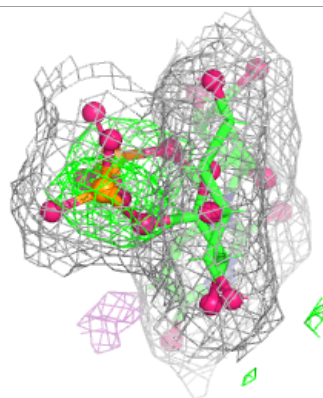
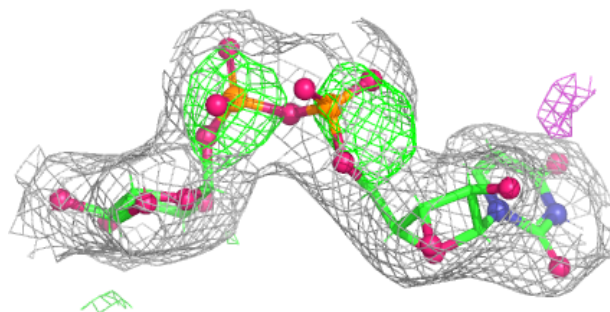
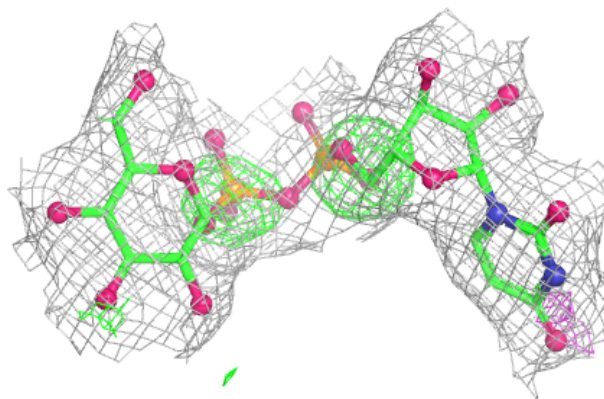
**Electron density around UPG F 301:**

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

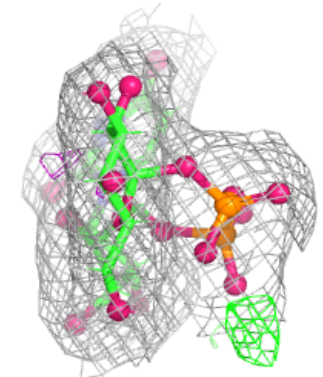
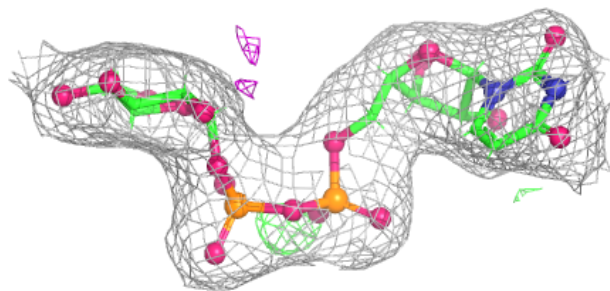
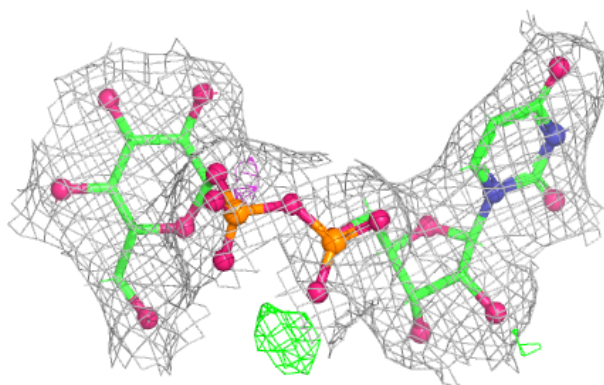


Electron density around UPG B 301:

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

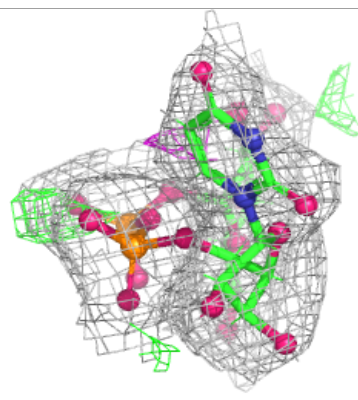
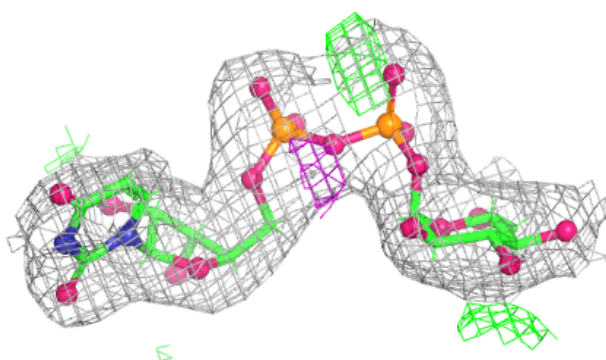
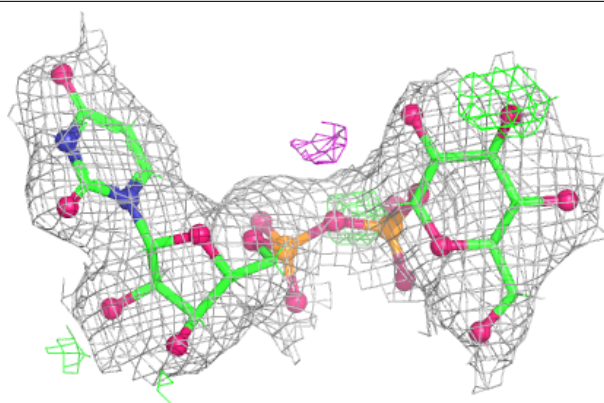
**Electron density around UPG C 301:**

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

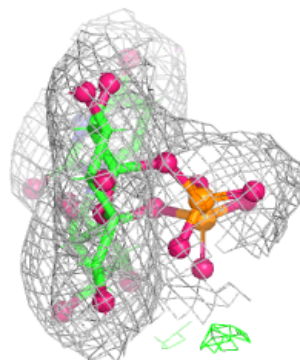
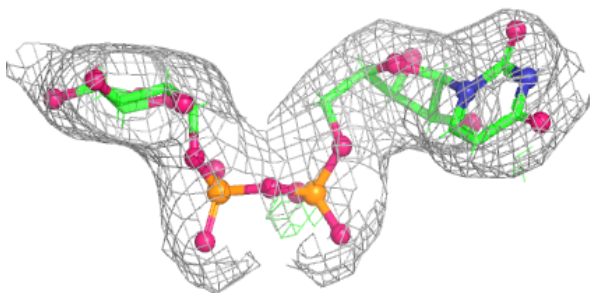
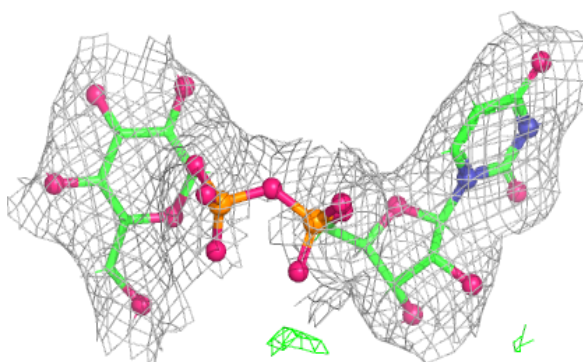


Electron density around UPG A 301:

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

**Electron density around UPG D 301:**

$2mF_o-DF_c$ (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.