



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:04 AM UTC

PDB ID : 8WWV / pdb_00008wwv
Title : 1-naphthylamine GS in complex with ADP and MetSox-P
Authors : Zhang, S.T.; Zhou, N.Y.
Deposited on : 2023-10-26
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

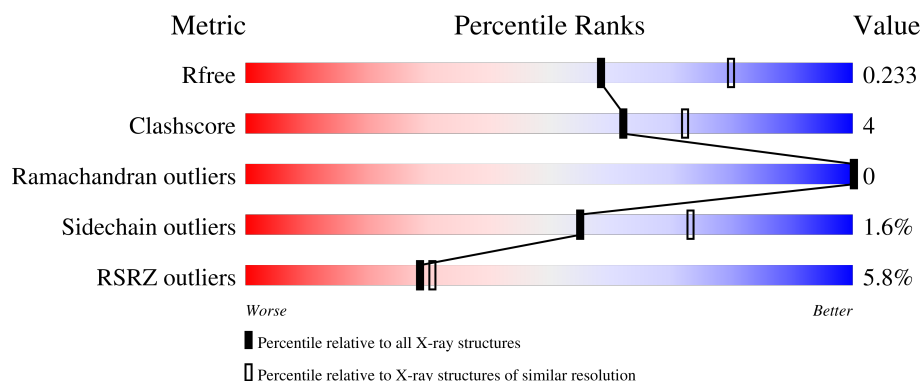
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	510	7% 85% 11% .
1	B	510	8% 85% 11% .
1	C	510	5% 86% 9% ..
1	D	510	5% 87% 9% ..
1	E	510	4% 87% 8% .

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Mol	Chain	Length	Quality of chain
1	F	510	4% 84% 12% ••

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 46408 atoms, of which 22530 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 Glutamine synthetase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	490	Total	C	H	N	O	S	0	0	0
			7565	2439	3737	643	728	18			
1	B	488	Total	C	H	N	O	S	0	0	0
			7524	2434	3705	641	726	18			
1	C	490	Total	C	H	N	O	S	0	0	0
			7560	2439	3732	643	728	18			
1	D	490	Total	C	H	N	O	S	0	0	0
			7552	2439	3724	643	728	18			
1	E	489	Total	C	H	N	O	S	0	0	0
			7571	2435	3751	642	725	18			
1	F	490	Total	C	H	N	O	S	0	0	0
			7583	2439	3755	643	728	18			

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
A	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
A	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
A	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
A	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1
A	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
A	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
A	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
A	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
A	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
A	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
A	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
A	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
A	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
A	0	GLY	-	expression tag	UNP A0A7Y1Q2L1
A	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
A	69	ALA	THR	conflict	UNP A0A7Y1Q2L1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
A	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
A	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
A	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
A	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
A	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
A	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
A	?	-	GLY	deletion	UNP A0A7Y1Q2L1
A	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
A	99	ASN	SER	conflict	UNP A0A7Y1Q2L1
A	101	THR	SER	conflict	UNP A0A7Y1Q2L1
A	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
A	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
A	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
A	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1
B	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
B	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
B	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
B	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
B	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1
B	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
B	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
B	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
B	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
B	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
B	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
B	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
B	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
B	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
B	0	GLY	-	expression tag	UNP A0A7Y1Q2L1
B	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
B	69	ALA	THR	conflict	UNP A0A7Y1Q2L1
B	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
B	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
B	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
B	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
B	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
B	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
B	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
B	?	-	GLY	deletion	UNP A0A7Y1Q2L1
B	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
B	99	ASN	SER	conflict	UNP A0A7Y1Q2L1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	101	THR	SER	conflict	UNP A0A7Y1Q2L1
B	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
B	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
B	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
B	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1
C	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
C	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
C	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
C	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
C	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1
C	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
C	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
C	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
C	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
C	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
C	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
C	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
C	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
C	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
C	0	GLY	-	expression tag	UNP A0A7Y1Q2L1
C	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
C	69	ALA	THR	conflict	UNP A0A7Y1Q2L1
C	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
C	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
C	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
C	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
C	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
C	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
C	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
C	?	-	GLY	deletion	UNP A0A7Y1Q2L1
C	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
C	99	ASN	SER	conflict	UNP A0A7Y1Q2L1
C	101	THR	SER	conflict	UNP A0A7Y1Q2L1
C	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
C	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
C	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
C	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1
D	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
D	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
D	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
D	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
D	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
D	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
D	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
D	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
D	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
D	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
D	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
D	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
D	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
D	0	GLY	-	expression tag	UNP A0A7Y1Q2L1
D	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
D	69	ALA	THR	conflict	UNP A0A7Y1Q2L1
D	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
D	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
D	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
D	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
D	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
D	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
D	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
D	?	-	GLY	deletion	UNP A0A7Y1Q2L1
D	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
D	99	ASN	SER	conflict	UNP A0A7Y1Q2L1
D	101	THR	SER	conflict	UNP A0A7Y1Q2L1
D	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
D	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
D	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
D	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1
E	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
E	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
E	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
E	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
E	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1
E	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
E	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
E	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
E	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
E	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
E	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
E	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
E	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
E	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
E	0	GLY	-	expression tag	UNP A0A7Y1Q2L1

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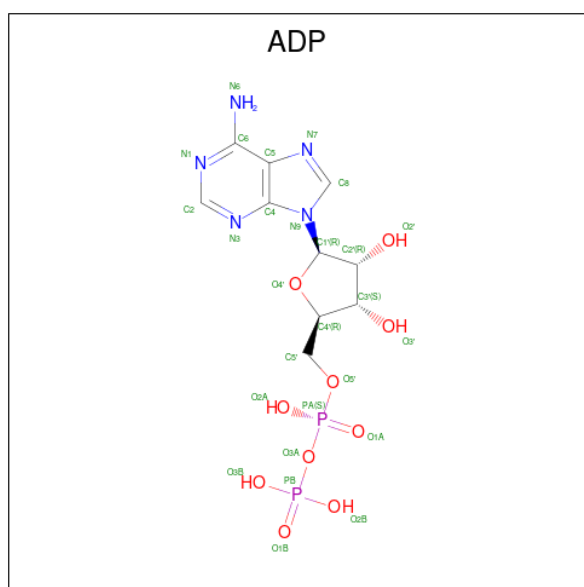
Chain	Residue	Modelled	Actual	Comment	Reference
E	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
E	69	ALA	THR	conflict	UNP A0A7Y1Q2L1
E	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
E	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
E	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
E	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
E	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
E	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
E	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
E	?	-	GLY	deletion	UNP A0A7Y1Q2L1
E	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
E	99	ASN	SER	conflict	UNP A0A7Y1Q2L1
E	101	THR	SER	conflict	UNP A0A7Y1Q2L1
E	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
E	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
E	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
E	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1
F	-14	TRP	-	expression tag	UNP A0A7Y1Q2L1
F	-13	SER	-	expression tag	UNP A0A7Y1Q2L1
F	-12	HIS	-	expression tag	UNP A0A7Y1Q2L1
F	-11	PRO	-	expression tag	UNP A0A7Y1Q2L1
F	-10	GLN	-	expression tag	UNP A0A7Y1Q2L1
F	-9	PHE	-	expression tag	UNP A0A7Y1Q2L1
F	-8	GLU	-	expression tag	UNP A0A7Y1Q2L1
F	-7	LYS	-	expression tag	UNP A0A7Y1Q2L1
F	-6	GLU	-	expression tag	UNP A0A7Y1Q2L1
F	-5	ASN	-	expression tag	UNP A0A7Y1Q2L1
F	-4	LEU	-	expression tag	UNP A0A7Y1Q2L1
F	-3	TYR	-	expression tag	UNP A0A7Y1Q2L1
F	-2	PHE	-	expression tag	UNP A0A7Y1Q2L1
F	-1	GLN	-	expression tag	UNP A0A7Y1Q2L1
F	0	GLY	-	expression tag	UNP A0A7Y1Q2L1
F	3	ARG	GLN	conflict	UNP A0A7Y1Q2L1
F	69	ALA	THR	conflict	UNP A0A7Y1Q2L1
F	70	ILE	MET	conflict	UNP A0A7Y1Q2L1
F	71	GLY	ALA	conflict	UNP A0A7Y1Q2L1
F	81	MET	ILE	conflict	UNP A0A7Y1Q2L1
F	85	LEU	PRO	conflict	UNP A0A7Y1Q2L1
F	88	THR	ALA	conflict	UNP A0A7Y1Q2L1
F	89	ALA	GLY	conflict	UNP A0A7Y1Q2L1
F	91	ASP	GLY	conflict	UNP A0A7Y1Q2L1
F	?	-	GLY	deletion	UNP A0A7Y1Q2L1

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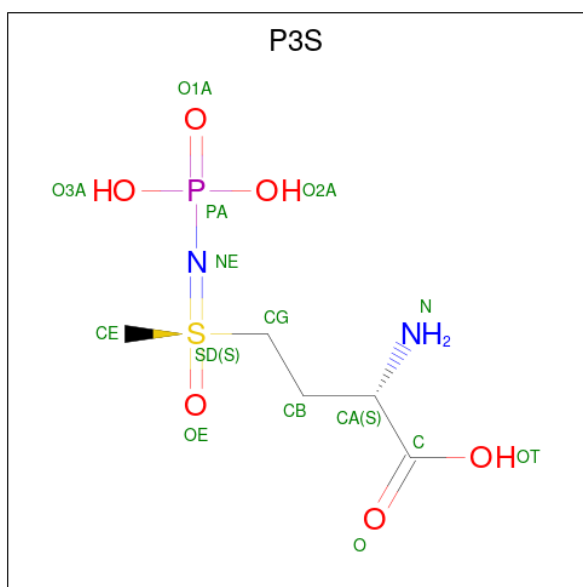
Chain	Residue	Modelled	Actual	Comment	Reference
F	93	GLU	ILE	conflict	UNP A0A7Y1Q2L1
F	99	ASN	SER	conflict	UNP A0A7Y1Q2L1
F	101	THR	SER	conflict	UNP A0A7Y1Q2L1
F	201	VAL	TYR	conflict	UNP A0A7Y1Q2L1
F	314	LEU	VAL	conflict	UNP A0A7Y1Q2L1
F	363	LYS	ARG	conflict	UNP A0A7Y1Q2L1
F	370	VAL	ILE	conflict	UNP A0A7Y1Q2L1

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 3 is L-METHIONINE-S-SULFOXIMINE PHOSPHATE (CCD ID: P3S) (formula: $C_5H_{13}N_2O_6PS$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		
3	B	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		
3	C	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		
3	D	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		
3	E	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		
3	F	1	Total	C	H	N	O	P	S	0	1
			25	5	10	2	6	1	1		

- Molecule 4 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Mn	0	0
			2	2		
4	B	2	Total	Mn	0	0
			2	2		
4	C	2	Total	Mn	0	0
			2	2		
4	D	2	Total	Mn	0	0
			2	2		
4	E	2	Total	Mn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	2	Total	Mn	0	0
			2	2		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

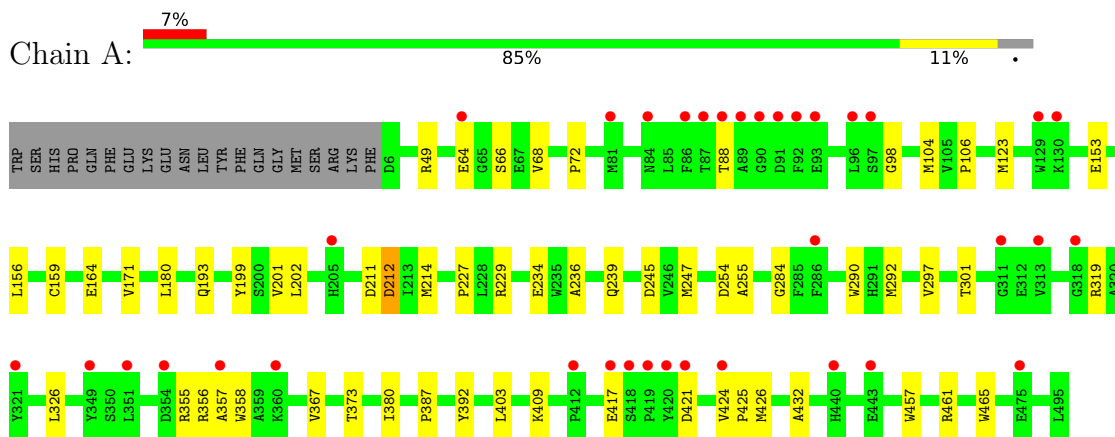
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	103	Total	O	0	0
			103	103		
6	B	107	Total	O	0	0
			107	107		
6	C	110	Total	O	0	0
			110	110		
6	D	111	Total	O	0	0
			111	111		
6	E	120	Total	O	0	0
			120	120		
6	F	106	Total	O	0	0
			106	106		

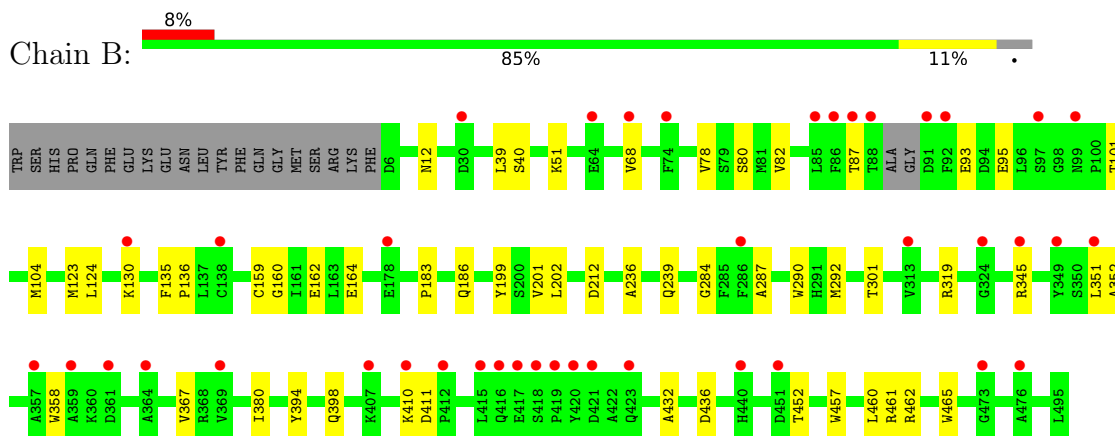
3 Residue-property plots [i](#)

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.

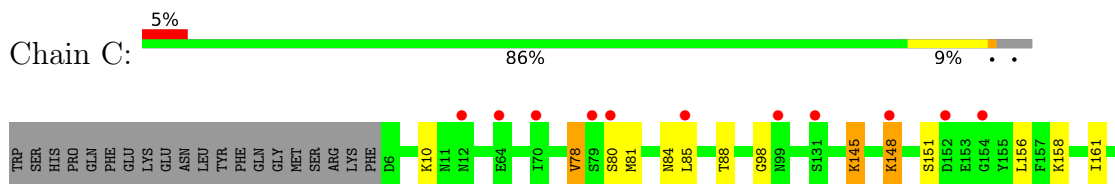
- Molecule 1: Glutamine synthetase

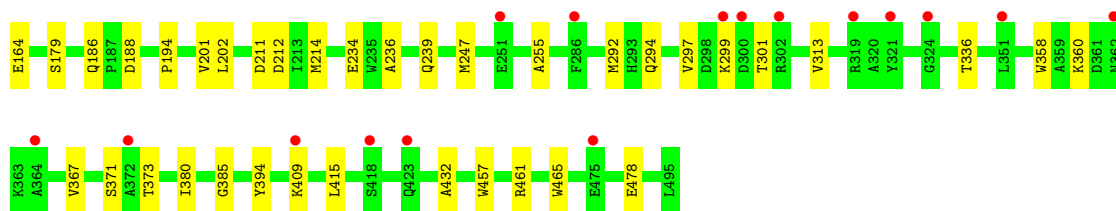


- Molecule 1: Glutamine synthetase

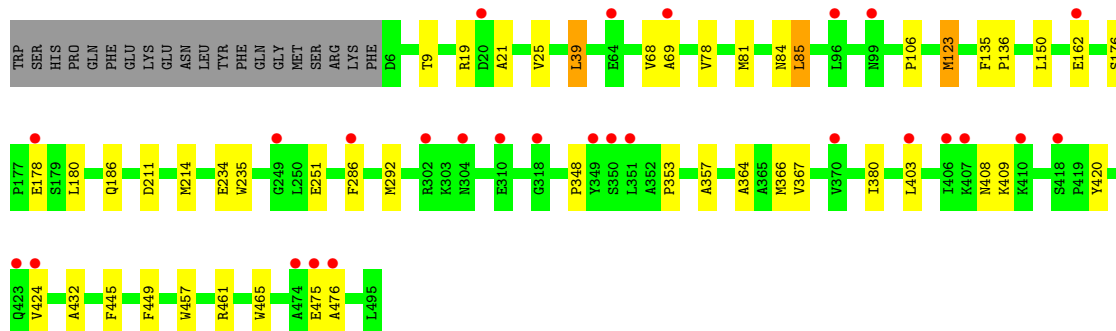
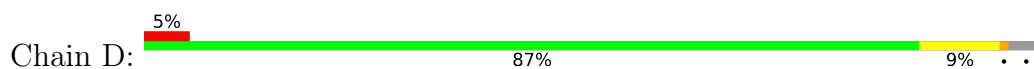


- Molecule 1: Glutamine synthetase

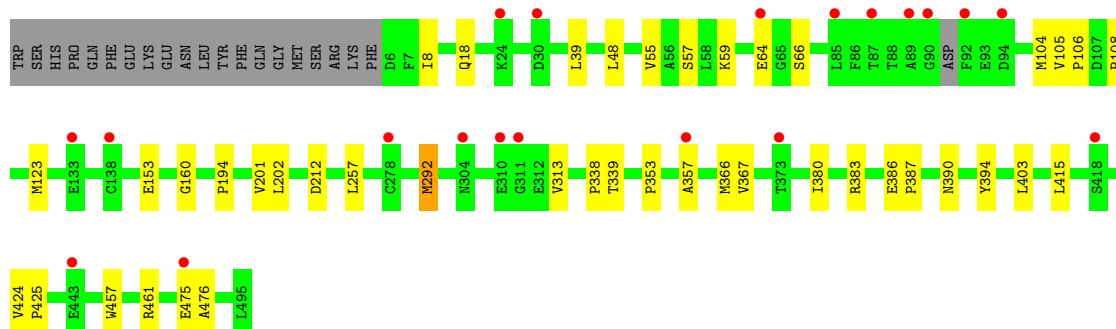
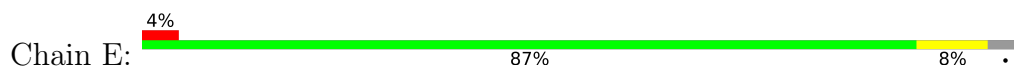




