



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2023 – 01:08 pm GMT

PDB ID : 8CDB
Title : Proulilysin E229A structure
Authors : Rodriguez-Banqueri, A.; Eckhard, U.; Gomis-Ruth, F.X.
Deposited on : 2023-01-30
Resolution : 4.50 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.32.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.32.1

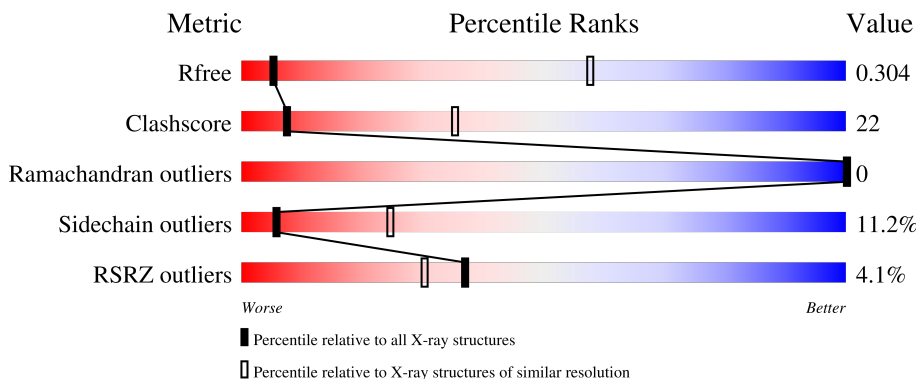
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1055 (5.20-3.80)
Clashscore	141614	1123 (5.20-3.80)
Ramachandran outliers	138981	1069 (5.20-3.80)
Sidechain outliers	138945	1050 (5.20-3.80)
RSRZ outliers	127900	1101 (5.30-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	360	
1	B	360	
1	C	360	
1	D	360	
1	E	360	

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Mol	Chain	Length	Quality of chain
1	F	360	<p>4% 43% 38% 15%</p>
1	G	360	<p>4% 44% 37% 15%</p>
1	H	360	<p>3% 44% 38% 15%</p>
1	I	360	<p>5% 45% 36% 15%</p>
1	J	360	<p>2% 45% 36% 15%</p>
1	K	360	<p>6% 45% 37% 15%</p>
1	L	360	<p>2% 46% 36% 15%</p>
1	M	360	<p>6% 45% 36% 15%</p>
1	N	360	<p>2% 44% 37% 15%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 33614 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Ulilysin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	306	2399	1502	419	464	14	0	0	0
1	B	306	2399	1502	419	464	14	0	0	0
1	C	306	2399	1502	419	464	14	0	0	0
1	D	306	2399	1502	419	464	14	0	0	0
1	E	306	2399	1502	419	464	14	0	0	0
1	F	306	2399	1502	419	464	14	0	0	0
1	G	306	2399	1502	419	464	14	0	0	0
1	H	306	2399	1502	419	464	14	0	0	0
1	I	306	2399	1502	419	464	14	0	0	0
1	J	306	2399	1502	419	464	14	0	0	0
1	K	306	2399	1502	419	464	14	0	0	0
1	L	306	2399	1502	419	464	14	0	0	0
1	M	306	2399	1502	419	464	14	0	0	0
1	N	306	2399	1502	419	464	14	0	0	0

