



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 8, 2026 – 10:16 AM JST

PDB ID : 9UVL / pdb_00009uvl
Title : Crystal structure of AGAP1-GDP complex
Authors : Zhang, S.; Zhang, Y.; Cheng, N.; Wang, J.; Li, Z.; Meng, G.; Zhang, H.
Deposited on : 2025-05-10
Resolution : 2.49 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 : 1.8.5 (274361), CSD as541be (2020)
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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

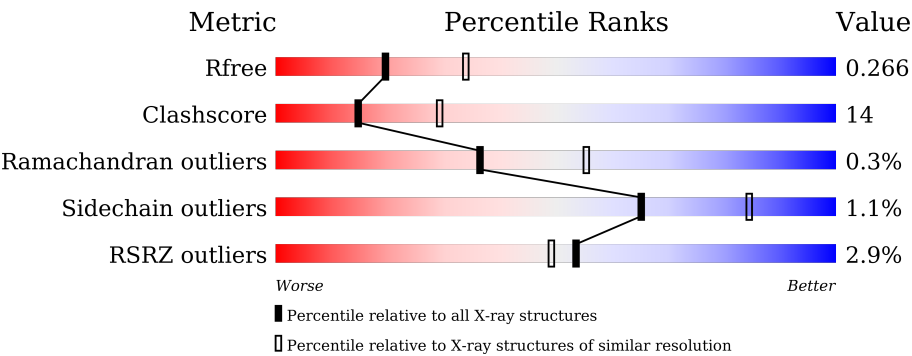
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	172	2% 77%17%••
1	B	172	% 70%24%•5%
1	C	172	% 80%13%•5%
1	D	172	2% 74%21%•5%
1	E	172	2% 82%14%•
1	F	172	3% 70%23%••5%

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Mol	Chain	Length	Quality of chain
1	G	172	7% 69% 27% 5%
1	H	172	3% 71% 24% • 5%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 21066 atoms, of which 10350 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 Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	166	Total	C	H	N	O	S	0	0	0
			2604	831	1292	224	253	4			
1	B	164	Total	C	H	N	O	S	0	0	0
			2581	823	1283	222	249	4			
1	C	164	Total	C	H	N	O	S	0	0	0
			2581	823	1283	222	249	4			
1	D	164	Total	C	H	N	O	S	0	0	0
			2581	823	1283	222	249	4			
1	E	165	Total	C	H	N	O	S	0	0	0
			2595	828	1288	223	252	4			
1	F	164	Total	C	H	N	O	S	0	0	0
			2581	823	1283	222	249	4			
1	G	164	Total	C	H	N	O	S	0	0	0
			2580	823	1282	222	249	4			
1	H	164	Total	C	H	N	O	S	0	0	0
			2581	823	1283	222	249	4			

There are 48 discrepancies between the modelled and reference sequences:

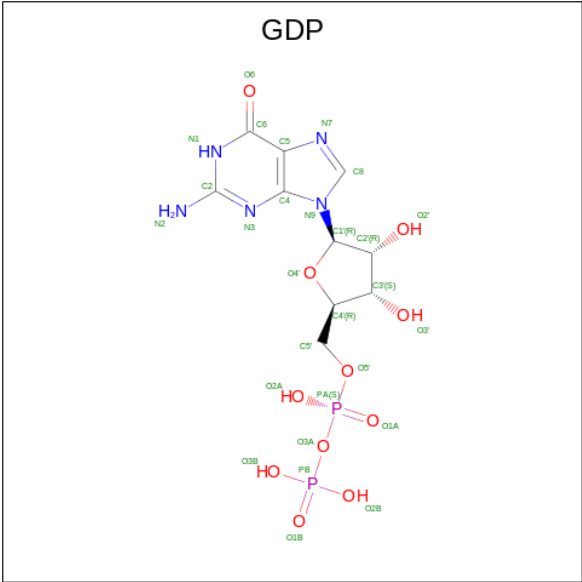
Chain	Residue	Modelled	Actual	Comment	Reference
A	64	GLY	-	expression tag	UNP Q9UPQ3
A	65	SER	-	expression tag	UNP Q9UPQ3
A	66	HIS	-	expression tag	UNP Q9UPQ3
A	67	MET	-	expression tag	UNP Q9UPQ3
A	68	LEU	-	expression tag	UNP Q9UPQ3
A	69	GLU	-	expression tag	UNP Q9UPQ3
B	64	GLY	-	expression tag	UNP Q9UPQ3
B	65	SER	-	expression tag	UNP Q9UPQ3
B	66	HIS	-	expression tag	UNP Q9UPQ3
B	67	MET	-	expression tag	UNP Q9UPQ3
B	68	LEU	-	expression tag	UNP Q9UPQ3
B	69	GLU	-	expression tag	UNP Q9UPQ3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	64	GLY	-	expression tag	UNP Q9UPQ3
C	65	SER	-	expression tag	UNP Q9UPQ3
C	66	HIS	-	expression tag	UNP Q9UPQ3
C	67	MET	-	expression tag	UNP Q9UPQ3
C	68	LEU	-	expression tag	UNP Q9UPQ3
C	69	GLU	-	expression tag	UNP Q9UPQ3
D	64	GLY	-	expression tag	UNP Q9UPQ3
D	65	SER	-	expression tag	UNP Q9UPQ3
D	66	HIS	-	expression tag	UNP Q9UPQ3
D	67	MET	-	expression tag	UNP Q9UPQ3
D	68	LEU	-	expression tag	UNP Q9UPQ3
D	69	GLU	-	expression tag	UNP Q9UPQ3
E	64	GLY	-	expression tag	UNP Q9UPQ3
E	65	SER	-	expression tag	UNP Q9UPQ3
E	66	HIS	-	expression tag	UNP Q9UPQ3
E	67	MET	-	expression tag	UNP Q9UPQ3
E	68	LEU	-	expression tag	UNP Q9UPQ3
E	69	GLU	-	expression tag	UNP Q9UPQ3
F	64	GLY	-	expression tag	UNP Q9UPQ3
F	65	SER	-	expression tag	UNP Q9UPQ3
F	66	HIS	-	expression tag	UNP Q9UPQ3
F	67	MET	-	expression tag	UNP Q9UPQ3
F	68	LEU	-	expression tag	UNP Q9UPQ3
F	69	GLU	-	expression tag	UNP Q9UPQ3
G	64	GLY	-	expression tag	UNP Q9UPQ3
G	65	SER	-	expression tag	UNP Q9UPQ3
G	66	HIS	-	expression tag	UNP Q9UPQ3
G	67	MET	-	expression tag	UNP Q9UPQ3
G	68	LEU	-	expression tag	UNP Q9UPQ3
G	69	GLU	-	expression tag	UNP Q9UPQ3
H	64	GLY	-	expression tag	UNP Q9UPQ3
H	65	SER	-	expression tag	UNP Q9UPQ3
H	66	HIS	-	expression tag	UNP Q9UPQ3
H	67	MET	-	expression tag	UNP Q9UPQ3
H	68	LEU	-	expression tag	UNP Q9UPQ3
H	69	GLU	-	expression tag	UNP Q9UPQ3

- Molecule 2 is GUANOSINE-5'-DIPHOSPHATE (CCD ID: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	B	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	B	1	Total	C	H	N	O	P	0	0
			38	10	10	5	11	2		
2	D	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	D	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	E	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	E	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	F	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		
2	G	1	Total	C	H	N	O	P	0	0
			37	10	9	5	11	2		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	5	Total	Mg	0	0
			5	5		
3	D	5	Total	Mg	0	0
			5	5		
3	E	6	Total	Mg	0	0
			6	6		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	2	Total	Mg	0	0
			2	2		
3	G	3	Total	Mg	0	0
			3	3		

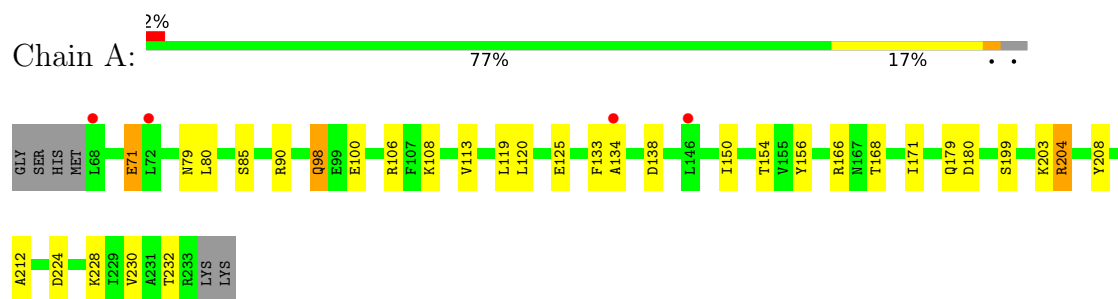
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		
4	B	9	Total	O	0	0
			9	9		
4	C	5	Total	O	0	0
			5	5		
4	D	7	Total	O	0	0
			7	7		
4	E	17	Total	O	0	0
			17	17		
4	F	9	Total	O	0	0
			9	9		
4	G	9	Total	O	0	0
			9	9		
4	H	1	Total	O	0	0
			1	1		

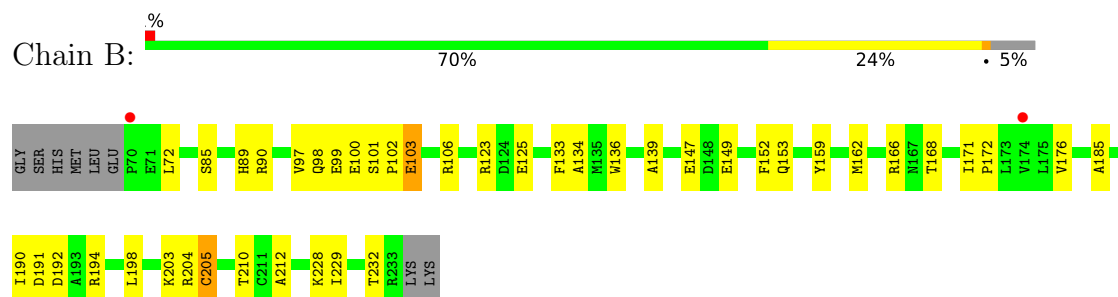
3 Residue-property plots [i](#)

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.

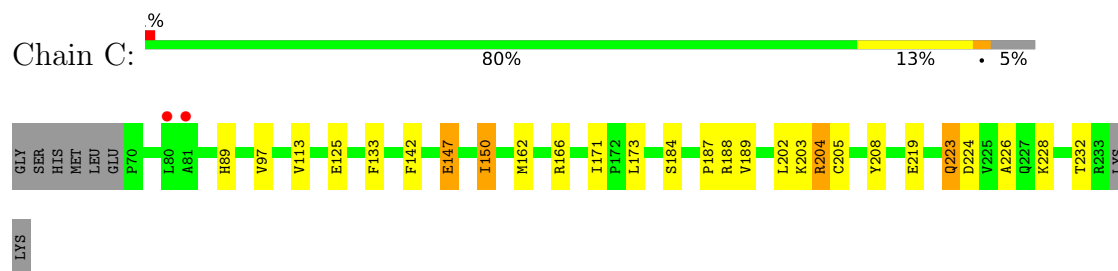
- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



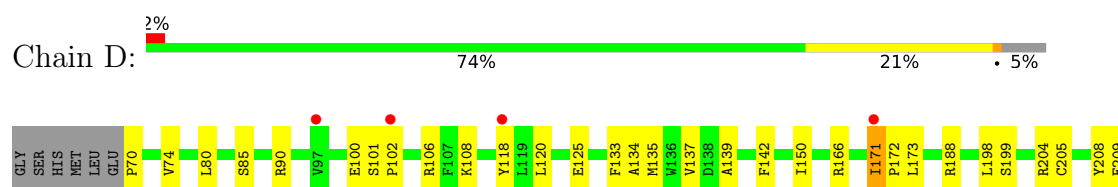
- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1

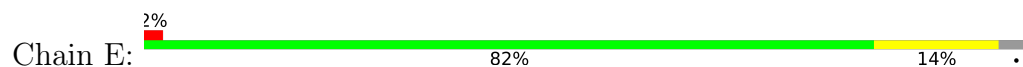


- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1

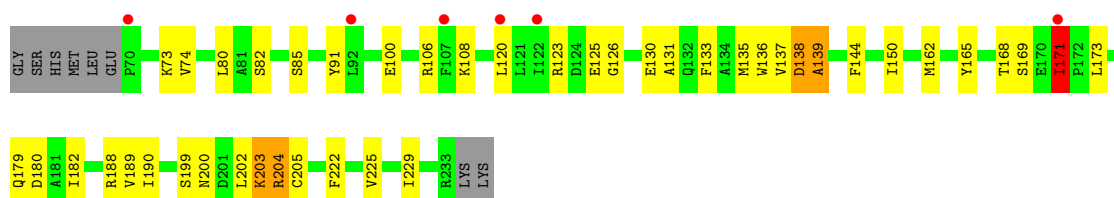




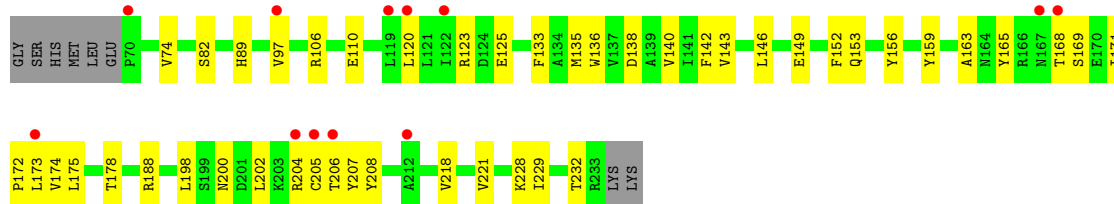
- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



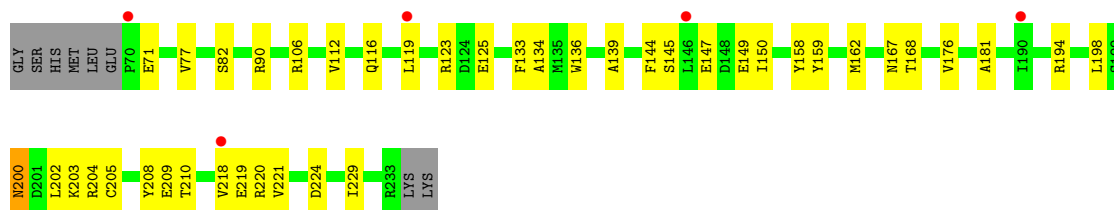
- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



- Molecule 1: Arf-GAP with GTPase, ANK repeat and PH domain-containing protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.41Å 124.68Å 127.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.17 – 2.49 35.17 – 2.49	Depositor EDS
% Data completeness (in resolution range)	95.8 (35.17-2.49) 96.2 (35.17-2.49)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.48Å)	Xtriage
Refinement program	PHENIX (1.14 _3260: ???)	Depositor
R, R_{free}	0.219 , 0.268 0.228 , 0.266	Depositor DCC
R_{free} test set	3145 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	67.1	Xtriage
Anisotropy	0.140	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.025 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21066	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.71	0/1337	0.89	1/1813 (0.1%)
1	B	0.75	0/1323	0.93	3/1793 (0.2%)
1	C	0.79	1/1323 (0.1%)	0.96	1/1793 (0.1%)
1	D	0.69	0/1323	0.93	1/1793 (0.1%)
1	E	0.65	0/1332	0.89	3/1806 (0.2%)
1	F	0.69	0/1323	0.97	3/1793 (0.2%)
1	G	0.63	0/1323	0.91	0/1793
1	H	0.71	1/1323 (0.1%)	0.97	3/1793 (0.2%)
All	All	0.71	2/10607 (0.0%)	0.93	15/14377 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	150	ILE	CG1-CD1	6.52	1.77	1.51
1	H	200	ASN	CG-ND2	5.19	1.44	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	171	ILE	N-CA-C	8.59	121.28	107.71
1	E	205	CYS	N-CA-C	6.92	120.79	107.44
1	B	205	CYS	N-CA-C	6.62	119.80	108.02
1	F	138	ASP	CB-CA-C	-6.03	101.71	111.06
1	D	171	ILE	N-CA-C	5.89	121.61	108.88
1	H	71	GLU	CB-CA-C	-5.67	99.96	109.48
1	H	150	ILE	CA-CB-CG1	-5.60	100.88	110.40
1	A	98	GLN	CA-CB-CG	-5.50	103.10	114.10
1	E	228	LYS	CG-CD-CE	5.50	123.94	111.30
1	B	103	GLU	CA-CB-CG	5.47	125.04	114.10
1	F	106	ARG	CA-CB-CG	5.40	124.91	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	186	ASN	N-CA-CB	-5.26	103.48	111.21
1	C	184	SER	CB-CA-C	5.19	120.64	110.67
1	B	203	LYS	CB-CG-CD	5.15	123.15	111.30
1	H	159	TYR	N-CA-C	-5.09	105.64	111.14

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	1312	1292	1290	30	0
1	B	1298	1283	1283	41	0
1	C	1298	1283	1283	25	0
1	D	1298	1283	1283	33	0
1	E	1307	1288	1287	16	0
1	F	1298	1283	1283	41	0
1	G	1298	1282	1281	60	0
1	H	1298	1283	1283	40	0
2	B	56	19	22	6	0
2	D	56	18	23	6	0
2	E	56	18	22	3	0
2	F	28	9	10	0	0
2	G	28	9	11	3	0
3	B	5	0	0	0	0
3	D	5	0	0	0	0
3	E	6	0	0	0	0
3	F	2	0	0	0	0
3	G	3	0	0	0	0
4	A	7	0	0	4	0
4	B	9	0	0	6	0
4	C	5	0	0	2	0
4	D	7	0	0	2	0
4	E	17	0	0	2	0
4	F	9	0	0	3	0
4	G	9	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	1	0	0	1	0
All	All	10716	10350	10361	291	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 14.