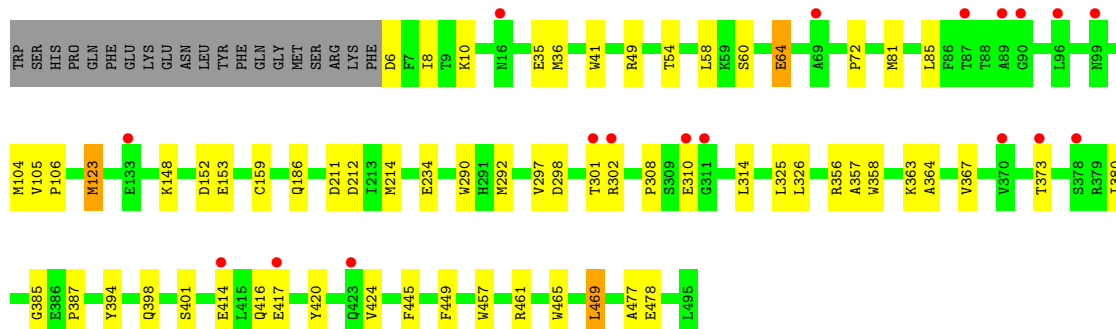
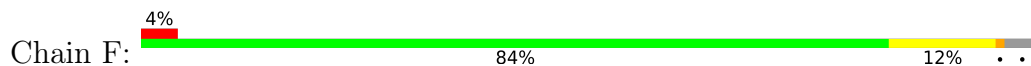
• Molecule 1: Glutamine synthetase



• Molecule 1: Glutamine synthetase



• Molecule 1: Glutamine synthetase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	123.06Å 140.48Å 216.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.77 – 2.30 43.77 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.77-2.30) 99.9 (43.77-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.29Å)	Xtrriage
Refinement program	PHENIX (1.20.1_4487: 000)	Depositor
R, R_{free}	0.212 , 0.232 0.213 , 0.233	Depositor DCC
R_{free} test set	8314 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	35.3	Xtrriage
Anisotropy	0.846	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	46408	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.08 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8856e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: P3S, MN, MG, ADP

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	0.41	2/3922 (0.1%)	0.68	3/5322 (0.1%)
1	B	0.41	2/3912 (0.1%)	0.60	1/5307 (0.0%)
1	C	0.34	0/3922	0.56	0/5322
1	D	0.45	0/3922	0.70	1/5322 (0.0%)
1	E	0.41	5/3913 (0.1%)	0.65	0/5308
1	F	0.32	1/3922 (0.0%)	0.54	0/5322
All	All	0.39	10/23513 (0.0%)	0.62	5/31903 (0.0%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	159	CYS	C-N	-10.63	1.28	1.33
1	A	159	CYS	C-N	-6.72	1.30	1.33
1	F	159	CYS	C-N	-6.42	1.30	1.33
1	E	425	PRO	C-O	-5.96	1.16	1.23
1	E	390	ASN	C-O	-5.56	1.17	1.24
1	E	338	PRO	C-O	-5.52	1.17	1.24
1	E	383	ARG	C-O	-5.44	1.16	1.24
1	A	425	PRO	C-O	-5.34	1.17	1.23
1	B	160	GLY	CA-C	5.11	1.56	1.51
1	E	387	PRO	C-O	-5.02	1.18	1.24

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	GLY	CA-C-O	-6.06	117.00	122.57
1	A	171	VAL	N-CA-C	-5.82	106.80	111.81
1	B	160	GLY	O-C-N	5.57	126.86	123.35
1	D	84	ASN	CA-C-O	-5.44	115.03	120.96
1	A	421	ASP	N-CA-C	-5.25	104.62	112.54

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	3828	3737	3755	33	0
1	B	3819	3705	3746	35	0
1	C	3828	3732	3755	31	0
1	D	3828	3724	3755	28	0
1	E	3820	3751	3750	21	0
1	F	3828	3755	3755	41	0
2	A	27	11	12	0	0
2	B	27	11	12	0	0
2	C	27	11	12	0	0
2	D	27	11	12	0	0
2	E	27	11	12	1	0
2	F	27	11	12	0	0
3	A	15	10	0	0	0
3	B	15	10	0	0	0
3	C	15	10	0	0	0
3	D	15	10	0	0	0
3	E	15	10	0	0	0
3	F	15	10	0	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
4	E	2	0	0	0	0
4	F	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
6	A	103	0	0	1	0
6	B	107	0	0	0	0
6	C	110	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	D	111	0	0	0	0
6	E	120	0	0	0	0
6	F	106	0	0	1	0
All	All	23878	22530	22588	184	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 4.

All (184) 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:F:36:MET:HE1	1:F:54:THR:HG22	1.40	1.04
1:F:298:ASP:HB3	1:F:301:THR:HG22	1.52	0.91
1:E:8:ILE:CD1	1:E:105:VAL:HG13	2.15	0.76
1:A:355:ARG:NH1	1:A:426:MET:SD	2.59	0.74
1:A:367:VAL:HG13	1:A:380:ILE:HG23	1.70	0.72
1:C:313:VAL:HG11	1:C:415:LEU:HD11	1.71	0.72
1:C:367:VAL:HG13	1:C:380:ILE:HG23	1.71	0.71
1:F:41:TRP:CD1	1:F:104:MET:HE3	2.25	0.71
1:F:58:LEU:HD11	1:F:123:MET:HE1	1.73	0.71
1:B:51:LYS:HE3	1:B:68:VAL:HG12	1.73	0.70
1:F:298:ASP:CB	1:F:301:THR:HG22	2.21	0.70
1:D:367:VAL:HG13	1:D:380:ILE:HG23	1.74	0.70
1:E:8:ILE:HD12	1:E:105:VAL:HG13	1.73	0.69
1:F:310:GLU:HA	1:F:373:THR:HG21	1.75	0.69
1:F:36:MET:HE1	1:F:54:THR:CG2	2.21	0.68
1:B:39:LEU:HD23	1:B:104:MET:HE2	1.75	0.67
1:C:151:SER:OG	1:C:299:LYS:HE2	1.95	0.67
1:B:367:VAL:HG13	1:B:380:ILE:HG23	1.77	0.65
1:F:41:TRP:HD1	1:F:104:MET:HE3	1.61	0.65
1:A:357:ALA:HB2	1:A:424:VAL:HG21	1.78	0.65
1:F:36:MET:HA	1:F:36:MET:HE2	1.79	0.65
1:F:148:LYS:NZ	1:F:152:ASP:OD1	2.26	0.65
1:A:68:VAL:HG13	1:A:104:MET:HE2	1.78	0.64
1:B:87:THR:HG23	1:B:87:THR:O	1.98	0.63
1:D:39:LEU:HD21	1:D:106:PRO:HG3	1.80	0.62
1:E:367:VAL:HG13	1:E:380:ILE:HG23	1.81	0.62
1:C:78:VAL:CG1	1:C:78:VAL:O	2.47	0.62
1:E:39:LEU:HD23	1:E:104:MET:HE3	1.84	0.60
1:F:36:MET:CE	1:F:54:THR:HG22	2.26	0.59
1:C:145:LYS:O	1:C:148:LYS:HG3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:ILE:HD12	1:C:247:MET:CE	2.34	0.58
1:F:297:VAL:CG1	1:F:302:ARG:HA	2.34	0.58
1:B:39:LEU:HD12	1:B:123:MET:HG2	1.86	0.58
1:F:367:VAL:HG13	1:F:380:ILE:HG23	1.85	0.57
1:A:326:LEU:O	1:A:356:ARG:HD2	2.04	0.56
1:C:78:VAL:O	1:C:78:VAL:HG12	2.04	0.56
1:F:326:LEU:HD11	1:F:358:TRP:HB3	1.86	0.56
1:C:88:THR:HB	1:F:417:GLU:HG3	1.87	0.55
1:B:78:VAL:HG11	1:B:186:GLN:HG2	1.89	0.54
1:E:357:ALA:HB2	1:E:424:VAL:HG21	1.89	0.54
1:B:436:ASP:OD1	1:B:462:ARG:NH2	2.41	0.54
1:F:6:ASP:O	1:F:10:LYS:HD3	2.08	0.54
1:D:353:PRO:HB3	1:D:366:MET:HA	1.89	0.54
1:C:360:LYS:NZ	1:C:371:SER:O	2.36	0.53
1:A:227:PRO:HB2	1:A:245:ASP:HB2	1.91	0.53
1:C:247:MET:HE1	1:C:255:ALA:CB	2.39	0.53
1:D:180:LEU:HD11	1:E:194:PRO:HG2	1.90	0.53
1:D:234:GLU:HG3	1:D:235:TRP:H	1.74	0.52
1:F:212:ASP:N	1:F:212:ASP:OD1	2.41	0.52
1:E:55:VAL:O	1:E:59:LYS:HG2	2.09	0.52
1:E:475:GLU:HG2	1:E:476:ALA:N	2.24	0.52
1:D:475:GLU:HG2	1:D:476:ALA:N	2.25	0.52
1:E:313:VAL:HG11	1:E:415:LEU:HD11	1.91	0.52
1:F:292:MET:HE3	1:F:398:GLN:HG3	1.92	0.52
1:F:357:ALA:HB2	1:F:424:VAL:HG21	1.91	0.52
1:A:68:VAL:CG1	1:A:104:MET:HE2	2.40	0.51
1:B:39:LEU:HD23	1:B:104:MET:CE	2.41	0.51
1:B:199:TYR:OH	1:B:284:GLY:O	2.27	0.51
1:A:326:LEU:HD11	1:A:358:TRP:HB3	1.92	0.51
1:E:48:LEU:HD21	1:E:257:LEU:HD21	1.93	0.51
1:A:106:PRO:HB3	1:A:123:MET:HE2	1.92	0.50
1:F:297:VAL:HG13	1:F:302:ARG:HA	1.93	0.50
1:C:212:ASP:N	1:C:212:ASP:OD1	2.45	0.50
1:D:21:ALA:O	1:D:25:VAL:HG23	2.12	0.50
1:F:457:TRP:O	1:F:461:ARG:HG2	2.11	0.50
1:D:68:VAL:HG22	1:D:69:ALA:N	2.27	0.49
1:F:35:GLU:O	1:F:36:MET:HE2	2.13	0.49
1:A:49:ARG:HD2	1:A:72:PRO:HB3	1.95	0.49
1:E:201:VAL:HG22	1:E:202:LEU:HG	1.95	0.49
1:B:212:ASP:N	1:B:212:ASP:OD1	2.45	0.49
1:C:358:TRP:CE2	1:C:415:LEU:HD23	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:211:ASP:HA	1:D:214:MET:HB2	1.95	0.48
1:A:212:ASP:OD1	1:A:212:ASP:N	2.43	0.48
1:C:457:TRP:O	1:C:461:ARG:HG2	2.14	0.48
1:A:457:TRP:O	1:A:461:ARG:HG2	2.14	0.47
1:E:39:LEU:HD21	1:E:106:PRO:HG3	1.96	0.47
1:A:229:ARG:HG3	1:E:57:SER:OG	2.13	0.47
1:B:93:GLU:HB3	1:B:452:THR:HG21	1.96	0.47
1:C:85:LEU:HD23	1:C:98:GLY:O	2.15	0.47
1:E:212:ASP:OD1	1:E:212:ASP:N	2.47	0.47
1:E:313:VAL:HG11	1:E:415:LEU:CD1	2.45	0.47
1:F:186:GLN:HG3	6:F:659:HOH:O	2.15	0.47
1:F:81:MET:HE2	1:F:85:LEU:HD12	1.96	0.46
1:B:12:ASN:O	1:B:12:ASN:CG	2.58	0.46
1:D:364:ALA:HB1	1:D:420:TYR:CE2	2.51	0.46
1:F:445:PHE:O	1:F:449:PHE:HB2	2.17	0.45
1:A:156:LEU:HB2	1:A:297:VAL:CG2	2.47	0.45
1:A:247:MET:HE1	1:A:255:ALA:CB	2.47	0.45
1:C:78:VAL:HG13	6:C:644:HOH:O	2.15	0.45
1:A:201:VAL:HG22	1:A:202:LEU:HG	1.97	0.45
1:D:78:VAL:HG11	1:D:186:GLN:HG2	1.98	0.45
1:B:290:TRP:CD1	1:B:290:TRP:C	2.94	0.45
1:D:348:PRO:O	1:D:420:TYR:CG	2.69	0.45
1:A:290:TRP:HB2	1:A:387:PRO:HA	1.98	0.45
1:B:410:LYS:HG3	1:B:411:ASP:N	2.32	0.45
1:D:9:THR:HG22	1:D:19:ARG:NH2	2.31	0.45
1:D:432:ALA:HB2	1:D:465:TRP:CE2	2.51	0.45
1:A:66:SER:HB3	1:A:104:MET:HE3	1.99	0.45
1:F:414:GLU:HG3	1:F:416:GLN:NE2	2.32	0.44
1:D:457:TRP:O	1:D:461:ARG:HG2	2.17	0.44
1:A:236:ALA:HB3	1:A:239:GLN:HB2	2.00	0.44
1:C:247:MET:HE1	1:C:255:ALA:HB3	2.00	0.44
1:F:49:ARG:HD2	1:F:72:PRO:HB3	2.00	0.44
1:A:254:ASP:OD1	1:A:392:TYR:OH	2.27	0.44
1:F:465:TRP:NE1	1:F:469:LEU:HD12	2.32	0.44
1:D:81:MET:HE2	1:D:85:LEU:HD12	1.99	0.44
1:E:160:GLY:HA3	2:E:501:ADP:H1'	1.98	0.44
1:B:87:THR:O	1:B:87:THR:CG2	2.65	0.44
1:B:457:TRP:O	1:B:461:ARG:HG2	2.18	0.44
1:C:81:MET:HE3	1:C:84:ASN:HA	2.00	0.43
1:F:211:ASP:HA	1:F:214:MET:HB2	2.00	0.43
1:B:319:ARG:HG3	1:B:358:TRP:CH2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:ALA:HB3	1:B:239:GLN:HB2	2.01	0.43
1:D:135:PHE:CD1	1:D:136:PRO:HD2	2.53	0.43
1:A:180:LEU:HD11	1:C:194:PRO:HG2	2.00	0.43
1:C:78:VAL:HG21	1:C:186:GLN:HG2	2.00	0.43
1:D:106:PRO:HB3	1:D:123:MET:CE	2.49	0.43
1:F:8:ILE:HG12	1:F:105:VAL:HG13	2.01	0.43
1:F:326:LEU:O	1:F:356:ARG:HD2	2.18	0.43
1:C:236:ALA:HB3	1:C:239:GLN:HB2	2.01	0.43
1:E:18:GLN:O	1:E:108:PRO:HG2	2.18	0.43
1:A:153:GLU:HG3	1:A:403:LEU:HD11	1.99	0.43
1:A:432:ALA:HB2	1:A:465:TRP:CE2	2.53	0.43
1:B:201:VAL:HG22	1:B:202:LEU:HG	1.99	0.43
1:B:39:LEU:HB3	1:B:104:MET:HE1	1.99	0.43
1:D:357:ALA:HB2	1:D:424:VAL:HG21	2.00	0.43
1:C:336:THR:HG22	1:C:394:TYR:CD1	2.54	0.43
1:E:153:GLU:HG3	1:E:403:LEU:HD11	2.00	0.43
1:F:290:TRP:HB2	1:F:387:PRO:HA	2.01	0.42
1:F:364:ALA:HB1	1:F:420:TYR:CE2	2.53	0.42
1:B:394:TYR:O	1:B:398:GLN:HG2	2.18	0.42
1:C:201:VAL:HG22	1:C:202:LEU:HG	2.00	0.42
1:B:410:LYS:HG3	1:B:411:ASP:H	1.85	0.42
1:C:179:SER:HB3	1:C:188:ASP:HB2	2.01	0.42
1:F:308:PRO:HB3	1:F:314:LEU:CA	2.49	0.42
1:E:292:MET:HE3	1:E:394:TYR:CE2	2.54	0.42
1:F:106:PRO:HB3	1:F:123:MET:HE3	2.01	0.42
1:D:150:LEU:HD12	1:D:403:LEU:HD13	2.01	0.42
1:A:247:MET:HE1	1:A:255:ALA:HB1	2.02	0.42
1:D:408:ASN:O	1:D:409:LYS:C	2.63	0.42
1:B:101:THR:HB	1:B:130:LYS:CE	2.50	0.42
1:D:68:VAL:HG22	1:D:69:ALA:O	2.20	0.42
1:D:176:SER:HB3	1:D:178:GLU:CD	2.45	0.42
1:A:409:LYS:HE3	1:A:409:LYS:HA	2.02	0.41
1:B:135:PHE:HA	1:B:136:PRO:HD3	1.96	0.41
1:B:432:ALA:HB2	1:B:465:TRP:CE2	2.55	0.41
1:C:385:GLY:HA3	1:C:394:TYR:CZ	2.55	0.41
1:D:106:PRO:HB3	1:D:123:MET:HE3	2.02	0.41
1:B:40:SER:O	1:B:124:LEU:HA	2.20	0.41
1:E:353:PRO:HB3	1:E:366:MET:HA	2.02	0.41
1:F:477:ALA:O	1:F:478:GLU:C	2.63	0.41
1:A:153:GLU:HA	1:A:153:GLU:OE1	2.21	0.41
1:A:199:TYR:OH	1:A:284:GLY:O	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:ASP:HA	1:C:214:MET:HB2	2.03	0.41
1:C:432:ALA:HB2	1:C:465:TRP:CE2	2.56	0.41
1:B:287:ALA:HB1	1:B:345:ARG:HG3	2.03	0.41
1:B:345:ARG:HE	1:B:352:ALA:HB2	1.86	0.41
1:C:156:LEU:HB2	1:C:297:VAL:HG22	2.02	0.41
1:F:60:SER:O	1:F:64:GLU:HG3	2.21	0.41
1:F:325:LEU:HD23	1:F:401:SER:HB2	2.03	0.41
1:F:385:GLY:HA3	1:F:394:TYR:CZ	2.56	0.41
1:C:158:LYS:O	1:C:294:GLN:HA	2.22	0.41
1:E:457:TRP:O	1:E:461:ARG:HG2	2.20	0.41
1:A:290:TRP:CD1	1:A:290:TRP:C	2.98	0.40
1:A:211:ASP:HA	1:A:214:MET:HB2	2.03	0.40
1:B:183:PRO:CG	1:D:286:PHE:HE2	2.35	0.40
1:D:68:VAL:CG2	1:D:69:ALA:N	2.84	0.40
1:A:193:GLN:HG2	6:A:686:HOH:O	2.20	0.40
1:A:319:ARG:HG3	1:A:358:TRP:CH2	2.57	0.40
1:D:445:PHE:O	1:D:449:PHE:HB2	2.21	0.40
1:F:297:VAL:CG1	1:F:298:ASP:N	2.84	0.40
1:B:82:VAL:HG21	1:B:460:LEU:HD12	2.03	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	488/510 (96%)	486 (100%)	2 (0%)	0	100	100
1	B	484/510 (95%)	481 (99%)	3 (1%)	0	100	100
1	C	488/510 (96%)	485 (99%)	3 (1%)	0	100	100
1	D	488/510 (96%)	485 (99%)	3 (1%)	0	100	100
1	E	485/510 (95%)	481 (99%)	4 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	488/510 (96%)	485 (99%)	3 (1%)	0	100	100
All	All	2921/3060 (96%)	2903 (99%)	18 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	408/427 (96%)	401 (98%)	7 (2%)	53	72
1	B	408/427 (96%)	404 (99%)	4 (1%)	68	82
1	C	408/427 (96%)	398 (98%)	10 (2%)	42	60
1	D	408/427 (96%)	402 (98%)	6 (2%)	57	75
1	E	407/427 (95%)	401 (98%)	6 (2%)	57	75
1	F	408/427 (96%)	403 (99%)	5 (1%)	63	79
All	All	2447/2562 (96%)	2409 (98%)	38 (2%)	55	73