There are 266 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	GLY	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	SER	-	expression tag	UNP Q8TL28
A	-15	SER	-	expression tag	UNP Q8TL28
A	-14	HIS	-	expression tag	UNP Q8TL28
A	-13	HIS	-	expression tag	UNP Q8TL28
A	-12	HIS	-	expression tag	UNP Q8TL28
A	-11	HIS	-	expression tag	UNP Q8TL28
A	-10	HIS	-	expression tag	UNP Q8TL28
A	-9	HIS	-	expression tag	UNP Q8TL28
A	-8	SER	-	expression tag	UNP Q8TL28
A	-7	SER	-	expression tag	UNP Q8TL28
A	-6	GLY	-	expression tag	UNP Q8TL28
A	-5	LEU	-	expression tag	UNP Q8TL28
A	-4	VAL	-	expression tag	UNP Q8TL28
A	-3	PRO	-	expression tag	UNP Q8TL28
A	-2	ARG	-	expression tag	UNP Q8TL28
A	-1	GLY	-	expression tag	UNP Q8TL28
A	0	SER	-	expression tag	UNP Q8TL28
A	229	ALA	GLU	variant	UNP Q8TL28
B	-17	GLY	-	expression tag	UNP Q8TL28
B	-16	SER	-	expression tag	UNP Q8TL28
B	-15	SER	-	expression tag	UNP Q8TL28
B	-14	HIS	-	expression tag	UNP Q8TL28
B	-13	HIS	-	expression tag	UNP Q8TL28
B	-12	HIS	-	expression tag	UNP Q8TL28
B	-11	HIS	-	expression tag	UNP Q8TL28
B	-10	HIS	-	expression tag	UNP Q8TL28
B	-9	HIS	-	expression tag	UNP Q8TL28
B	-8	SER	-	expression tag	UNP Q8TL28
B	-7	SER	-	expression tag	UNP Q8TL28
B	-6	GLY	-	expression tag	UNP Q8TL28
B	-5	LEU	-	expression tag	UNP Q8TL28
B	-4	VAL	-	expression tag	UNP Q8TL28
B	-3	PRO	-	expression tag	UNP Q8TL28
B	-2	ARG	-	expression tag	UNP Q8TL28
B	-1	GLY	-	expression tag	UNP Q8TL28
B	0	SER	-	expression tag	UNP Q8TL28
B	229	ALA	GLU	variant	UNP Q8TL28
C	-17	GLY	-	expression tag	UNP Q8TL28
C	-16	SER	-	expression tag	UNP Q8TL28
C	-15	SER	-	expression tag	UNP Q8TL28
C	-14	HIS	-	expression tag	UNP Q8TL28
C	-13	HIS	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-12	HIS	-	expression tag	UNP Q8TL28
C	-11	HIS	-	expression tag	UNP Q8TL28
C	-10	HIS	-	expression tag	UNP Q8TL28
C	-9	HIS	-	expression tag	UNP Q8TL28
C	-8	SER	-	expression tag	UNP Q8TL28
C	-7	SER	-	expression tag	UNP Q8TL28
C	-6	GLY	-	expression tag	UNP Q8TL28
C	-5	LEU	-	expression tag	UNP Q8TL28
C	-4	VAL	-	expression tag	UNP Q8TL28
C	-3	PRO	-	expression tag	UNP Q8TL28
C	-2	ARG	-	expression tag	UNP Q8TL28
C	-1	GLY	-	expression tag	UNP Q8TL28
C	0	SER	-	expression tag	UNP Q8TL28
C	229	ALA	GLU	variant	UNP Q8TL28
D	-17	GLY	-	expression tag	UNP Q8TL28
D	-16	SER	-	expression tag	UNP Q8TL28
D	-15	SER	-	expression tag	UNP Q8TL28
D	-14	HIS	-	expression tag	UNP Q8TL28
D	-13	HIS	-	expression tag	UNP Q8TL28
D	-12	HIS	-	expression tag	UNP Q8TL28
D	-11	HIS	-	expression tag	UNP Q8TL28
D	-10	HIS	-	expression tag	UNP Q8TL28
D	-9	HIS	-	expression tag	UNP Q8TL28
D	-8	SER	-	expression tag	UNP Q8TL28
D	-7	SER	-	expression tag	UNP Q8TL28
D	-6	GLY	-	expression tag	UNP Q8TL28
D	-5	LEU	-	expression tag	UNP Q8TL28
D	-4	VAL	-	expression tag	UNP Q8TL28
D	-3	PRO	-	expression tag	UNP Q8TL28
D	-2	ARG	-	expression tag	UNP Q8TL28
D	-1	GLY	-	expression tag	UNP Q8TL28
D	0	SER	-	expression tag	UNP Q8TL28
D	229	ALA	GLU	variant	UNP Q8TL28
E	-17	GLY	-	expression tag	UNP Q8TL28
E	-16	SER	-	expression tag	UNP Q8TL28
E	-15	SER	-	expression tag	UNP Q8TL28
E	-14	HIS	-	expression tag	UNP Q8TL28
E	-13	HIS	-	expression tag	UNP Q8TL28
E	-12	HIS	-	expression tag	UNP Q8TL28
E	-11	HIS	-	expression tag	UNP Q8TL28
E	-10	HIS	-	expression tag	UNP Q8TL28
E	-9	HIS	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-8	SER	-	expression tag	UNP Q8TL28
E	-7	SER	-	expression tag	UNP Q8TL28
E	-6	GLY	-	expression tag	UNP Q8TL28
E	-5	LEU	-	expression tag	UNP Q8TL28
E	-4	VAL	-	expression tag	UNP Q8TL28
E	-3	PRO	-	expression tag	UNP Q8TL28
E	-2	ARG	-	expression tag	UNP Q8TL28
E	-1	GLY	-	expression tag	UNP Q8TL28
E	0	SER	-	expression tag	UNP Q8TL28
E	229	ALA	GLU	variant	UNP Q8TL28
F	-17	GLY	-	expression tag	UNP Q8TL28
F	-16	SER	-	expression tag	UNP Q8TL28
F	-15	SER	-	expression tag	UNP Q8TL28
F	-14	HIS	-	expression tag	UNP Q8TL28
F	-13	HIS	-	expression tag	UNP Q8TL28
F	-12	HIS	-	expression tag	UNP Q8TL28
F	-11	HIS	-	expression tag	UNP Q8TL28
F	-10	HIS	-	expression tag	UNP Q8TL28
F	-9	HIS	-	expression tag	UNP Q8TL28
F	-8	SER	-	expression tag	UNP Q8TL28
F	-7	SER	-	expression tag	UNP Q8TL28
F	-6	GLY	-	expression tag	UNP Q8TL28
F	-5	LEU	-	expression tag	UNP Q8TL28
F	-4	VAL	-	expression tag	UNP Q8TL28
F	-3	PRO	-	expression tag	UNP Q8TL28
F	-2	ARG	-	expression tag	UNP Q8TL28
F	-1	GLY	-	expression tag	UNP Q8TL28
F	0	SER	-	expression tag	UNP Q8TL28
F	229	ALA	GLU	variant	UNP Q8TL28
G	-17	GLY	-	expression tag	UNP Q8TL28
G	-16	SER	-	expression tag	UNP Q8TL28
G	-15	SER	-	expression tag	UNP Q8TL28
G	-14	HIS	-	expression tag	UNP Q8TL28
G	-13	HIS	-	expression tag	UNP Q8TL28
G	-12	HIS	-	expression tag	UNP Q8TL28
G	-11	HIS	-	expression tag	UNP Q8TL28
G	-10	HIS	-	expression tag	UNP Q8TL28
G	-9	HIS	-	expression tag	UNP Q8TL28
G	-8	SER	-	expression tag	UNP Q8TL28
G	-7	SER	-	expression tag	UNP Q8TL28
G	-6	GLY	-	expression tag	UNP Q8TL28
G	-5	LEU	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	VAL	-	expression tag	UNP Q8TL28
G	-3	PRO	-	expression tag	UNP Q8TL28
G	-2	ARG	-	expression tag	UNP Q8TL28
G	-1	GLY	-	expression tag	UNP Q8TL28
G	0	SER	-	expression tag	UNP Q8TL28
G	229	ALA	GLU	variant	UNP Q8TL28
H	-17	GLY	-	expression tag	UNP Q8TL28
H	-16	SER	-	expression tag	UNP Q8TL28
H	-15	SER	-	expression tag	UNP Q8TL28
H	-14	HIS	-	expression tag	UNP Q8TL28
H	-13	HIS	-	expression tag	UNP Q8TL28
H	-12	HIS	-	expression tag	UNP Q8TL28
H	-11	HIS	-	expression tag	UNP Q8TL28
H	-10	HIS	-	expression tag	UNP Q8TL28
H	-9	HIS	-	expression tag	UNP Q8TL28
H	-8	SER	-	expression tag	UNP Q8TL28
H	-7	SER	-	expression tag	UNP Q8TL28
H	-6	GLY	-	expression tag	UNP Q8TL28
H	-5	LEU	-	expression tag	UNP Q8TL28
H	-4	VAL	-	expression tag	UNP Q8TL28
H	-3	PRO	-	expression tag	UNP Q8TL28
H	-2	ARG	-	expression tag	UNP Q8TL28
H	-1	GLY	-	expression tag	UNP Q8TL28
H	0	SER	-	expression tag	UNP Q8TL28
H	229	ALA	GLU	variant	UNP Q8TL28
I	-17	GLY	-	expression tag	UNP Q8TL28
I	-16	SER	-	expression tag	UNP Q8TL28
I	-15	SER	-	expression tag	UNP Q8TL28
I	-14	HIS	-	expression tag	UNP Q8TL28
I	-13	HIS	-	expression tag	UNP Q8TL28
I	-12	HIS	-	expression tag	UNP Q8TL28
I	-11	HIS	-	expression tag	UNP Q8TL28
I	-10	HIS	-	expression tag	UNP Q8TL28
I	-9	HIS	-	expression tag	UNP Q8TL28
I	-8	SER	-	expression tag	UNP Q8TL28
I	-7	SER	-	expression tag	UNP Q8TL28
I	-6	GLY	-	expression tag	UNP Q8TL28
I	-5	LEU	-	expression tag	UNP Q8TL28
I	-4	VAL	-	expression tag	UNP Q8TL28
I	-3	PRO	-	expression tag	UNP Q8TL28
I	-2	ARG	-	expression tag	UNP Q8TL28
I	-1	GLY	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
I	0	SER	-	expression tag	UNP Q8TL28
I	229	ALA	GLU	variant	UNP Q8TL28
J	-17	GLY	-	expression tag	UNP Q8TL28
J	-16	SER	-	expression tag	UNP Q8TL28
J	-15	SER	-	expression tag	UNP Q8TL28
J	-14	HIS	-	expression tag	UNP Q8TL28
J	-13	HIS	-	expression tag	UNP Q8TL28
J	-12	HIS	-	expression tag	UNP Q8TL28
J	-11	HIS	-	expression tag	UNP Q8TL28
J	-10	HIS	-	expression tag	UNP Q8TL28
J	-9	HIS	-	expression tag	UNP Q8TL28
J	-8	SER	-	expression tag	UNP Q8TL28
J	-7	SER	-	expression tag	UNP Q8TL28
J	-6	GLY	-	expression tag	UNP Q8TL28
J	-5	LEU	-	expression tag	UNP Q8TL28
J	-4	VAL	-	expression tag	UNP Q8TL28
J	-3	PRO	-	expression tag	UNP Q8TL28
J	-2	ARG	-	expression tag	UNP Q8TL28
J	-1	GLY	-	expression tag	UNP Q8TL28
J	0	SER	-	expression tag	UNP Q8TL28
J	229	ALA	GLU	variant	UNP Q8TL28
K	-17	GLY	-	expression tag	UNP Q8TL28
K	-16	SER	-	expression tag	UNP Q8TL28
K	-15	SER	-	expression tag	UNP Q8TL28
K	-14	HIS	-	expression tag	UNP Q8TL28
K	-13	HIS	-	expression tag	UNP Q8TL28
K	-12	HIS	-	expression tag	UNP Q8TL28
K	-11	HIS	-	expression tag	UNP Q8TL28
K	-10	HIS	-	expression tag	UNP Q8TL28
K	-9	HIS	-	expression tag	UNP Q8TL28
K	-8	SER	-	expression tag	UNP Q8TL28
K	-7	SER	-	expression tag	UNP Q8TL28
K	-6	GLY	-	expression tag	UNP Q8TL28
K	-5	LEU	-	expression tag	UNP Q8TL28
K	-4	VAL	-	expression tag	UNP Q8TL28
K	-3	PRO	-	expression tag	UNP Q8TL28
K	-2	ARG	-	expression tag	UNP Q8TL28
K	-1	GLY	-	expression tag	UNP Q8TL28
K	0	SER	-	expression tag	UNP Q8TL28
K	229	ALA	GLU	variant	UNP Q8TL28
L	-17	GLY	-	expression tag	UNP Q8TL28
L	-16	SER	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
L	-15	SER	-	expression tag	UNP Q8TL28
L	-14	HIS	-	expression tag	UNP Q8TL28
L	-13	HIS	-	expression tag	UNP Q8TL28
L	-12	HIS	-	expression tag	UNP Q8TL28
L	-11	HIS	-	expression tag	UNP Q8TL28
L	-10	HIS	-	expression tag	UNP Q8TL28
L	-9	HIS	-	expression tag	UNP Q8TL28
L	-8	SER	-	expression tag	UNP Q8TL28
L	-7	SER	-	expression tag	UNP Q8TL28
L	-6	GLY	-	expression tag	UNP Q8TL28
L	-5	LEU	-	expression tag	UNP Q8TL28
L	-4	VAL	-	expression tag	UNP Q8TL28
L	-3	PRO	-	expression tag	UNP Q8TL28
L	-2	ARG	-	expression tag	UNP Q8TL28
L	-1	GLY	-	expression tag	UNP Q8TL28
L	0	SER	-	expression tag	UNP Q8TL28
L	229	ALA	GLU	variant	UNP Q8TL28
M	-17	GLY	-	expression tag	UNP Q8TL28
M	-16	SER	-	expression tag	UNP Q8TL28
M	-15	SER	-	expression tag	UNP Q8TL28
M	-14	HIS	-	expression tag	UNP Q8TL28
M	-13	HIS	-	expression tag	UNP Q8TL28
M	-12	HIS	-	expression tag	UNP Q8TL28
M	-11	HIS	-	expression tag	UNP Q8TL28
M	-10	HIS	-	expression tag	UNP Q8TL28
M	-9	HIS	-	expression tag	UNP Q8TL28
M	-8	SER	-	expression tag	UNP Q8TL28
M	-7	SER	-	expression tag	UNP Q8TL28
M	-6	GLY	-	expression tag	UNP Q8TL28
M	-5	LEU	-	expression tag	UNP Q8TL28
M	-4	VAL	-	expression tag	UNP Q8TL28
M	-3	PRO	-	expression tag	UNP Q8TL28
M	-2	ARG	-	expression tag	UNP Q8TL28
M	-1	GLY	-	expression tag	UNP Q8TL28
M	0	SER	-	expression tag	UNP Q8TL28
M	229	ALA	GLU	variant	UNP Q8TL28
N	-17	GLY	-	expression tag	UNP Q8TL28
N	-16	SER	-	expression tag	UNP Q8TL28
N	-15	SER	-	expression tag	UNP Q8TL28
N	-14	HIS	-	expression tag	UNP Q8TL28
N	-13	HIS	-	expression tag	UNP Q8TL28
N	-12	HIS	-	expression tag	UNP Q8TL28

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Chain	Residue	Modelled	Actual	Comment	Reference
N	-11	HIS	-	expression tag	UNP Q8TL28
N	-10	HIS	-	expression tag	UNP Q8TL28
N	-9	HIS	-	expression tag	UNP Q8TL28
N	-8	SER	-	expression tag	UNP Q8TL28
N	-7	SER	-	expression tag	UNP Q8TL28
N	-6	GLY	-	expression tag	UNP Q8TL28
N	-5	LEU	-	expression tag	UNP Q8TL28
N	-4	VAL	-	expression tag	UNP Q8TL28
N	-3	PRO	-	expression tag	UNP Q8TL28
N	-2	ARG	-	expression tag	UNP Q8TL28
N	-1	GLY	-	expression tag	UNP Q8TL28
N	0	SER	-	expression tag	UNP Q8TL28
N	229	ALA	GLU	variant	UNP Q8TL28

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0
2	I	1	Total Zn 1 1	0	0
2	J	1	Total Zn 1 1	0	0
2	K	1	Total Zn 1 1	0	0
2	L	1	Total Zn 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	M	1	Total 1	Zn 1	0	0
2	N	1	Total 1	Zn 1	0	0

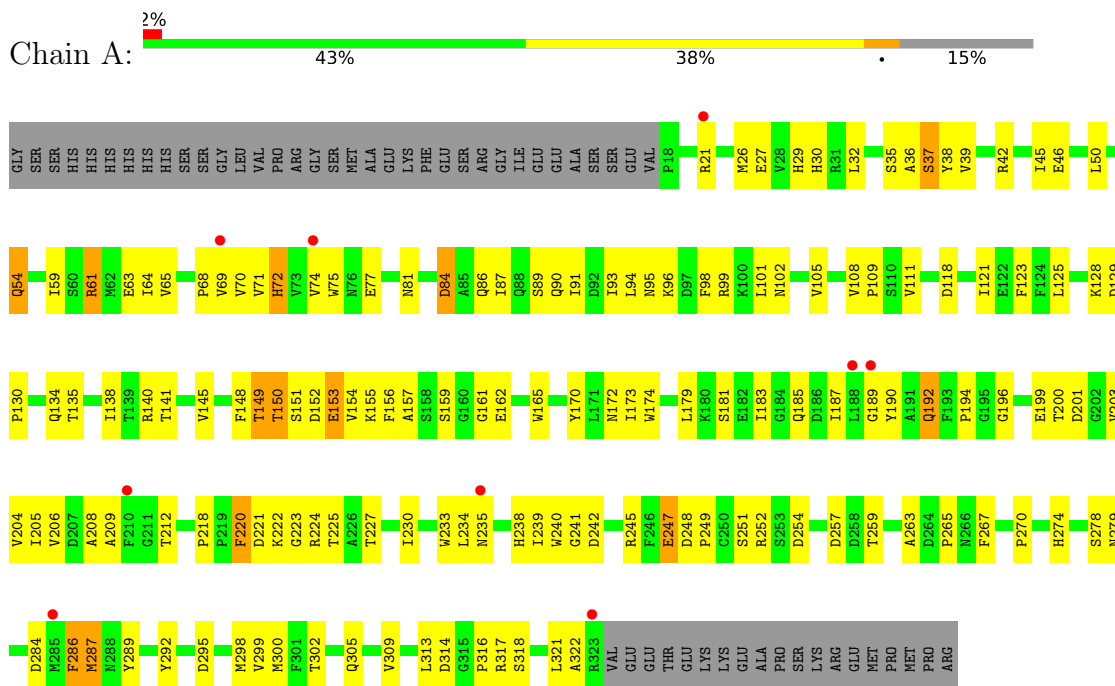
- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Ca 1	0	0
3	B	1	Total 1	Ca 1	0	0
3	C	1	Total 1	Ca 1	0	0
3	D	1	Total 1	Ca 1	0	0
3	E	1	Total 1	Ca 1	0	0
3	F	1	Total 1	Ca 1	0	0
3	G	1	Total 1	Ca 1	0	0
3	H	1	Total 1	Ca 1	0	0
3	I	1	Total 1	Ca 1	0	0
3	J	1	Total 1	Ca 1	0	0
3	K	1	Total 1	Ca 1	0	0
3	L	1	Total 1	Ca 1	0	0
3	M	1	Total 1	Ca 1	0	0
3	N	1	Total 1	Ca 1	0	0