All (291) 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:150:ILE:CD1	1:C:150:ILE:CG1	1.77	1.57
1:G:173:LEU:N	1:G:204:ARG:O	1.58	1.33
1:B:185:ALA:O	1:C:204:ARG:NH1	1.63	1.30
1:G:174:VAL:C	1:G:205:CYS:SG	2.28	1.15
1:G:173:LEU:HB2	1:G:205:CYS:HB2	1.25	1.15
1:G:175:LEU:N	1:G:205:CYS:SG	2.22	1.12
1:D:135:MET:HA	1:D:166:ARG:HH21	1.19	1.04
1:G:173:LEU:CB	1:G:205:CYS:HB2	1.89	1.01
2:B:301:GDP:H5''	2:B:302:GDP:HN21	1.25	0.99
1:G:174:VAL:CA	1:G:205:CYS:SG	2.52	0.97
1:E:166:ARG:NE	1:E:171:ILE:HD11	1.83	0.94
1:G:173:LEU:CA	1:G:204:ARG:O	2.18	0.92
1:H:198:LEU:HD12	1:H:202:LEU:HD13	1.50	0.91
1:E:69:GLU:N	4:E:401:HOH:O	2.02	0.91
1:D:135:MET:HA	1:D:166:ARG:NH2	1.87	0.89
2:D:301:GDP:C2	2:D:302:GDP:C2	2.59	0.89
1:H:106:ARG:HD2	1:H:123:ARG:CZ	2.03	0.88
2:E:301:GDP:H5''	2:E:302:GDP:HN21	1.39	0.87
1:F:139:ALA:N	1:F:171:ILE:HD11	1.91	0.85
1:G:146:LEU:HD22	1:G:188:ARG:NH2	1.93	0.82
1:B:204:ARG:HH12	1:B:232:THR:HG21	1.43	0.81
1:F:135:MET:HE2	1:F:165:TYR:O	1.80	0.81
1:B:103:GLU:OE1	4:B:401:HOH:O	1.97	0.79
1:A:138:ASP:C	1:A:171:ILE:HD11	2.08	0.79
1:A:134:ALA:O	1:A:166:ARG:HG2	1.83	0.79
1:F:108:LYS:HG2	2:G:401:GDP:O3'	1.84	0.77
2:D:301:GDP:N2	2:D:302:GDP:N2	2.32	0.77
1:A:106:ARG:NH2	4:A:302:HOH:O	2.16	0.76
1:G:163:ALA:HB2	1:G:168:THR:HG21	1.66	0.76
1:H:198:LEU:HD12	1:H:202:LEU:CD1	2.15	0.75
1:G:174:VAL:HA	1:G:205:CYS:SG	2.27	0.74
1:G:169:SER:O	1:G:204:ARG:NE	2.21	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:ASP:OD2	4:A:301:HOH:O	2.05	0.74
1:H:168:THR:OG1	4:H:301:HOH:O	2.05	0.74
2:B:301:GDP:C5'	2:B:302:GDP:HN21	2.01	0.74
1:G:140:VAL:HB	1:G:173:LEU:CD2	2.19	0.73
1:F:168:THR:OG1	4:F:401:HOH:O	2.07	0.72
2:D:301:GDP:N2	2:D:302:GDP:C2	2.58	0.72
1:G:110:GLU:O	4:G:501:HOH:O	2.07	0.71
1:B:168:THR:O	1:B:171:ILE:HG22	1.91	0.71
2:B:301:GDP:H5''	2:B:302:GDP:N2	2.04	0.70
1:B:204:ARG:HH12	1:B:232:THR:CG2	2.04	0.70
1:G:106:ARG:HD2	1:G:123:ARG:NH2	2.06	0.70
1:B:205:CYS:O	4:B:402:HOH:O	2.09	0.70
1:B:204:ARG:HH22	1:B:232:THR:HG23	1.55	0.70
1:C:89:HIS:ND1	1:C:97:VAL:HG12	2.07	0.70
1:E:167:ASN:OD1	1:E:169:SER:HB2	1.91	0.70
1:D:134:ALA:O	1:D:166:ARG:HG3	1.92	0.69
1:F:139:ALA:H	1:F:171:ILE:HD11	1.56	0.69
1:B:101:SER:C	4:B:401:HOH:O	2.35	0.69
1:B:147:GLU:O	1:C:203:LYS:NZ	2.27	0.68
1:D:173:LEU:HB2	1:D:205:CYS:HB3	1.75	0.68
1:F:168:THR:O	1:F:171:ILE:HG22	1.93	0.68
1:C:150:ILE:CD1	1:C:150:ILE:CB	2.71	0.67
1:H:200:ASN:HB3	1:H:204:ARG:HE	1.60	0.67
1:H:202:LEU:O	1:H:203:LYS:HB2	1.93	0.66
1:A:108:LYS:HE3	4:A:302:HOH:O	1.95	0.66
1:B:103:GLU:CD	4:B:401:HOH:O	2.39	0.66
1:A:208:TYR:OH	1:A:224:ASP:OD2	2.14	0.66
1:A:134:ALA:O	1:A:166:ARG:CG	2.43	0.65
1:H:198:LEU:O	1:H:202:LEU:HD13	1.97	0.65
1:B:176:VAL:HG13	1:B:210:THR:HG23	1.78	0.65
1:A:138:ASP:O	1:A:171:ILE:HD11	1.97	0.65
2:B:302:GDP:H5'	2:B:302:GDP:O2B	1.97	0.65
1:H:198:LEU:CD1	1:H:202:LEU:CD1	2.76	0.64
1:D:204:ARG:O	1:D:204:ARG:HG3	1.98	0.64
1:H:106:ARG:HD2	1:H:123:ARG:NH1	2.13	0.64
1:F:202:LEU:O	4:F:402:HOH:O	2.15	0.64
1:G:135:MET:HE1	1:G:165:TYR:O	1.99	0.63
1:G:172:PRO:HA	1:G:204:ARG:HB3	1.79	0.63
1:C:166:ARG:HG2	1:C:171:ILE:HD11	1.80	0.63
1:H:106:ARG:CG	1:H:123:ARG:HG2	2.29	0.63
1:A:71:GLU:HB3	1:A:119:LEU:HB3	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:ALA:O	1:B:166:ARG:NH2	2.33	0.61
1:G:106:ARG:HE	1:G:123:ARG:NH1	1.99	0.61
1:D:171:ILE:HG23	1:D:171:ILE:O	2.00	0.61
1:D:137:VAL:HG23	1:D:171:ILE:CD1	2.31	0.61
1:A:199:SER:O	1:A:204:ARG:HA	2.00	0.60
1:B:85:SER:HB2	1:B:100:GLU:CD	2.26	0.60
1:D:139:ALA:HA	1:D:171:ILE:HG13	1.83	0.60
1:H:198:LEU:CD1	1:H:202:LEU:HD13	2.29	0.60
1:A:168:THR:O	1:A:171:ILE:HG22	2.02	0.60
1:F:80:LEU:HD11	1:F:126:GLY:O	2.02	0.60
1:E:166:ARG:CZ	1:E:171:ILE:HD11	2.30	0.60
1:G:82:SER:HB2	1:G:143:VAL:HG12	1.84	0.60
1:F:222:PHE:O	1:F:225:VAL:HG12	2.01	0.59
1:F:108:LYS:HE2	2:G:401:GDP:O3'	2.02	0.59
1:G:89:HIS:HD2	1:G:97:VAL:HB	1.67	0.59
1:G:163:ALA:HB2	1:G:168:THR:CG2	2.33	0.59
1:D:80:LEU:HD21	1:D:102:PRO:HA	1.85	0.58
2:E:301:GDP:C5'	2:E:302:GDP:HN21	2.14	0.58
1:F:138:ASP:O	1:F:229:ILE:HG12	2.04	0.58
1:G:159:TYR:OH	1:G:168:THR:HB	2.02	0.58
1:A:138:ASP:O	1:A:171:ILE:CD1	2.51	0.58
1:G:173:LEU:CB	1:G:204:ARG:O	2.51	0.58
1:G:156:TYR:CE2	1:G:198:LEU:HD13	2.39	0.58
1:C:147:GLU:C	1:C:147:GLU:OE1	2.46	0.58
1:G:159:TYR:CZ	1:G:168:THR:HB	2.38	0.58
1:H:106:ARG:HG2	1:H:123:ARG:HG2	1.86	0.57
1:C:219:GLU:HG3	1:C:223:GLN:NE2	2.19	0.57
1:F:85:SER:N	1:F:100:GLU:OE2	2.32	0.57
1:B:106:ARG:HD3	1:B:123:ARG:NH1	2.20	0.57
1:D:135:MET:CA	1:D:166:ARG:NH2	2.64	0.57
1:G:146:LEU:HD22	1:G:188:ARG:CZ	2.34	0.57
1:G:174:VAL:C	1:G:205:CYS:HG	2.10	0.57
1:F:171:ILE:HG23	1:F:171:ILE:O	2.04	0.57
1:G:174:VAL:N	1:G:205:CYS:SG	2.78	0.57
1:F:137:VAL:HG23	1:F:171:ILE:HD13	1.86	0.57
1:B:185:ALA:O	1:C:204:ARG:CZ	2.46	0.56
1:B:204:ARG:HH22	1:B:232:THR:CG2	2.19	0.56
1:G:140:VAL:HB	1:G:173:LEU:HD22	1.87	0.56
1:D:137:VAL:HG23	1:D:171:ILE:HD12	1.87	0.56
1:C:89:HIS:ND1	1:C:97:VAL:CG1	2.68	0.56
1:D:227:GLN:O	1:D:230:VAL:HG22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:123:ARG:HG3	1:B:136:TRP:CH2	2.41	0.55
2:B:302:GDP:O2B	2:B:302:GDP:C5'	2.54	0.55
1:B:89:HIS:ND1	1:B:97:VAL:HG12	2.22	0.55
1:H:200:ASN:OD1	1:H:204:ARG:HD3	2.07	0.55
1:E:82:SER:HB2	1:E:143:VAL:HG12	1.88	0.55
1:F:173:LEU:O	1:F:205:CYS:HB2	2.07	0.55
1:F:200:ASN:HA	1:F:204:ARG:HA	1.89	0.55
1:B:147:GLU:OE2	1:C:203:LYS:NZ	2.39	0.55
1:C:228:LYS:O	1:C:232:THR:HG23	2.08	0.54
1:D:139:ALA:HA	1:D:171:ILE:CG1	2.37	0.54
1:G:135:MET:CE	1:G:165:TYR:O	2.56	0.54
1:H:200:ASN:CB	1:H:204:ARG:HE	2.19	0.54
1:D:142:PHE:HE2	1:D:173:LEU:HD22	1.72	0.54
1:F:182:ILE:HD13	1:F:188:ARG:HB2	1.90	0.54
1:C:187:PRO:O	1:C:189:VAL:HG13	2.07	0.54
1:H:198:LEU:CD1	1:H:202:LEU:HD11	2.38	0.54
1:B:101:SER:HB3	4:B:401:HOH:O	2.07	0.54
1:D:125:GLU:HG3	1:D:133:PHE:CD1	2.42	0.53
1:D:188:ARG:NH2	1:D:209:GLU:OE2	2.41	0.53
1:G:173:LEU:HB2	1:G:205:CYS:CB	2.17	0.53
1:A:228:LYS:O	1:A:232:THR:HG23	2.08	0.53
1:G:146:LEU:HD22	1:G:188:ARG:HH21	1.73	0.53
1:A:113:VAL:HG11	1:A:230:VAL:HG21	1.90	0.53
1:H:176:VAL:HG13	1:H:210:THR:HG23	1.91	0.53
1:F:139:ALA:HA	1:F:171:ILE:CG1	2.39	0.53
1:G:142:PHE:HE2	1:G:173:LEU:HD22	1.73	0.53
1:F:130:GLU:HA	1:F:130:GLU:OE1	2.08	0.52
1:D:90:ARG:HG2	1:D:218:VAL:HG21	1.91	0.52
1:G:163:ALA:HA	1:G:168:THR:HG23	1.91	0.52
1:G:173:LEU:C	1:G:205:CYS:SG	2.93	0.52
1:D:172:PRO:HB3	1:D:228:LYS:HB3	1.92	0.52
1:G:146:LEU:CD2	1:G:188:ARG:NH2	2.69	0.52
1:G:200:ASN:O	1:G:202:LEU:O	2.28	0.51
1:H:147:GLU:HG2	1:H:181:ALA:HB3	1.92	0.51
1:F:135:MET:CE	1:F:165:TYR:O	2.54	0.51
1:A:125:GLU:HG3	1:A:133:PHE:CD1	2.46	0.51
1:D:198:LEU:O	1:D:199:SER:C	2.54	0.51
1:G:152:PHE:CE1	1:G:198:LEU:HD22	2.45	0.51
1:C:113:VAL:HG21	1:C:226:ALA:HB1	1.93	0.51
1:F:85:SER:HB2	1:F:100:GLU:OE2	2.11	0.51
1:C:203:LYS:HB2	4:C:303:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:GLN:O	1:A:98:GLN:HG2	2.11	0.50
1:D:108:LYS:NZ	4:D:402:HOH:O	2.44	0.50
1:D:135:MET:C	1:D:166:ARG:HH22	2.20	0.50
1:B:159:TYR:CD2	1:F:150:ILE:HG13	2.47	0.50
1:F:125:GLU:HG3	1:F:133:PHE:CG	2.47	0.50
1:B:90:ARG:HD3	1:B:212:ALA:O	2.12	0.49
1:B:172:PRO:HB3	1:B:228:LYS:HB3	1.93	0.49
2:B:301:GDP:C5'	2:B:302:GDP:N2	2.68	0.49
1:E:110:GLU:CD	4:E:406:HOH:O	2.54	0.49
1:A:79:ASN:OD1	1:A:79:ASN:C	2.55	0.49
1:G:149:GLU:OE2	1:G:153:GLN:NE2	2.45	0.49
1:H:139:ALA:HB3	1:H:229:ILE:HD11	1.94	0.49
1:G:138:ASP:O	1:G:172:PRO:HD2	2.12	0.49
1:G:146:LEU:HD12	1:G:178:THR:O	2.12	0.49
1:H:147:GLU:CG	1:H:181:ALA:HB3	2.41	0.49
1:H:149:GLU:CD	1:H:194:ARG:HH12	2.21	0.49
1:H:82:SER:HB2	1:H:144:PHE:HA	1.95	0.49
1:C:142:PHE:HE2	1:C:173:LEU:HD22	1.78	0.49
1:H:112:VAL:HA	1:H:116:GLN:O	2.12	0.49
1:H:134:ALA:HA	1:H:162:MET:CE	2.43	0.49
1:G:175:LEU:HD21	1:G:198:LEU:HD23	1.95	0.48
1:H:106:ARG:HG3	1:H:123:ARG:HG2	1.94	0.48
1:B:192:ASP:OD1	1:B:192:ASP:N	2.46	0.48
1:C:97:VAL:HG13	1:C:97:VAL:O	2.13	0.48
1:F:189:VAL:HG23	1:F:190:ILE:CD1	2.42	0.48
1:F:137:VAL:HG23	1:F:171:ILE:CD1	2.43	0.48
2:E:301:GDP:H5''	2:E:302:GDP:N2	2.20	0.48
1:A:156:TYR:OH	1:E:150:ILE:HD11	2.14	0.48
1:D:135:MET:C	1:D:166:ARG:NH2	2.72	0.47
1:G:171:ILE:HG23	1:G:171:ILE:O	2.14	0.47
1:B:152:PHE:CE1	1:B:198:LEU:HD22	2.49	0.47
1:G:74:VAL:CG2	1:G:120:LEU:HD11	2.44	0.47
1:F:179:GLN:O	1:F:180:ASP:C	2.55	0.47
1:A:138:ASP:CA	1:A:171:ILE:HD11	2.44	0.47
1:D:205:CYS:HA	4:D:405:HOH:O	2.15	0.47
1:G:206:THR:HG22	1:G:207:TYR:N	2.29	0.47
1:H:208:TYR:CE1	1:H:220:ARG:HD3	2.49	0.47
1:H:125:GLU:HG3	1:H:133:PHE:CD1	2.50	0.47
1:B:149:GLU:OE2	1:B:153:GLN:NE2	2.47	0.47
1:F:82:SER:HB2	1:F:144:PHE:HA	1.97	0.47
1:H:77:VAL:HG12	1:H:158:TYR:HD2	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:140:VAL:HB	1:G:173:LEU:HD23	1.97	0.46
1:E:166:ARG:HE	1:E:171:ILE:HD11	1.73	0.46
1:B:139:ALA:HB3	1:B:229:ILE:HD11	1.97	0.46
1:F:203:LYS:NZ	4:F:401:HOH:O	2.48	0.46
1:D:188:ARG:HH22	1:D:209:GLU:CD	2.23	0.46
1:H:198:LEU:HD12	1:H:198:LEU:O	2.15	0.46
1:B:229:ILE:O	1:B:232:THR:OG1	2.34	0.46
1:F:74:VAL:HG21	1:F:120:LEU:HD11	1.98	0.46
1:G:89:HIS:CD2	1:G:97:VAL:HB	2.50	0.46
1:C:173:LEU:O	1:C:205:CYS:HB2	2.16	0.45
1:C:188:ARG:NH1	4:C:301:HOH:O	2.29	0.45
1:F:125:GLU:HG3	1:F:133:PHE:CD1	2.51	0.45
1:B:125:GLU:HG3	1:B:133:PHE:CD1	2.51	0.45
1:D:227:GLN:O	1:D:230:VAL:CG2	2.65	0.45
1:A:90:ARG:HD3	1:A:212:ALA:O	2.17	0.45
1:A:166:ARG:HD2	4:A:304:HOH:O	2.17	0.45
1:D:106:ARG:HB2	2:D:302:GDP:C8	2.51	0.45
1:G:125:GLU:HG3	1:G:133:PHE:CD1	2.52	0.45
1:H:145:SER:OG	1:H:147:GLU:HG3	2.16	0.45
1:C:219:GLU:HG3	1:C:223:GLN:HE22	1.81	0.45
1:C:125:GLU:HG3	1:C:133:PHE:CD1	2.52	0.45
1:D:85:SER:HB2	1:D:100:GLU:CD	2.41	0.45
1:G:169:SER:O	1:G:204:ARG:CD	2.65	0.45
1:A:125:GLU:CG	1:A:133:PHE:CD1	3.00	0.45
1:F:131:ALA:HB2	1:F:165:TYR:CD2	2.51	0.45
1:G:172:PRO:HA	1:G:204:ARG:CB	2.45	0.44
1:B:106:ARG:HG2	1:B:123:ARG:HG2	1.98	0.44
1:E:145:SER:OG	1:E:147:GLU:HG3	2.18	0.44
1:C:202:LEU:O	1:C:203:LYS:HB2	2.17	0.44
1:G:106:ARG:HD2	1:G:123:ARG:CZ	2.47	0.44
1:C:162:MET:HE3	1:C:162:MET:HB3	1.90	0.44
1:D:208:TYR:OH	1:D:224:ASP:OD2	2.31	0.44
1:G:228:LYS:O	1:G:232:THR:HG23	2.17	0.44
1:G:106:ARG:HG2	1:G:123:ARG:HG2	1.98	0.44
1:G:138:ASP:HB3	1:G:229:ILE:HD13	1.99	0.44
1:B:72:LEU:HD22	1:B:229:ILE:CG2	2.48	0.44
1:G:171:ILE:O	1:G:173:LEU:HG	2.18	0.44
1:F:139:ALA:HA	1:F:171:ILE:HG13	1.99	0.43
1:G:74:VAL:HG23	1:G:120:LEU:HD11	1.99	0.43
1:B:106:ARG:HD3	1:B:123:ARG:HH11	1.83	0.43
1:B:152:PHE:CG	1:B:190:ILE:HD12	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:89:HIS:CD2	1:G:97:VAL:CG2	3.01	0.43
1:B:191:ASP:OD1	1:B:194:ARG:HG3	2.18	0.43
1:G:123:ARG:HD2	1:G:136:TRP:CZ2	2.54	0.43
2:D:301:GDP:HN22	2:D:302:GDP:N2	2.12	0.43
1:E:72:LEU:HD12	1:E:111:ILE:HD11	2.01	0.43
2:D:301:GDP:C2	2:D:302:GDP:N1	2.87	0.43
1:B:152:PHE:CE1	1:B:198:LEU:CD2	3.02	0.43
1:H:208:TYR:HE1	1:H:220:ARG:HD3	1.84	0.43
1:F:91:TYR:HB2	1:F:222:PHE:CE2	2.54	0.42
1:G:208:TYR:CD2	1:G:221:VAL:HG22	2.54	0.42
1:A:179:GLN:O	1:A:180:ASP:C	2.63	0.42
1:B:102:PRO:N	4:B:401:HOH:O	2.51	0.42
1:G:172:PRO:C	1:G:173:LEU:HG	2.44	0.42
1:G:198:LEU:O	1:G:202:LEU:HD12	2.20	0.42
1:H:208:TYR:OH	1:H:224:ASP:OD2	2.26	0.42
1:H:202:LEU:O	1:H:203:LYS:CB	2.66	0.42
1:H:219:GLU:HG3	1:H:219:GLU:O	2.19	0.42
1:C:208:TYR:OH	1:C:224:ASP:OD2	2.30	0.42
1:F:162:MET:HE2	1:F:162:MET:HB3	1.86	0.42
1:B:134:ALA:HA	1:B:162:MET:HE3	2.01	0.42
1:A:85:SER:HB2	1:A:100:GLU:CD	2.45	0.42
1:E:163:ALA:HA	1:E:168:THR:HG23	2.01	0.42
1:A:108:LYS:HA	1:A:120:LEU:O	2.20	0.42
1:F:169:SER:OG	1:F:203:LYS:NZ	2.51	0.42
1:F:199:SER:O	1:F:204:ARG:HA	2.20	0.42
1:D:74:VAL:HG21	1:D:120:LEU:HD11	2.02	0.41
1:A:150:ILE:O	1:A:154:THR:OG1	2.33	0.41
1:A:166:ARG:HD2	1:A:166:ARG:HA	1.84	0.41
1:F:125:GLU:CG	1:F:133:PHE:CD1	3.03	0.41
1:B:159:TYR:CZ	1:B:168:THR:HB	2.56	0.41
1:H:200:ASN:CG	1:H:204:ARG:NE	2.79	0.41
1:H:198:LEU:HD11	1:H:202:LEU:HD11	2.02	0.41
1:A:113:VAL:CG1	1:A:230:VAL:HG21	2.51	0.41
1:C:142:PHE:CE2	1:C:173:LEU:HD22	2.55	0.41
1:F:108:LYS:CG	2:G:401:GDP:O3'	2.62	0.41
1:B:159:TYR:HD2	1:F:150:ILE:HG13	1.86	0.41
1:D:224:ASP:OD1	1:D:224:ASP:C	2.63	0.41
1:H:123:ARG:HD2	1:H:136:TRP:CZ2	2.56	0.41
1:H:147:GLU:HG2	1:H:181:ALA:CB	2.51	0.41
1:H:167:ASN:OD1	1:H:167:ASN:C	2.64	0.41
1:E:174:VAL:HG11	1:E:221:VAL:HG13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:90:ARG:HG2	1:H:218:VAL:HG21	2.03	0.40
1:A:171:ILE:O	1:A:171:ILE:HG23	2.22	0.40
1:A:203:LYS:HE3	1:E:148:ASP:OD1	2.21	0.40
1:B:72:LEU:HD22	1:B:229:ILE:HG21	2.04	0.40
1:F:123:ARG:HD2	1:F:136:TRP:CZ3	2.57	0.40
1:E:71:GLU:HG2	1:E:119:LEU:HD23	2.03	0.40
1:D:70:PRO:HG2	1:D:118:TYR:CE1	2.56	0.40
1:D:214:TYR:CE2	1:H:209:GLU:HG3	2.57	0.40
1:E:121:LEU:HD23	1:E:121:LEU:HA	1.91	0.40
1:E:167:ASN:C	1:E:169:SER:H	2.29	0.40
1:B:98:GLN:O	1:B:99:GLU:C	2.64	0.40
1:D:173:LEU:O	1:D:205:CYS:HB2	2.22	0.40
1:F:73:LYS:HG2	1:F:136:TRP:CZ2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	164/172 (95%)	160 (98%)	3 (2%)	1 (1%)	21	38
1	B	162/172 (94%)	160 (99%)	2 (1%)	0	100	100
1	C	162/172 (94%)	158 (98%)	4 (2%)	0	100	100
1	D	162/172 (94%)	155 (96%)	7 (4%)	0	100	100
1	E	163/172 (95%)	159 (98%)	4 (2%)	0	100	100
1	F	162/172 (94%)	156 (96%)	4 (2%)	2 (1%)	10	20
1	G	162/172 (94%)	159 (98%)	3 (2%)	0	100	100
1	H	162/172 (94%)	155 (96%)	6 (4%)	1 (1%)	21	38
All	All	1299/1376 (94%)	1262 (97%)	33 (2%)	4 (0%)	36	55

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	204	ARG
1	F	139	ALA
1	F	204	ARG
1	H	205	CYS

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	141/147 (96%)	139 (99%)	2 (1%)	59	81
1	B	140/147 (95%)	140 (100%)	0	100	100
1	C	140/147 (95%)	137 (98%)	3 (2%)	47	74
1	D	140/147 (95%)	138 (99%)	2 (1%)	59	81
1	E	141/147 (96%)	141 (100%)	0	100	100
1	F	140/147 (95%)	138 (99%)	2 (1%)	59	81
1	G	140/147 (95%)	139 (99%)	1 (1%)	76	89
1	H	140/147 (95%)	138 (99%)	2 (1%)	59	81
All	All	1122/1176 (95%)	1110 (99%)	12 (1%)	65	84

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	GLU
1	A	80	LEU
1	C	147	GLU
1	C	204	ARG
1	C	223	GLN
1	D	101	SER
1	D	150	ILE
1	F	171	ILE
1	F	203	LYS
1	G	218	VAL