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	GLU
1	A	88	THR
1	A	212	ASP
1	A	292	MET
1	A	301	THR
1	A	373	THR
1	A	417	GLU
1	B	80	SER
1	B	95	GLU
1	B	292	MET
1	B	301	THR
1	C	10	LYS
1	C	78	VAL

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Mol	Chain	Res	Type
1	C	80	SER
1	C	145	LYS
1	C	148	LYS
1	C	292	MET
1	C	301	THR
1	C	373	THR
1	C	409	LYS
1	C	478	GLU
1	D	39	LEU
1	D	85	LEU
1	D	123	MET
1	D	162	GLU
1	D	251	GLU
1	D	292	MET
1	E	64	GLU
1	E	66	SER
1	E	123	MET
1	E	292	MET
1	E	339	THR
1	E	386	GLU
1	F	64	GLU
1	F	123	MET
1	F	153	GLU
1	F	234	GLU
1	F	469	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	99	ASN
1	A	209	GLN
1	A	408	ASN
1	B	128	HIS
1	B	270	HIS
1	B	408	ASN
1	B	416	GLN
1	C	84	ASN
1	C	186	GLN
1	C	205	HIS
1	C	209	GLN
1	D	193	GLN

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Mol	Chain	Res	Type
1	D	270	HIS
1	E	128	HIS
1	E	186	GLN
1	E	191	GLN
1	E	205	HIS
1	F	84	ASN
1	F	347	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 36 ligands modelled in this entry, 18 are monoatomic - leaving 18 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	ADP	C	501	5,4	28,29,29	1.39	6 (21%)	43,45,45	2.03	12 (27%)
2	ADP	B	501	5,4	28,29,29	1.39	5 (17%)	43,45,45	2.03	11 (25%)
2	ADP	D	501	5,4	28,29,29	1.34	4 (14%)	43,45,45	2.03	13 (30%)
2	ADP	A	501	5,4	28,29,29	1.39	6 (21%)	43,45,45	2.00	12 (27%)
2	ADP	E	501	5,4	28,29,29	1.43	5 (17%)	43,45,45	2.00	12 (27%)
2	ADP	F	501	5,4	28,29,29	1.38	6 (21%)	43,45,45	2.01	13 (30%)

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	ADP	C	501	5,4	-	1/16/32/32	0/3/3/3
2	ADP	B	501	5,4	-	4/16/32/32	0/3/3/3
2	ADP	D	501	5,4	-	2/16/32/32	0/3/3/3
2	ADP	A	501	5,4	-	1/16/32/32	0/3/3/3
2	ADP	E	501	5,4	-	2/16/32/32	0/3/3/3
2	ADP	F	501	5,4	-	4/16/32/32	0/3/3/3

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	501	ADP	C5-C4	4.18	1.46	1.39
2	F	501	ADP	C5-C4	4.12	1.46	1.39
2	A	501	ADP	C5-C4	4.12	1.46	1.39
2	C	501	ADP	C5-C4	4.12	1.46	1.39
2	B	501	ADP	C5-C4	4.12	1.46	1.39
2	D	501	ADP	C5-C4	4.06	1.46	1.39
2	E	501	ADP	C5-C6	2.85	1.48	1.41
2	B	501	ADP	C5-C6	2.79	1.48	1.41
2	D	501	ADP	C5-C6	2.77	1.48	1.41
2	F	501	ADP	C5-C6	2.73	1.48	1.41
2	E	501	ADP	C8-N7	2.73	1.36	1.31
2	C	501	ADP	C5-C6	2.72	1.48	1.41
2	F	501	ADP	C8-N7	2.72	1.36	1.31
2	A	501	ADP	C5-C6	2.66	1.48	1.41
2	B	501	ADP	C8-N7	2.59	1.36	1.31
2	C	501	ADP	C8-N7	2.55	1.36	1.31
2	D	501	ADP	C8-N7	2.54	1.36	1.31
2	E	501	ADP	PA-O3A	2.52	1.62	1.59
2	A	501	ADP	C8-N7	2.47	1.36	1.31
2	A	501	ADP	PA-O3A	2.37	1.62	1.59
2	C	501	ADP	PA-O3A	2.34	1.62	1.59
2	E	501	ADP	C4-N9	-2.30	1.32	1.37
2	B	501	ADP	PA-O3A	2.23	1.61	1.59
2	A	501	ADP	C5-N7	-2.21	1.35	1.39
2	F	501	ADP	C4-N9	-2.16	1.33	1.37
2	C	501	ADP	C4-N9	-2.13	1.33	1.37
2	A	501	ADP	C4-N9	-2.09	1.33	1.37
2	C	501	ADP	C5-N7	-2.05	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	501	ADP	PA-O3A	2.05	1.61	1.59
2	B	501	ADP	C5-N7	-2.05	1.35	1.39
2	D	501	ADP	C5-N7	-2.04	1.35	1.39
2	F	501	ADP	C5-N7	-2.03	1.35	1.39

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	ADP	C5-C4-N3	-5.57	119.05	126.72
2	D	501	ADP	C5-C4-N3	-5.31	119.40	126.72
2	F	501	ADP	C5-C4-N3	-5.20	119.55	126.72
2	A	501	ADP	C5-C4-N3	-5.19	119.58	126.72
2	C	501	ADP	C5-C4-N3	-5.02	119.80	126.72
2	E	501	ADP	C5-C4-N3	-4.98	119.86	126.72
2	B	501	ADP	N3-C4-N9	4.24	134.37	127.17
2	D	501	ADP	N3-C4-N9	4.20	134.32	127.17
2	C	501	ADP	N3-C2-N1	-4.18	122.25	128.58
2	F	501	ADP	N3-C2-N1	-4.12	122.34	128.58
2	B	501	ADP	C2-N3-C4	4.10	121.85	111.83
2	A	501	ADP	N3-C2-N1	-4.08	122.40	128.58
2	A	501	ADP	N3-C4-N9	4.08	134.10	127.17
2	B	501	ADP	N3-C2-N1	-4.07	122.42	128.58
2	D	501	ADP	N3-C2-N1	-4.04	122.46	128.58
2	E	501	ADP	N3-C2-N1	-4.04	122.46	128.58
2	E	501	ADP	C4-C5-N7	-4.03	105.98	110.58
2	B	501	ADP	C4-C5-N7	-4.00	106.00	110.58
2	F	501	ADP	C2-N3-C4	3.97	121.53	111.83
2	D	501	ADP	C2-N3-C4	3.94	121.46	111.83
2	C	501	ADP	N3-C4-N9	3.92	133.83	127.17
2	A	501	ADP	C2-N3-C4	3.90	121.35	111.83
2	C	501	ADP	C4-C5-N7	-3.89	106.13	110.58
2	C	501	ADP	C2-N3-C4	3.88	121.31	111.83
2	E	501	ADP	C2-N3-C4	3.88	121.30	111.83
2	D	501	ADP	C4-C5-N7	-3.87	106.16	110.58
2	F	501	ADP	C4-C5-N7	-3.85	106.18	110.58
2	F	501	ADP	N3-C4-N9	3.84	133.69	127.17
2	C	501	ADP	C4-N9-C8	3.76	109.69	105.74
2	D	501	ADP	C4-N9-C8	3.75	109.68	105.74
2	A	501	ADP	C4-C5-N7	-3.72	106.33	110.58
2	E	501	ADP	N3-C4-N9	3.58	133.26	127.17
2	A	501	ADP	O4'-C1'-N9	3.56	114.93	108.09
2	E	501	ADP	O4'-C1'-N9	3.51	114.84	108.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	ADP	C4-N9-C8	3.51	109.43	105.74
2	E	501	ADP	C4-N9-C8	3.40	109.31	105.74
2	B	501	ADP	C4-N9-C8	3.40	109.31	105.74
2	F	501	ADP	C4-N9-C8	3.33	109.23	105.74
2	E	501	ADP	C6-C5-N7	3.12	138.11	132.09
2	E	501	ADP	C4-N9-C1'	-3.10	119.39	126.63
2	B	501	ADP	C5-N7-C8	3.09	108.31	103.45
2	B	501	ADP	O4'-C1'-N9	3.09	114.02	108.09
2	D	501	ADP	C5-N7-C8	3.07	108.27	103.45
2	C	501	ADP	N9-C8-N7	-3.05	109.60	113.94
2	D	501	ADP	N9-C8-N7	-3.05	109.61	113.94
2	C	501	ADP	C5-N7-C8	3.03	108.20	103.45
2	F	501	ADP	C6-C5-N7	3.00	137.88	132.09
2	C	501	ADP	C6-C5-N7	2.99	137.86	132.09
2	E	501	ADP	N9-C8-N7	-2.97	109.72	113.94
2	E	501	ADP	C5-N7-C8	2.97	108.11	103.45
2	F	501	ADP	O4'-C1'-N9	2.95	113.75	108.09
2	A	501	ADP	C5-N7-C8	2.93	108.05	103.45
2	B	501	ADP	C6-C5-N7	2.92	137.72	132.09
2	B	501	ADP	N9-C8-N7	-2.92	109.80	113.94
2	F	501	ADP	C5-N7-C8	2.89	107.99	103.45
2	F	501	ADP	N9-C8-N7	-2.88	109.84	113.94
2	C	501	ADP	C4-N9-C1'	-2.87	119.92	126.63
2	A	501	ADP	N9-C8-N7	-2.87	109.87	113.94
2	F	501	ADP	C4-N9-C1'	-2.86	119.93	126.63
2	D	501	ADP	C6-C5-N7	2.82	137.52	132.09
2	A	501	ADP	C6-C5-N7	2.79	137.48	132.09
2	D	501	ADP	C4-N9-C1'	-2.73	120.26	126.63
2	B	501	ADP	C4-N9-C1'	-2.48	120.83	126.63
2	C	501	ADP	O4'-C1'-N9	2.44	112.78	108.09
2	C	501	ADP	C2-N1-C6	2.30	122.50	118.73
2	D	501	ADP	C2'-C1'-N9	2.21	118.80	113.30
2	A	501	ADP	C2-N1-C6	2.20	122.34	118.73
2	A	501	ADP	C4-N9-C1'	-2.19	121.51	126.63
2	D	501	ADP	O4'-C1'-N9	2.19	112.29	108.09
2	F	501	ADP	O2A-PA-O3A	2.15	113.08	107.27
2	F	501	ADP	C2-N1-C6	2.14	122.24	118.73
2	E	501	ADP	C2-N1-C6	2.09	122.16	118.73
2	D	501	ADP	C2-N1-C6	2.03	122.06	118.73

There are no chirality outliers.

All (14) torsion outliers are listed below:

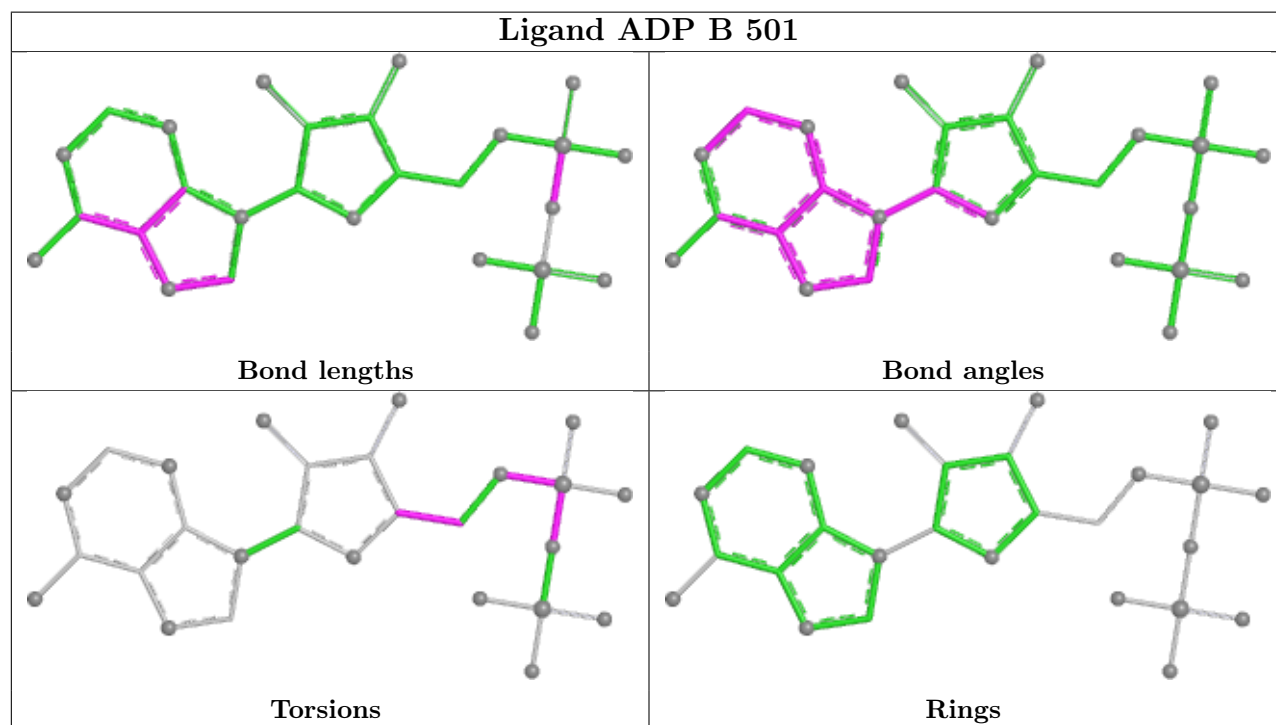
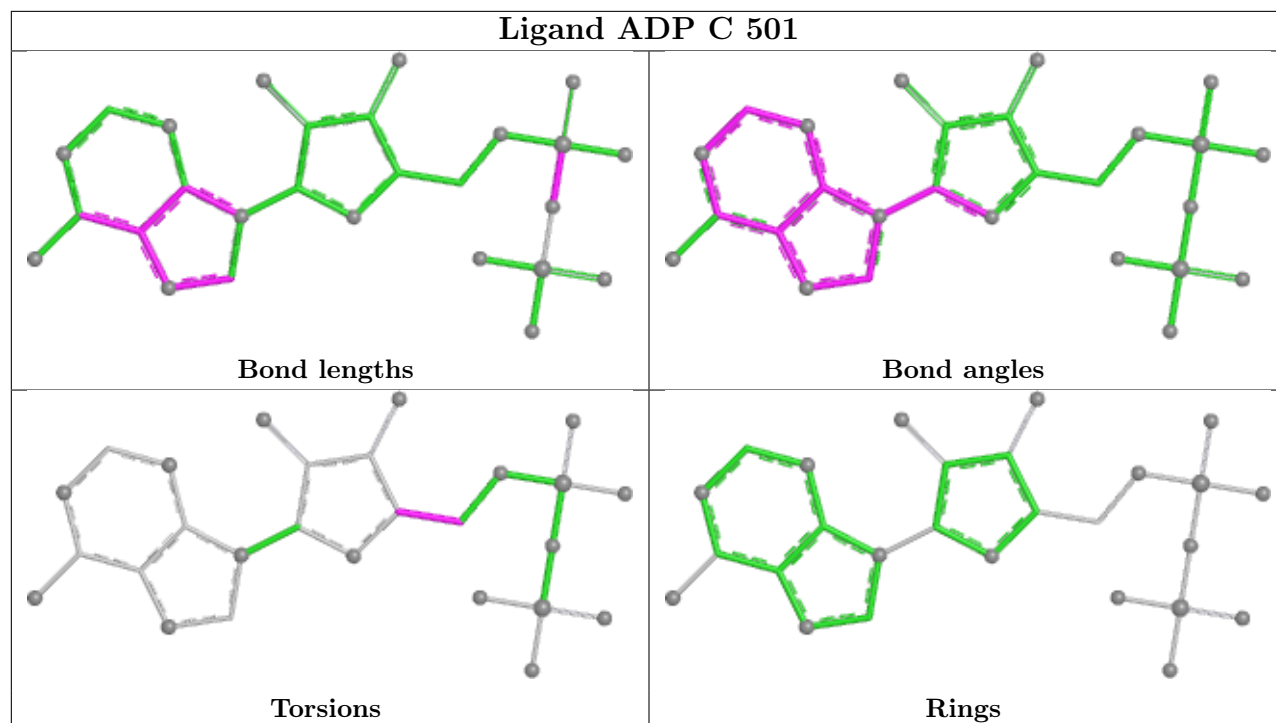
Mol	Chain	Res	Type	Atoms
2	F	501	ADP	PA-O3A-PB-O3B
2	F	501	ADP	O4'-C4'-C5'-O5'
2	E	501	ADP	PA-O3A-PB-O2B
2	D	501	ADP	O4'-C4'-C5'-O5'
2	B	501	ADP	C5'-O5'-PA-O1A
2	C	501	ADP	O4'-C4'-C5'-O5'
2	B	501	ADP	O4'-C4'-C5'-O5'
2	E	501	ADP	O4'-C4'-C5'-O5'
2	F	501	ADP	PA-O3A-PB-O1B
2	D	501	ADP	PA-O3A-PB-O2B
2	A	501	ADP	O4'-C4'-C5'-O5'
2	B	501	ADP	PB-O3A-PA-O2A
2	F	501	ADP	C3'-C4'-C5'-O5'
2	B	501	ADP	PB-O3A-PA-O1A