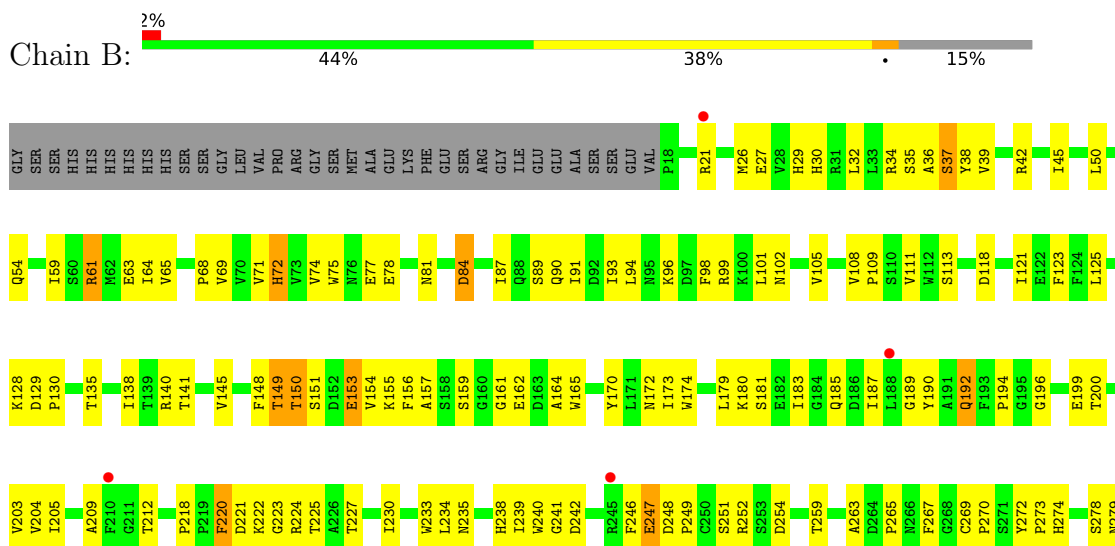
3 Residue-property plots [i](#)