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Mol	Chain	Res	Type
1	H	119	LEU
1	H	221	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	GLN
1	C	223	GLN
1	D	223	GLN
1	E	200	ASN
1	E	227	GLN
1	F	200	ASN
1	F	223	GLN
1	G	89	HIS
1	G	153	GLN
1	G	167	ASN
1	G	223	GLN
1	H	153	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 21 are monoatomic - leaving 8 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	GDP	B	302	3	28,30,30	1.93	10 (35%)	44,47,47	3.18	18 (40%)
2	GDP	D	301	3	28,30,30	1.13	3 (10%)	44,47,47	1.91	8 (18%)
2	GDP	E	301	3	28,30,30	1.13	3 (10%)	44,47,47	1.91	8 (18%)
2	GDP	F	301	3	28,30,30	1.62	5 (17%)	44,47,47	2.89	17 (38%)
2	GDP	D	302	3	28,30,30	1.12	3 (10%)	44,47,47	1.90	8 (18%)
2	GDP	G	401	3	28,30,30	1.56	6 (21%)	44,47,47	2.99	22 (50%)
2	GDP	E	302	3	28,30,30	1.82	7 (25%)	44,47,47	3.39	17 (38%)
2	GDP	B	301	3	28,30,30	1.13	3 (10%)	44,47,47	1.91	8 (18%)

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	GDP	B	302	3	-	5/16/32/32	0/3/3/3
2	GDP	D	301	3	-	3/16/32/32	0/3/3/3
2	GDP	E	301	3	-	4/16/32/32	0/3/3/3
2	GDP	F	301	3	-	8/16/32/32	0/3/3/3
2	GDP	D	302	3	-	3/16/32/32	0/3/3/3
2	GDP	G	401	3	-	5/16/32/32	0/3/3/3
2	GDP	E	302	3	-	7/16/32/32	0/3/3/3
2	GDP	B	301	3	-	4/16/32/32	0/3/3/3

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	302	GDP	O4'-C1'	5.29	1.54	1.42
2	F	301	GDP	O4'-C1'	5.18	1.54	1.42
2	E	302	GDP	C5-N7	-4.11	1.31	1.39
2	G	401	GDP	O4'-C1'	4.02	1.51	1.42
2	E	302	GDP	O4'-C1'	3.93	1.51	1.42
2	B	302	GDP	C5-N7	-3.35	1.32	1.39
2	F	301	GDP	C6-N1	-3.26	1.32	1.38
2	B	302	GDP	C6-N1	-3.25	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	302	GDP	C6-N1	-3.22	1.32	1.38
2	E	302	GDP	C5-C4	3.11	1.47	1.38
2	F	301	GDP	C5-C4	3.11	1.47	1.38
2	D	301	GDP	C5-C4	3.08	1.47	1.38
2	B	301	GDP	C5-C4	3.06	1.47	1.38
2	G	401	GDP	C6-N1	-3.05	1.33	1.38
2	E	301	GDP	C5-C4	3.05	1.47	1.38
2	D	302	GDP	C5-C4	3.04	1.47	1.38
2	G	401	GDP	C5-N7	-3.01	1.33	1.39
2	E	302	GDP	PA-O5'	2.77	1.70	1.59
2	G	401	GDP	C5-C4	2.74	1.46	1.38
2	E	302	GDP	PA-O1A	2.69	1.60	1.50
2	B	302	GDP	C5-C4	2.67	1.46	1.38
2	F	301	GDP	C5-N7	-2.67	1.33	1.39
2	B	302	GDP	PA-O5'	2.53	1.69	1.59
2	B	302	GDP	PA-O1A	2.46	1.59	1.50
2	B	301	GDP	C6-N1	-2.42	1.34	1.38
2	D	301	GDP	C6-N1	-2.40	1.34	1.38
2	B	302	GDP	C4-N9	-2.39	1.31	1.38
2	E	301	GDP	C6-N1	-2.39	1.34	1.38
2	D	302	GDP	C6-N1	-2.38	1.34	1.38
2	E	302	GDP	C4-N9	-2.23	1.32	1.38
2	B	302	GDP	O4'-C4'	2.22	1.50	1.45
2	B	301	GDP	C5-N7	-2.18	1.34	1.39
2	G	401	GDP	C2'-C1'	-2.13	1.46	1.53
2	D	302	GDP	C5-N7	-2.13	1.35	1.39
2	E	301	GDP	C5-N7	-2.12	1.35	1.39
2	D	301	GDP	C5-N7	-2.10	1.35	1.39
2	G	401	GDP	C4-N9	-2.10	1.32	1.38
2	B	302	GDP	C2-N1	-2.03	1.32	1.37
2	B	302	GDP	C2'-C3'	2.01	1.58	1.53
2	F	301	GDP	O6-C6	2.01	1.27	1.23

All (106) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	302	GDP	C2'-C1'-N9	12.59	148.90	113.22
2	B	302	GDP	C2'-C1'-N9	9.64	140.54	113.22
2	E	302	GDP	O4'-C4'-C5'	9.27	139.87	109.37
2	F	301	GDP	C2'-C3'-C4'	8.36	118.89	102.64
2	F	301	GDP	C2'-C1'-N9	7.85	135.46	113.22
2	G	401	GDP	C2'-C1'-N9	7.51	134.50	113.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	GDP	O4'-C4'-C5'	6.44	130.57	109.37
2	B	302	GDP	C2'-C3'-C4'	6.44	115.15	102.64
2	G	401	GDP	C2'-C3'-C4'	6.41	115.09	102.64
2	D	301	GDP	C5-C4-N3	-6.19	118.42	128.46
2	B	301	GDP	C5-C4-N3	-6.17	118.44	128.46
2	D	302	GDP	C5-C4-N3	-6.17	118.45	128.46
2	E	301	GDP	C5-C4-N3	-6.16	118.46	128.46
2	E	302	GDP	O4'-C4'-C3'	-6.14	92.96	105.11
2	F	301	GDP	C5-C4-N3	-6.13	118.51	128.46
2	B	302	GDP	C5-C4-N3	-6.08	118.60	128.46
2	B	302	GDP	O4'-C4'-C5'	6.03	129.21	109.37
2	E	302	GDP	C5-C4-N3	-5.76	119.12	128.46
2	B	302	GDP	O4'-C4'-C3'	-5.45	94.33	105.11
2	G	401	GDP	C5-C4-N3	-5.34	119.81	128.46
2	B	302	GDP	C1'-N9-C8	-5.28	111.68	126.70
2	E	302	GDP	C2'-C3'-C4'	5.07	112.50	102.64
2	F	301	GDP	N9-C4-N3	4.97	135.92	125.94
2	B	302	GDP	C1'-N9-C4	4.97	141.26	126.50
2	E	302	GDP	C1'-N9-C8	-4.97	112.56	126.70
2	B	302	GDP	N9-C4-N3	4.95	135.89	125.94
2	G	401	GDP	N9-C4-N3	4.92	135.82	125.94
2	B	301	GDP	C2-N3-C4	4.90	121.04	112.30
2	D	302	GDP	C2-N3-C4	4.90	121.04	112.30
2	D	301	GDP	C2-N3-C4	4.90	121.03	112.30
2	E	301	GDP	C2-N3-C4	4.88	121.00	112.30
2	F	301	GDP	C1'-N9-C8	-4.82	112.98	126.70
2	D	301	GDP	N9-C4-N3	4.68	135.33	125.94
2	B	301	GDP	N9-C4-N3	4.66	135.29	125.94
2	E	301	GDP	N9-C4-N3	4.62	135.21	125.94
2	D	302	GDP	N9-C4-N3	4.61	135.20	125.94
2	E	302	GDP	N9-C4-N3	4.56	135.09	125.94
2	E	302	GDP	C1'-N9-C4	4.42	139.62	126.50
2	G	401	GDP	O3'-C3'-C2'	-4.41	97.56	111.82
2	G	401	GDP	C1'-N9-C8	-4.34	114.35	126.70
2	G	401	GDP	C5'-C4'-C3'	-4.00	100.17	115.18
2	B	302	GDP	C4'-O4'-C1'	3.98	118.27	109.47
2	B	302	GDP	O5'-PA-O1A	3.95	124.50	109.07
2	F	301	GDP	C1'-N9-C4	3.93	138.19	126.50
2	G	401	GDP	O6-C6-C5	-3.90	116.26	126.60
2	B	302	GDP	C2-N3-C4	3.88	119.21	112.30
2	B	302	GDP	O2A-PA-O1A	-3.80	93.44	112.24
2	F	301	GDP	C2-N3-C4	3.80	119.06	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	GDP	O4'-C4'-C3'	-3.79	97.62	105.11
2	F	301	GDP	O3'-C3'-C4'	-3.69	100.38	111.05
2	E	301	GDP	PA-O3A-PB	-3.58	120.53	132.83
2	B	301	GDP	PA-O3A-PB	-3.57	120.57	132.83
2	G	401	GDP	C2-N3-C4	3.57	118.65	112.30
2	G	401	GDP	O2'-C2'-C1'	-3.56	98.13	110.02
2	D	301	GDP	PA-O3A-PB	-3.55	120.64	132.83
2	F	301	GDP	O4'-C4'-C3'	-3.55	98.09	105.11
2	F	301	GDP	O4'-C4'-C5'	3.53	120.97	109.37
2	D	302	GDP	PA-O3A-PB	-3.50	120.83	132.83
2	G	401	GDP	C4'-O4'-C1'	3.37	116.91	109.47
2	B	302	GDP	O6-C6-C5	-3.22	118.05	126.60
2	G	401	GDP	O5'-PA-O1A	3.19	121.53	109.07
2	E	301	GDP	C6-C5-N7	3.18	136.17	130.25
2	D	302	GDP	C6-C5-N7	3.18	136.16	130.25
2	B	301	GDP	C6-C5-N7	3.17	136.15	130.25
2	D	301	GDP	C6-C5-N7	3.15	136.10	130.25
2	E	302	GDP	C2-N3-C4	3.09	117.81	112.30
2	G	401	GDP	C1'-N9-C4	3.08	135.65	126.50
2	B	302	GDP	O3B-PB-O3A	2.96	114.55	104.64
2	G	401	GDP	O2B-PB-O3A	2.88	114.29	104.64
2	F	301	GDP	N2-C2-N3	-2.82	114.24	119.73
2	E	302	GDP	C4'-O4'-C1'	2.80	115.66	109.47
2	D	302	GDP	C4-C5-N7	-2.68	106.48	110.72
2	E	301	GDP	C4-C5-N7	-2.68	106.48	110.72
2	B	302	GDP	O6-C6-N1	2.68	125.25	120.12
2	D	301	GDP	C4-C5-N7	-2.66	106.52	110.72
2	B	301	GDP	C4-C5-N7	-2.64	106.53	110.72
2	E	302	GDP	C2-N1-C6	-2.63	120.31	125.10
2	E	302	GDP	O2B-PB-O3A	2.61	113.39	104.64
2	G	401	GDP	C2-N1-C6	-2.52	120.50	125.10
2	F	301	GDP	O6-C6-C5	-2.51	119.94	126.60
2	E	302	GDP	O2A-PA-O5'	2.46	119.19	107.75
2	D	301	GDP	C3'-C2'-C1'	2.46	106.11	101.43
2	E	301	GDP	C3'-C2'-C1'	2.45	106.08	101.43
2	D	302	GDP	C3'-C2'-C1'	2.44	106.07	101.43
2	B	301	GDP	C3'-C2'-C1'	2.44	106.06	101.43
2	E	302	GDP	C5'-C4'-C3'	-2.38	106.25	115.18
2	G	401	GDP	C5-C6-N1	2.35	119.16	113.19
2	B	302	GDP	C2-N1-C6	-2.35	120.81	125.10
2	G	401	GDP	O6-C6-N1	2.32	124.57	120.12
2	G	401	GDP	O3'-C3'-C4'	2.30	117.70	111.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	302	GDP	O5'-C5'-C4'	-2.27	101.16	108.99
2	G	401	GDP	N2-C2-N3	-2.26	115.33	119.73
2	F	301	GDP	C5'-C4'-C3'	2.17	123.30	115.18
2	B	301	GDP	O6-C6-C5	-2.16	120.87	126.60
2	E	301	GDP	O6-C6-C5	-2.16	120.88	126.60
2	D	302	GDP	O6-C6-C5	-2.14	120.92	126.60
2	D	301	GDP	O6-C6-C5	-2.14	120.92	126.60
2	F	301	GDP	C3'-C2'-C1'	-2.14	97.37	101.43
2	B	302	GDP	O2'-C2'-C1'	-2.11	102.97	110.02
2	F	301	GDP	O2A-PA-O5'	2.09	117.44	107.75
2	E	302	GDP	N1-C2-N3	2.08	127.20	123.32
2	B	302	GDP	O2B-PB-O1B	-2.07	102.56	110.68
2	F	301	GDP	O3'-C3'-C2'	-2.06	105.17	111.82
2	F	301	GDP	O5'-C5'-C4'	-2.03	101.99	108.99
2	G	401	GDP	C6-C5-N7	2.02	134.01	130.25
2	E	302	GDP	O2'-C2'-C3'	-2.01	105.32	111.82

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301	GDP	C5'-O5'-PA-O2A
2	B	302	GDP	C5'-O5'-PA-O3A
2	E	301	GDP	C5'-O5'-PA-O1A
2	E	301	GDP	C5'-O5'-PA-O2A
2	E	302	GDP	PA-O3A-PB-O3B
2	E	302	GDP	C5'-O5'-PA-O3A
2	F	301	GDP	PA-O3A-PB-O3B
2	F	301	GDP	C5'-O5'-PA-O3A
2	F	301	GDP	C3'-C4'-C5'-O5'
2	G	401	GDP	C5'-O5'-PA-O3A
2	G	401	GDP	C5'-O5'-PA-O2A
2	B	302	GDP	O4'-C4'-C5'-O5'
2	E	302	GDP	O4'-C4'-C5'-O5'
2	E	302	GDP	C3'-C4'-C5'-O5'
2	G	401	GDP	O4'-C4'-C5'-O5'
2	G	401	GDP	C3'-C4'-C5'-O5'
2	B	302	GDP	C3'-C4'-C5'-O5'
2	F	301	GDP	O4'-C4'-C5'-O5'
2	F	301	GDP	O4'-C1'-N9-C4
2	B	302	GDP	PB-O3A-PA-O5'
2	B	301	GDP	C5'-O5'-PA-O3A

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Mol	Chain	Res	Type	Atoms
2	B	301	GDP	C5'-O5'-PA-O1A
2	B	302	GDP	C5'-O5'-PA-O2A
2	E	302	GDP	C5'-O5'-PA-O2A
2	F	301	GDP	C5'-O5'-PA-O1A
2	F	301	GDP	C5'-O5'-PA-O2A
2	G	401	GDP	C5'-O5'-PA-O1A
2	D	302	GDP	C3'-C4'-C5'-O5'
2	D	301	GDP	C4'-C5'-O5'-PA
2	B	301	GDP	O4'-C4'-C5'-O5'
2	E	301	GDP	O4'-C4'-C5'-O5'
2	E	302	GDP	PA-O3A-PB-O1B
2	D	302	GDP	PB-O3A-PA-O1A
2	D	301	GDP	O4'-C4'-C5'-O5'
2	D	301	GDP	PA-O3A-PB-O2B
2	F	301	GDP	PA-O3A-PB-O2B
2	E	301	GDP	C5'-O5'-PA-O3A
2	D	302	GDP	PB-O3A-PA-O2A
2	E	302	GDP	C5'-O5'-PA-O1A

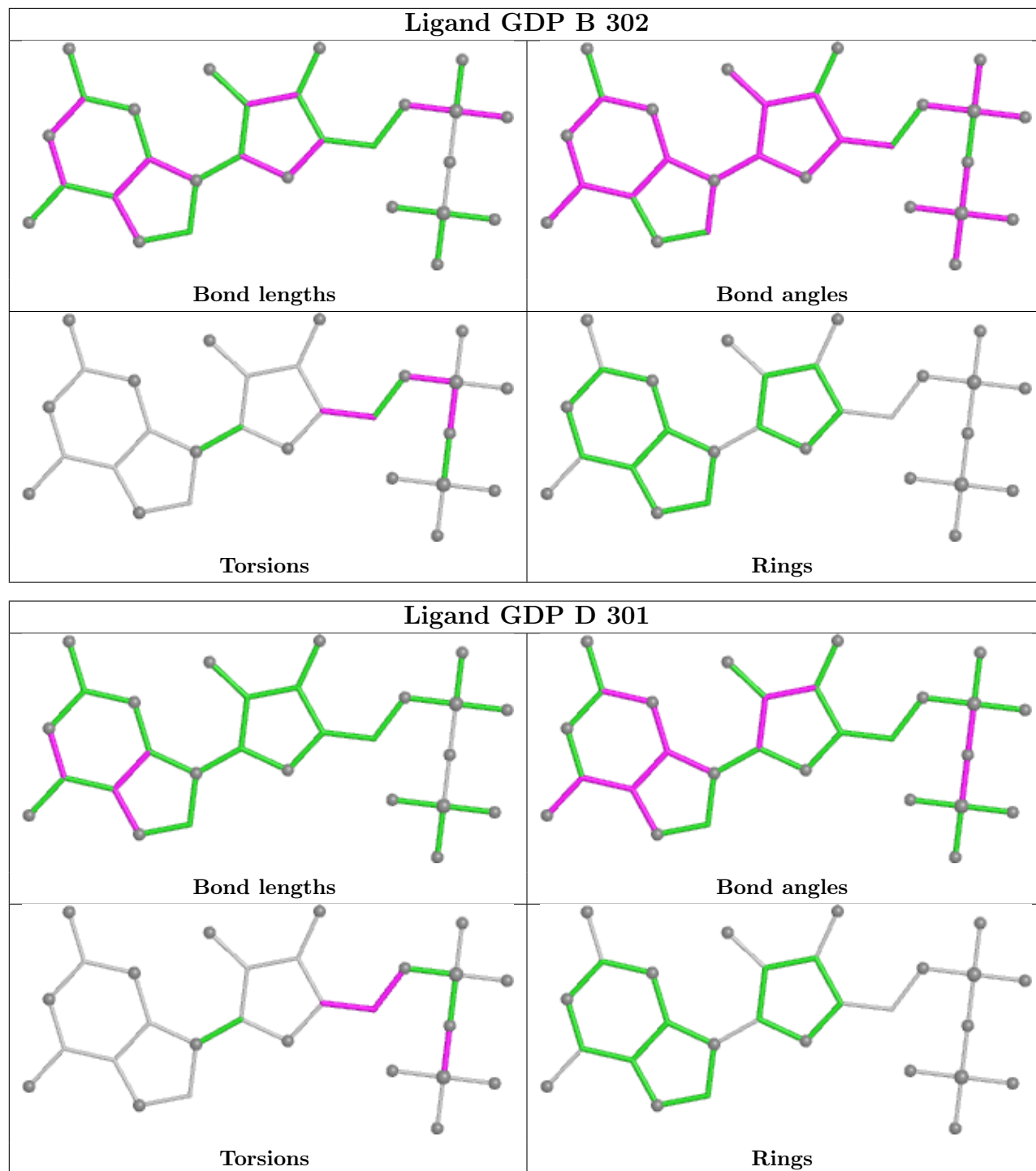
There are no ring outliers.

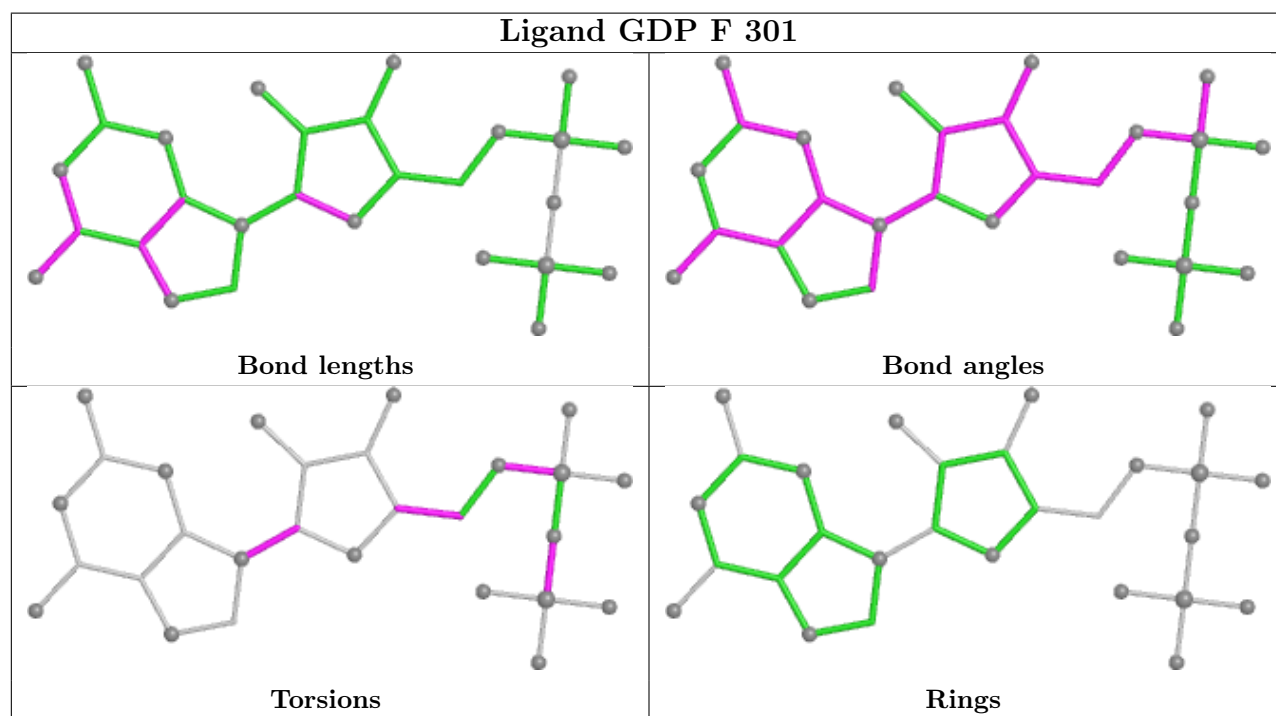
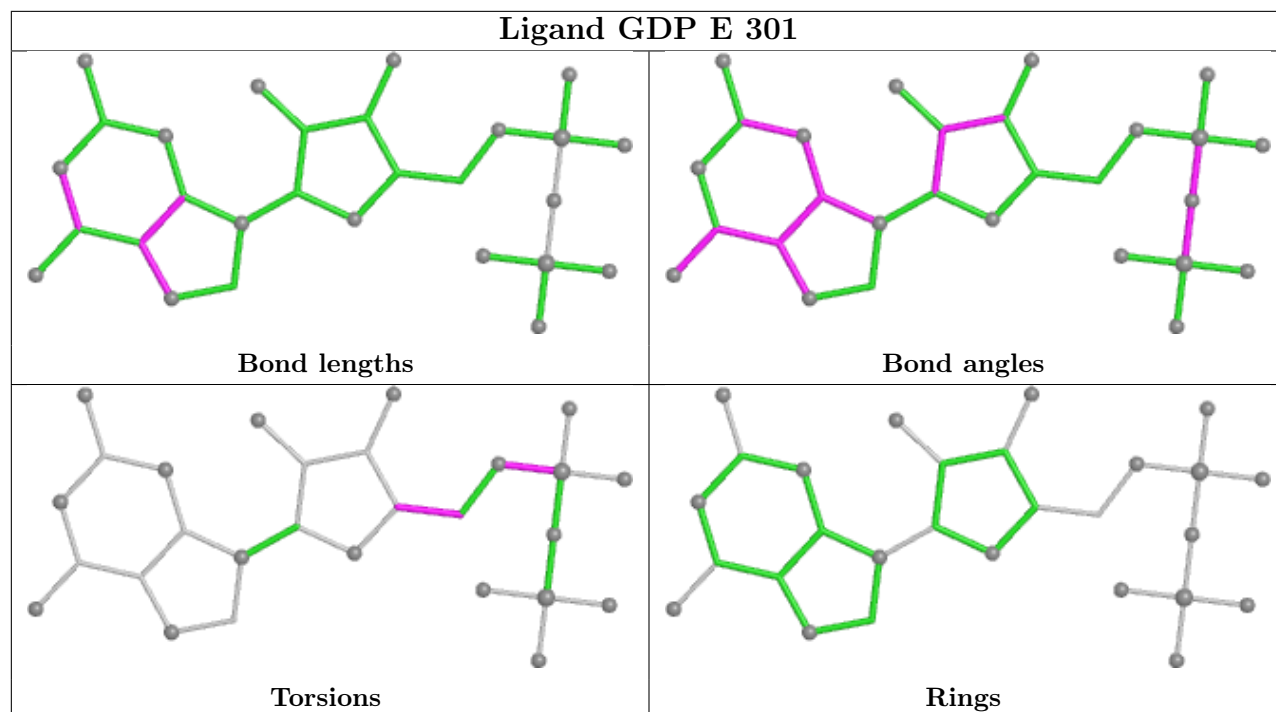
7 monomers are involved in 18 short contacts:

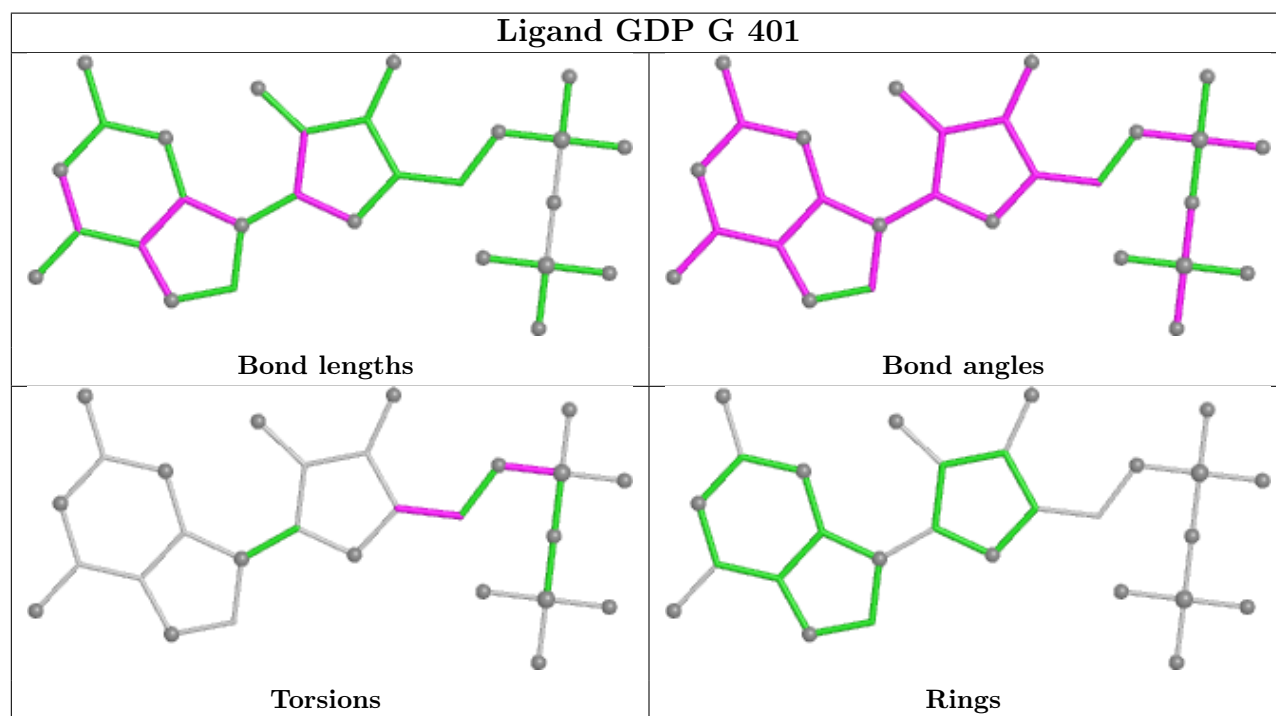
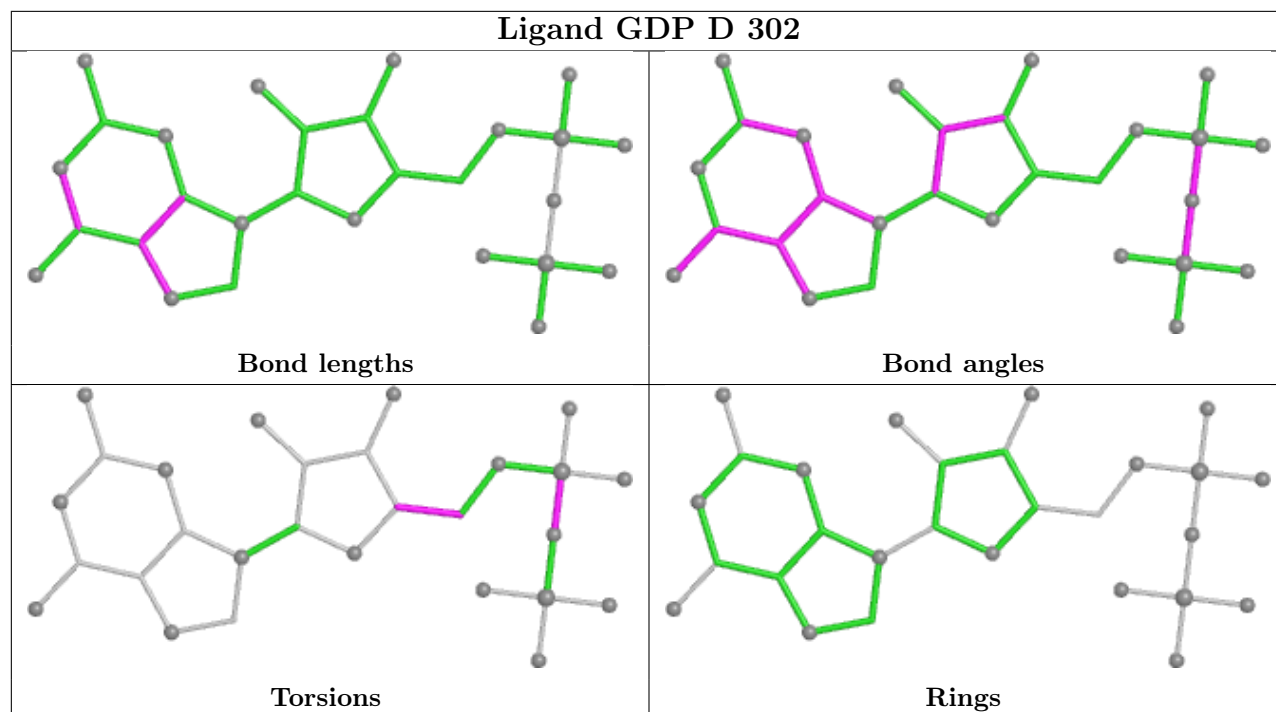
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	302	GDP	6	0
2	D	301	GDP	5	0
2	E	301	GDP	3	0
2	D	302	GDP	6	0
2	G	401	GDP	3	0
2	E	302	GDP	3	0
2	B	301	GDP	4	0

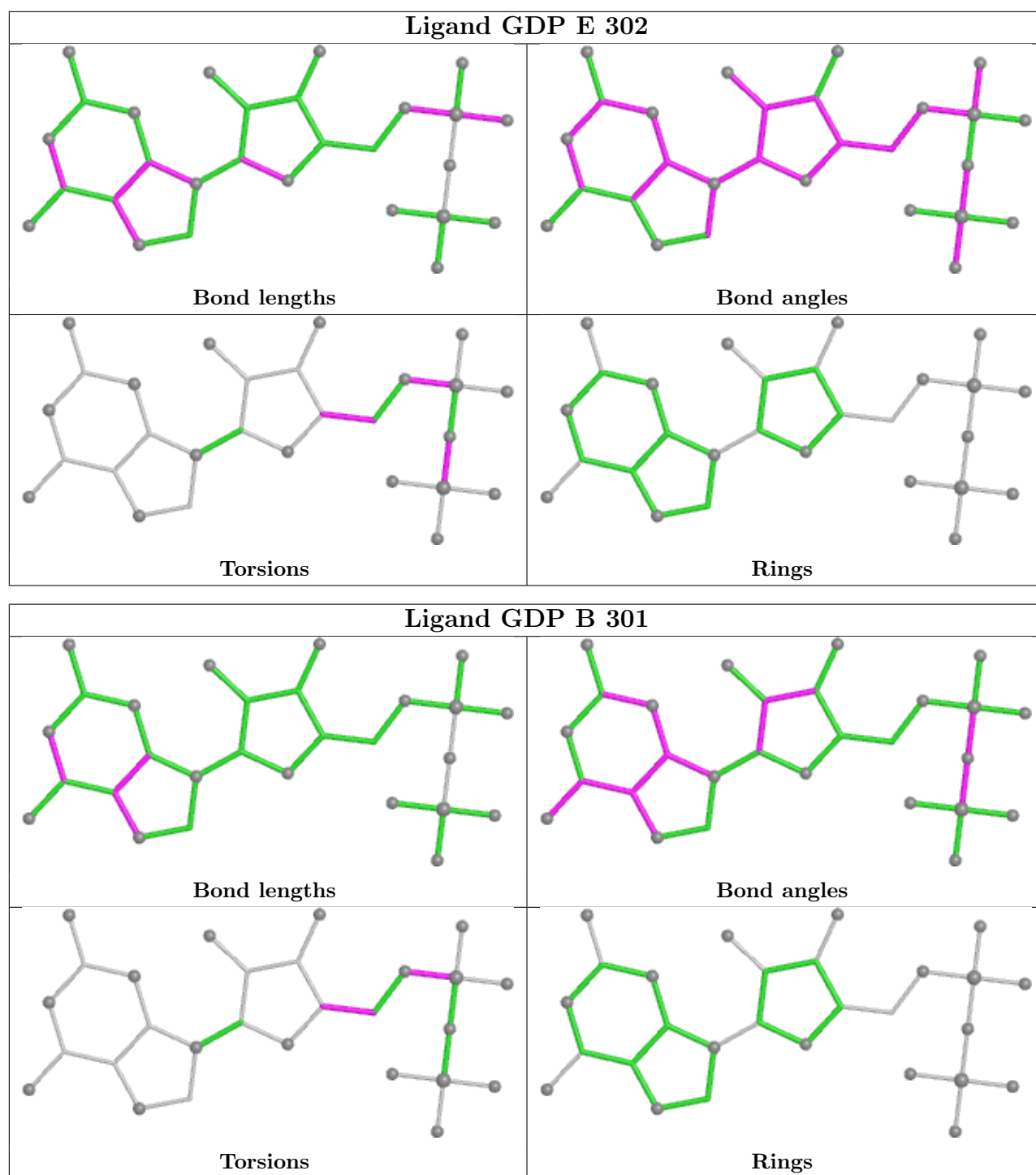
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	166/172 (96%)	0.17	4 (2%) 59 55	55, 77, 118, 135	0
1	B	164/172 (95%)	0.03	2 (1%) 76 73	52, 75, 110, 145	0
1	C	164/172 (95%)	0.22	2 (1%) 76 73	53, 78, 118, 146	0
1	D	164/172 (95%)	0.16	4 (2%) 59 55	55, 81, 126, 142	0
1	E	165/172 (95%)	0.35	3 (1%) 67 64	59, 82, 120, 139	0
1	F	164/172 (95%)	0.33	6 (3%) 45 40	57, 82, 125, 140	0
1	G	164/172 (95%)	0.54	12 (7%) 21 19	60, 88, 132, 152	0
1	H	164/172 (95%)	0.40	5 (3%) 52 48	59, 86, 132, 151	0
All	All	1315/1376 (95%)	0.28	38 (2%) 53 49	52, 81, 126, 152	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	205	CYS	5.8
1	H	146	LEU	3.9
1	C	80	LEU	3.7
1	G	168	THR	3.6
1	E	214	TYR	3.6
1	H	70	PRO	3.5
1	F	171	ILE	3.5
1	A	146	LEU	3.4
1	F	120	LEU	3.1
1	F	122	ILE	3.0
1	F	92	LEU	2.9
1	D	102	PRO	2.8
1	H	218	VAL	2.8
1	B	70	PRO	2.8
1	G	122	ILE	2.7
1	D	118	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	68	LEU	2.7
1	B	174	VAL	2.7
1	C	81	ALA	2.7
1	A	134	ALA	2.7
1	E	96	TYR	2.6
1	G	206	THR	2.5
1	H	119	LEU	2.5
1	G	173	LEU	2.4
1	F	107	PHE	2.4
1	G	120	LEU	2.3
1	E	156	TYR	2.3
1	G	70	PRO	2.3
1	G	212	ALA	2.3
1	D	171	ILE	2.2
1	F	70	PRO	2.2
1	H	190	ILE	2.2
1	G	119	LEU	2.2
1	A	72	LEU	2.1
1	G	167	ASN	2.1
1	G	97	VAL	2.1
1	G	204	ARG	2.1
1	D	97	VAL	2.1

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

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

6.3 Carbohydrates [\(i\)](#)

There are no oligosaccharides in this entry.

6.4 Ligands [\(i\)](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GDP	E	301	28/28	0.58	0.18	54,64,92,94	0

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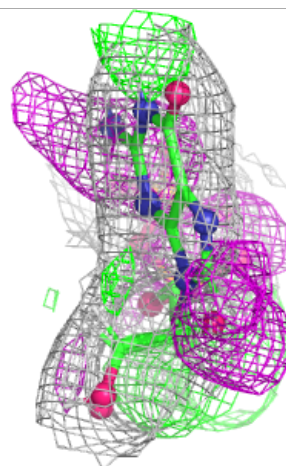
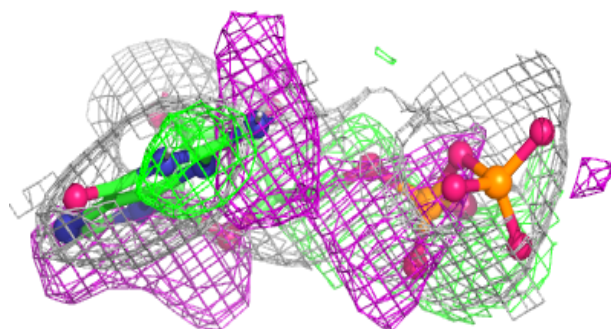
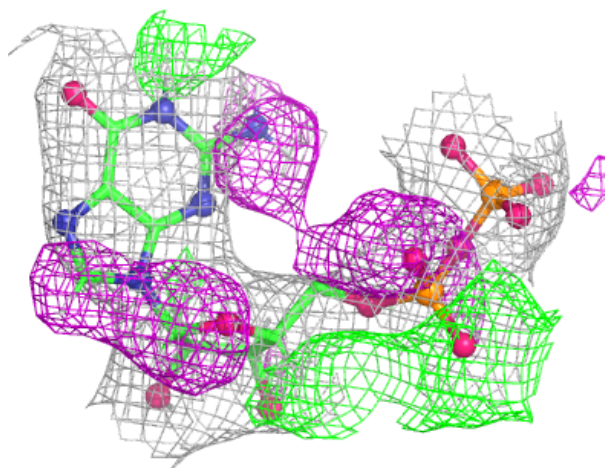
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GDP	B	301	28/28	0.59	0.16	61,70,89,105	0
2	GDP	D	302	28/28	0.61	0.16	62,70,85,93	0
2	GDP	D	301	28/28	0.65	0.16	59,68,82,90	0
2	GDP	E	302	28/28	0.78	0.13	54,65,94,98	0
2	GDP	B	302	28/28	0.79	0.13	54,67,85,95	0
3	MG	G	403	1/1	0.80	0.22	82,82,82,82	0
2	GDP	F	301	28/28	0.81	0.11	59,78,100,103	0
2	GDP	G	401	28/28	0.84	0.10	62,76,96,107	0
3	MG	E	307	1/1	0.87	0.10	54,54,54,54	0
3	MG	D	304	1/1	0.87	0.25	57,57,57,57	0
3	MG	E	303	1/1	0.88	0.08	54,54,54,54	0
3	MG	G	402	1/1	0.89	0.07	60,60,60,60	0
3	MG	D	305	1/1	0.89	0.08	61,61,61,61	0
3	MG	D	307	1/1	0.90	0.15	69,69,69,69	0
3	MG	B	306	1/1	0.91	0.10	66,66,66,66	0
3	MG	E	308	1/1	0.91	0.09	58,58,58,58	0
3	MG	F	303	1/1	0.91	0.07	52,52,52,52	0
3	MG	D	303	1/1	0.91	0.07	50,50,50,50	0
3	MG	B	303	1/1	0.91	0.08	53,53,53,53	0
3	MG	D	306	1/1	0.92	0.15	54,54,54,54	0
3	MG	E	304	1/1	0.92	0.06	52,52,52,52	0
3	MG	E	305	1/1	0.92	0.09	58,58,58,58	0
3	MG	B	307	1/1	0.92	0.20	62,62,62,62	0
3	MG	G	404	1/1	0.94	0.05	50,50,50,50	0
3	MG	B	304	1/1	0.95	0.07	55,55,55,55	0
3	MG	F	302	1/1	0.95	0.08	56,56,56,56	0
3	MG	B	305	1/1	0.96	0.10	52,52,52,52	0
3	MG	E	306	1/1	0.97	0.10	61,61,61,61	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.

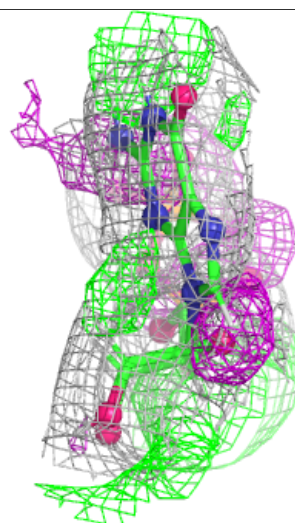
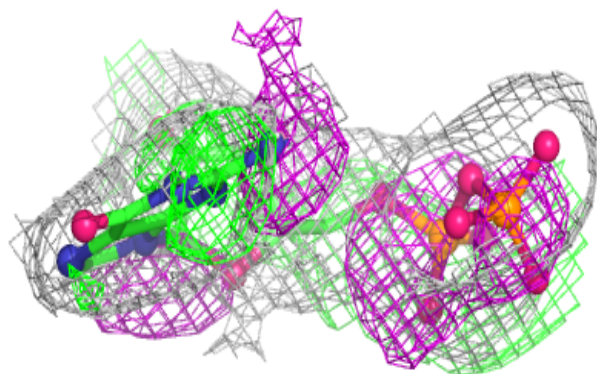
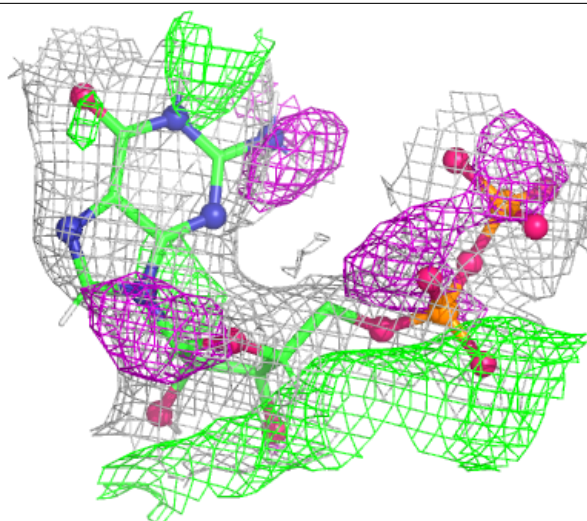
Electron density around GDP E 301:

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



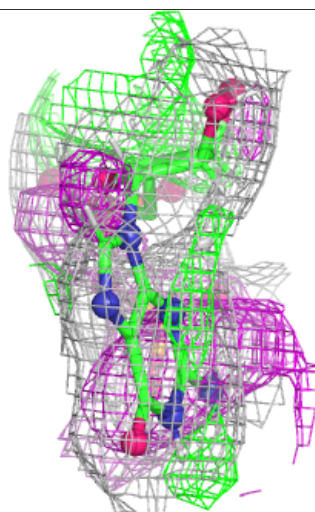
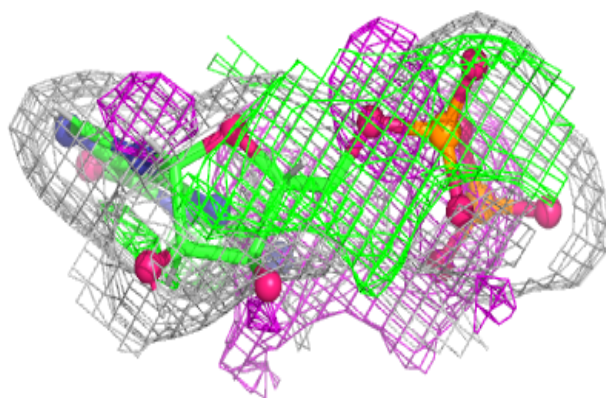
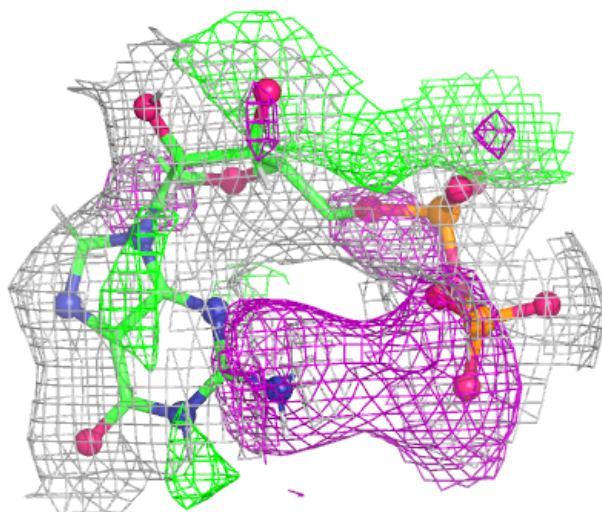
Electron density around GDP B 301:

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



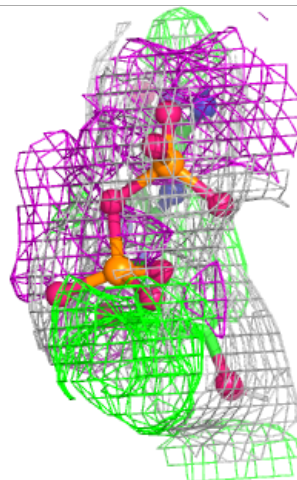
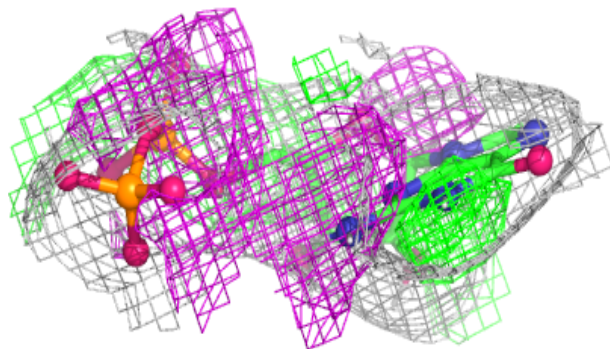
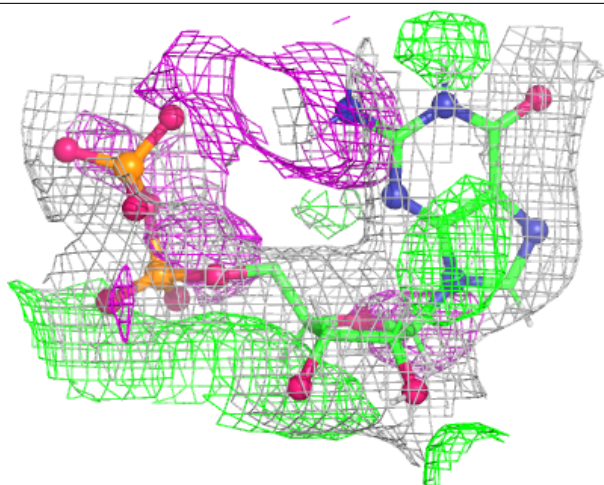
Electron density around GDP D 302:

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and green (positive)



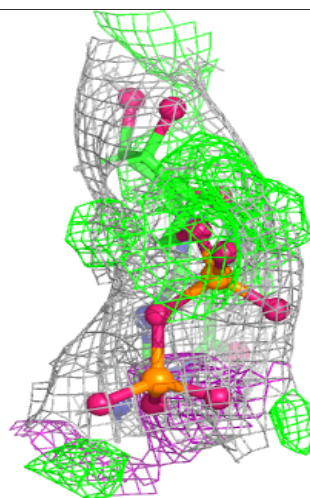
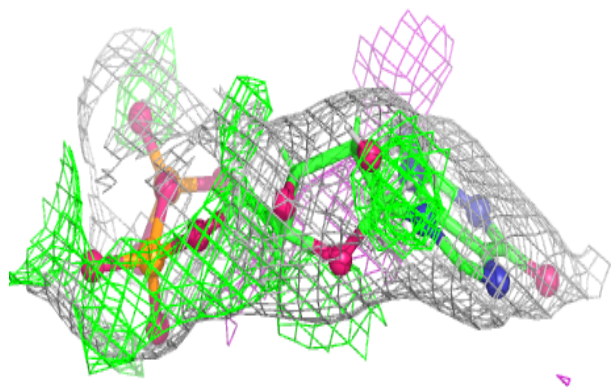
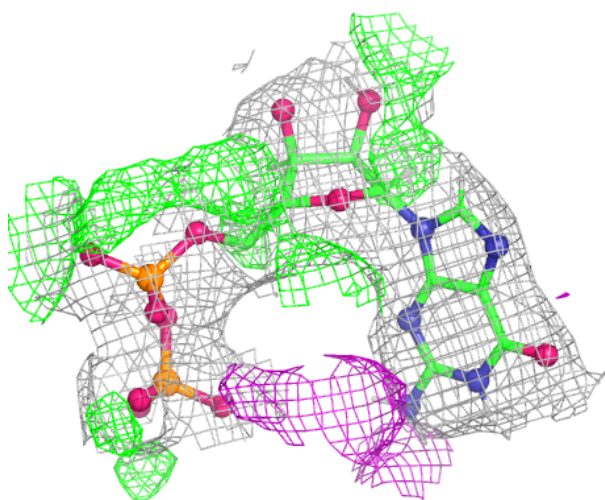
Electron density around GDP D 301:

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



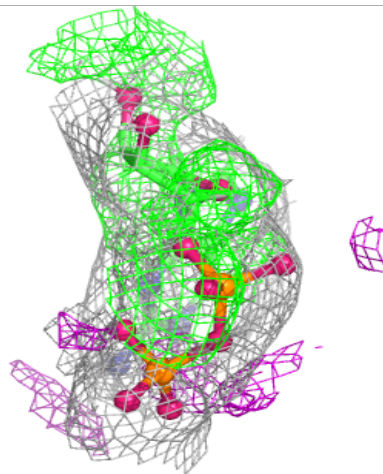
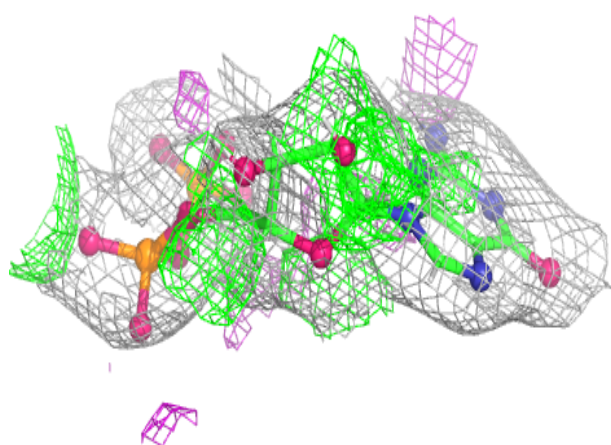
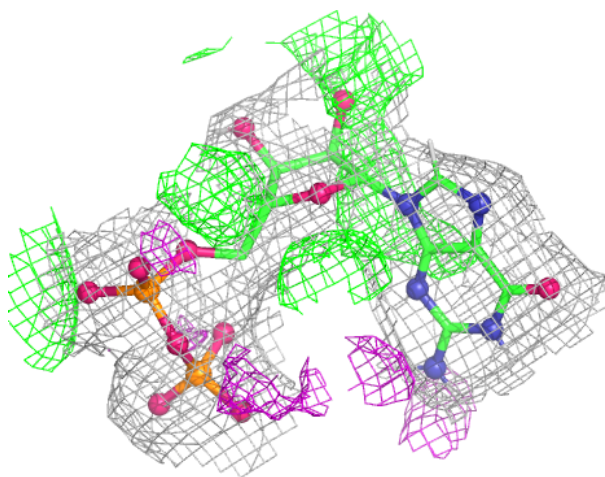
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and green (positive)



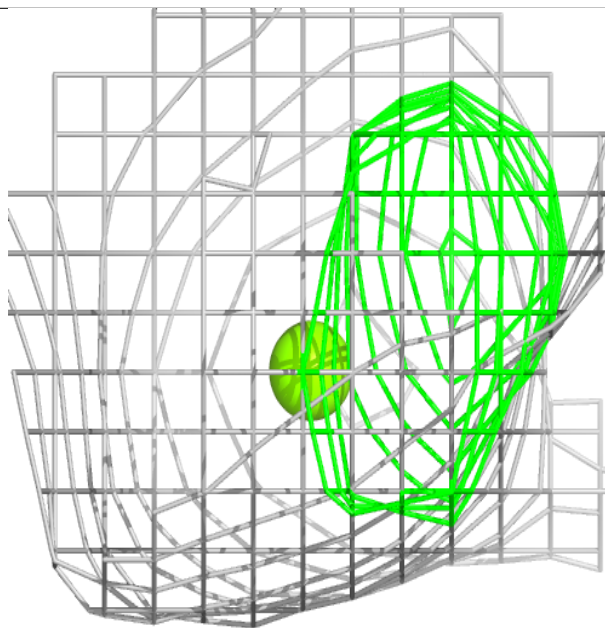
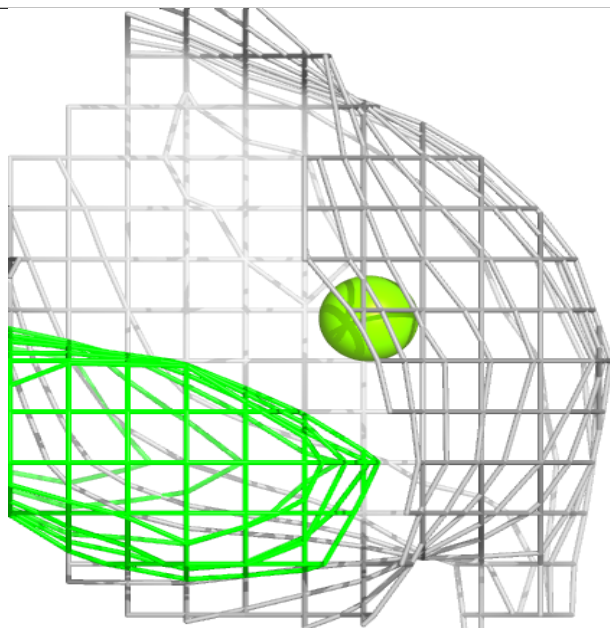
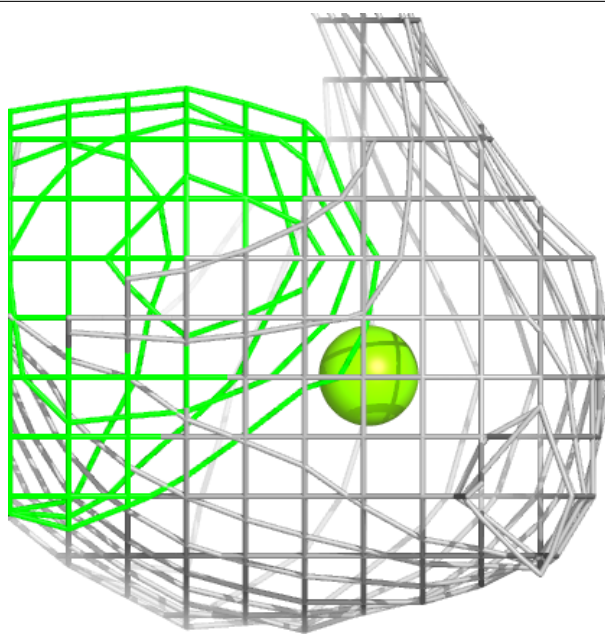
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and green (positive)



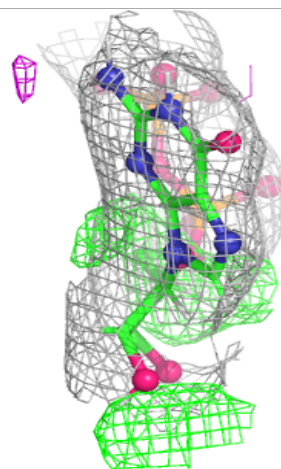
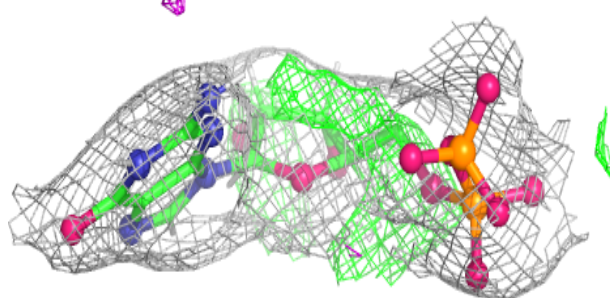
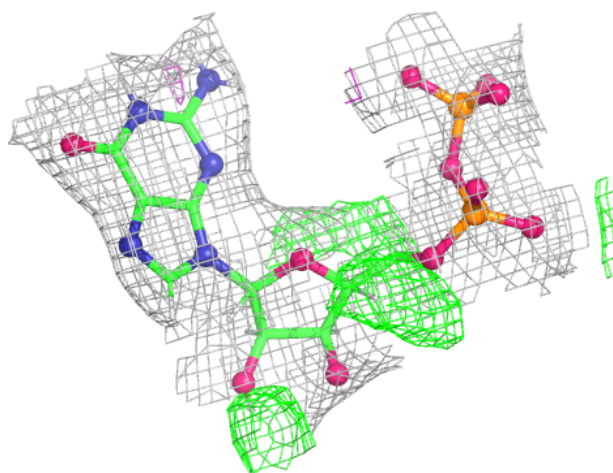
Electron density around MG G 403:

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



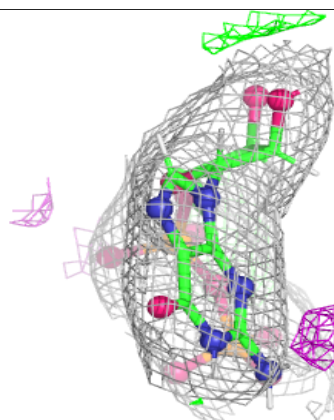
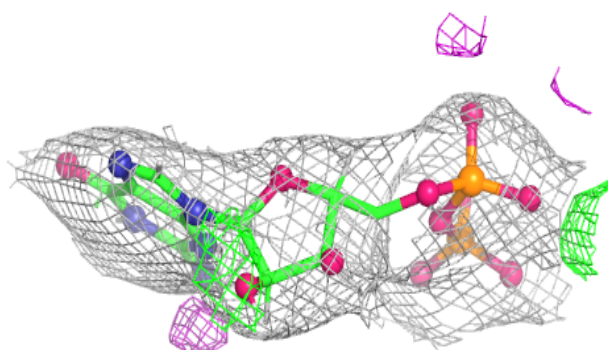
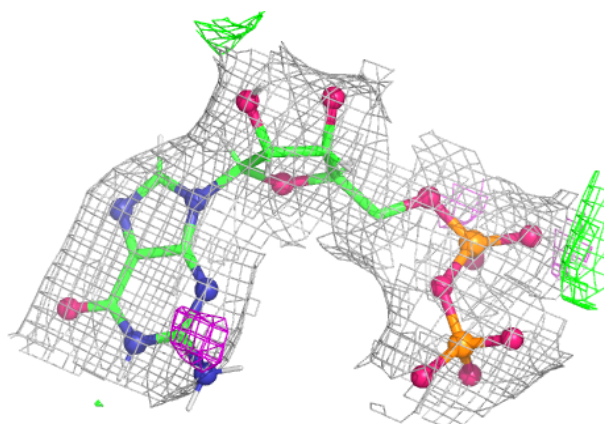
Electron density around GDP F 301:

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and green (positive)



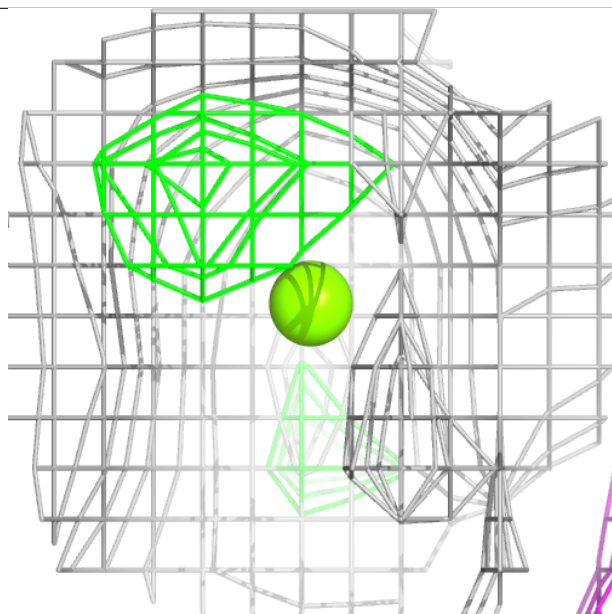
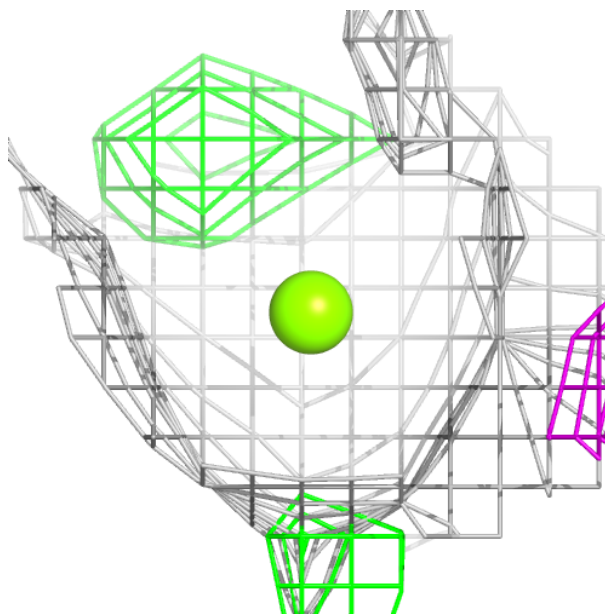
Electron density around GDP G 401:

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and green (positive)



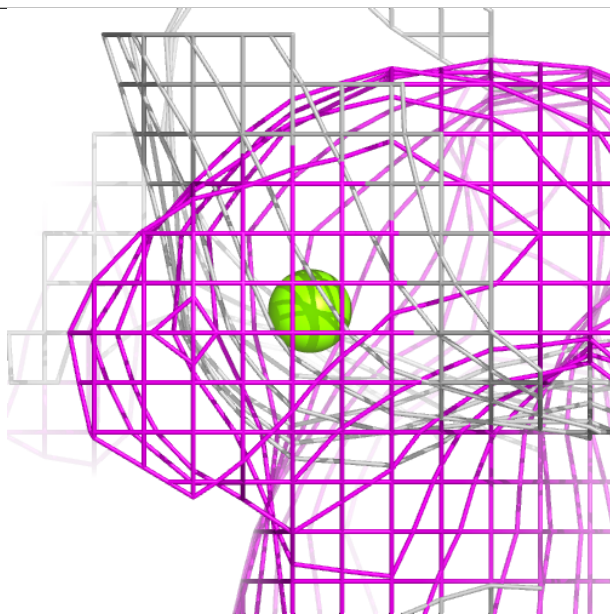
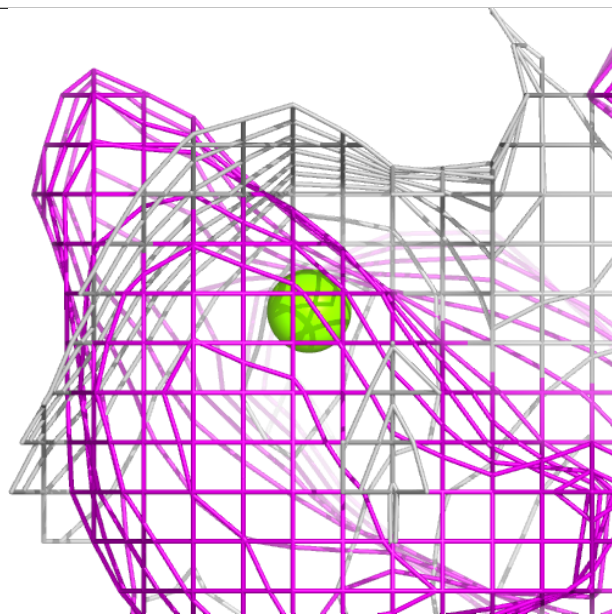
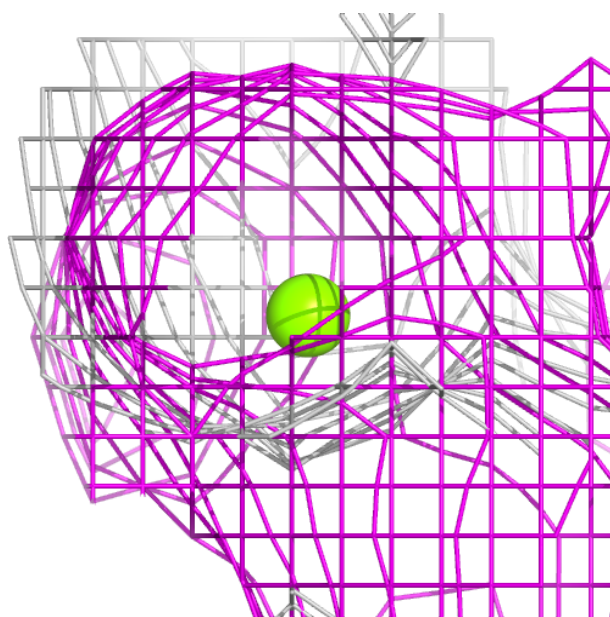
Electron density around MG E 307:

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and green (positive)