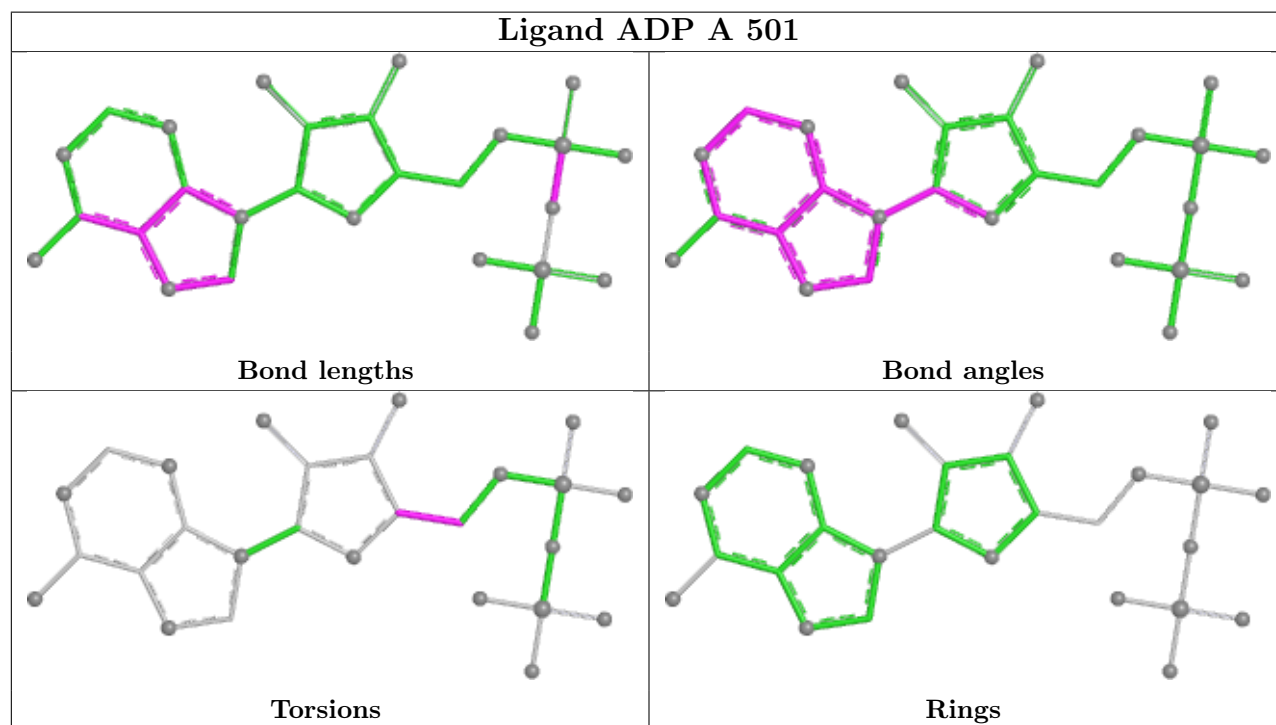
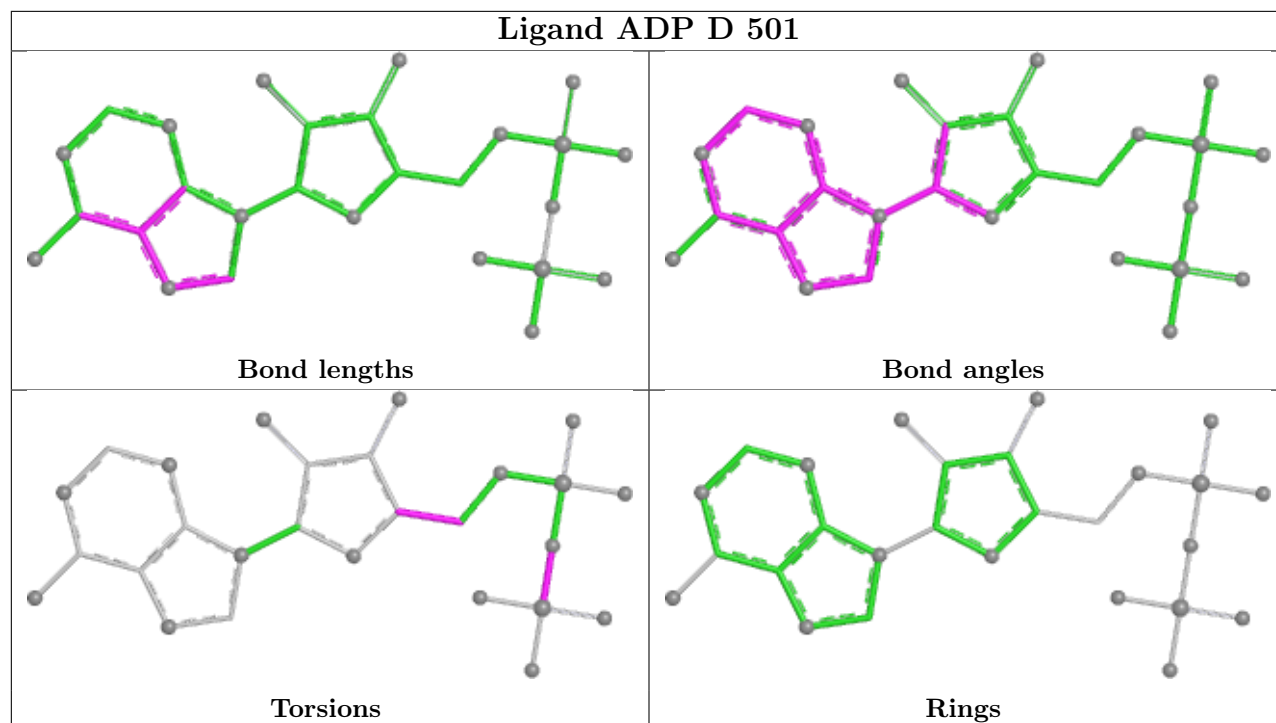
There are no ring outliers.

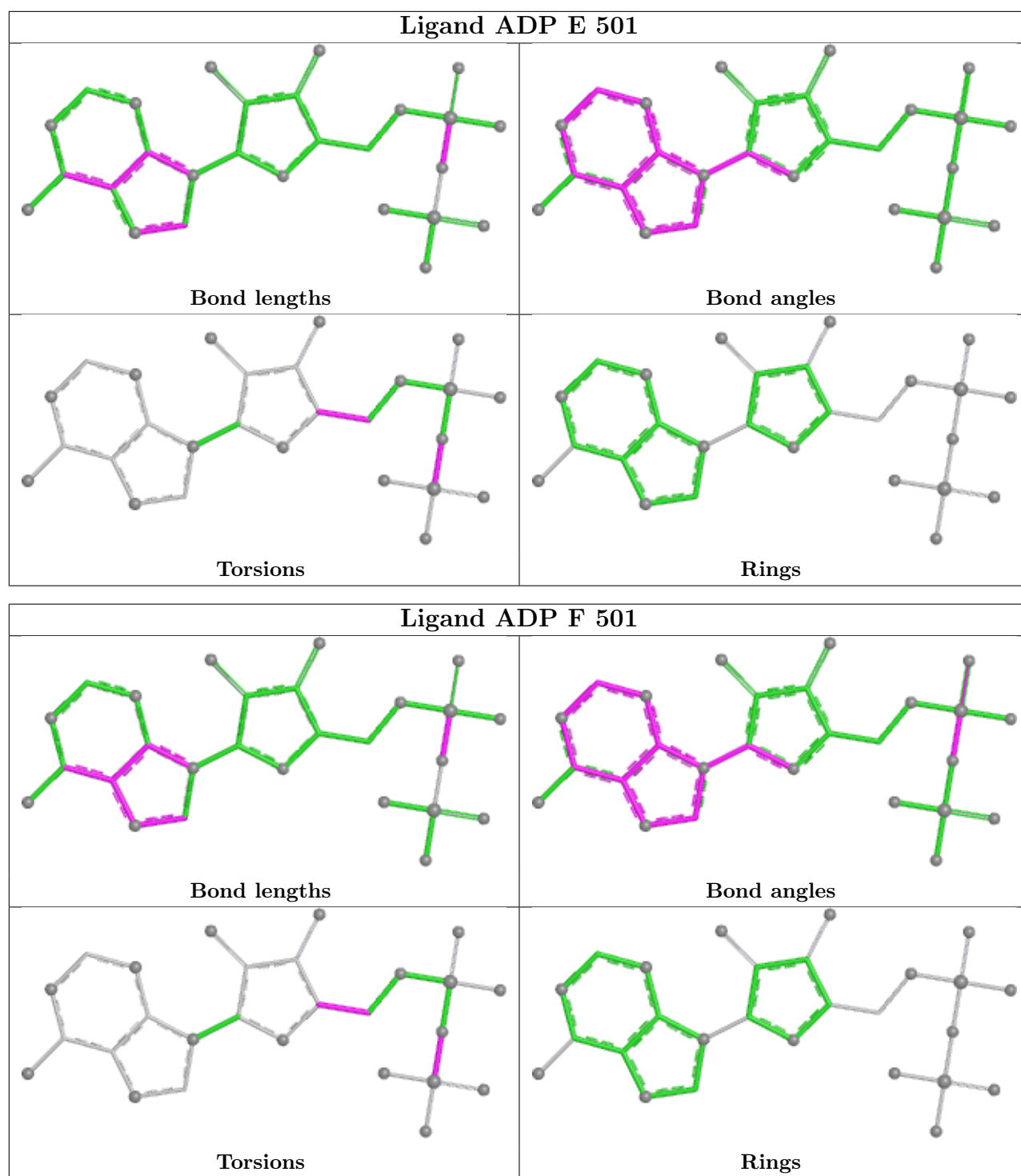
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	501	ADP	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	490/510 (96%)	0.57	36 (7%)	21	23	30, 53, 89, 114	0
1	B	488/510 (95%)	0.57	41 (8%)	17	18	32, 53, 86, 114	0
1	C	490/510 (96%)	0.60	27 (5%)	30	32	35, 51, 82, 103	0
1	D	490/510 (96%)	0.52	27 (5%)	30	32	32, 50, 78, 97	0
1	E	489/510 (95%)	0.44	20 (4%)	41	43	33, 49, 78, 103	0
1	F	490/510 (96%)	0.47	18 (3%)	45	48	33, 50, 82, 105	0
All	All	2937/3060 (95%)	0.53	169 (5%)	29	31	30, 51, 83, 114	0

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	92	PHE	6.9
1	A	420	TYR	6.1
1	E	90	GLY	5.5
1	A	64	GLU	5.2
1	D	351	LEU	5.1
1	B	88	THR	4.7
1	C	423	GLN	4.4
1	B	345	ARG	4.4
1	A	419	PRO	4.1
1	C	351	LEU	4.1
1	E	92	PHE	3.8
1	D	410	LYS	3.8
1	A	88	THR	3.7
1	E	278	CYS	3.6
1	B	440	HIS	3.6
1	C	85	LEU	3.5
1	C	148	LYS	3.4
1	E	85	LEU	3.3
1	B	92	PHE	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	351	LEU	3.3
1	E	418	SER	3.3
1	A	89	ALA	3.2
1	D	475	GLU	3.2
1	A	349	TYR	3.2
1	B	357	ALA	3.2
1	C	70	ILE	3.2
1	C	302	ARG	3.2
1	C	364	ALA	3.1
1	F	423	GLN	3.1
1	C	154	GLY	3.1
1	C	79	SER	3.0
1	E	311	GLY	3.0
1	B	417	GLU	3.0
1	C	131	SER	3.0
1	B	86	PHE	2.9
1	B	421	ASP	2.9
1	D	407	LYS	2.9
1	E	443	GLU	2.9
1	B	351	LEU	2.9
1	B	130	LYS	2.8
1	A	313	VAL	2.8
1	E	87	THR	2.8
1	B	91	ASP	2.8
1	C	418	SER	2.8
1	D	318	GLY	2.8
1	D	286	PHE	2.8
1	A	424	VAL	2.8
1	C	409	LYS	2.8
1	A	354	ASP	2.8
1	C	80	SER	2.8
1	C	321	TYR	2.8
1	D	20	ASP	2.7
1	E	94	ASP	2.7
1	F	311	GLY	2.7
1	A	357	ALA	2.7
1	B	359	ALA	2.7
1	C	286	PHE	2.7
1	D	350	SER	2.7
1	D	370	VAL	2.7
1	C	64	GLU	2.7
1	E	138	CYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	349	TYR	2.6
1	B	99	ASN	2.6
1	B	87	THR	2.6
1	B	476	ALA	2.6
1	D	476	ALA	2.6
1	A	443	GLU	2.6
1	B	74	PHE	2.6
1	A	311	GLY	2.6
1	F	87	THR	2.6
1	B	415	LEU	2.6
1	A	91	ASP	2.6
1	A	90	GLY	2.6
1	B	97	SER	2.6
1	D	423	GLN	2.6
1	C	99	ASN	2.6
1	B	412	PRO	2.6
1	B	473	GLY	2.6
1	F	373	THR	2.6
1	B	313	VAL	2.5
1	E	64	GLU	2.5
1	A	421	ASP	2.5
1	A	87	THR	2.5
1	A	440	HIS	2.5
1	B	286	PHE	2.5
1	D	178	GLU	2.5
1	A	360	LYS	2.5
1	E	373	THR	2.5
1	A	286	PHE	2.5
1	F	378	SER	2.5
1	B	361	ASP	2.5
1	C	300	ASP	2.5
1	D	99	ASN	2.5
1	A	205	HIS	2.5
1	F	310	GLU	2.4
1	D	406	ILE	2.4
1	A	412	PRO	2.4
1	C	12	ASN	2.4
1	A	475	GLU	2.4
1	F	414	GLU	2.4
1	E	357	ALA	2.4
1	F	370	VAL	2.4
1	D	302	ARG	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	321	TYR	2.3
1	F	99	ASN	2.3
1	A	86	PHE	2.3
1	A	129	TRP	2.3
1	B	420	TYR	2.3
1	C	324	GLY	2.3
1	F	90	GLY	2.3
1	A	81	MET	2.3
1	D	64	GLU	2.3
1	E	133	GLU	2.3
1	B	418	SER	2.3
1	B	451	ASP	2.2
1	E	30	ASP	2.2
1	C	251	GLU	2.2
1	C	372	ALA	2.2
1	D	69	ALA	2.2
1	B	419	PRO	2.2
1	B	410	LYS	2.2
1	A	418	SER	2.2
1	D	349	TYR	2.2
1	C	362	ASN	2.2
1	F	417	GLU	2.2
1	B	416	GLN	2.2
1	B	138	CYS	2.2
1	A	93	GLU	2.1
1	B	64	GLU	2.1
1	B	178	GLU	2.1
1	D	162	GLU	2.1
1	F	133	GLU	2.1
1	B	364	ALA	2.1
1	E	89	ALA	2.1
1	F	89	ALA	2.1
1	F	302	ARG	2.1
1	A	130	LYS	2.1
1	E	24	LYS	2.1
1	B	369	VAL	2.1
1	D	424	VAL	2.1
1	D	249	GLY	2.1
1	D	418	SER	2.1
1	F	69	ALA	2.1
1	E	475	GLU	2.1
1	F	96	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	152	ASP	2.1
1	A	97	SER	2.1
1	B	407	LYS	2.1
1	C	475	GLU	2.1
1	E	310	GLU	2.1
1	B	30	ASP	2.1
1	E	304	ASN	2.1
1	F	16	ASN	2.1
1	B	68	VAL	2.1
1	B	324	GLY	2.1
1	C	299	LYS	2.0
1	F	301	THR	2.0
1	A	417	GLU	2.0
1	D	310	GLU	2.0
1	A	84	ASN	2.0
1	A	96	LEU	2.0
1	D	96	LEU	2.0
1	D	403	LEU	2.0
1	C	319	ARG	2.0
1	A	318	GLY	2.0
1	D	474	ALA	2.0
1	B	423	GLN	2.0
1	B	85	LEU	2.0
1	D	304	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