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

- Molecule 1: Ulilysin

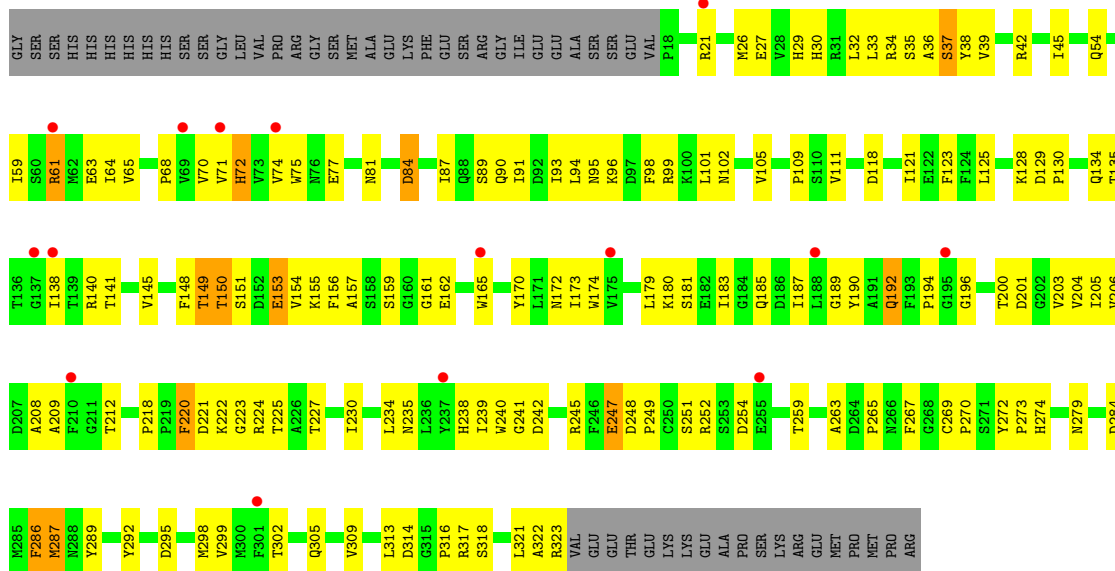


- Molecule 1: Ulilysin

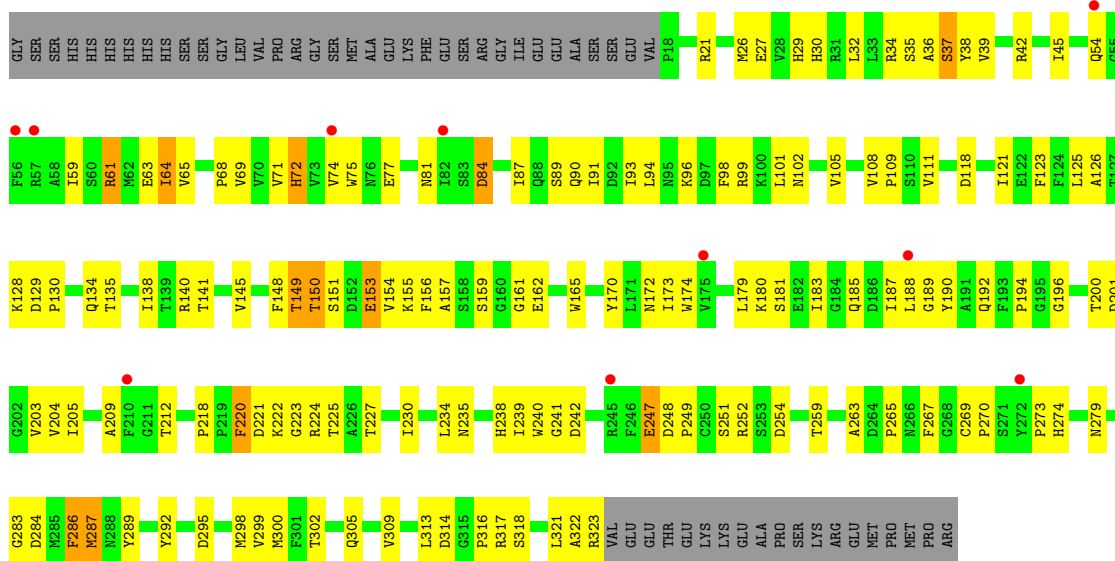




- Molecule 1: Ulilysin

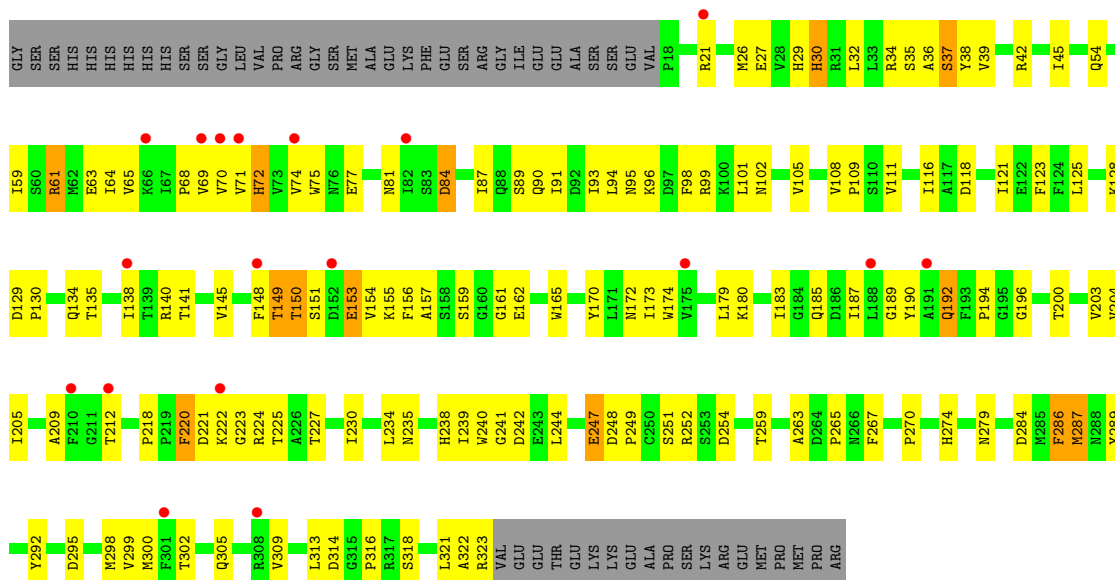


- Molecule 1: Ulilysin

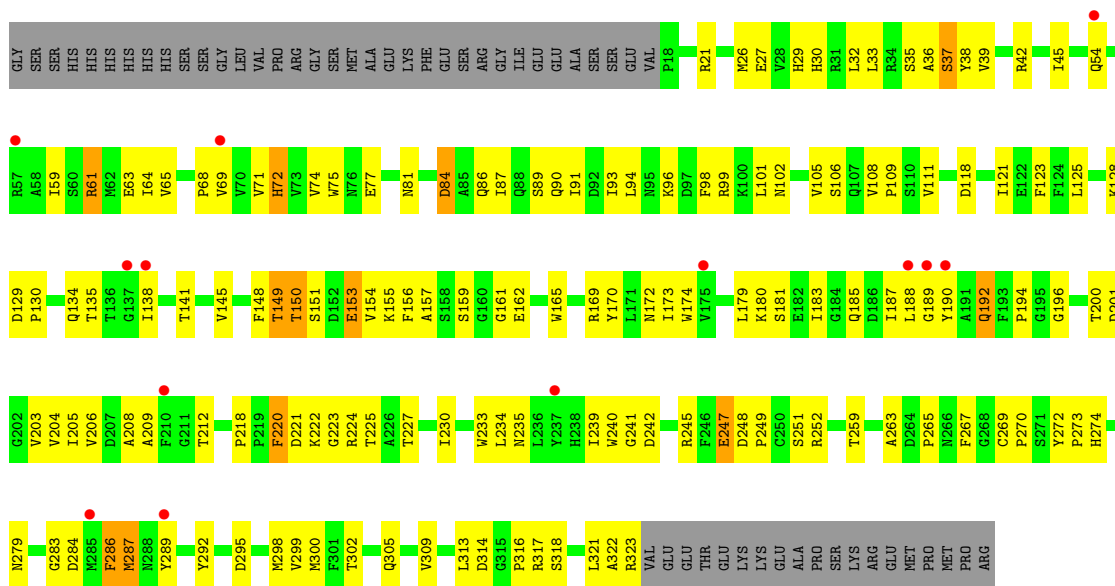


- Molecule 1: Ulilysin

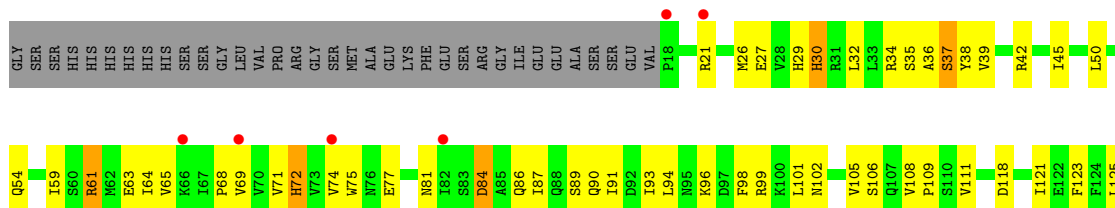


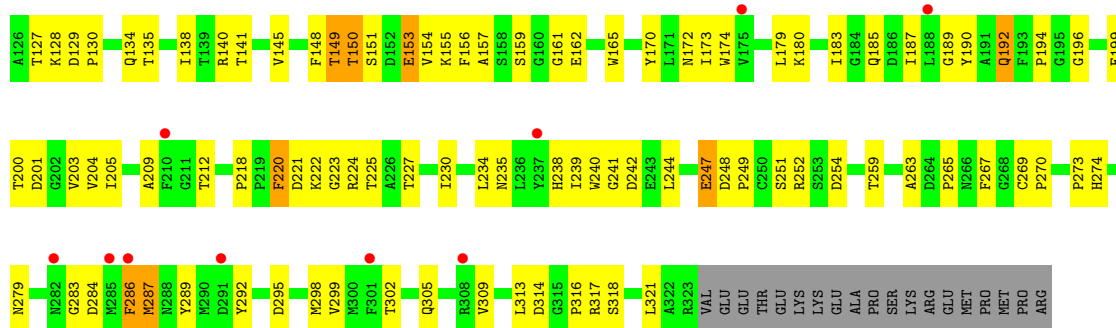


• Molecule 1: Ulilycin

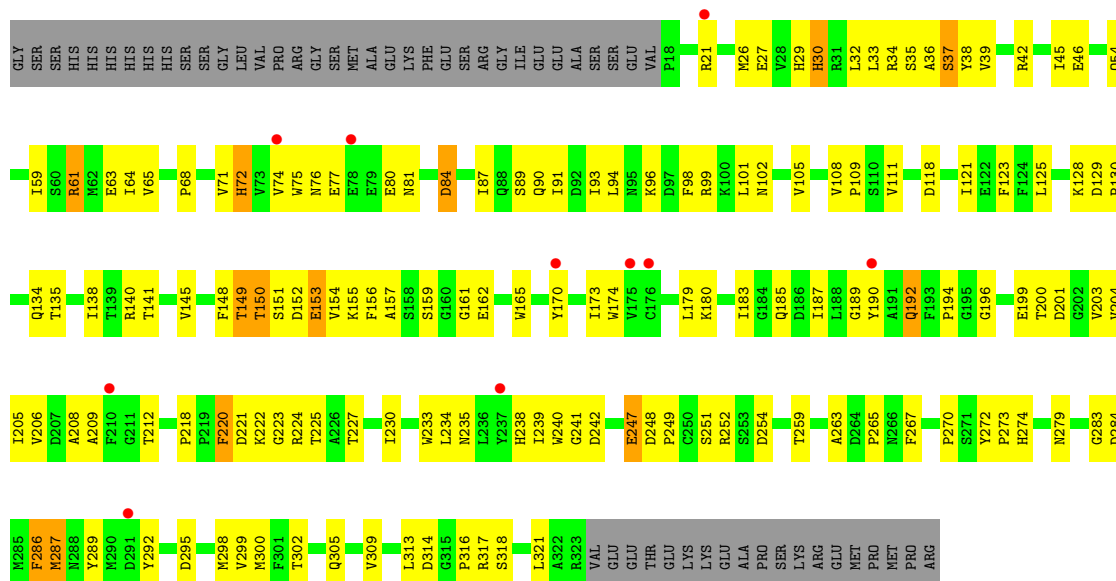
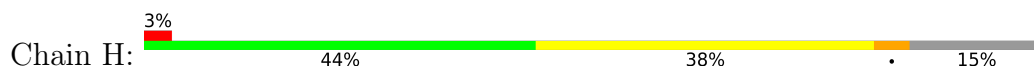


• Molecule 1: Ulilycin

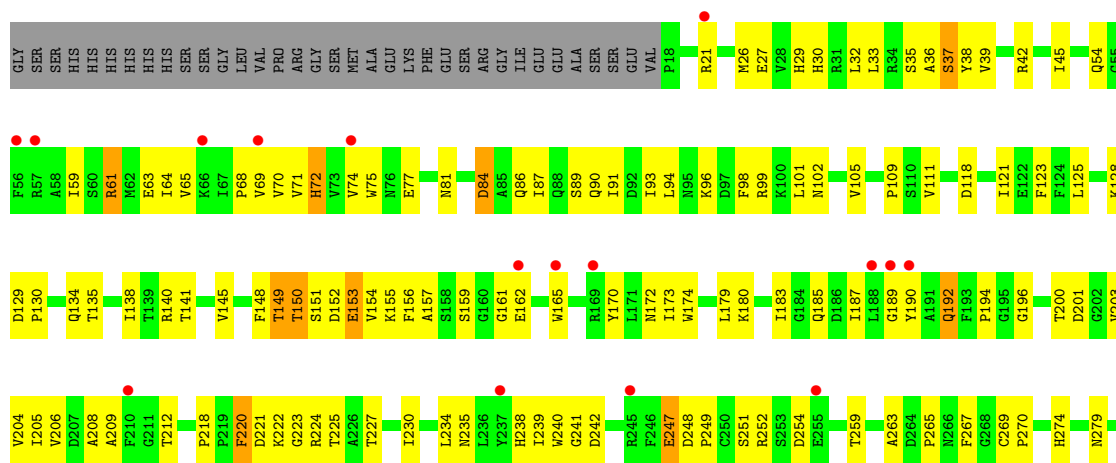
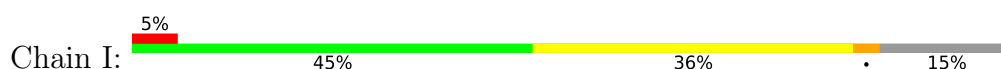