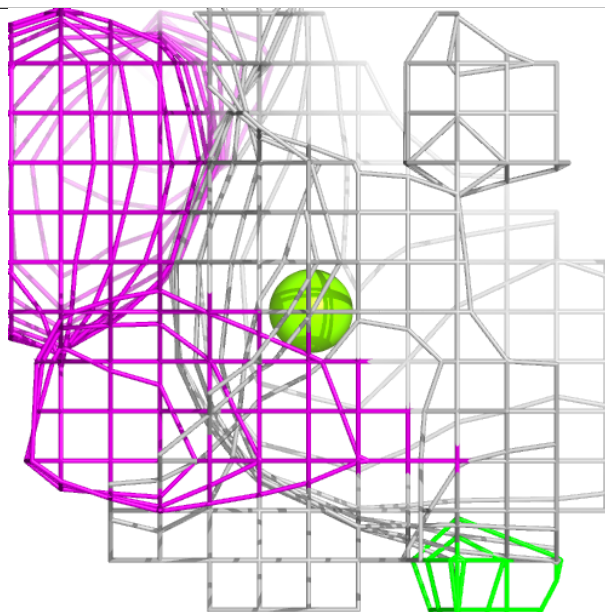
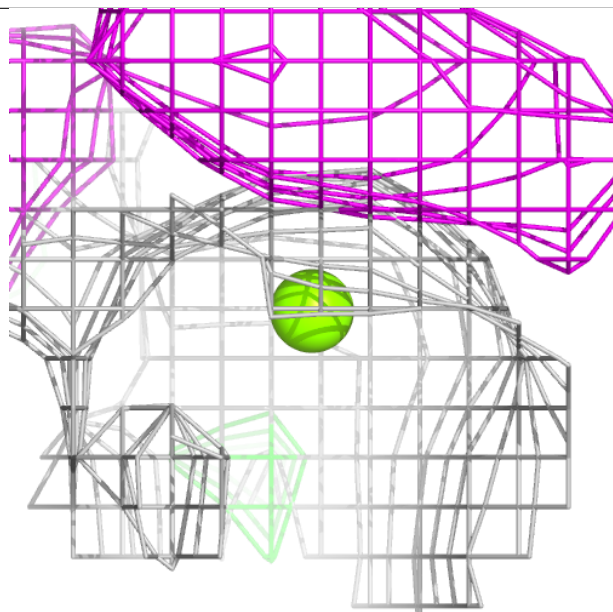
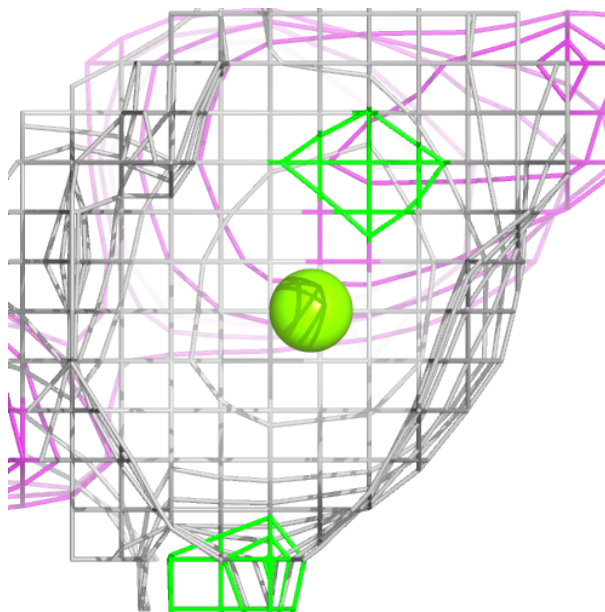
Electron density around MG D 304:

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and green (positive)



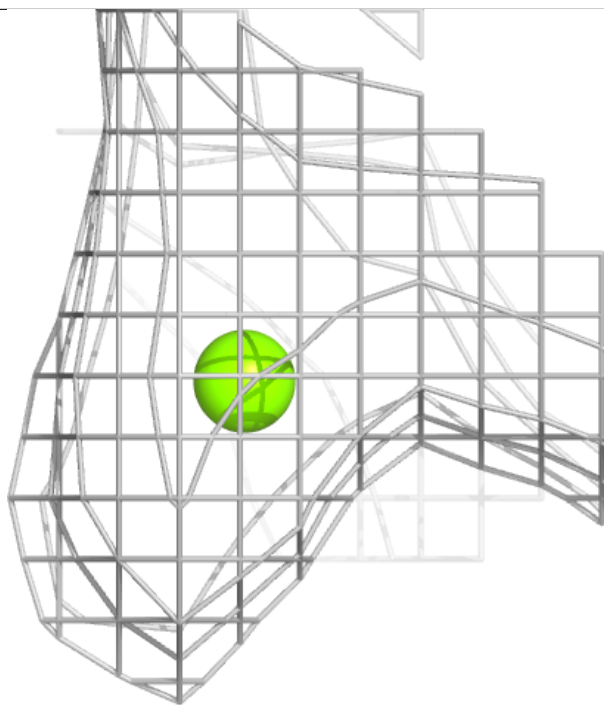
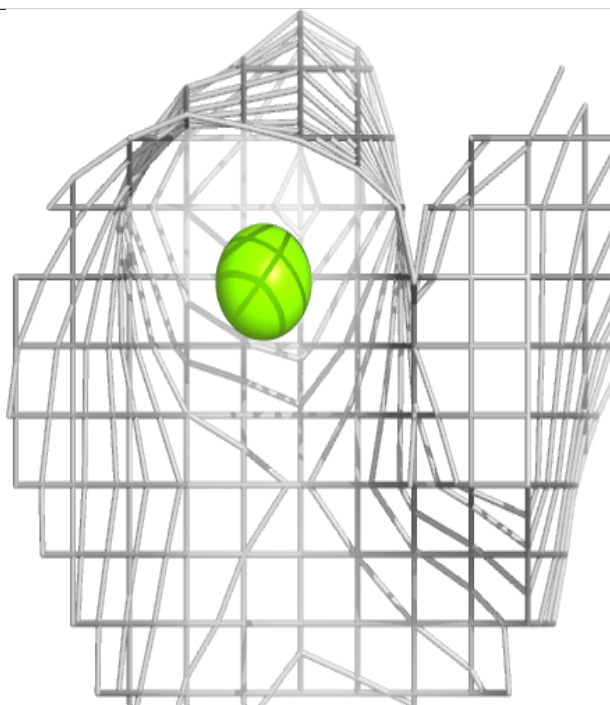
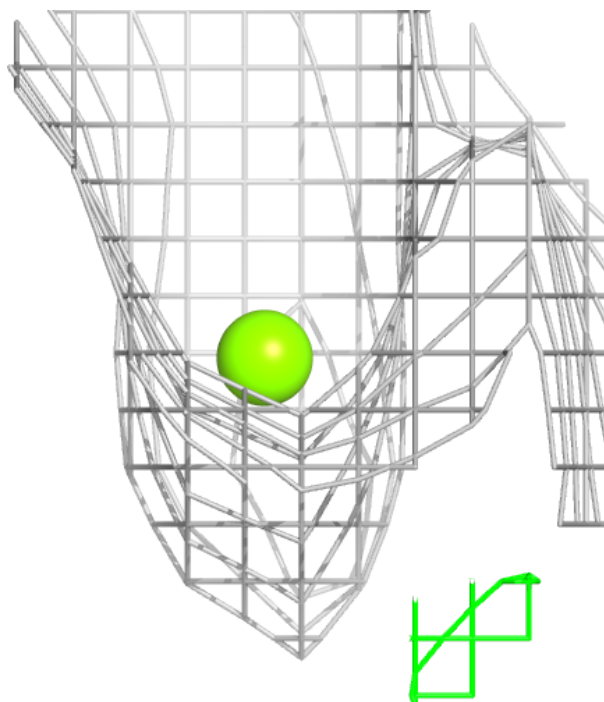
Electron density around MG E 303:

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and green (positive)



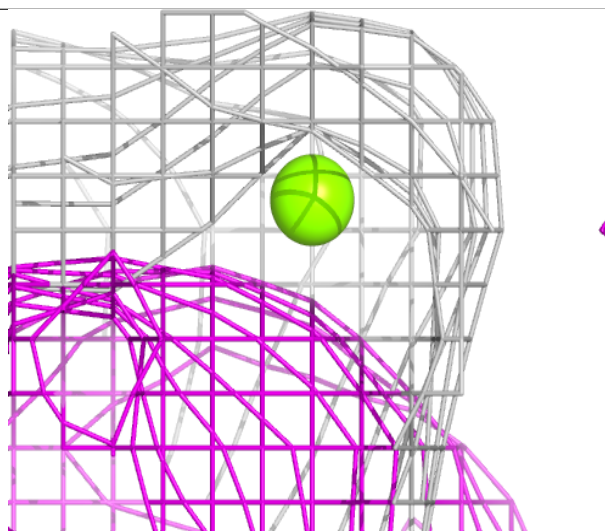
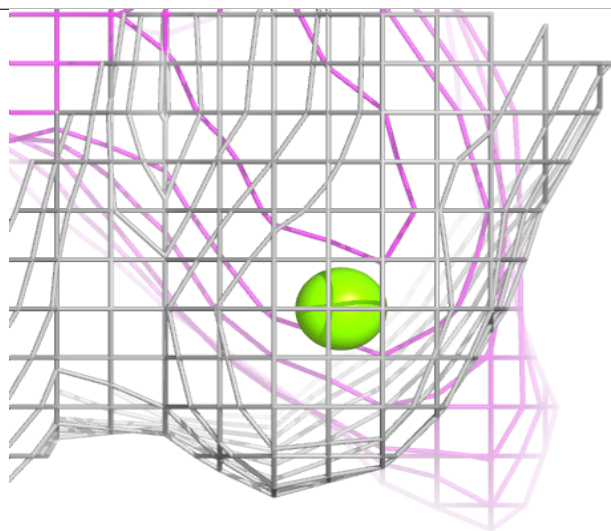
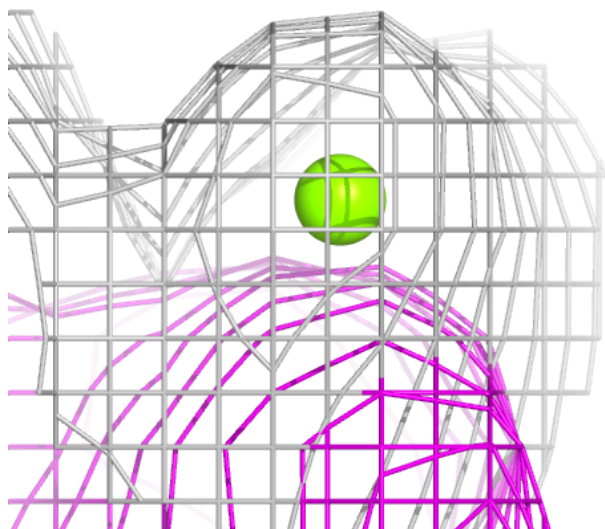
Electron density around MG G 402:

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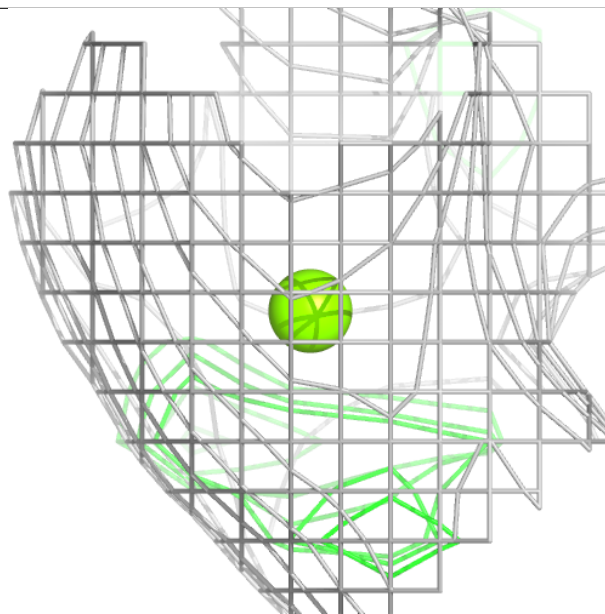
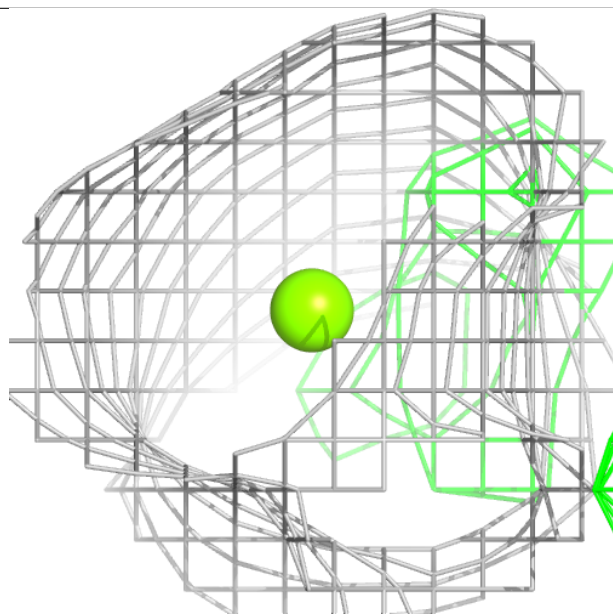
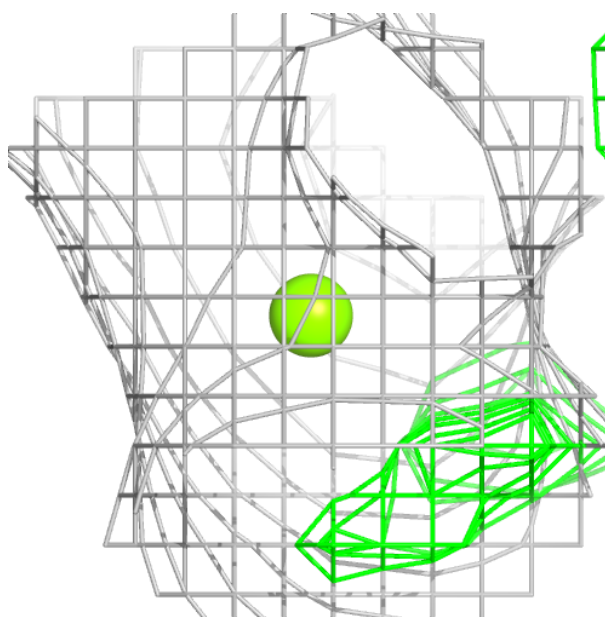
Electron density around MG D 305:

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and green (positive)



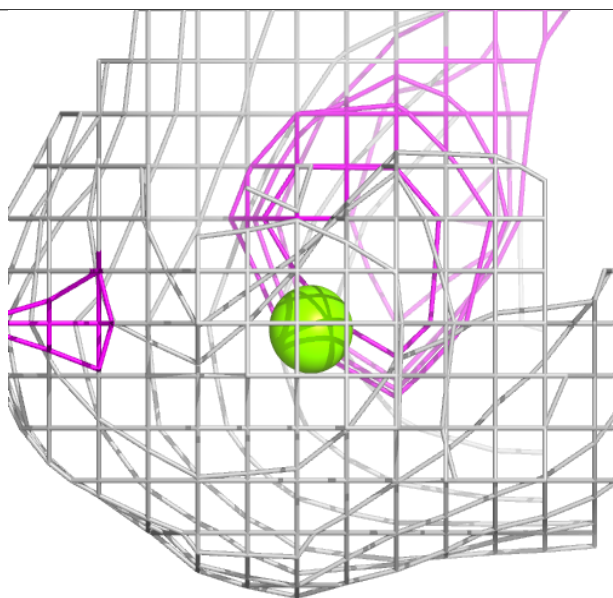
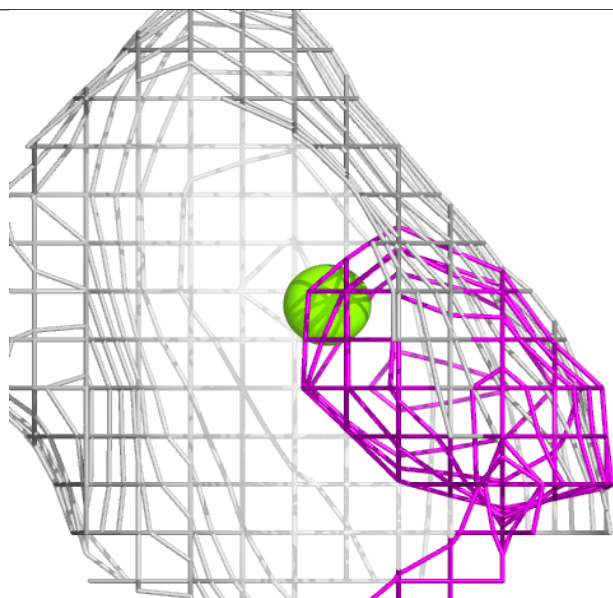
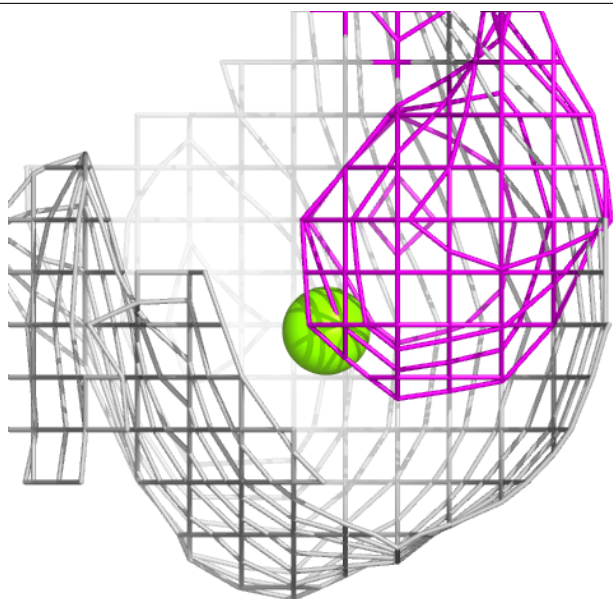
Electron density around MG D 307:

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and green (positive)



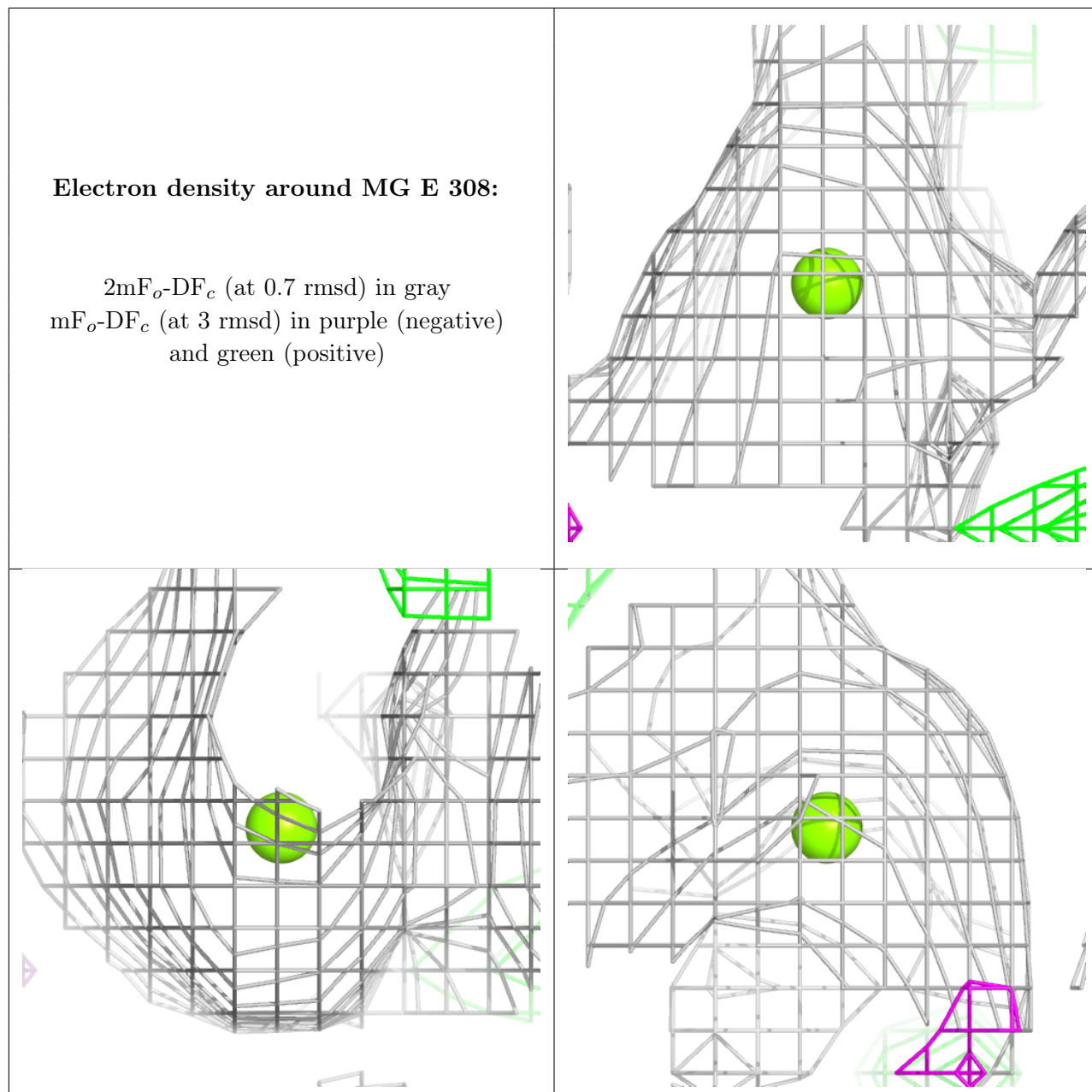
Electron density around MG B 306:

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and green (positive)