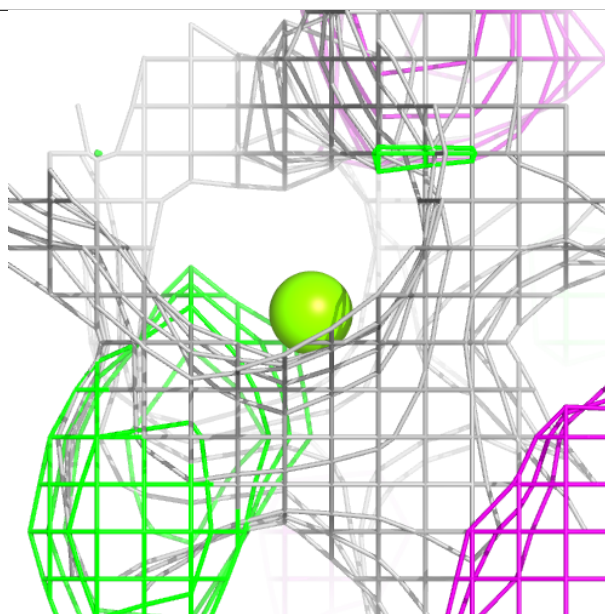
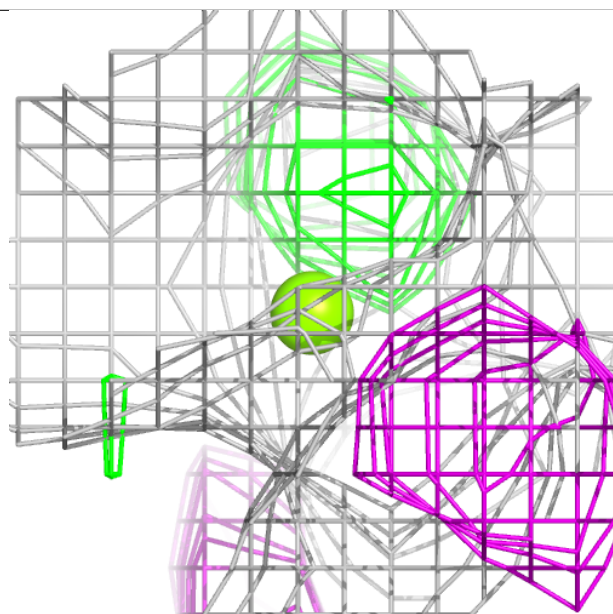
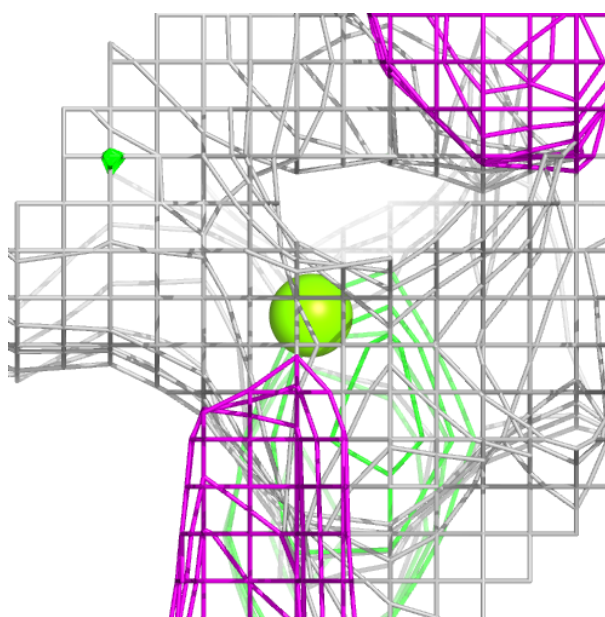
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	P3S	B	502[A]	15/15	0.78	0.17	54,72,80,81	1
3	P3S	B	502[B]	15/15	0.78	0.17	54,72,80,81	1
3	P3S	C	502[A]	15/15	0.78	0.16	52,79,95,103	1
3	P3S	C	502[B]	15/15	0.78	0.16	52,79,95,103	1
5	MG	A	505	1/1	0.78	0.14	53,53,53,53	0
5	MG	D	505	1/1	0.81	0.16	55,55,55,55	0
3	P3S	D	502[A]	15/15	0.82	0.16	47,72,83,88	1
5	MG	C	505	1/1	0.82	0.10	64,64,64,64	0
3	P3S	D	502[B]	15/15	0.82	0.16	47,72,83,88	1
5	MG	B	505	1/1	0.83	0.17	54,54,54,54	0
4	MN	D	504	1/1	0.84	0.13	79,79,79,79	0
3	P3S	A	502[B]	15/15	0.85	0.13	50,63,73,77	1
3	P3S	A	502[A]	15/15	0.85	0.13	50,63,73,77	1
4	MN	F	504	1/1	0.88	0.13	63,63,63,63	0
5	MG	F	505	1/1	0.89	0.16	45,45,45,45	0
2	ADP	D	501	27/27	0.90	0.11	41,50,56,57	0
4	MN	B	504	1/1	0.91	0.14	80,80,80,80	0
2	ADP	A	501	27/27	0.91	0.10	40,49,59,63	0
3	P3S	F	502[A]	15/15	0.91	0.14	43,52,55,56	1
5	MG	E	505	1/1	0.91	0.09	41,41,41,41	0
3	P3S	F	502[B]	15/15	0.91	0.14	43,52,55,56	1
2	ADP	B	501	27/27	0.92	0.10	42,51,60,62	0
2	ADP	C	501	27/27	0.92	0.10	47,56,68,72	0
4	MN	C	503	1/1	0.93	0.09	59,59,59,59	0
2	ADP	F	501	27/27	0.94	0.09	41,45,56,58	0
4	MN	C	504	1/1	0.94	0.14	85,85,85,85	0
2	ADP	E	501	27/27	0.95	0.09	38,44,55,56	0
4	MN	A	504	1/1	0.96	0.13	65,65,65,65	0
3	P3S	E	502[A]	15/15	0.96	0.08	41,44,52,52	1
3	P3S	E	502[B]	15/15	0.96	0.08	41,44,52,52	1
4	MN	F	503	1/1	0.97	0.11	47,47,47,47	0
4	MN	E	503	1/1	0.97	0.13	61,61,61,61	0
4	MN	E	504	1/1	0.98	0.06	59,59,59,59	0
4	MN	A	503	1/1	0.98	0.11	54,54,54,54	0
4	MN	D	503	1/1	0.98	0.05	50,50,50,50	0
4	MN	B	503	1/1	0.99	0.06	54,54,54,54	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.

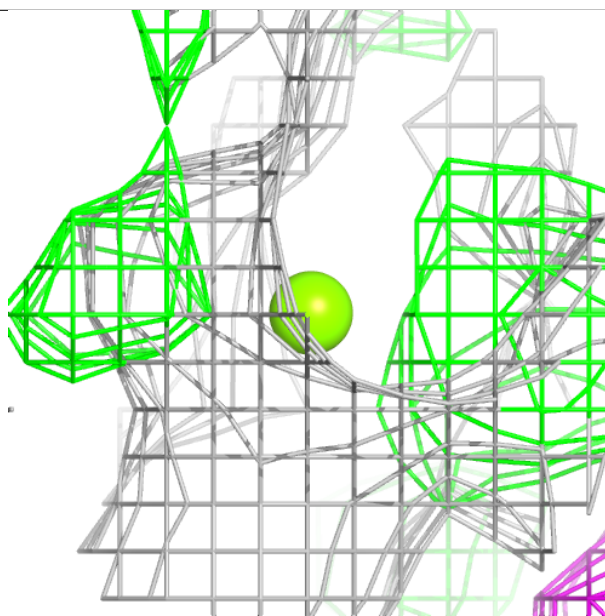
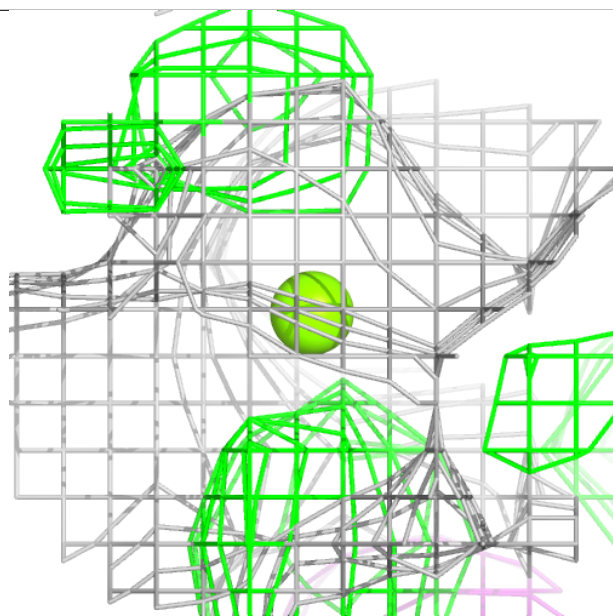
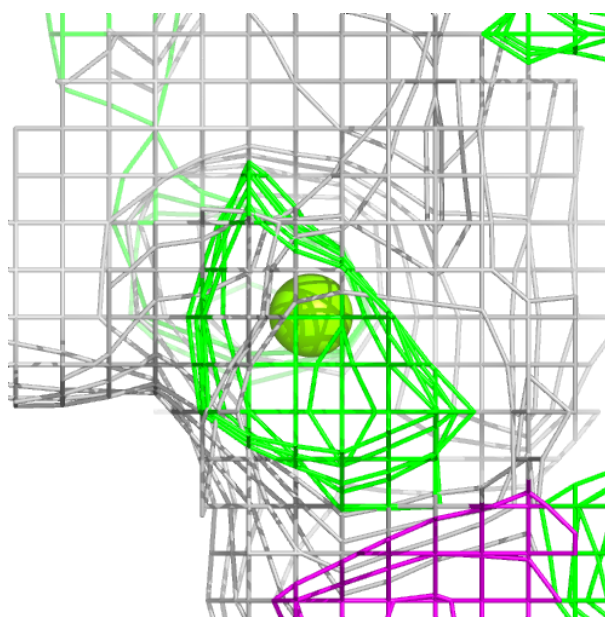
Electron density around MG A 505:

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



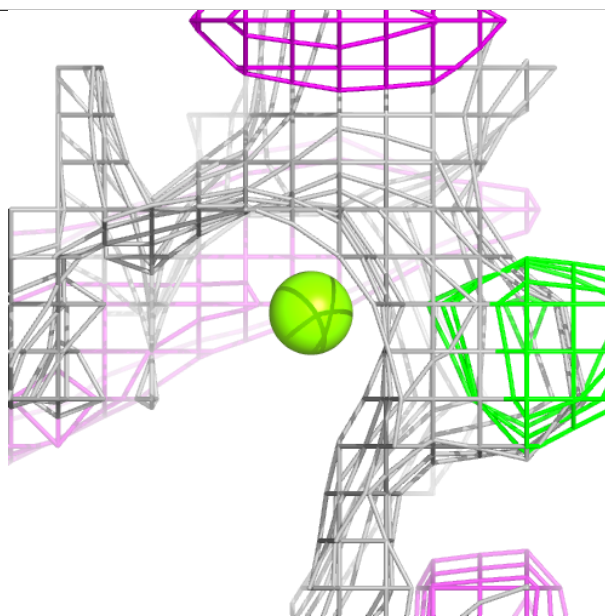
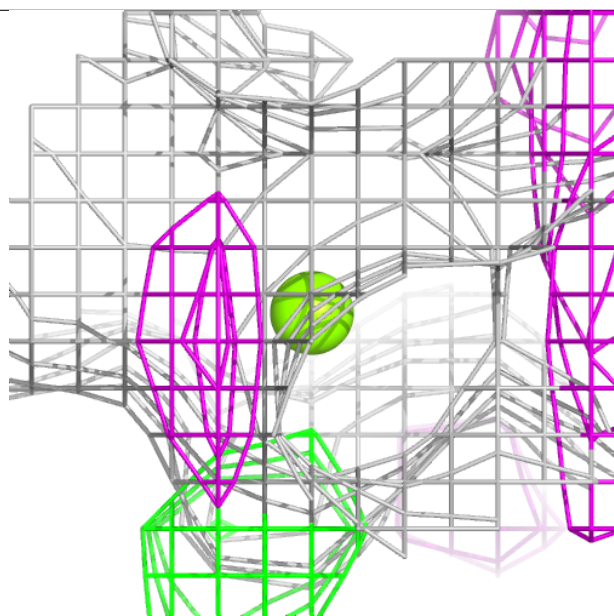
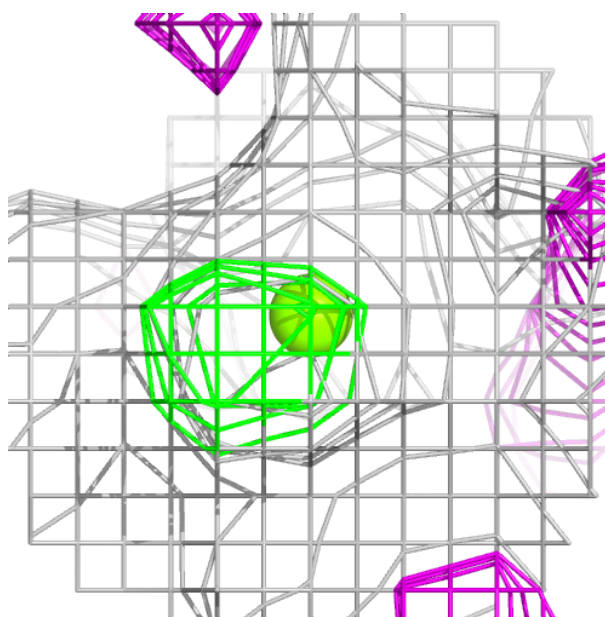
Electron density around MG D 505:

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



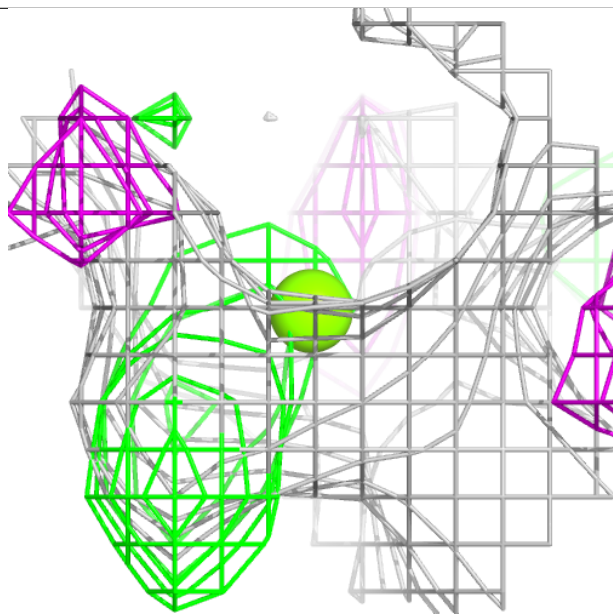
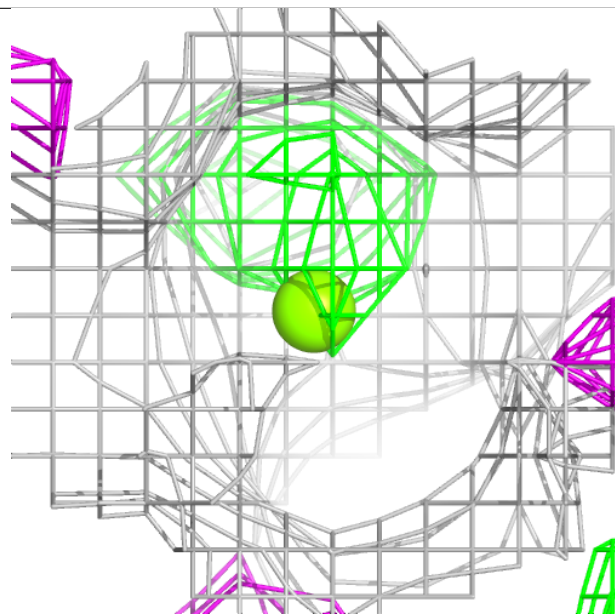
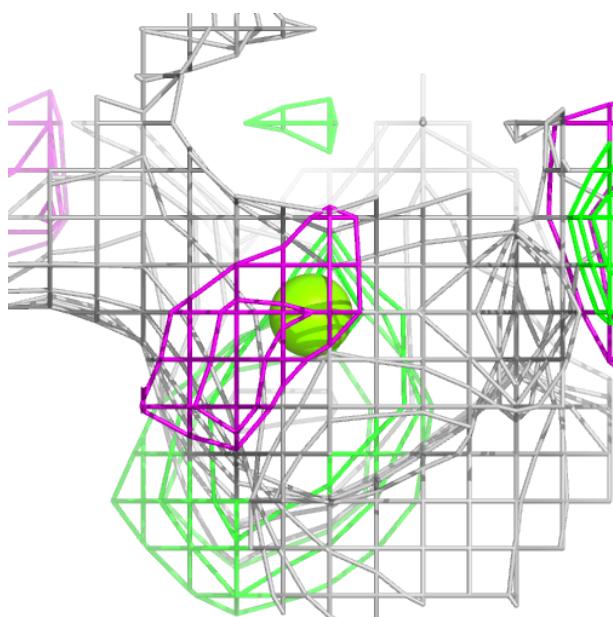
Electron density around MG C 505:

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



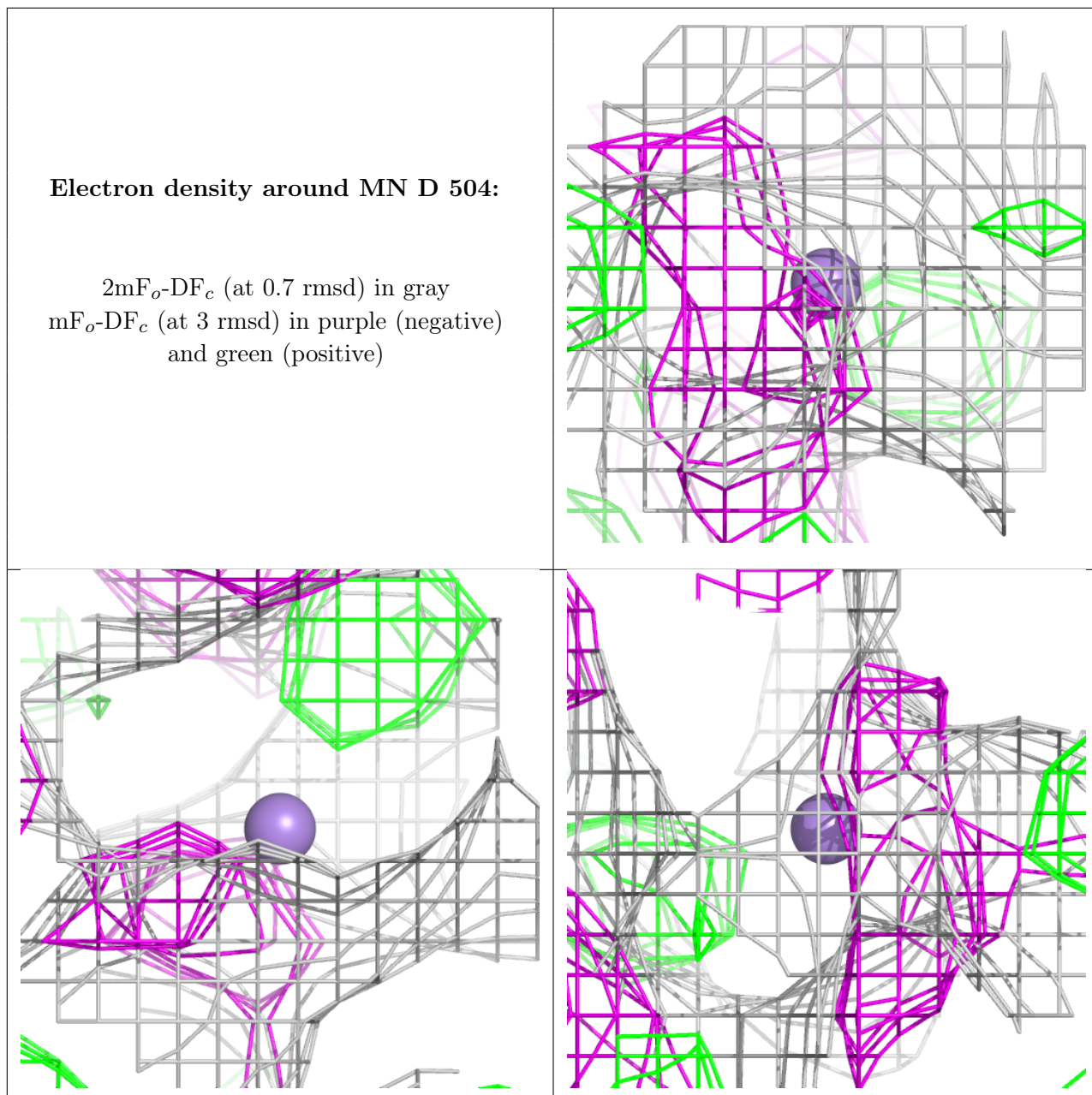
Electron density around MG B 505:

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



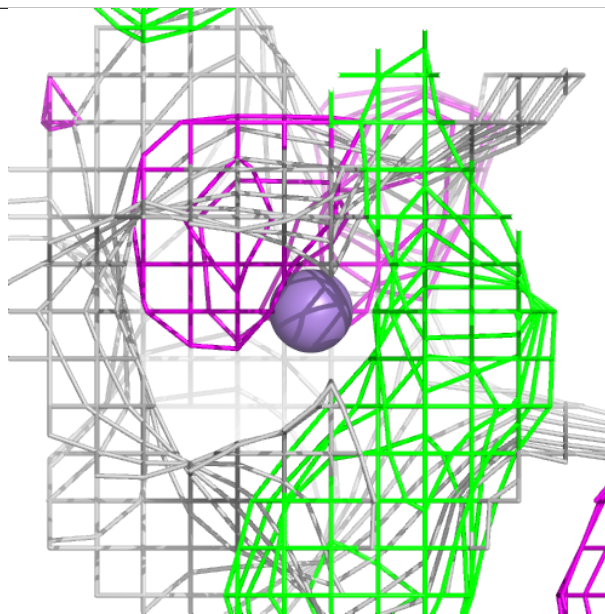
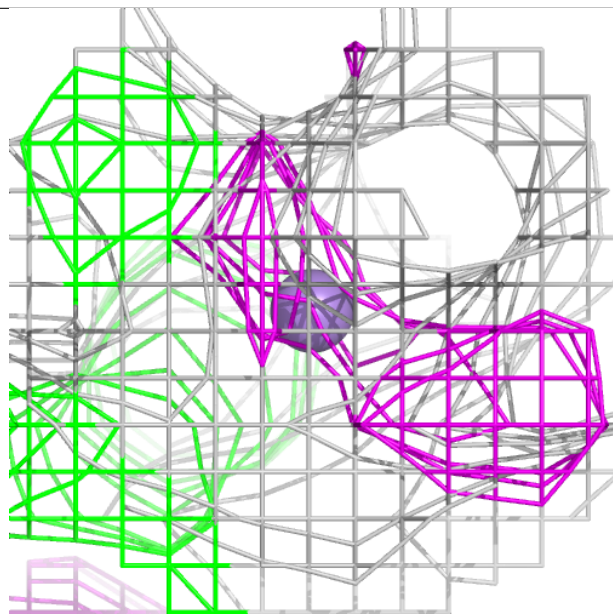
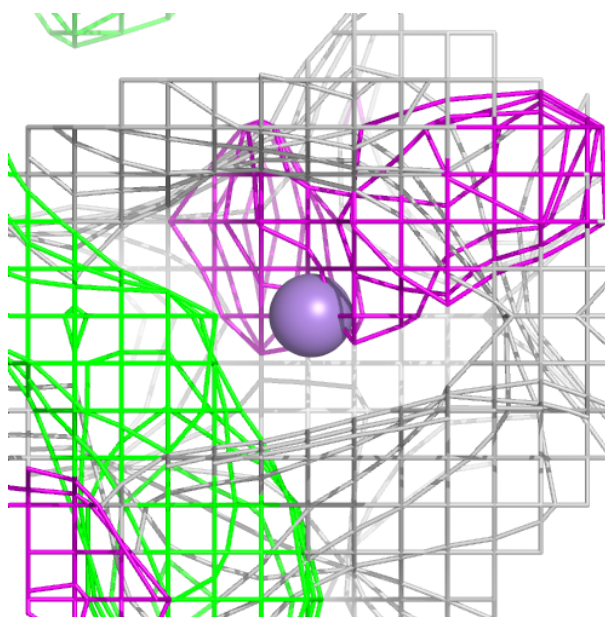
Electron density around MN D 504:

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



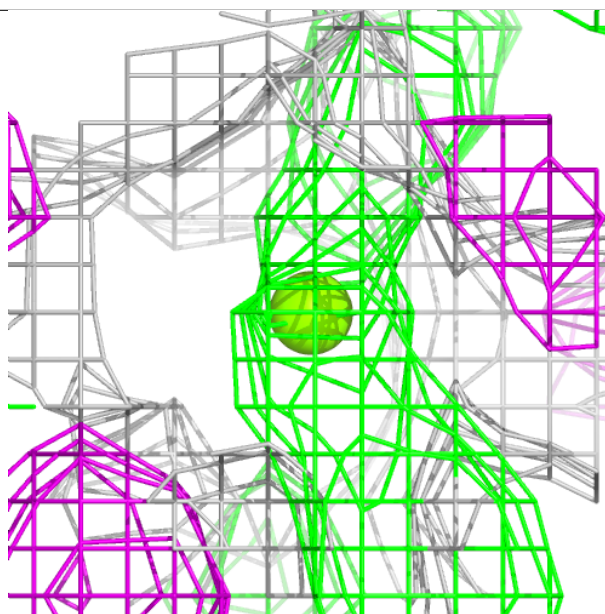
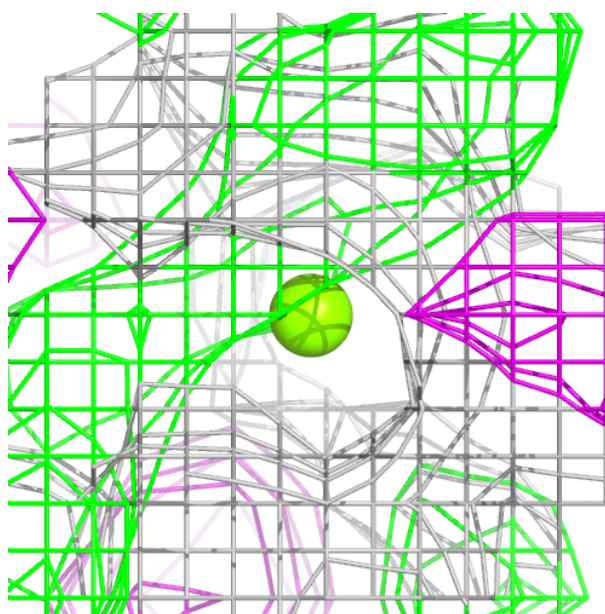
Electron density around MN F 504:

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



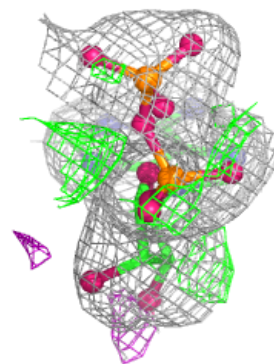
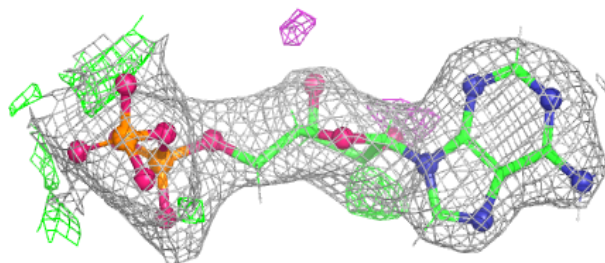
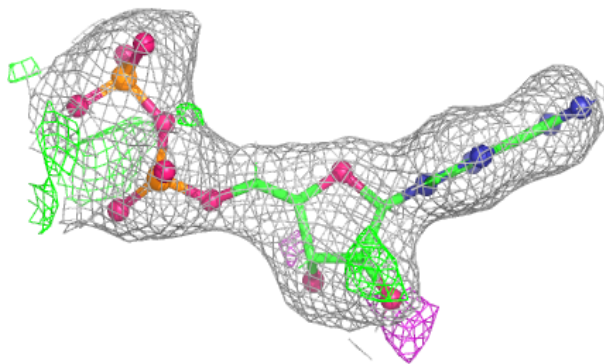
Electron density around MG F 505:

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



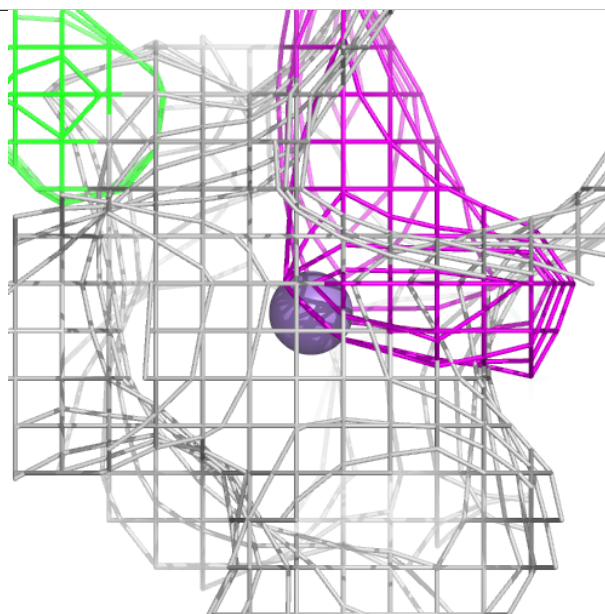
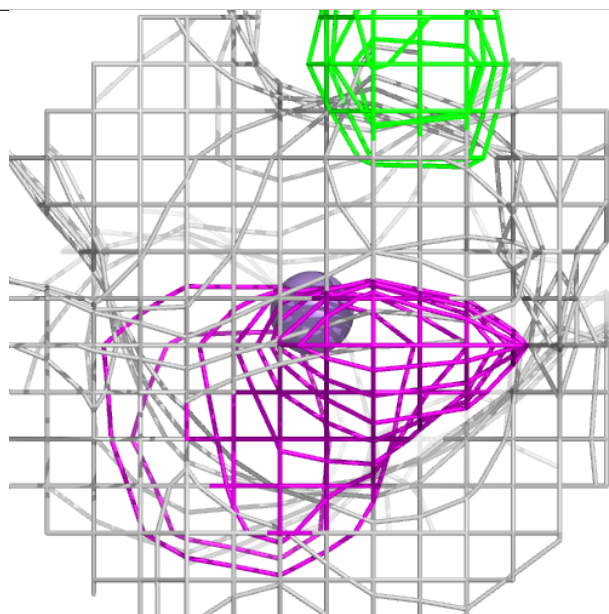
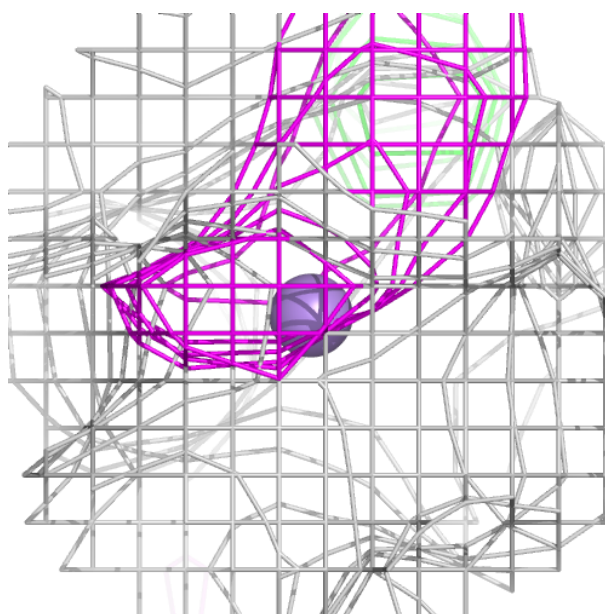
Electron density around ADP D 501:

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



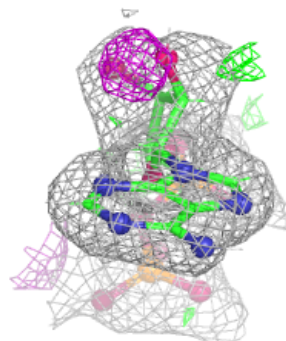
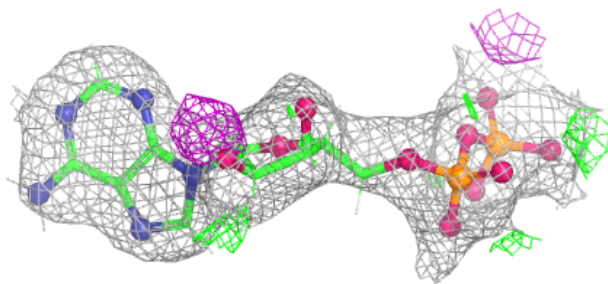
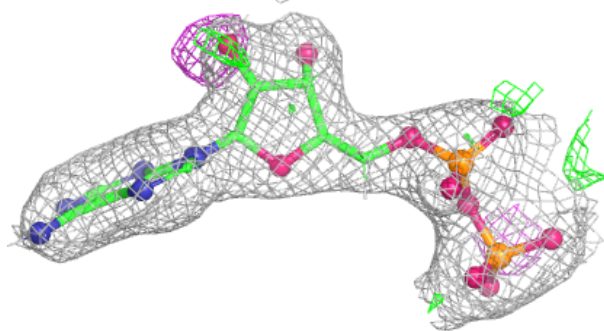
Electron density around MN B 504:

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



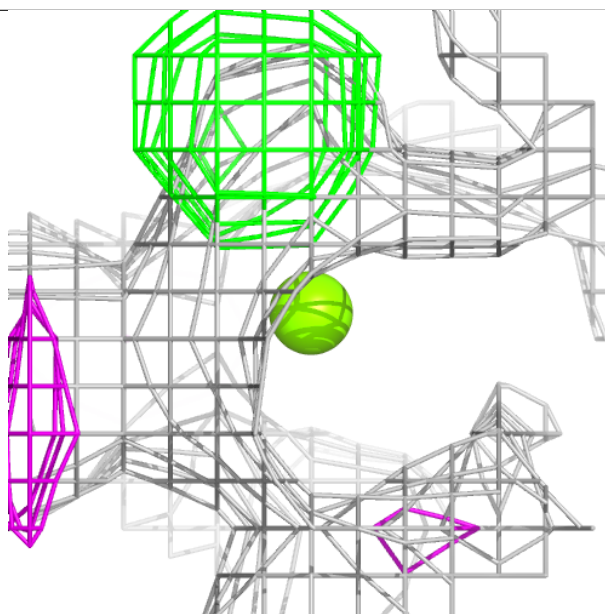
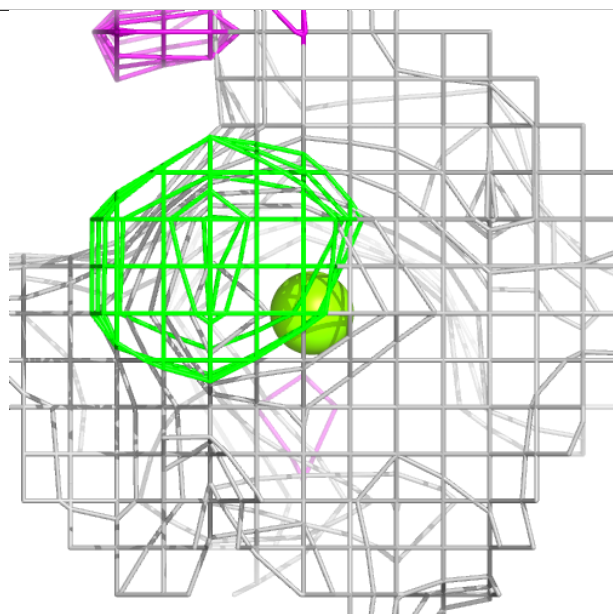
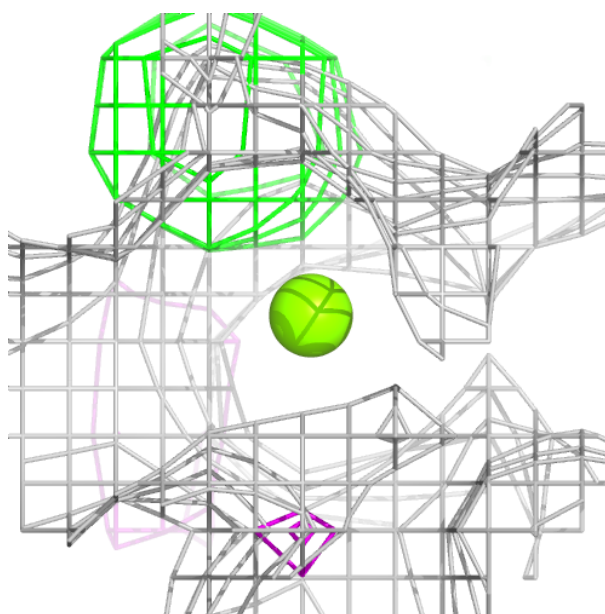
Electron density around ADP A 501:

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



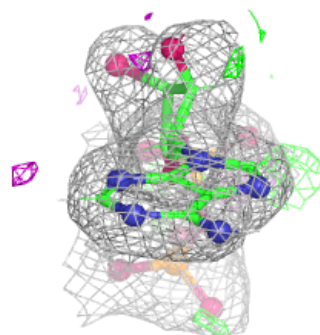
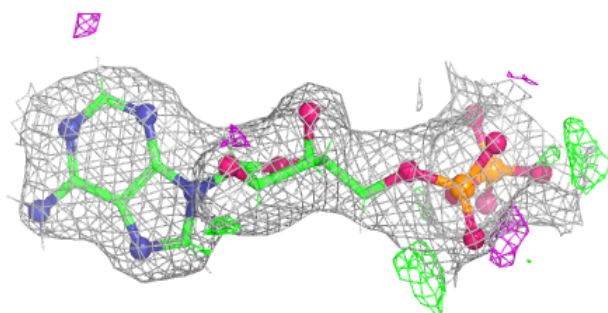
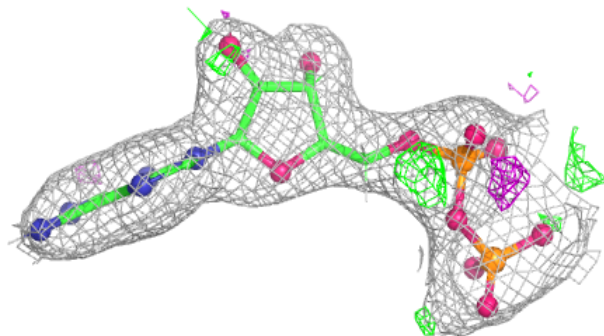
Electron density around MG E 505:

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and green (positive)