• Molecule 1: Ulilysin

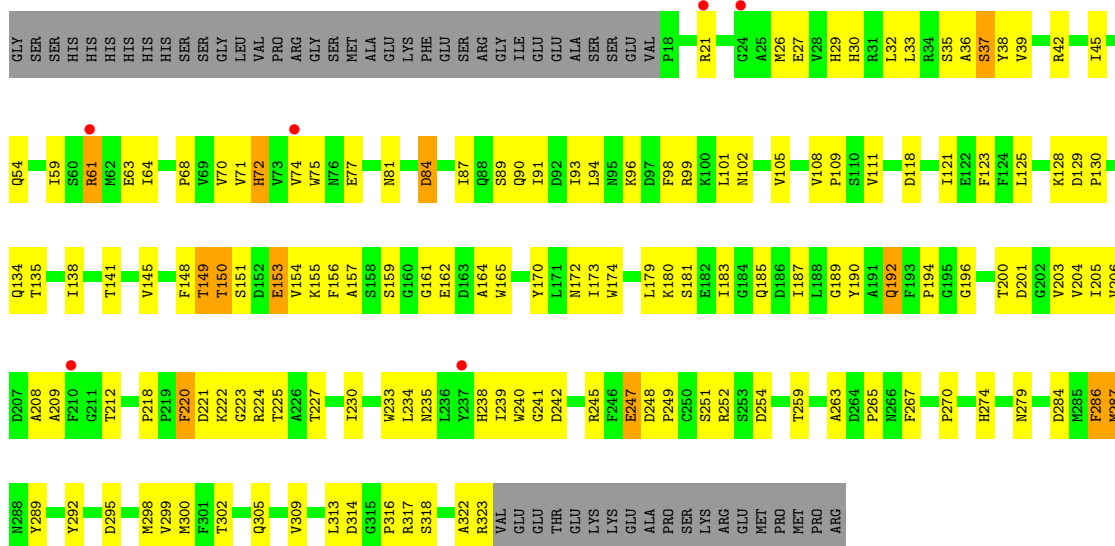


• Molecule 1: Ulilysin

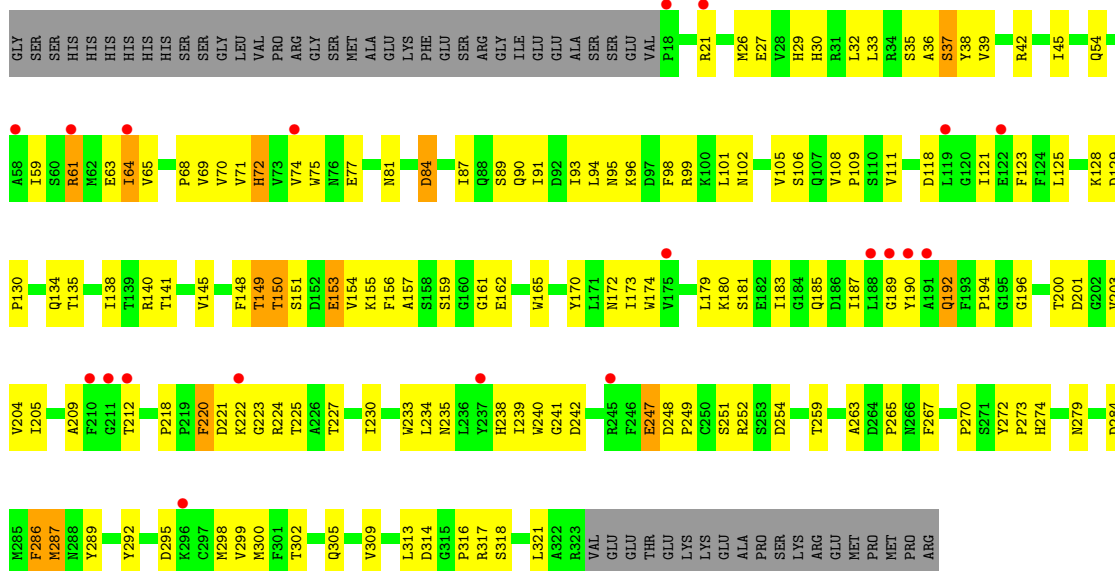




- Molecule 1: Uliysin

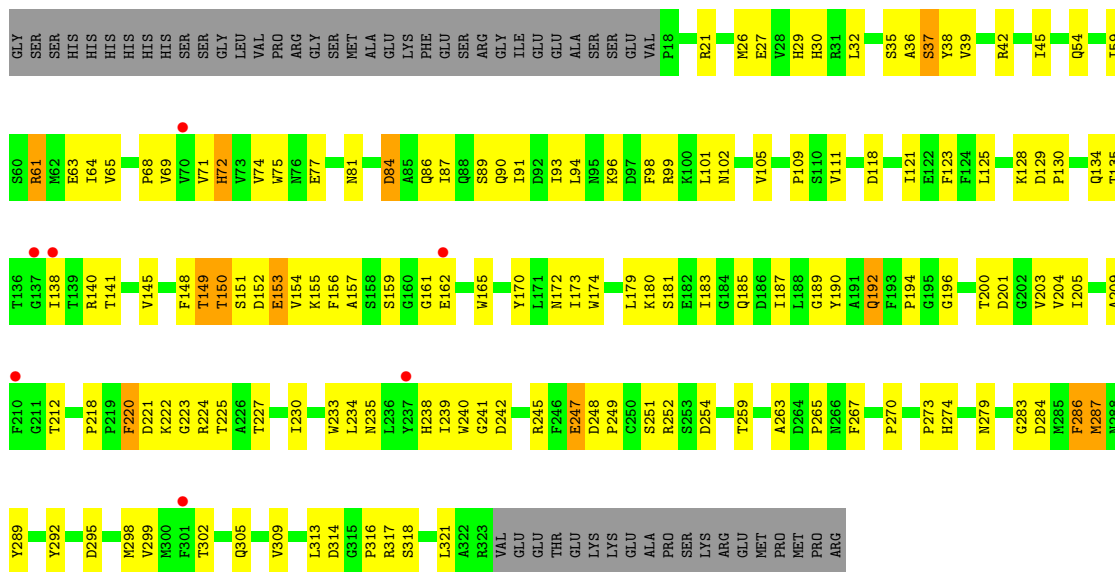


- Molecule 1: Uliysin

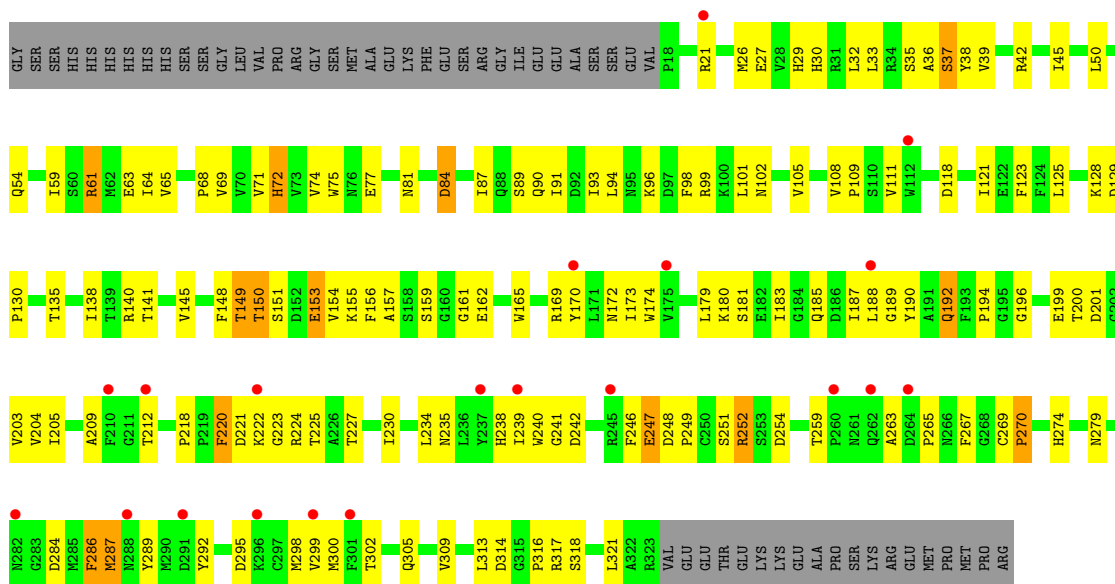


- Molecule 1: Uliysin

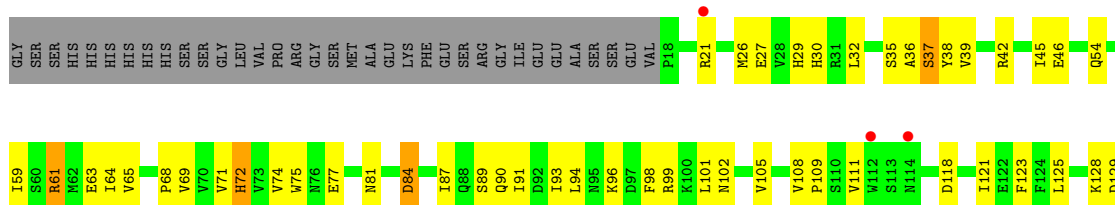


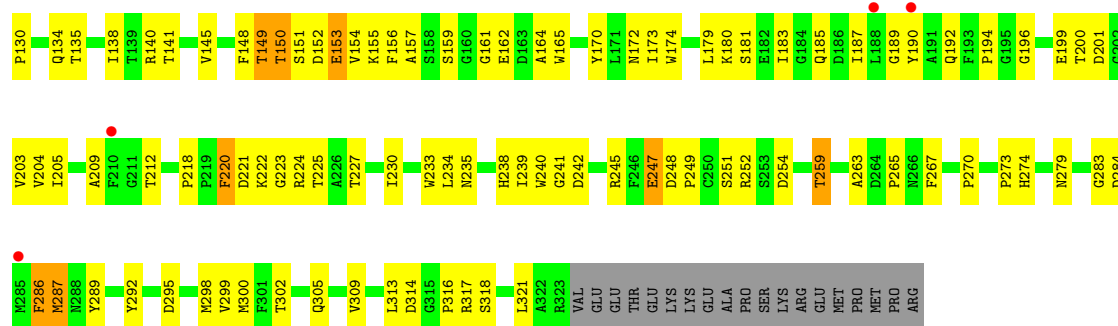


- Molecule 1: Ulilysin



- Molecule 1: Ulilysin





4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	192.10Å 544.40Å 185.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	94.72 – 4.50 94.72 – 4.50	Depositor EDS
% Data completeness (in resolution range)	98.6 (94.72-4.50) 99.6 (94.72-4.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.20 (at 4.47Å)	Xtrriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.289 , 0.302 0.293 , 0.304	Depositor DCC
R_{free} test set	845 reflections (1.46%)	wwPDB-VP
Wilson B-factor (Å ²)	131.9	Xtrriage
Anisotropy	1.029	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 151.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	33614	wwPDB-VP
Average B, all atoms (Å ²)	171.0	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 34.00 % 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 7.3091e-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: CA, ZN

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	0.38	0/2457	0.58	0/3340
1	B	0.42	0/2457	0.59	0/3340
1	C	0.39	0/2457	0.59	0/3340
1	D	0.40	0/2457	0.59	0/3340
1	E	0.41	0/2457	0.60	0/3340
1	F	0.41	0/2457	0.60	0/3340
1	G	0.43	0/2457	0.61	0/3340
1	H	0.47	1/2457 (0.0%)	0.62	0/3340
1	I	0.42	0/2457	0.59	0/3340
1	J	0.43	0/2457	0.60	0/3340
1	K	0.40	0/2457	0.59	0/3340
1	L	0.41	0/2457	0.59	0/3340
1	M	0.39	0/2457	0.59	0/3340
1	N	0.42	0/2457	0.60	0/3340
All	All	0.41	1/34398 (0.0%)	0.60	0/46760

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	30	HIS	CB-CG	5.32	1.59	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2399	0	2279	112	0
1	B	2399	0	2279	113	1
1	C	2399	0	2279	107	0
1	D	2399	0	2279	107	0
1	E	2399	0	2279	104	0
1	F	2399	0	2279	113	0
1	G	2399	0	2279	104	0
1	H	2399	0	2279	108	0
1	I	2399	0	2279	106	0
1	J	2399	0	2279	106	0
1	K	2399	0	2279	104	0
1	L	2399	0	2279	102	0
1	M	2399	0	2279	103	0
1	N	2399	0	2279	104	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0
2	N	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
3	M	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	N	1	0	0	0	0
All	All	33614	0	31906	1446	1

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 22.