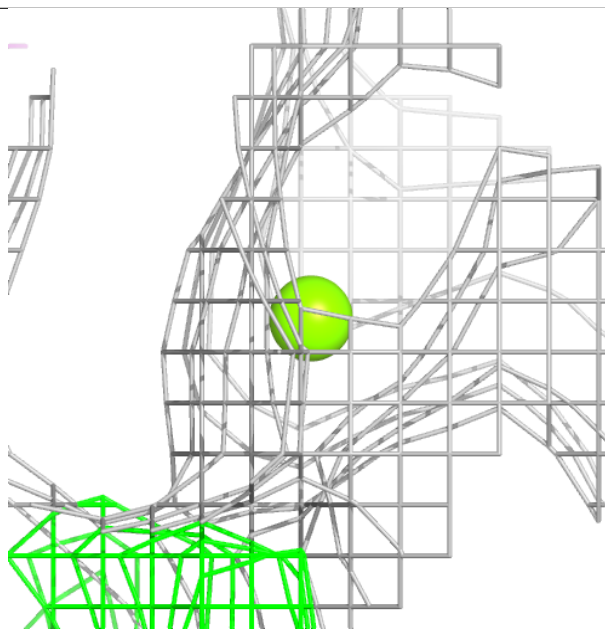
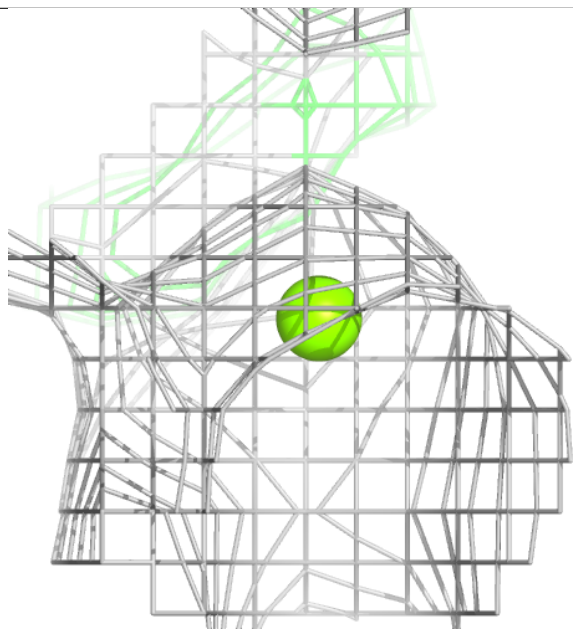
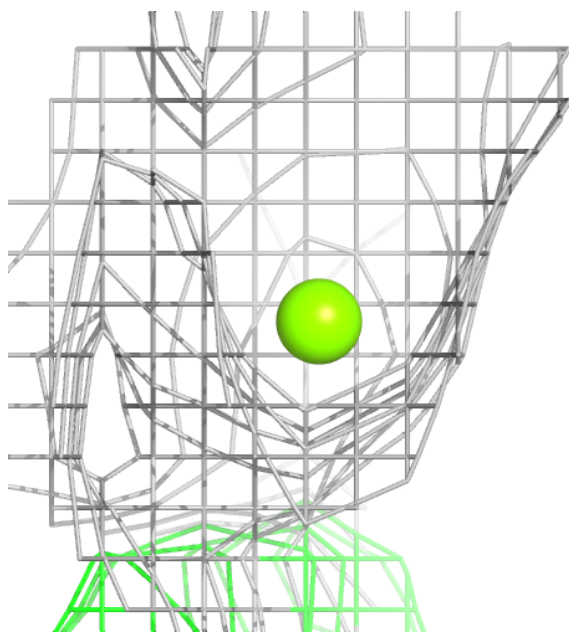
Electron density around MG E 308:

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



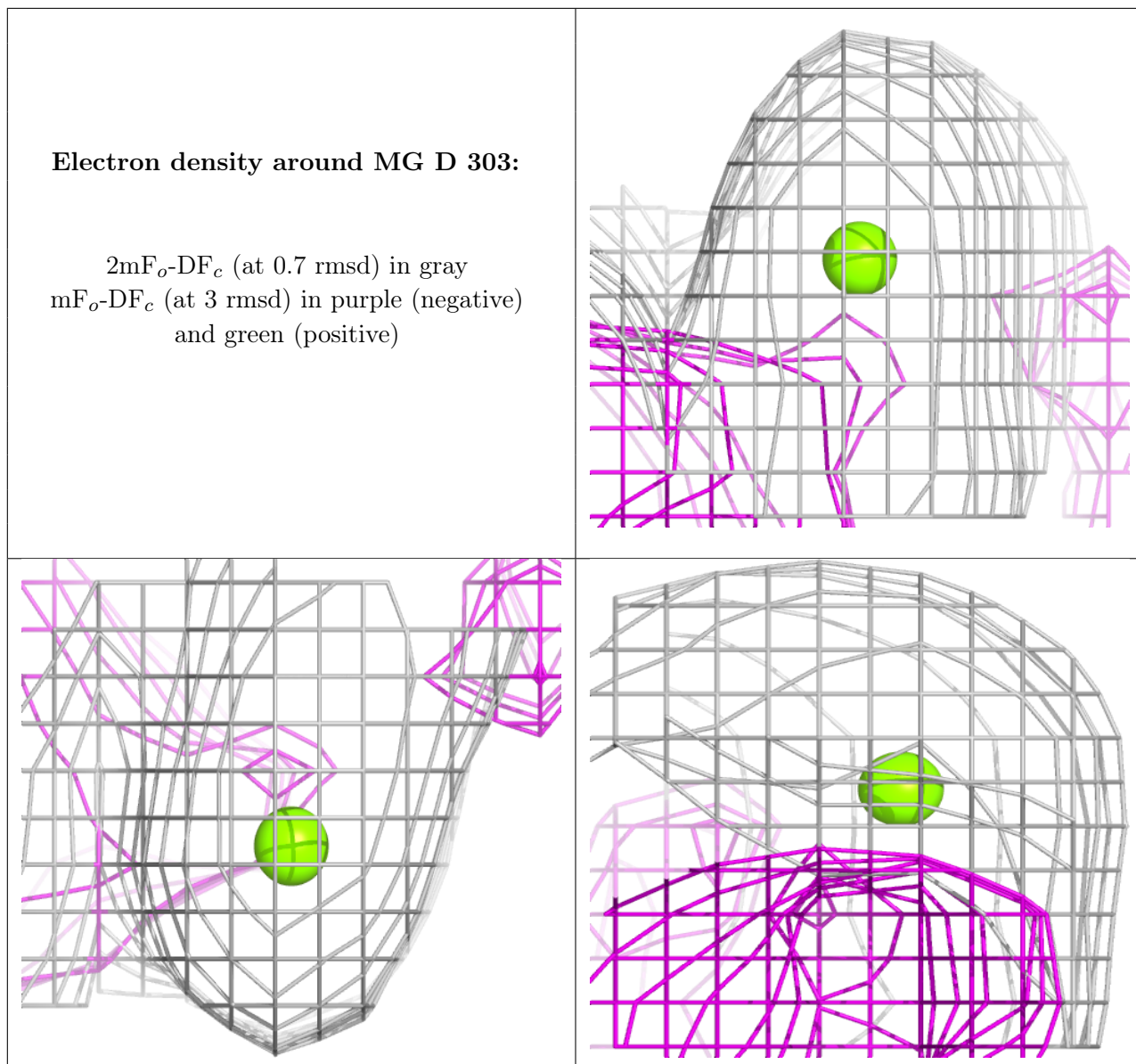
Electron density around MG F 303:

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



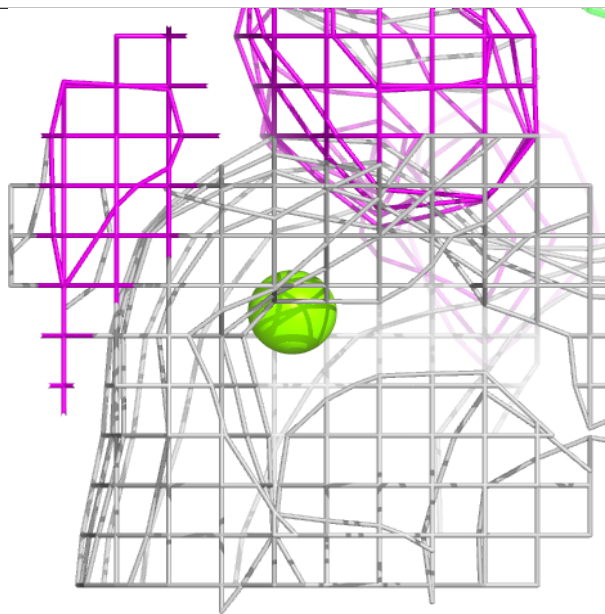
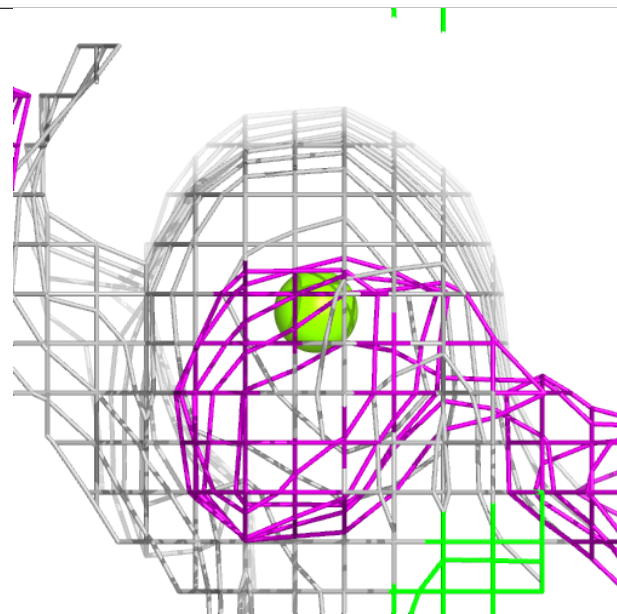
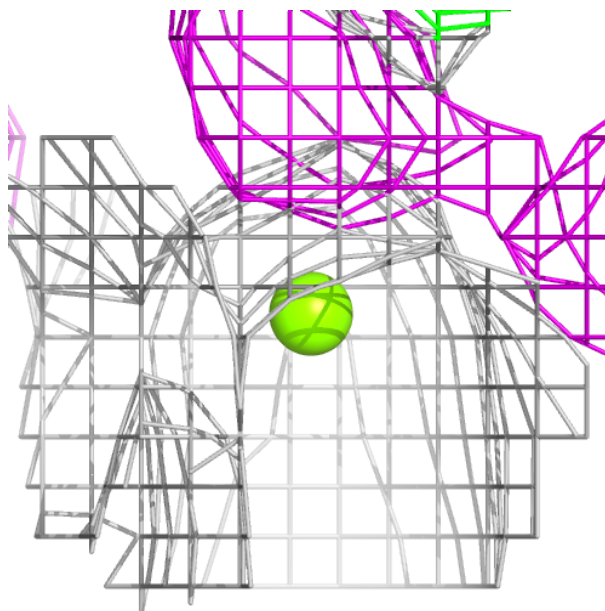
Electron density around MG D 303:

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



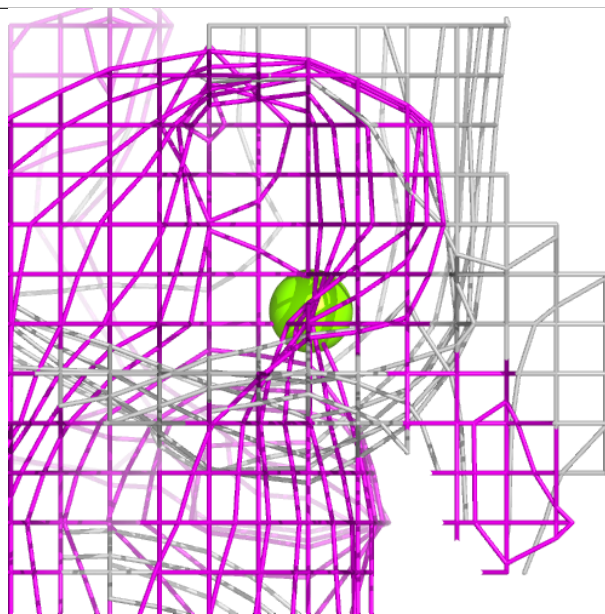
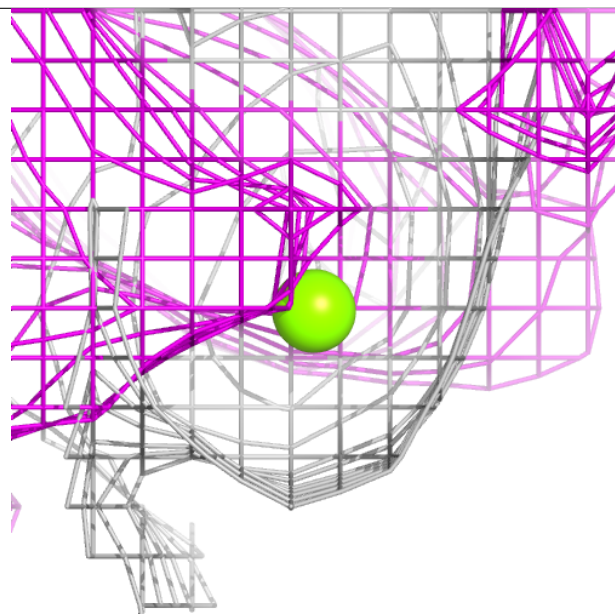
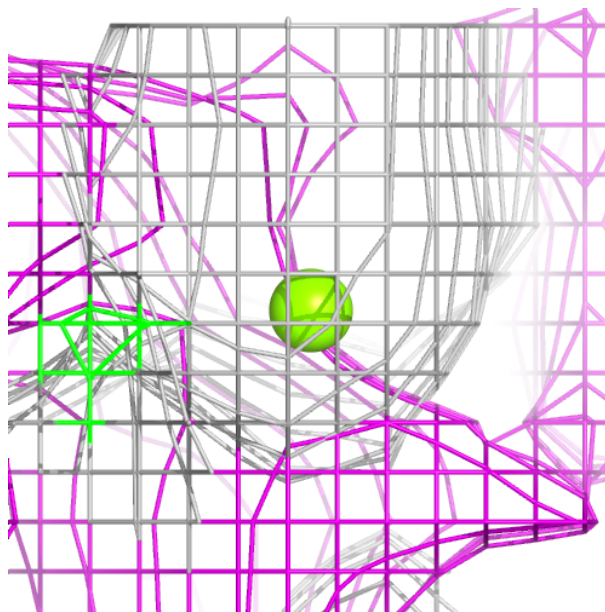
Electron density around MG B 303:

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



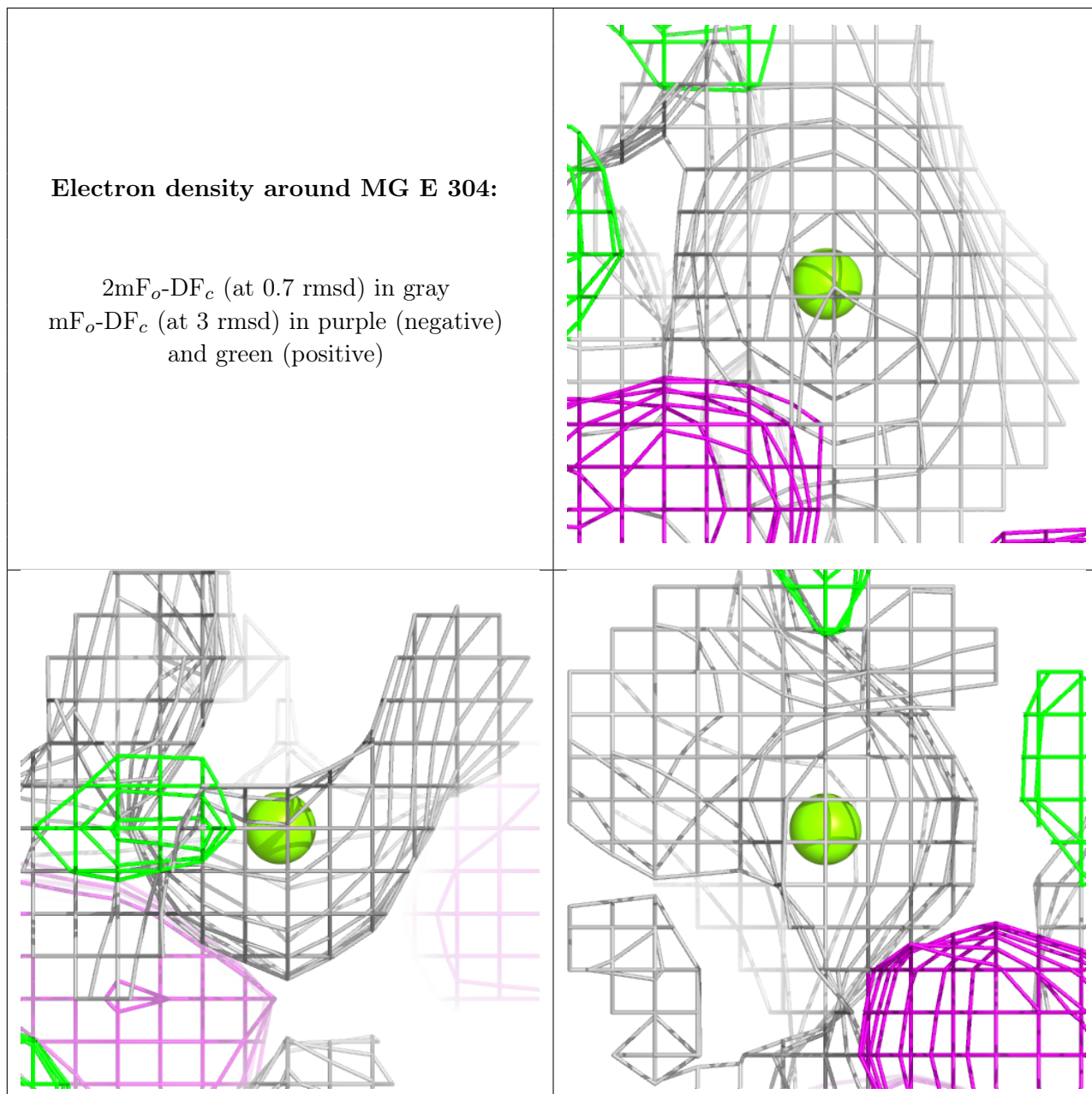
Electron density around MG D 306:

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



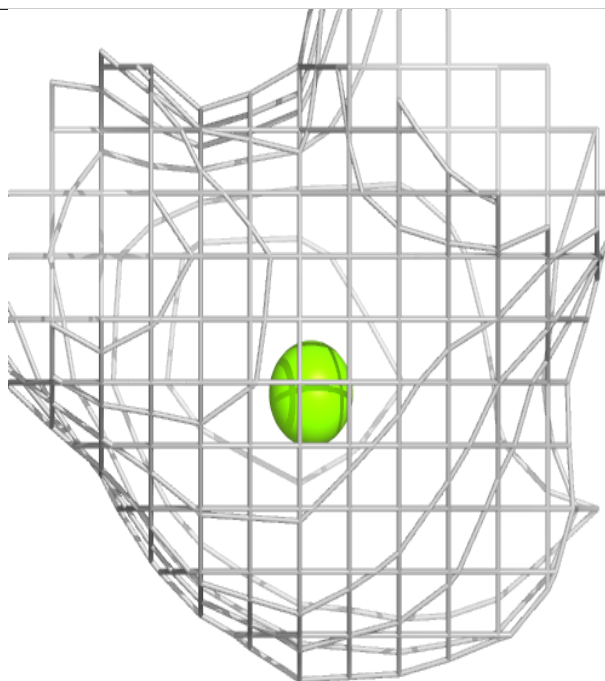
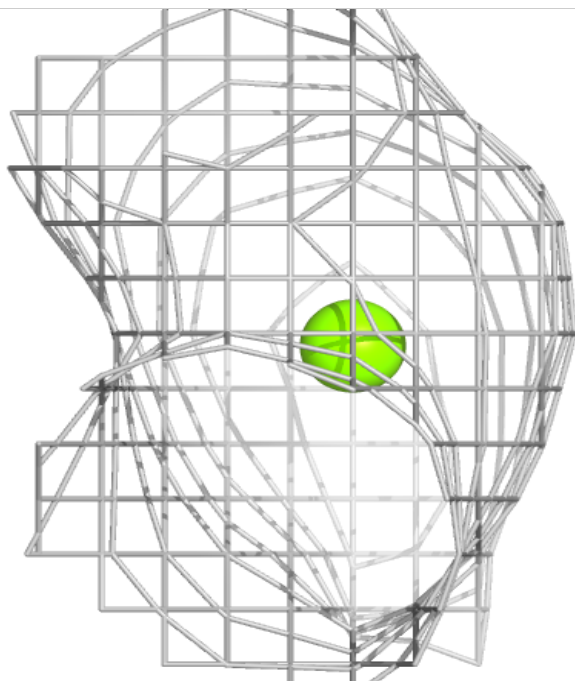
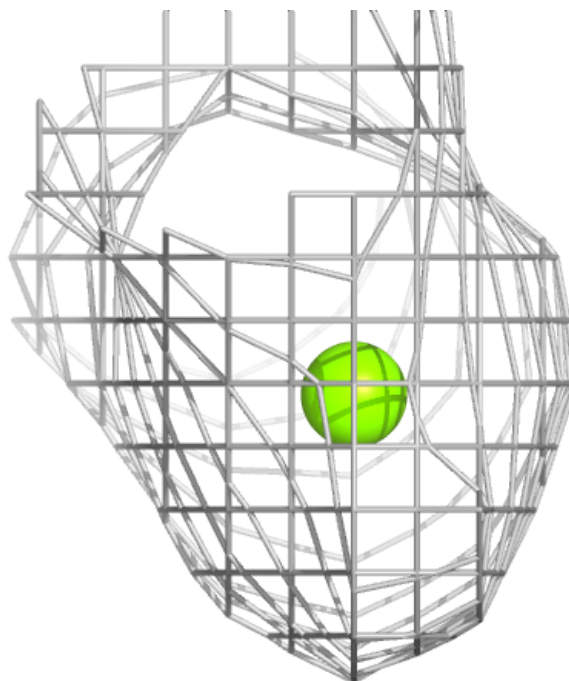
Electron density around MG E 304:

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



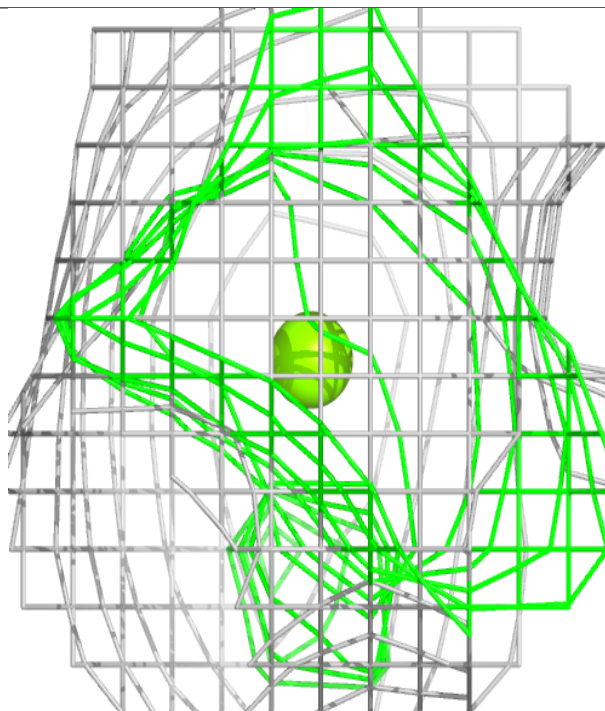
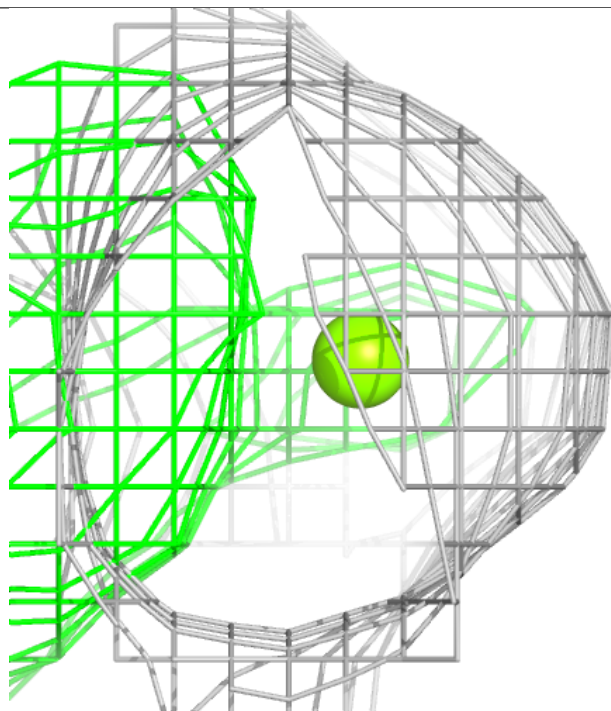
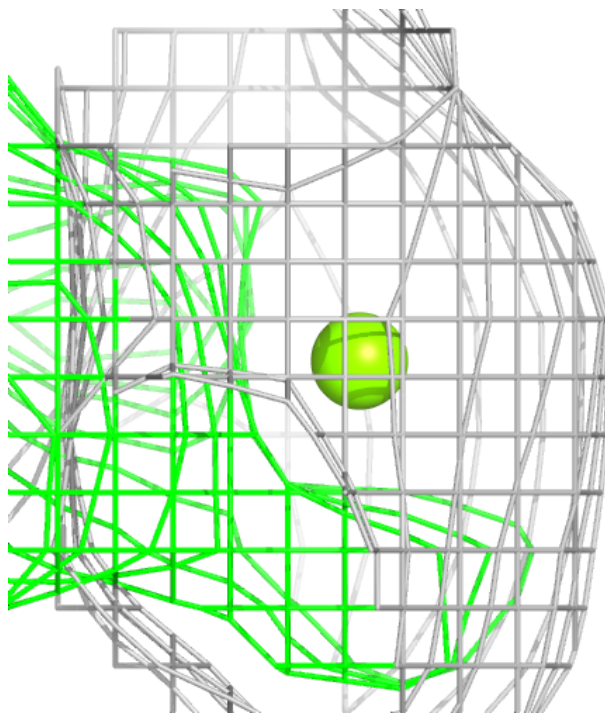
Electron density around MG E 305:

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



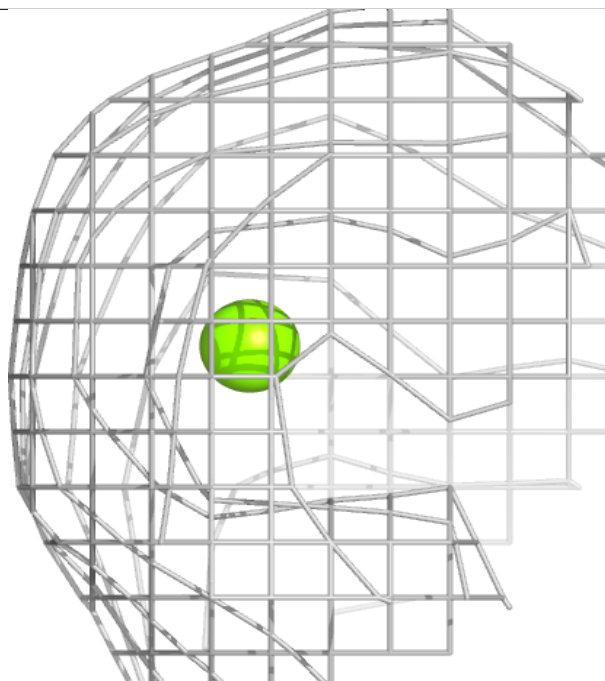
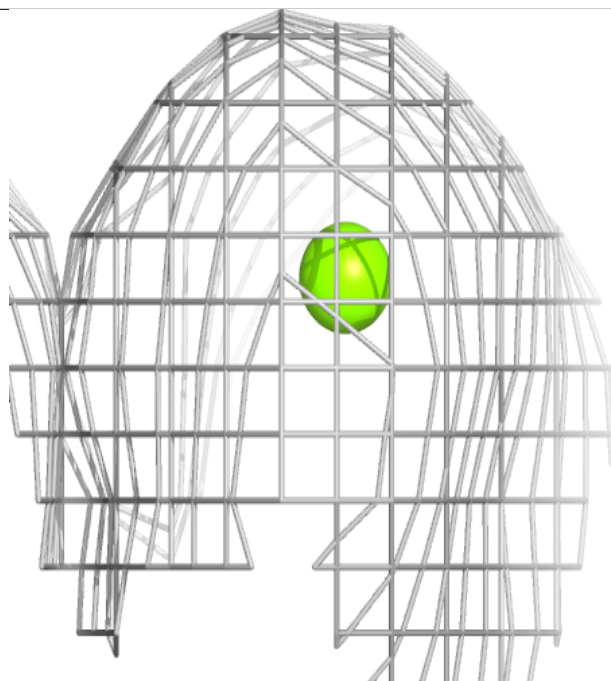
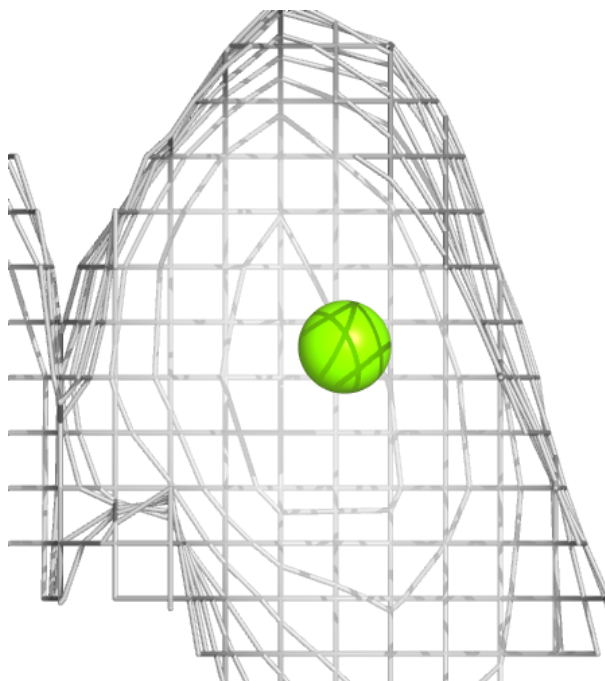
Electron density around MG B 307:

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



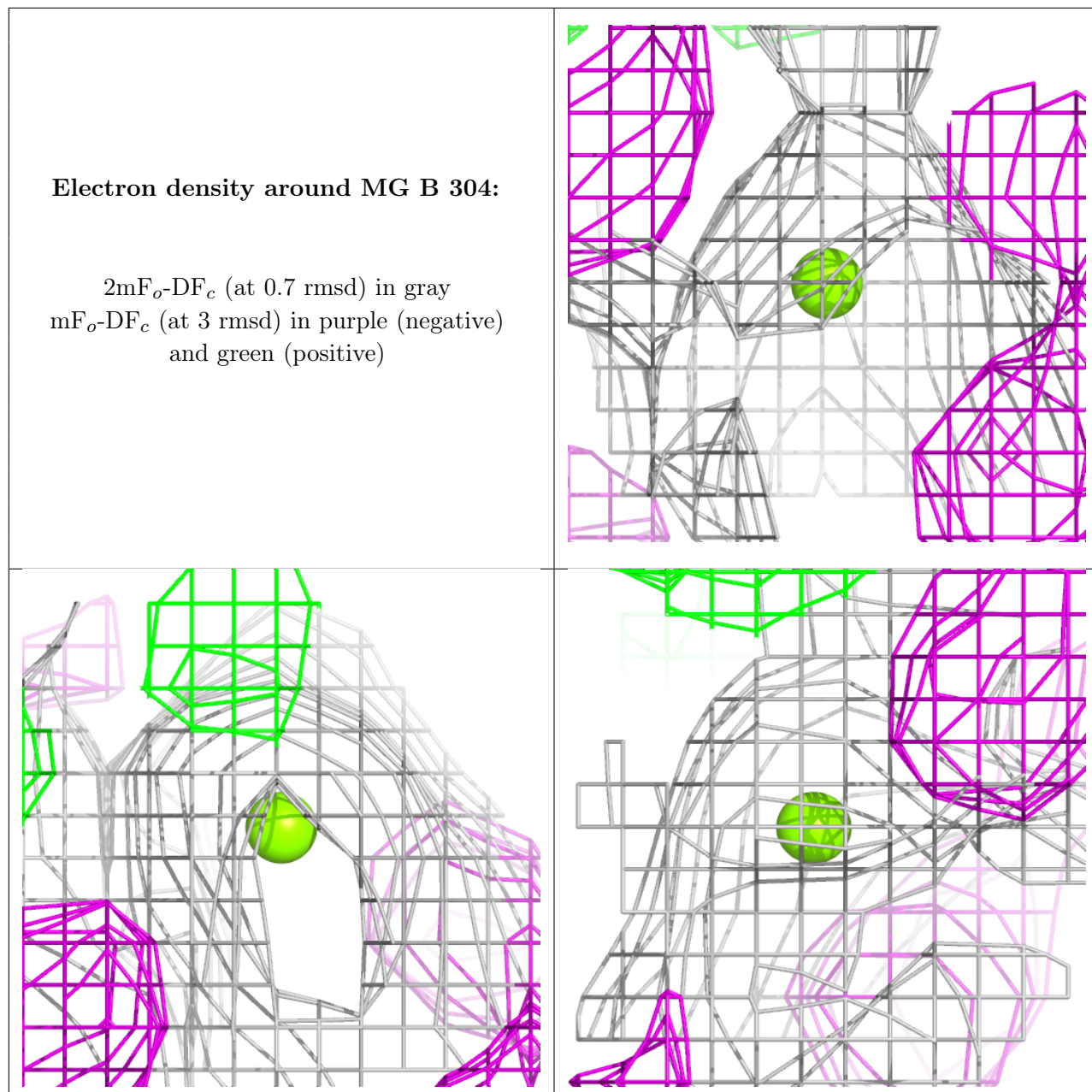
Electron density around MG G 404:

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



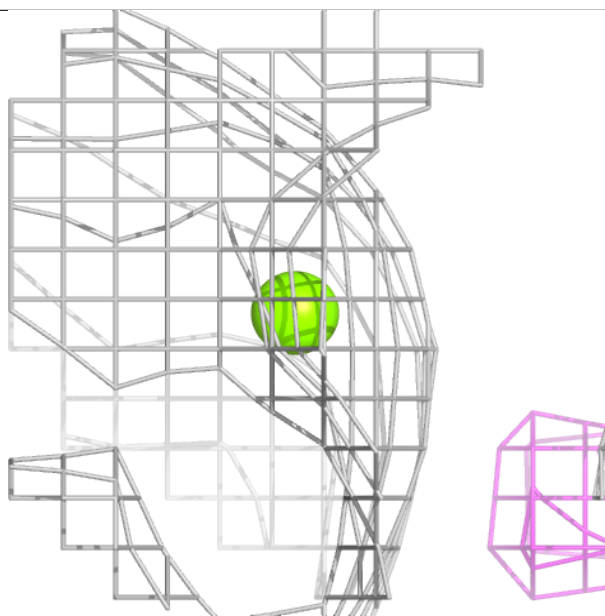
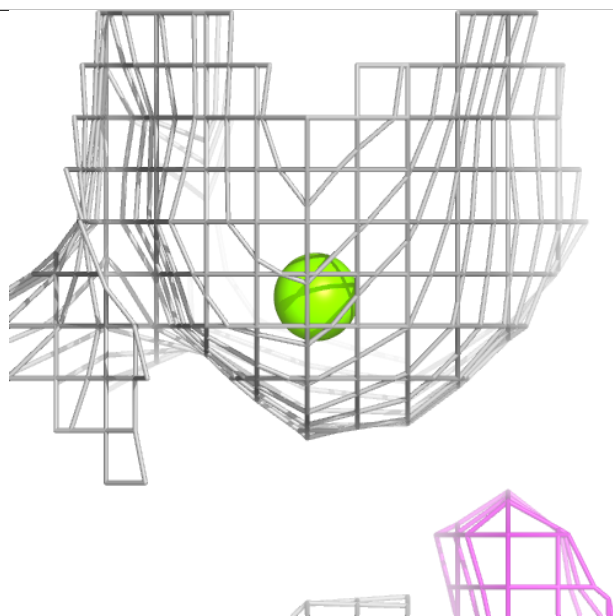
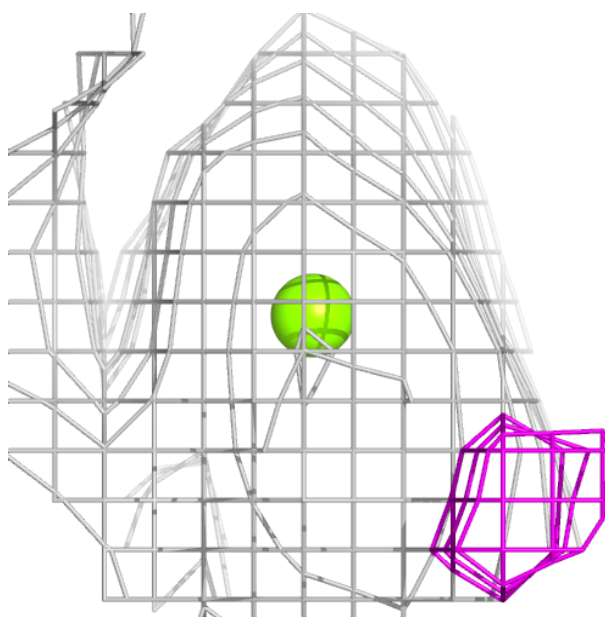
Electron density around MG B 304:

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



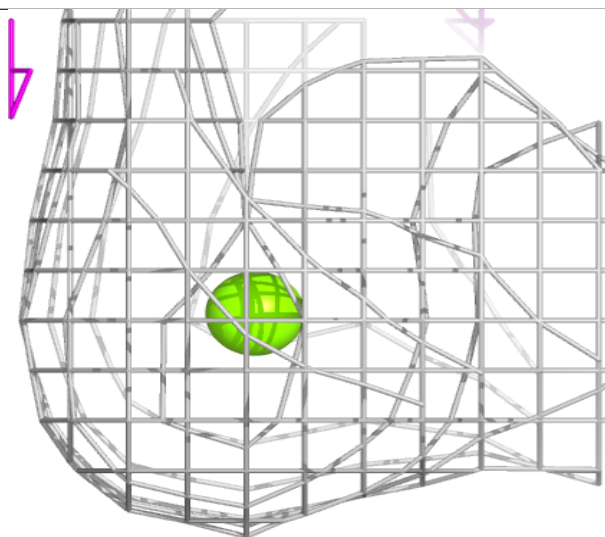
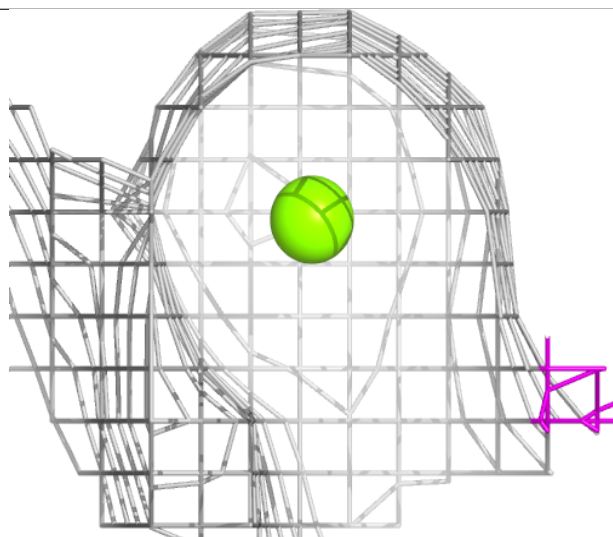
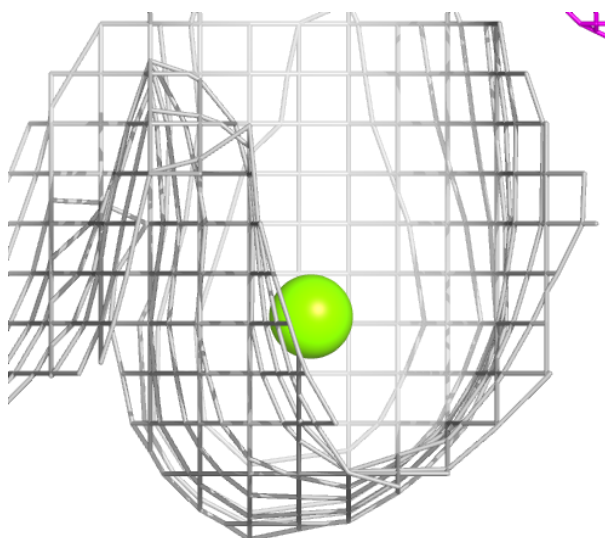
Electron density around MG F 302:

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



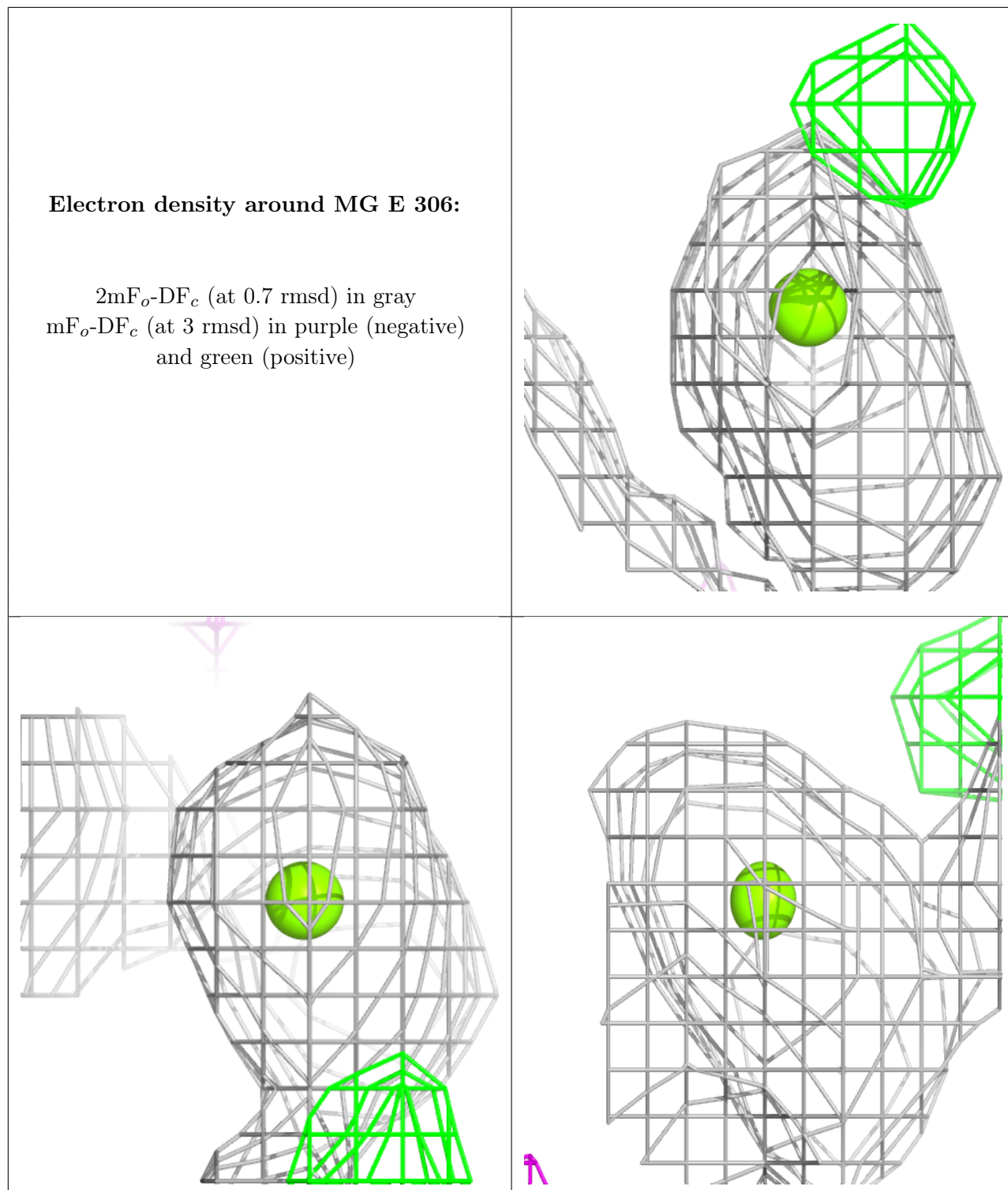
Electron density around MG B 305:

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



Electron density around MG E 306:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.