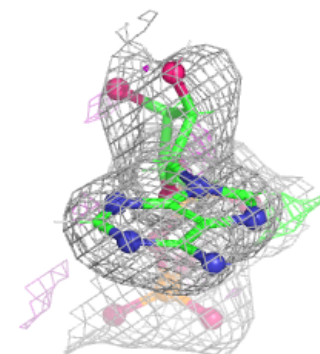
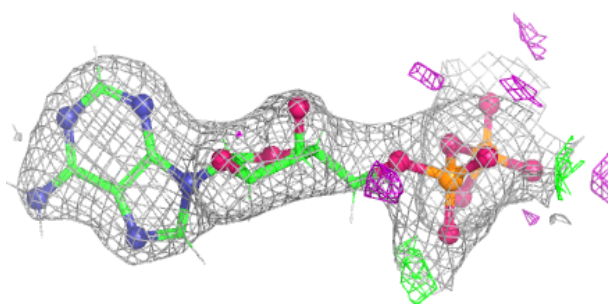
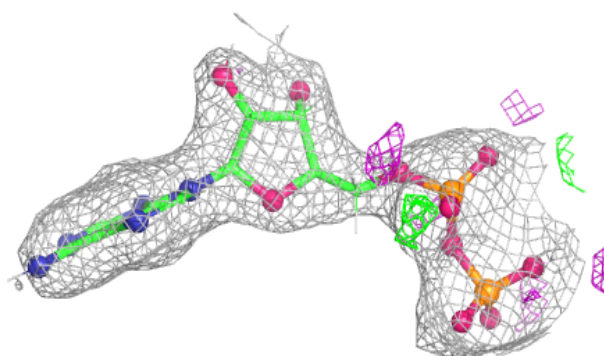


Electron density around ADP B 501:

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

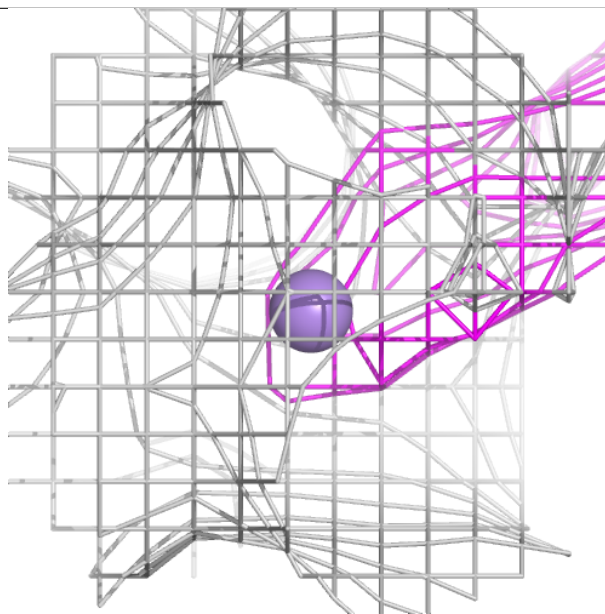
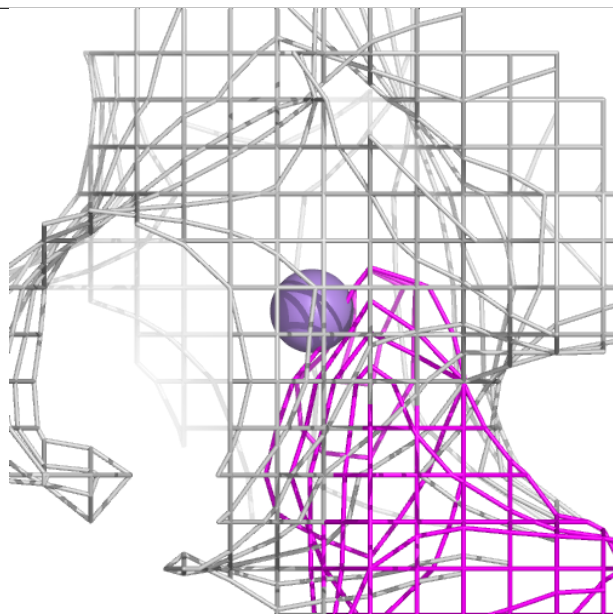
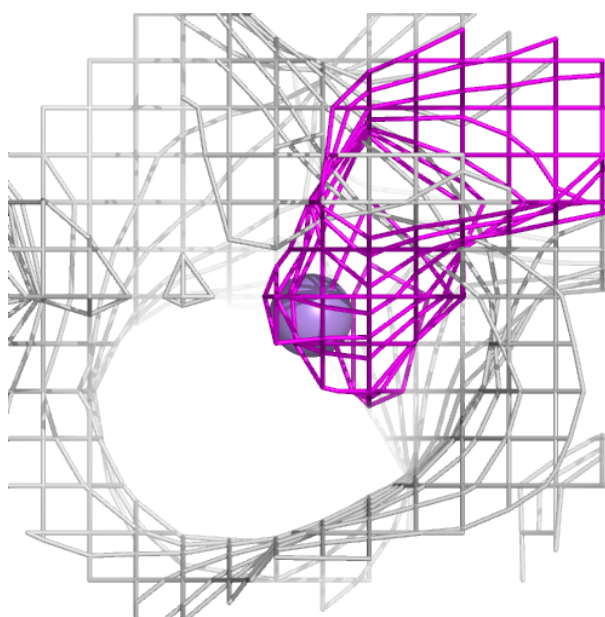
**Electron density around ADP C 501:**

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and green (positive)



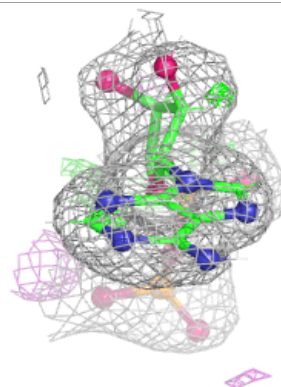
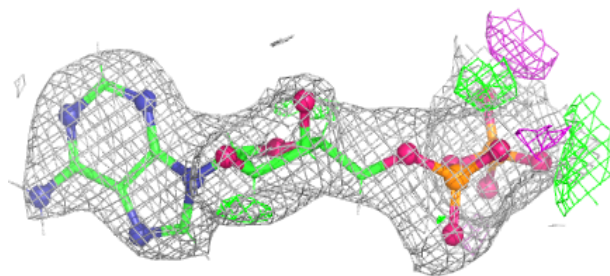
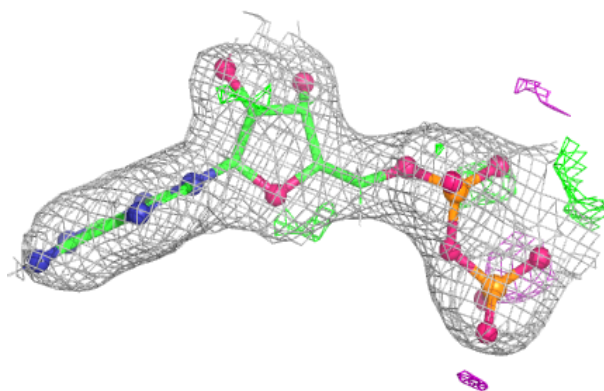
Electron density around MN C 503:

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



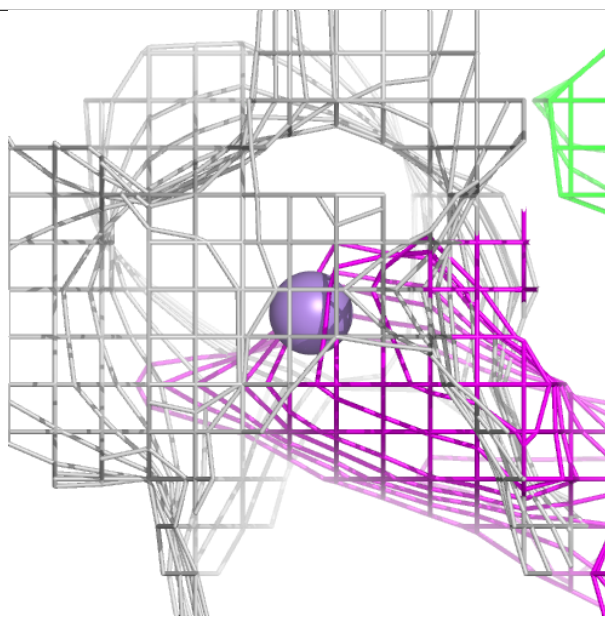
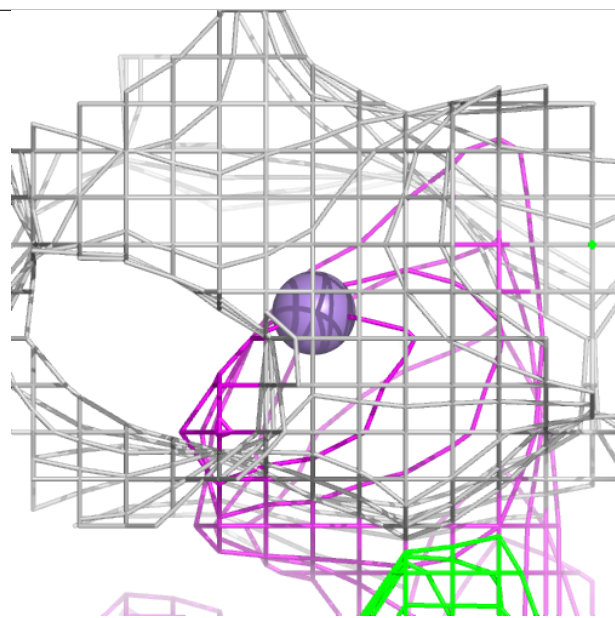
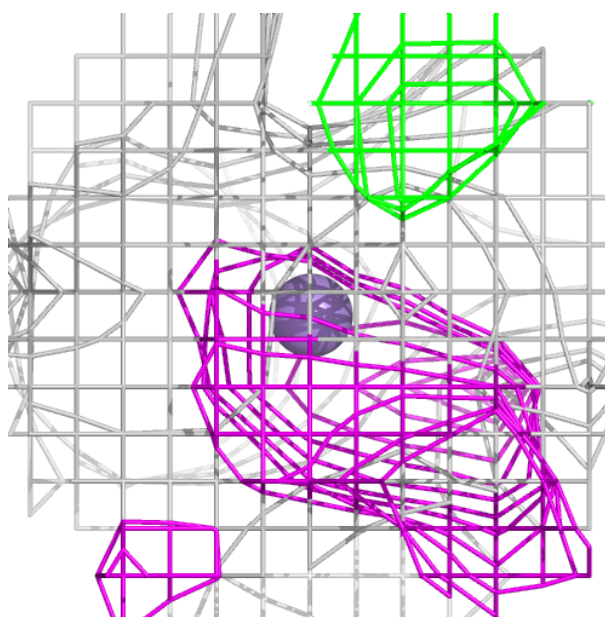
Electron density around ADP F 501:

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



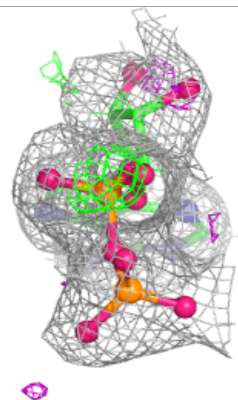
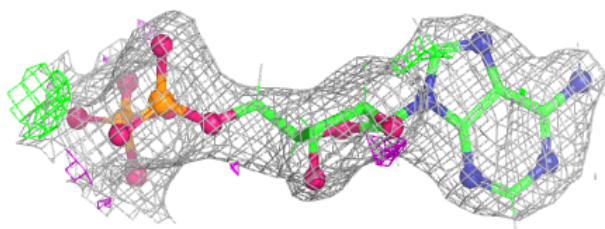
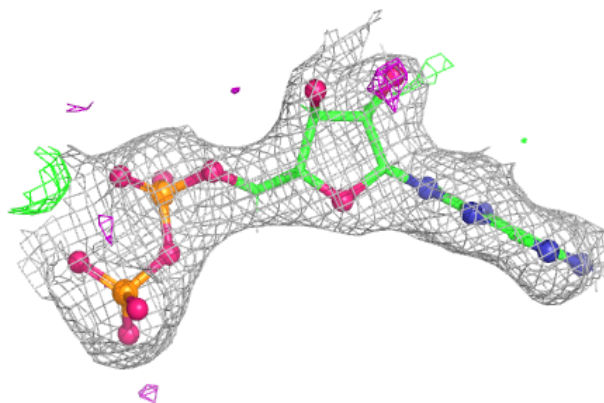
Electron density around MN C 504:

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



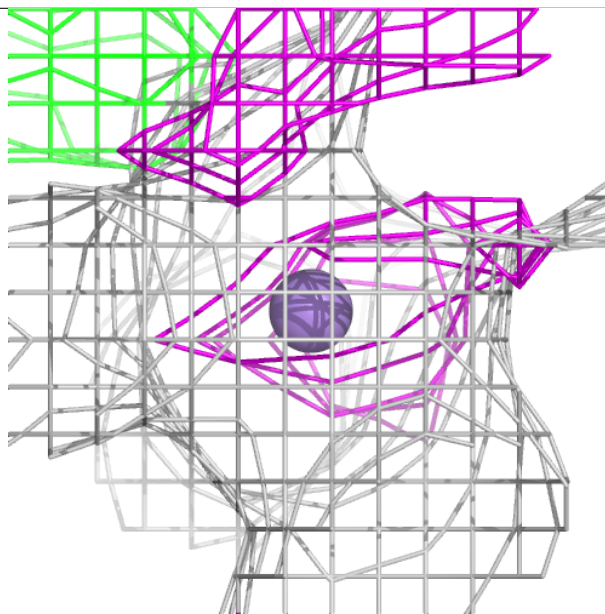
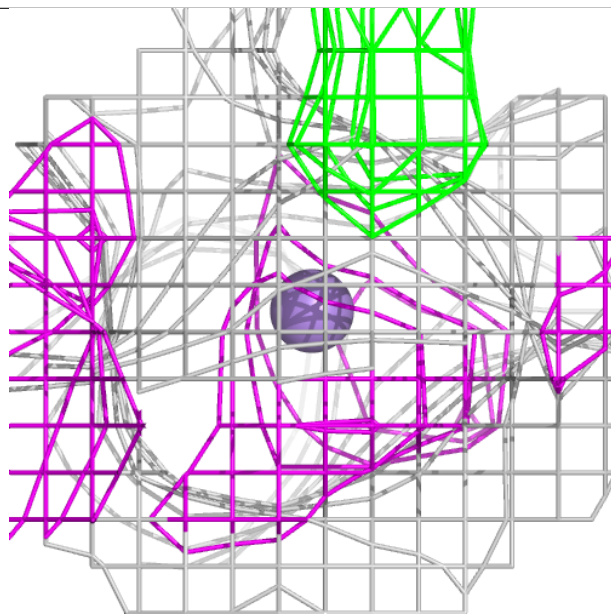
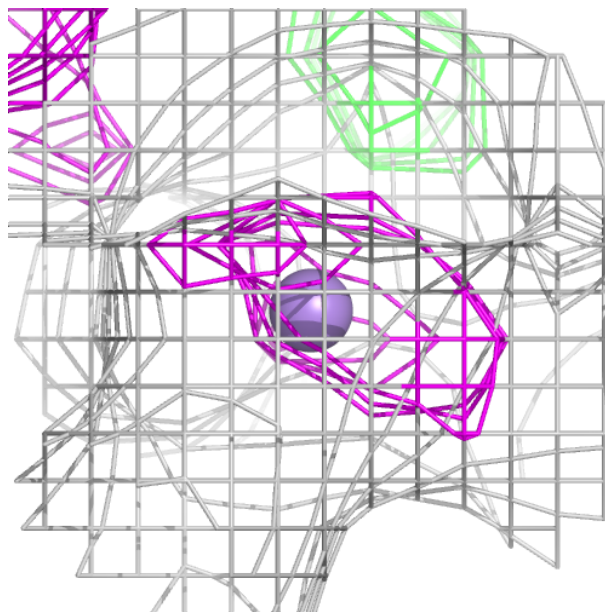
Electron density around ADP E 501:

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



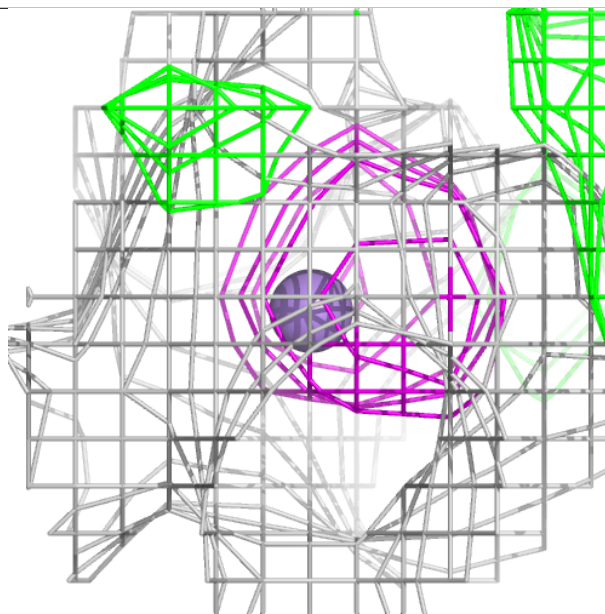
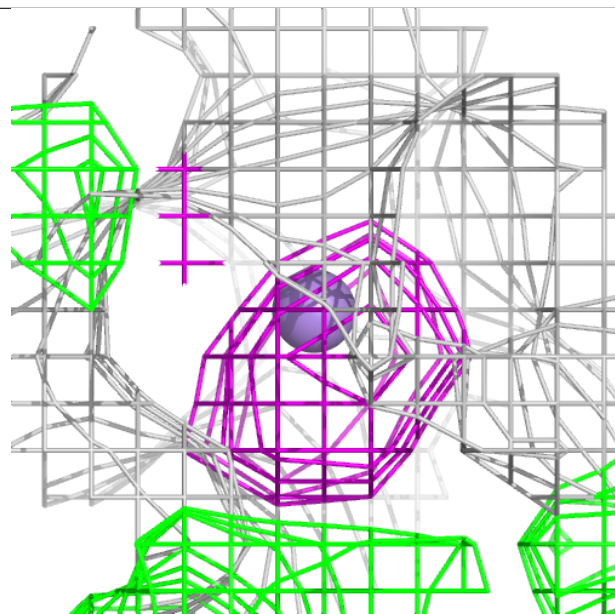
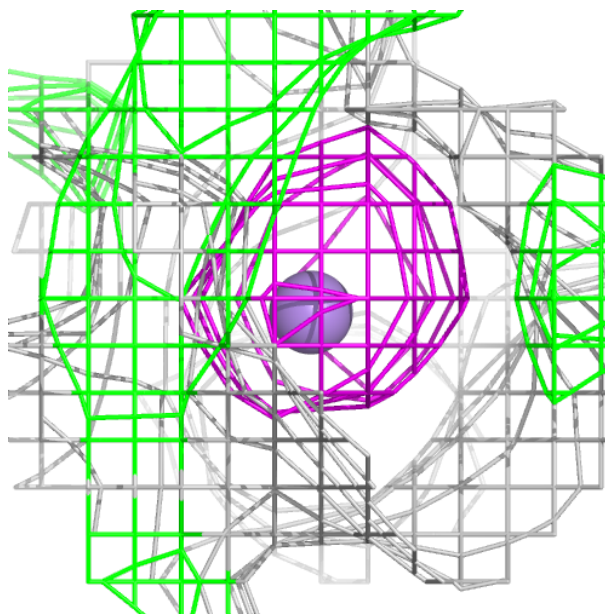
Electron density around MN A 504:

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



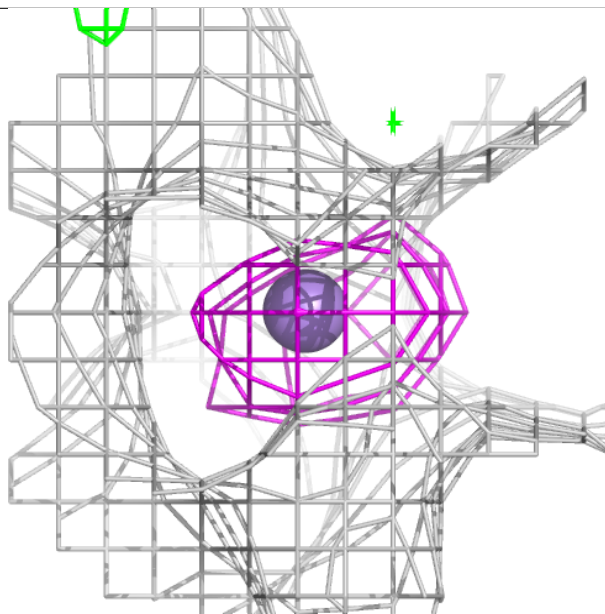
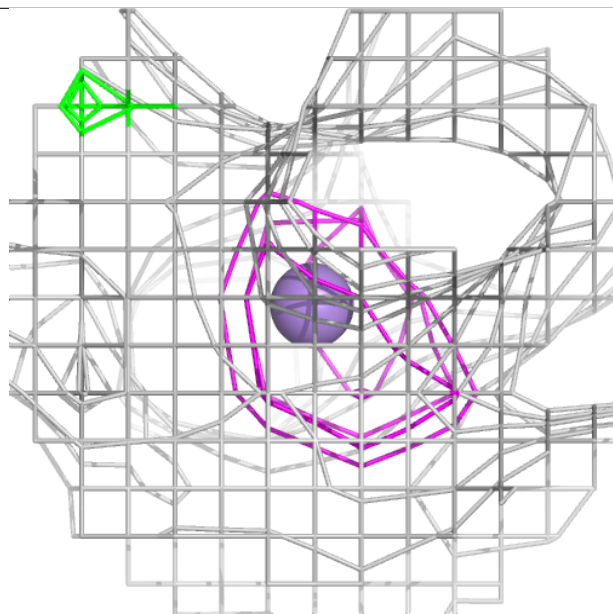
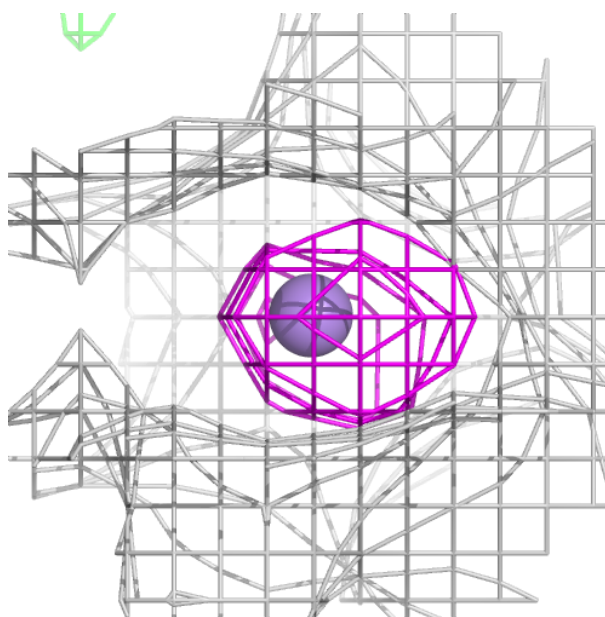
Electron density around MN F 503:

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



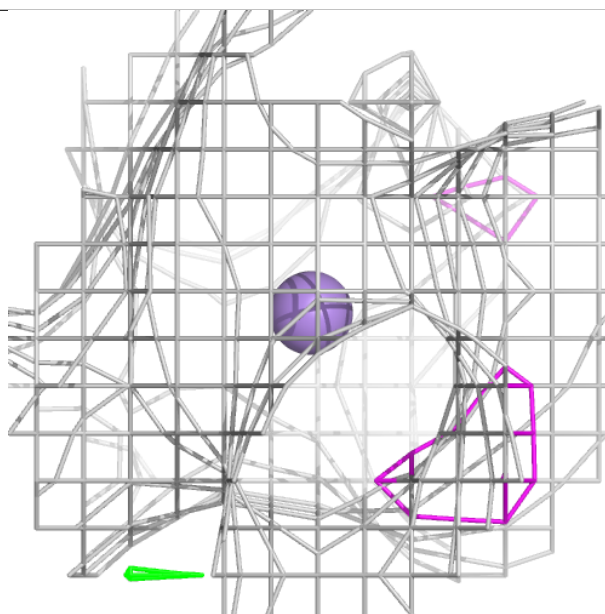
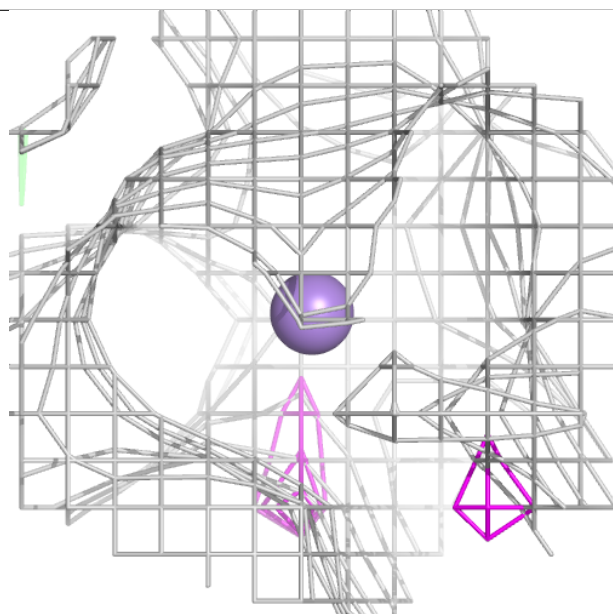
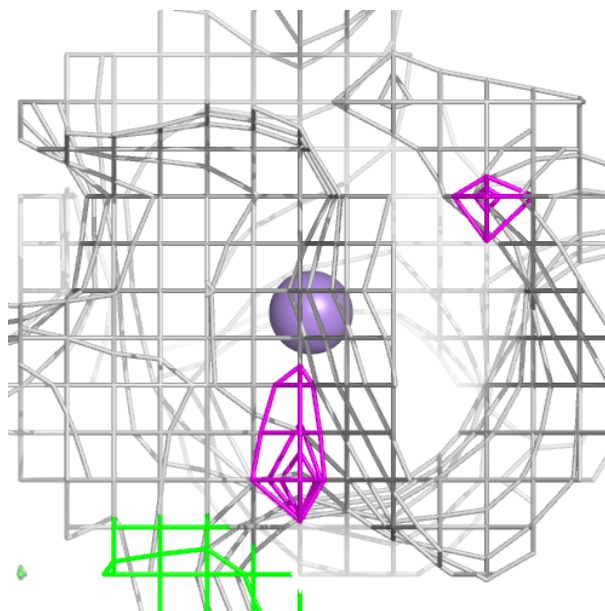
Electron density around MN E 503:

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



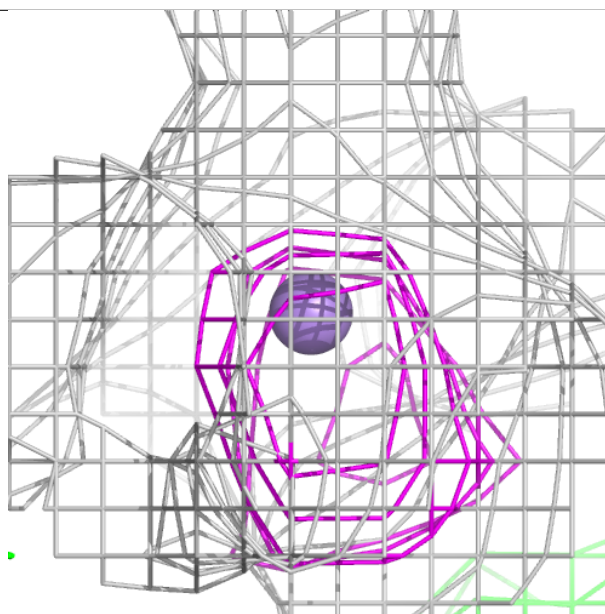
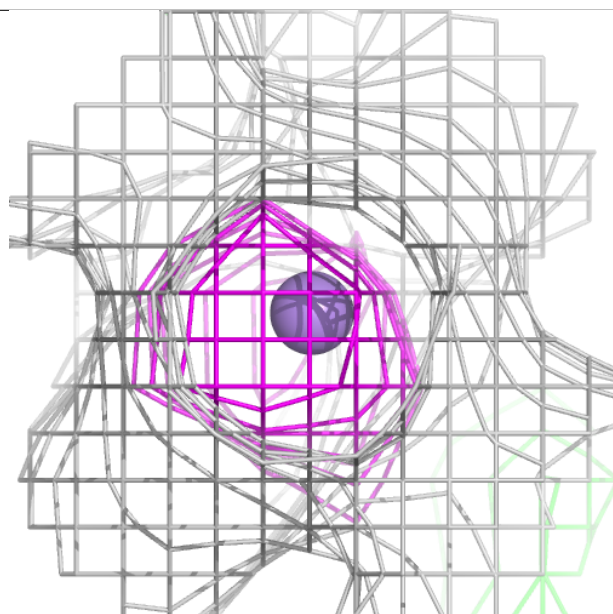
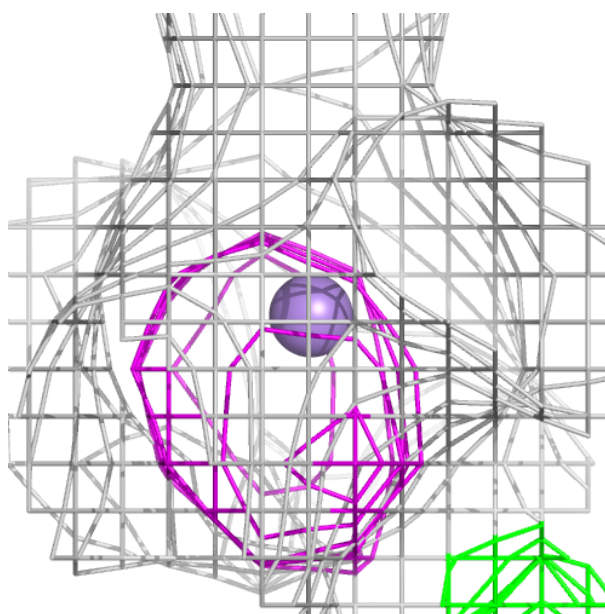
Electron density around MN E 504:

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



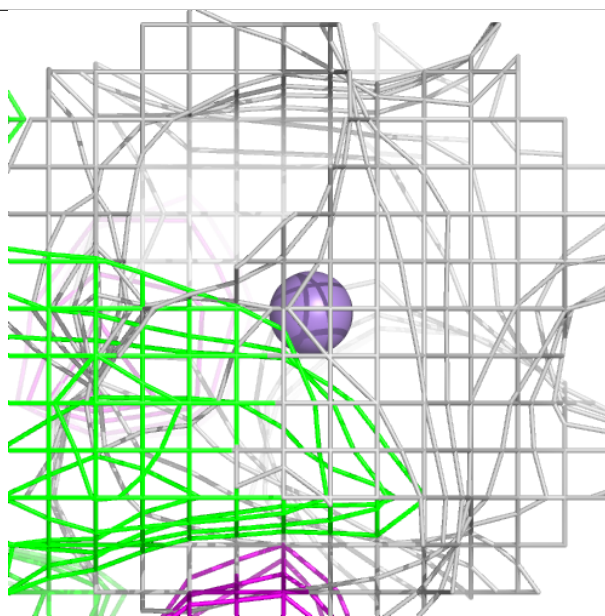
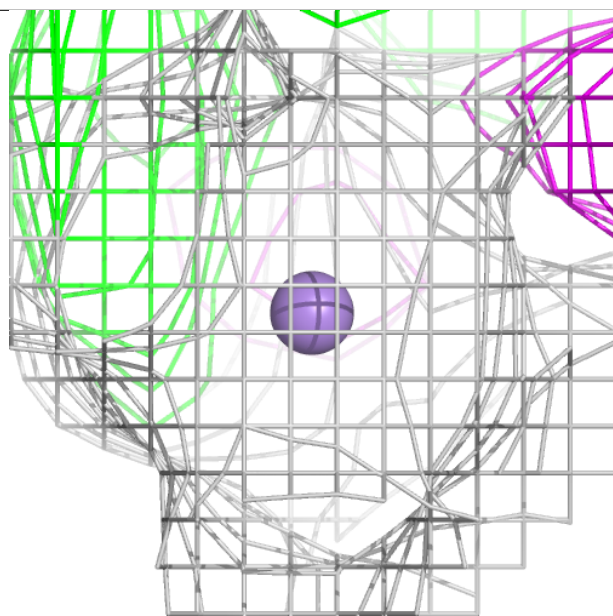
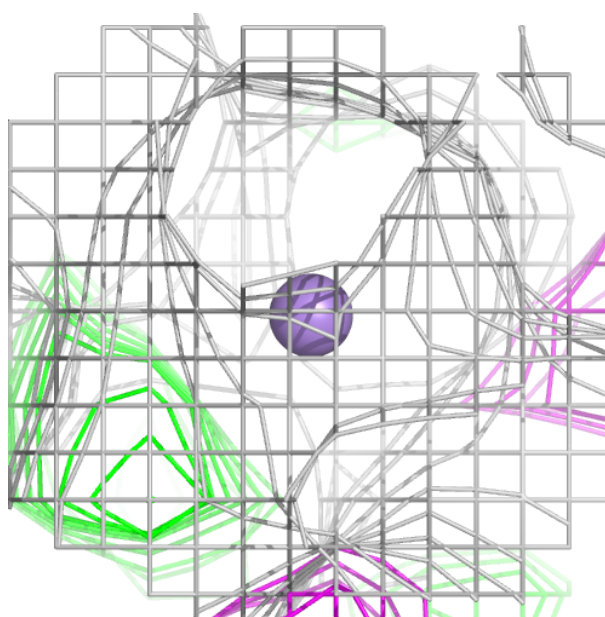
Electron density around MN A 503:

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



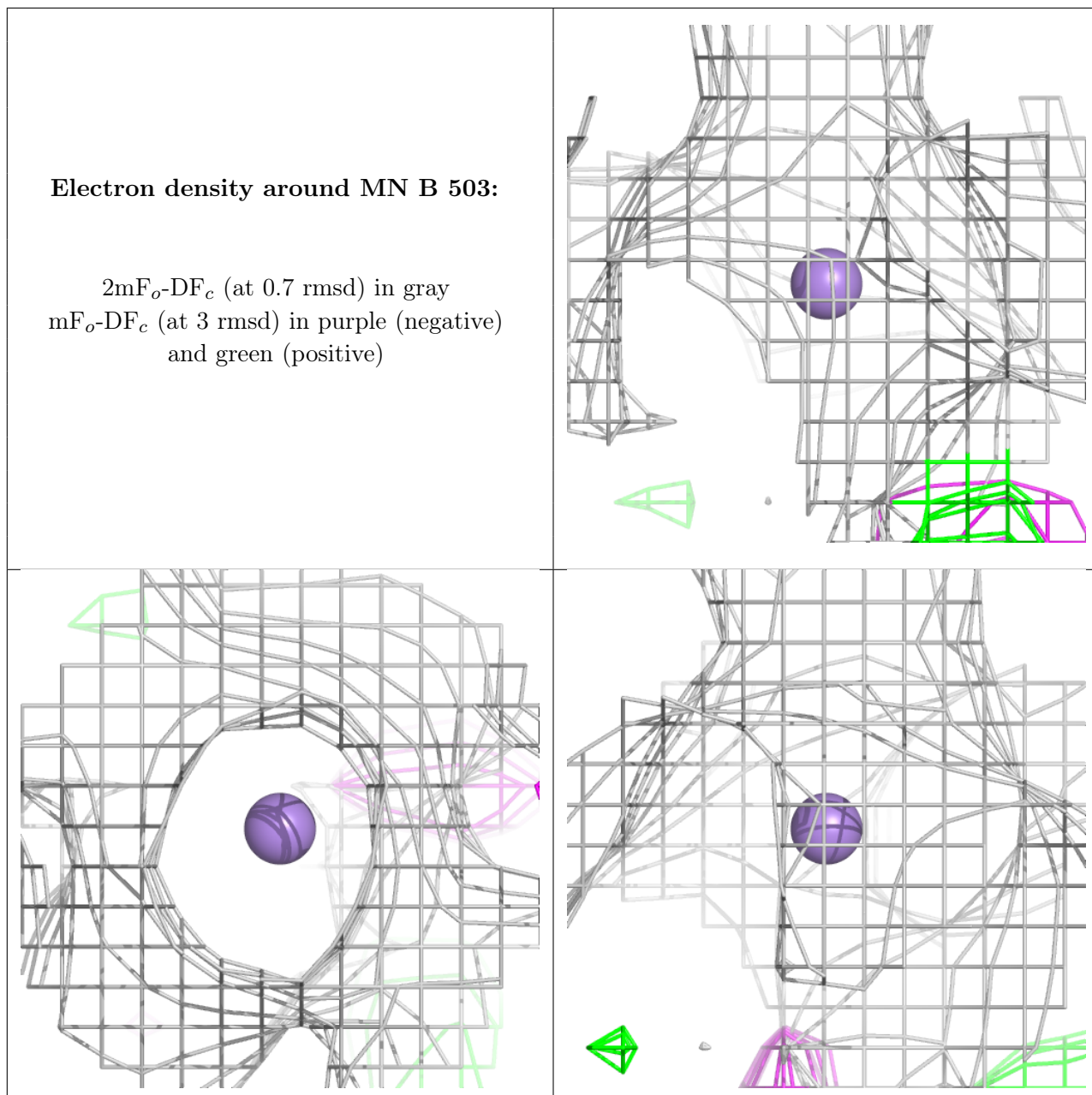
Electron density around MN D 503:

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



Electron density around MN B 503:

$2mF_o-DF_c$ (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.