All (1446) 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:N:174:TRP:HB2	1:N:204:VAL:HG22	1.61	0.82
1:H:174:TRP:HB2	1:H:204:VAL:HG22	1.62	0.82
1:I:174:TRP:HB2	1:I:204:VAL:HG22	1.62	0.81
1:B:174:TRP:HB2	1:B:204:VAL:HG22	1.62	0.80
1:G:174:TRP:HB2	1:G:204:VAL:HG22	1.64	0.79
1:K:174:TRP:HB2	1:K:204:VAL:HG22	1.65	0.79
1:D:174:TRP:HB2	1:D:204:VAL:HG22	1.64	0.79
1:L:174:TRP:HB2	1:L:204:VAL:HG22	1.64	0.79
1:M:174:TRP:HB2	1:M:204:VAL:HG22	1.66	0.78
1:F:174:TRP:HB2	1:F:204:VAL:HG22	1.64	0.78
1:N:29:HIS:HA	1:N:32:LEU:HD12	1.66	0.78
1:H:29:HIS:HA	1:H:32:LEU:HD12	1.65	0.77
1:J:174:TRP:HB2	1:J:204:VAL:HG22	1.66	0.77
1:E:29:HIS:HA	1:E:32:LEU:HD12	1.66	0.77
1:A:29:HIS:HA	1:A:32:LEU:HD12	1.66	0.76
1:K:224:ARG:HD2	1:K:289:TYR:HE1	1.50	0.76
1:A:174:TRP:HB2	1:A:204:VAL:HG22	1.67	0.76
1:J:29:HIS:HA	1:J:32:LEU:HD12	1.65	0.76
1:D:29:HIS:HA	1:D:32:LEU:HD12	1.68	0.76
1:E:174:TRP:HB2	1:E:204:VAL:HG22	1.66	0.76
1:K:29:HIS:HA	1:K:32:LEU:HD12	1.67	0.75
1:L:29:HIS:HA	1:L:32:LEU:HD12	1.68	0.75
1:A:279:ASN:HA	1:B:279:ASN:H	1.52	0.75
1:B:224:ARG:HD2	1:B:289:TYR:HE1	1.52	0.75
1:C:224:ARG:HD2	1:C:289:TYR:HE1	1.53	0.74
1:B:29:HIS:HA	1:B:32:LEU:HD12	1.69	0.74
1:C:174:TRP:HB2	1:C:204:VAL:HG22	1.68	0.74
1:I:29:HIS:HA	1:I:32:LEU:HD12	1.68	0.74
1:C:29:HIS:HA	1:C:32:LEU:HD12	1.70	0.74
1:E:224:ARG:HD2	1:E:289:TYR:HE1	1.52	0.74
1:M:29:HIS:HA	1:M:32:LEU:HD12	1.67	0.74
1:G:224:ARG:HD2	1:G:289:TYR:HE1	1.52	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:224:ARG:HD2	1:J:289:TYR:HE1	1.52	0.74
1:M:224:ARG:HD2	1:M:289:TYR:HE1	1.52	0.74
1:M:151:SER:HB2	1:M:153:GLU:HG3	1.70	0.74
1:F:249:PRO:HB2	1:F:263:ALA:HB1	1.71	0.73
1:H:224:ARG:HD2	1:H:289:TYR:HE1	1.53	0.73
1:L:234:LEU:HB3	1:L:309:VAL:HG13	1.70	0.73
1:F:224:ARG:HD2	1:F:289:TYR:HE1	1.53	0.73
1:J:45:ILE:HG12	1:J:316:PRO:HG3	1.69	0.73
1:M:45:ILE:HG12	1:M:316:PRO:HG3	1.70	0.73
1:A:224:ARG:HD2	1:A:289:TYR:HE1	1.53	0.72
1:A:279:ASN:H	1:B:279:ASN:HA	1.54	0.72
1:G:45:ILE:HG12	1:G:316:PRO:HG3	1.69	0.72
1:D:224:ARG:HD2	1:D:289:TYR:HE1	1.53	0.72
1:I:151:SER:HB2	1:I:153:GLU:HG3	1.70	0.72
1:G:249:PRO:HB2	1:G:263:ALA:HB1	1.72	0.72
1:I:249:PRO:HB2	1:I:263:ALA:HB1	1.70	0.72
1:G:29:HIS:HA	1:G:32:LEU:HD12	1.71	0.72
1:E:45:ILE:HG12	1:E:316:PRO:HG3	1.72	0.72
1:B:45:ILE:HG12	1:B:316:PRO:HG3	1.72	0.71
1:K:151:SER:HB2	1:K:153:GLU:HG3	1.71	0.71
1:K:249:PRO:HB2	1:K:263:ALA:HB1	1.72	0.71
1:I:224:ARG:HD2	1:I:289:TYR:HE1	1.55	0.71
1:J:151:SER:HB2	1:J:153:GLU:HG3	1.72	0.71
1:N:224:ARG:HD2	1:N:289:TYR:HE1	1.53	0.71
1:A:249:PRO:HB2	1:A:263:ALA:HB1	1.72	0.71
1:F:29:HIS:HA	1:F:32:LEU:HD12	1.72	0.71
1:E:151:SER:HB2	1:E:153:GLU:HG3	1.72	0.71
1:B:249:PRO:HB2	1:B:263:ALA:HB1	1.72	0.71
1:J:249:PRO:HB2	1:J:263:ALA:HB1	1.72	0.71
1:F:151:SER:HB2	1:F:153:GLU:HG3	1.73	0.71
1:F:222:LYS:HB2	1:F:295:ASP:HB3	1.71	0.71
1:C:151:SER:HB2	1:C:153:GLU:HG3	1.72	0.71
1:C:247:GLU:HG2	1:C:251:SER:HB2	1.72	0.71
1:H:249:PRO:HB2	1:H:263:ALA:HB1	1.72	0.71
1:K:45:ILE:HG12	1:K:316:PRO:HG3	1.73	0.71
1:L:224:ARG:HD2	1:L:289:TYR:HE1	1.55	0.71
1:N:249:PRO:HB2	1:N:263:ALA:HB1	1.72	0.71
1:D:151:SER:HB2	1:D:153:GLU:HG3	1.73	0.70
1:M:247:GLU:HG2	1:M:251:SER:HB2	1.73	0.70
1:A:151:SER:HB2	1:A:153:GLU:HG3	1.71	0.70
1:I:259:THR:HG22	1:I:302:THR:HG21	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:249:PRO:HB2	1:L:263:ALA:HB1	1.71	0.70
1:K:234:LEU:HB3	1:K:309:VAL:HG13	1.73	0.70
1:H:151:SER:HB2	1:H:153:GLU:HG3	1.73	0.70
1:J:247:GLU:HG2	1:J:251:SER:HB2	1.72	0.70
1:L:151:SER:HB2	1:L:153:GLU:HG3	1.72	0.70
1:G:151:SER:HB2	1:G:153:GLU:HG3	1.74	0.70
1:B:151:SER:HB2	1:B:153:GLU:HG3	1.74	0.70
1:C:222:LYS:HB2	1:C:295:ASP:HB3	1.74	0.70
1:C:249:PRO:HB2	1:C:263:ALA:HB1	1.73	0.69
1:L:259:THR:HG22	1:L:302:THR:HG21	1.74	0.69
1:G:247:GLU:HG2	1:G:251:SER:HB2	1.74	0.69
1:K:222:LYS:HB2	1:K:295:ASP:HB3	1.74	0.69
1:A:45:ILE:HG12	1:A:316:PRO:HG3	1.74	0.69
1:C:45:ILE:HG12	1:C:316:PRO:HG3	1.73	0.69
1:D:45:ILE:HG12	1:D:316:PRO:HG3	1.73	0.69
1:E:249:PRO:HB2	1:E:263:ALA:HB1	1.73	0.69
1:I:45:ILE:HG12	1:I:316:PRO:HG3	1.73	0.69
1:L:45:ILE:HG12	1:L:316:PRO:HG3	1.73	0.69
1:F:45:ILE:HG12	1:F:316:PRO:HG3	1.75	0.69
1:M:249:PRO:HB2	1:M:263:ALA:HB1	1.73	0.69
1:N:45:ILE:HG12	1:N:316:PRO:HG3	1.75	0.69
1:N:151:SER:HB2	1:N:153:GLU:HG3	1.75	0.69
1:D:249:PRO:HB2	1:D:263:ALA:HB1	1.74	0.69
1:B:234:LEU:HB3	1:B:309:VAL:HG13	1.73	0.69
1:E:247:GLU:HG2	1:E:251:SER:HB2	1.74	0.69
1:H:222:LYS:HB2	1:H:295:ASP:HB3	1.75	0.69
1:J:234:LEU:HB3	1:J:309:VAL:HG13	1.74	0.69
1:M:259:THR:HG22	1:M:302:THR:HG21	1.75	0.69
1:C:234:LEU:HB3	1:C:309:VAL:HG13	1.76	0.68
1:D:247:GLU:HG2	1:D:251:SER:HB2	1.75	0.68
1:H:45:ILE:HG12	1:H:316:PRO:HG3	1.74	0.68
1:J:222:LYS:HB2	1:J:295:ASP:HB3	1.75	0.68
1:A:234:LEU:HB3	1:A:309:VAL:HG13	1.76	0.68
1:G:259:THR:HG22	1:G:302:THR:HG21	1.76	0.68
1:I:222:LYS:HB2	1:I:295:ASP:HB3	1.76	0.68
1:K:247:GLU:HG2	1:K:251:SER:HB2	1.75	0.68
1:D:259:THR:HG22	1:D:302:THR:HG21	1.74	0.68
1:J:180:LYS:NZ	1:L:134:GLN:OE1	2.27	0.68
1:N:259:THR:HG22	1:N:302:THR:HG21	1.76	0.68
1:D:222:LYS:HB2	1:D:295:ASP:HB3	1.75	0.68
1:E:222:LYS:HB2	1:E:295:ASP:HB3	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:247:GLU:HG2	1:F:251:SER:HB2	1.75	0.68
1:N:222:LYS:HB2	1:N:295:ASP:HB3	1.76	0.68
1:A:222:LYS:HB2	1:A:295:ASP:HB3	1.76	0.67
1:H:247:GLU:HG2	1:H:251:SER:HB2	1.77	0.67
1:A:247:GLU:HG2	1:A:251:SER:HB2	1.77	0.67
1:F:259:THR:HG22	1:F:302:THR:HG21	1.77	0.67
1:G:222:LYS:HB2	1:G:295:ASP:HB3	1.77	0.67
1:G:234:LEU:HB3	1:G:309:VAL:HG13	1.75	0.67
1:I:87:ILE:HD13	1:I:138:ILE:HD13	1.77	0.67
1:L:222:LYS:HB2	1:L:295:ASP:HB3	1.75	0.67
1:I:234:LEU:HB3	1:I:309:VAL:HG13	1.77	0.66
1:J:259:THR:HG22	1:J:302:THR:HG21	1.77	0.66
1:K:259:THR:HG22	1:K:302:THR:HG21	1.77	0.66
1:M:222:LYS:HB2	1:M:295:ASP:HB3	1.77	0.66
1:B:259:THR:HG22	1:B:302:THR:HG21	1.77	0.66
1:H:234:LEU:HB3	1:H:309:VAL:HG13	1.76	0.66
1:B:222:LYS:HB2	1:B:295:ASP:HB3	1.78	0.66
1:N:247:GLU:HG2	1:N:251:SER:HB2	1.77	0.66
1:B:241:GLY:HA3	1:B:252:ARG:HB2	1.78	0.66
1:N:234:LEU:HB3	1:N:309:VAL:HG13	1.76	0.66
1:C:259:THR:HG22	1:C:302:THR:HG21	1.77	0.66
1:F:234:LEU:HB3	1:F:309:VAL:HG13	1.77	0.66
1:L:247:GLU:HG2	1:L:251:SER:HB2	1.77	0.66
1:D:87:ILE:HD13	1:D:138:ILE:HD13	1.78	0.65
1:B:247:GLU:HG2	1:B:251:SER:HB2	1.77	0.65
1:C:241:GLY:HA3	1:C:252:ARG:HB2	1.77	0.65
1:E:234:LEU:HB3	1:E:309:VAL:HG13	1.77	0.65
1:E:259:THR:HG22	1:E:302:THR:HG21	1.76	0.65
1:I:134:GLN:OE1	1:K:180:LYS:NZ	2.29	0.65
1:I:247:GLU:HG2	1:I:251:SER:HB2	1.79	0.65
1:G:241:GLY:HA3	1:G:252:ARG:HB2	1.79	0.65
1:D:234:LEU:HB3	1:D:309:VAL:HG13	1.77	0.65
1:I:74:VAL:HG11	1:I:161:GLY:HA2	1.78	0.65
1:F:87:ILE:HD13	1:F:138:ILE:HD13	1.78	0.65
1:M:234:LEU:HB3	1:M:309:VAL:HG13	1.79	0.65
1:J:74:VAL:HG11	1:J:161:GLY:HA2	1.79	0.64
1:B:180:LYS:NZ	1:D:134:GLN:OE1	2.31	0.64
1:C:74:VAL:HG11	1:C:161:GLY:HA2	1.80	0.64
1:D:241:GLY:HA3	1:D:252:ARG:HB2	1.78	0.64
1:H:259:THR:HG22	1:H:302:THR:HG21	1.80	0.64
1:H:21:ARG:HD2	1:H:292:TYR:HA	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:180:LYS:NZ	1:J:134:GLN:OE1	2.31	0.63
1:K:21:ARG:HD2	1:K:292:TYR:HA	1.79	0.63
1:E:89:SER:HB2	1:E:212:THR:HG23	1.81	0.63
1:G:21:ARG:HD2	1:G:292:TYR:HA	1.81	0.63
1:H:87:ILE:HD13	1:H:138:ILE:HD13	1.80	0.63
1:H:89:SER:HB2	1:H:212:THR:HG23	1.81	0.63
1:A:259:THR:HG22	1:A:302:THR:HG21	1.79	0.63
1:E:241:GLY:HA3	1:E:252:ARG:HB2	1.81	0.63
1:G:279:ASN:HA	1:H:279:ASN:H	1.64	0.63
1:H:190:TYR:O	1:H:203:VAL:HG13	1.99	0.63
1:A:72:HIS:HD2	1:A:162:GLU:HB3	1.64	0.63
1:G:218:PRO:HA	1:G:220:PHE:H	1.64	0.63
1:G:279:ASN:H	1:H:279:ASN:HA	1.63	0.63
1:I:148:PHE:CD1	1:I:154:VAL:HB	2.34	0.63
1:A:74:VAL:HG11	1:A:161:GLY:HA2	1.81	0.63
1:M:87:ILE:HD13	1:M:138:ILE:HD13	1.80	0.63
1:L:72:HIS:HD2	1:L:162:GLU:HB3	1.63	0.63
1:L:241:GLY:HA3	1:L:252:ARG:HB2	1.81	0.63
1:L:74:VAL:HG11	1:L:161:GLY:HA2	1.80	0.63
1:A:241:GLY:HA3	1:A:252:ARG:HB2	1.82	0.62
1:E:190:TYR:O	1:E:203:VAL:HG13	1.99	0.62
1:B:87:ILE:HD13	1:B:138:ILE:HD13	1.80	0.62
1:L:87:ILE:HD13	1:L:138:ILE:HD13	1.79	0.62
1:E:21:ARG:HD2	1:E:292:TYR:HA	1.82	0.62
1:N:241:GLY:HA3	1:N:252:ARG:HB2	1.79	0.62
1:G:87:ILE:HD13	1:G:138:ILE:HD13	1.82	0.62
1:I:72:HIS:HD2	1:I:162:GLU:HB3	1.65	0.62
1:J:21:ARG:HD2	1:J:292:TYR:HA	1.82	0.62
1:A:102:ASN:ND2	1:A:299:VAL:O	2.32	0.62
1:F:72:HIS:HD2	1:F:162:GLU:HB3	1.65	0.62
1:F:190:TYR:O	1:F:203:VAL:HG13	2.00	0.62
1:L:21:ARG:HD2	1:L:292:TYR:HA	1.82	0.62
1:E:148:PHE:CD1	1:E:154:VAL:HB	2.34	0.62
1:G:74:VAL:HG11	1:G:161:GLY:HA2	1.80	0.62
1:L:190:TYR:O	1:L:203:VAL:HG13	2.00	0.62
1:M:74:VAL:HG11	1:M:161:GLY:HA2	1.82	0.62
1:B:102:ASN:ND2	1:B:299:VAL:O	2.32	0.61
1:D:74:VAL:HG11	1:D:161:GLY:HA2	1.81	0.61
1:K:74:VAL:HG11	1:K:161:GLY:HA2	1.82	0.61
1:E:134:GLN:OE1	1:G:180:LYS:NZ	2.33	0.61
1:B:218:PRO:HA	1:B:220:PHE:H	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:190:TYR:O	1:J:203:VAL:HG13	2.00	0.61
1:E:87:ILE:HD13	1:E:138:ILE:HD13	1.82	0.61
1:H:241:GLY:HA3	1:H:252:ARG:HB2	1.82	0.61
1:B:154:VAL:O	1:B:161:GLY:HA3	2.00	0.61
1:D:102:ASN:ND2	1:D:299:VAL:O	2.31	0.61
1:B:21:ARG:HD2	1:B:292:TYR:HA	1.81	0.61
1:J:72:HIS:HD2	1:J:162:GLU:HB3	1.65	0.61
1:K:241:GLY:HA3	1:K:252:ARG:HB2	1.81	0.61
1:D:21:ARG:HD2	1:D:292:TYR:HA	1.82	0.61
1:E:74:VAL:HG11	1:E:161:GLY:HA2	1.81	0.61
1:K:72:HIS:HD2	1:K:162:GLU:HB3	1.66	0.61
1:N:87:ILE:HD13	1:N:138:ILE:HD13	1.82	0.61
1:A:87:ILE:HD13	1:A:138:ILE:HD13	1.83	0.61
1:I:241:GLY:HA3	1:I:252:ARG:HB2	1.83	0.61
1:J:179:LEU:HB3	1:J:187:ILE:HG13	1.83	0.61
1:N:74:VAL:HG11	1:N:161:GLY:HA2	1.83	0.61
1:N:179:LEU:HB3	1:N:187:ILE:HG13	1.83	0.61
1:A:179:LEU:HB3	1:A:187:ILE:HG13	1.83	0.61
1:A:21:ARG:HD2	1:A:292:TYR:HA	1.83	0.60
1:C:102:ASN:ND2	1:C:299:VAL:O	2.33	0.60
1:D:218:PRO:HA	1:D:220:PHE:H	1.66	0.60
1:K:87:ILE:HD13	1:K:138:ILE:HD13	1.82	0.60
1:K:148:PHE:CD1	1:K:154:VAL:HB	2.36	0.60
1:L:180:LYS:NZ	1:N:134:GLN:OE1	2.34	0.60
1:L:218:PRO:HA	1:L:220:PHE:H	1.66	0.60
1:A:90:GLN:O	1:A:94:LEU:HG	2.01	0.60
1:A:154:VAL:O	1:A:161:GLY:HA3	2.02	0.60
1:C:87:ILE:HD13	1:C:138:ILE:HD13	1.83	0.60
1:K:179:LEU:HB3	1:K:187:ILE:HG13	1.83	0.60
1:A:218:PRO:HA	1:A:220:PHE:H	1.67	0.60
1:D:180:LYS:NZ	1:F:134:GLN:OE1	2.34	0.60
1:H:74:VAL:HG11	1:H:161:GLY:HA2	1.82	0.60
1:I:190:TYR:O	1:I:203:VAL:HG13	2.00	0.60
1:J:241:GLY:HA3	1:J:252:ARG:HB2	1.82	0.60
1:C:21:ARG:HD2	1:C:292:TYR:HA	1.84	0.60
1:C:148:PHE:CD1	1:C:154:VAL:HB	2.36	0.60
1:C:218:PRO:HA	1:C:220:PHE:H	1.67	0.60
1:D:148:PHE:CD1	1:D:154:VAL:HB	2.37	0.60
1:E:102:ASN:ND2	1:E:299:VAL:O	2.34	0.60
1:F:74:VAL:HG11	1:F:161:GLY:HA2	1.82	0.60
1:E:72:HIS:HD2	1:E:162:GLU:HB3	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:179:LEU:HB3	1:F:187:ILE:HG13	1.83	0.60
1:F:218:PRO:HA	1:F:220:PHE:H	1.66	0.60
1:M:21:ARG:HD2	1:M:292:TYR:HA	1.82	0.60
1:A:190:TYR:O	1:A:203:VAL:HG13	2.02	0.60
1:G:72:HIS:CD2	1:G:162:GLU:HB3	2.37	0.60
1:I:72:HIS:CD2	1:I:162:GLU:HB3	2.37	0.60
1:N:90:GLN:O	1:N:94:LEU:HG	2.02	0.60
1:F:89:SER:HB2	1:F:212:THR:HG23	1.84	0.59
1:K:218:PRO:HA	1:K:220:PHE:H	1.67	0.59
1:K:279:ASN:H	1:L:279:ASN:HA	1.67	0.59
1:M:218:PRO:HA	1:M:220:PHE:H	1.67	0.59
1:M:241:GLY:HA3	1:M:252:ARG:HB2	1.82	0.59
1:B:227:THR:HA	1:B:230:ILE:HD12	1.84	0.59
1:C:190:TYR:O	1:C:203:VAL:HG13	2.02	0.59
1:F:241:GLY:HA3	1:F:252:ARG:HB2	1.83	0.59
1:G:102:ASN:ND2	1:G:299:VAL:O	2.34	0.59
1:J:89:SER:HB2	1:J:212:THR:HG23	1.84	0.59
1:L:102:ASN:ND2	1:L:299:VAL:O	2.34	0.59
1:G:190:TYR:O	1:G:203:VAL:HG13	2.02	0.59
1:J:218:PRO:HA	1:J:220:PHE:H	1.66	0.59
1:L:72:HIS:CD2	1:L:162:GLU:HB3	2.37	0.59
1:L:89:SER:HB2	1:L:212:THR:HG23	1.84	0.59
1:N:190:TYR:O	1:N:203:VAL:HG13	2.02	0.59
1:C:90:GLN:O	1:C:94:LEU:HG	2.02	0.59
1:J:148:PHE:CD1	1:J:154:VAL:HB	2.37	0.59
1:L:179:LEU:HB3	1:L:187:ILE:HG13	1.84	0.59
1:M:179:LEU:HB3	1:M:187:ILE:HG13	1.85	0.59
1:N:102:ASN:ND2	1:N:299:VAL:O	2.33	0.59
1:A:72:HIS:CD2	1:A:162:GLU:HB3	2.37	0.59
1:I:240:TRP:CZ3	1:I:265:PRO:HB3	2.36	0.59
1:A:96:LYS:HA	1:A:101:LEU:HB2	1.83	0.59
1:B:74:VAL:HG11	1:B:161:GLY:HA2	1.84	0.59
1:F:180:LYS:NZ	1:H:134:GLN:OE1	2.36	0.59
1:M:89:SER:HB2	1:M:212:THR:HG23	1.85	0.59
1:C:179:LEU:HB3	1:C:187:ILE:HG13	1.84	0.59
1:H:218:PRO:HA	1:H:220:PHE:H	1.66	0.59
1:M:190:TYR:O	1:M:203:VAL:HG13	2.02	0.59
1:B:96:LYS:HA	1:B:101:LEU:HB2	1.84	0.59
1:D:190:TYR:O	1:D:203:VAL:HG13	2.03	0.59
1:G:134:GLN:OE1	1:I:180:LYS:NZ	2.36	0.59
1:G:148:PHE:CD1	1:G:154:VAL:HB	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:21:ARG:HD2	1:I:292:TYR:HA	1.84	0.59
1:K:190:TYR:O	1:K:203:VAL:HG13	2.03	0.59
1:G:89:SER:HB2	1:G:212:THR:HG23	1.85	0.59
1:J:87:ILE:HD13	1:J:138:ILE:HD13	1.84	0.59
1:F:21:ARG:HD2	1:F:292:TYR:HA	1.85	0.58
1:I:96:LYS:HA	1:I:101:LEU:HB2	1.85	0.58
1:I:102:ASN:ND2	1:I:299:VAL:O	2.35	0.58
1:M:72:HIS:HD2	1:M:162:GLU:HB3	1.66	0.58
1:N:148:PHE:CD1	1:N:154:VAL:HB	2.38	0.58
1:F:72:HIS:CD2	1:F:162:GLU:HB3	2.38	0.58
1:J:72:HIS:CD2	1:J:162:GLU:HB3	2.38	0.58
1:B:148:PHE:CD1	1:B:154:VAL:HB	2.38	0.58
1:H:148:PHE:CD1	1:H:154:VAL:HB	2.39	0.58
1:N:89:SER:HB2	1:N:212:THR:HG23	1.83	0.58
1:N:227:THR:HA	1:N:230:ILE:HD12	1.84	0.58
1:C:89:SER:HB2	1:C:212:THR:HG23	1.85	0.58
1:C:96:LYS:HA	1:C:101:LEU:HB2	1.86	0.58
1:E:90:GLN:O	1:E:94:LEU:HG	2.04	0.58
1:K:102:ASN:ND2	1:K:299:VAL:O	2.33	0.58
1:N:218:PRO:HA	1:N:220:PHE:H	1.67	0.58
1:B:190:TYR:O	1:B:203:VAL:HG13	2.03	0.58
1:F:154:VAL:O	1:F:161:GLY:HA3	2.04	0.58
1:D:179:LEU:HB3	1:D:187:ILE:HG13	1.85	0.58
1:I:179:LEU:HB3	1:I:187:ILE:HG13	1.84	0.58
1:L:148:PHE:CD1	1:L:154:VAL:HB	2.38	0.58
1:M:279:ASN:HA	1:N:279:ASN:H	1.69	0.58
1:D:240:TRP:CZ3	1:D:265:PRO:HB3	2.39	0.58
1:H:192:GLN:OE1	1:H:200:THR:OG1	2.17	0.58
1:I:153:GLU:O	1:I:159:SER:HB2	2.04	0.58
1:I:218:PRO:HA	1:I:220:PHE:H	1.67	0.58
1:N:154:VAL:O	1:N:161:GLY:HA3	2.03	0.58
1:D:227:THR:HA	1:D:230:ILE:HD12	1.86	0.58
1:H:154:VAL:O	1:H:161:GLY:HA3	2.04	0.58
1:K:96:LYS:HA	1:K:101:LEU:HB2	1.84	0.58
1:M:148:PHE:CD1	1:M:154:VAL:HB	2.39	0.58
1:K:183:ILE:HG13	1:K:185:GLN:H	1.69	0.57
1:M:154:VAL:O	1:M:161:GLY:HA3	2.03	0.57
1:N:21:ARG:HD2	1:N:292:TYR:HA	1.86	0.57
1:D:89:SER:HB2	1:D:212:THR:HG23	1.85	0.57
1:A:71:VAL:HA	1:A:173:ILE:HB	1.86	0.57
1:A:148:PHE:CD1	1:A:154:VAL:HB	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:96:LYS:HA	1:E:101:LEU:HB2	1.85	0.57
1:K:72:HIS:CD2	1:K:162:GLU:HB3	2.39	0.57
1:L:96:LYS:HA	1:L:101:LEU:HB2	1.86	0.57
1:M:90:GLN:O	1:M:94:LEU:HG	2.03	0.57
1:F:227:THR:HA	1:F:230:ILE:HD12	1.86	0.57
1:I:89:SER:HB2	1:I:212:THR:HG23	1.85	0.57
1:B:179:LEU:HB3	1:B:187:ILE:HG13	1.85	0.57
1:C:279:ASN:H	1:D:279:ASN:HA	1.69	0.57
1:D:154:VAL:O	1:D:161:GLY:HA3	2.05	0.57
1:M:153:GLU:O	1:M:159:SER:HB2	2.05	0.57
1:M:279:ASN:H	1:N:279:ASN:HA	1.70	0.57
1:D:72:HIS:CD2	1:D:162:GLU:HB3	2.39	0.57
1:F:148:PHE:CD1	1:F:154:VAL:HB	2.39	0.57
1:G:72:HIS:HD2	1:G:162:GLU:HB3	1.69	0.57
1:A:89:SER:HB2	1:A:212:THR:HG23	1.87	0.57
1:I:154:VAL:O	1:I:161:GLY:HA3	2.04	0.57
1:K:279:ASN:HA	1:L:279:ASN:H	1.69	0.57
1:N:96:LYS:HA	1:N:101:LEU:HB2	1.87	0.57
1:J:96:LYS:HA	1:J:101:LEU:HB2	1.85	0.57
1:K:89:SER:HB2	1:K:212:THR:HG23	1.86	0.57
1:E:179:LEU:HB3	1:E:187:ILE:HG13	1.86	0.57
1:G:90:GLN:O	1:G:94:LEU:HG	2.04	0.57
1:I:183:ILE:HG13	1:I:185:GLN:H	1.69	0.57
1:M:72:HIS:CD2	1:M:162:GLU:HB3	2.40	0.57
1:M:227:THR:HA	1:M:230:ILE:HD12	1.86	0.57
1:B:90:GLN:O	1:B:94:LEU:HG	2.05	0.56
1:H:227:THR:HA	1:H:230:ILE:HD12	1.87	0.56
1:K:90:GLN:O	1:K:94:LEU:HG	2.04	0.56
1:L:71:VAL:HA	1:L:173:ILE:HB	1.87	0.56
1:A:227:THR:HA	1:A:230:ILE:HD12	1.86	0.56
1:H:102:ASN:ND2	1:H:299:VAL:O	2.37	0.56
1:J:240:TRP:CZ3	1:J:265:PRO:HB3	2.40	0.56
1:E:72:HIS:CD2	1:E:162:GLU:HB3	2.39	0.56
1:E:218:PRO:HA	1:E:220:PHE:H	1.69	0.56
1:G:240:TRP:CZ3	1:G:265:PRO:HB3	2.40	0.56
1:K:240:TRP:CZ3	1:K:265:PRO:HB3	2.41	0.56
1:L:90:GLN:O	1:L:94:LEU:HG	2.06	0.56
1:L:227:THR:HA	1:L:230:ILE:HD12	1.86	0.56
1:N:72:HIS:CD2	1:N:162:GLU:HB3	2.39	0.56
1:B:89:SER:HB2	1:B:212:THR:HG23	1.86	0.56
1:I:227:THR:HA	1:I:230:ILE:HD12	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:154:VAL:O	1:C:161:GLY:HA3	2.05	0.56
1:D:96:LYS:HA	1:D:101:LEU:HB2	1.88	0.56
1:J:71:VAL:HA	1:J:173:ILE:HB	1.87	0.56
1:J:90:GLN:O	1:J:94:LEU:HG	2.06	0.56
1:L:154:VAL:O	1:L:161:GLY:HA3	2.05	0.56
1:N:71:VAL:HA	1:N:173:ILE:HB	1.88	0.56
1:C:153:GLU:O	1:C:159:SER:HB2	2.06	0.56
1:H:179:LEU:HB3	1:H:187:ILE:HG13	1.87	0.56
1:A:153:GLU:O	1:A:159:SER:HB2	2.04	0.56
1:C:72:HIS:CD2	1:C:162:GLU:HB3	2.40	0.56
1:C:183:ILE:HG13	1:C:185:GLN:H	1.71	0.56
1:C:279:ASN:HA	1:D:279:ASN:H	1.70	0.56
1:B:72:HIS:CD2	1:B:162:GLU:HB3	2.40	0.56
1:D:153:GLU:O	1:D:159:SER:HB2	2.05	0.56
1:F:96:LYS:HA	1:F:101:LEU:HB2	1.88	0.56
1:H:90:GLN:O	1:H:94:LEU:HG	2.05	0.56
1:I:279:ASN:H	1:J:279:ASN:HA	1.71	0.56
1:M:96:LYS:HA	1:M:101:LEU:HB2	1.87	0.56
1:F:201:ASP:OD1	1:F:317:ARG:NH2	2.30	0.56
1:H:72:HIS:CD2	1:H:162:GLU:HB3	2.41	0.56
1:A:109:PRO:HB2	1:A:111:VAL:HG12	1.88	0.55
1:G:154:VAL:O	1:G:161:GLY:HA3	2.06	0.55
1:G:179:LEU:HB3	1:G:187:ILE:HG13	1.87	0.55
1:H:183:ILE:HG13	1:H:185:GLN:H	1.70	0.55
1:H:240:TRP:CZ3	1:H:265:PRO:HB3	2.41	0.55
1:L:153:GLU:O	1:L:159:SER:HB2	2.06	0.55
1:M:183:ILE:HG13	1:M:185:GLN:H	1.71	0.55
1:E:154:VAL:O	1:E:161:GLY:HA3	2.06	0.55
1:F:153:GLU:O	1:F:159:SER:HB2	2.07	0.55
1:K:154:VAL:O	1:K:161:GLY:HA3	2.05	0.55
1:M:240:TRP:CZ3	1:M:265:PRO:HB3	2.42	0.55
1:E:240:TRP:CZ3	1:E:265:PRO:HB3	2.41	0.55
1:F:183:ILE:HG13	1:F:185:GLN:H	1.70	0.55
1:I:279:ASN:HA	1:J:279:ASN:H	1.71	0.55
1:J:149:THR:HB	1:J:151:SER:H	1.71	0.55
1:K:153:GLU:O	1:K:159:SER:HB2	2.06	0.55
1:J:227:THR:HA	1:J:230:ILE:HD12	1.88	0.55
1:D:90:GLN:O	1:D:94:LEU:HG	2.07	0.55
1:E:227:THR:HA	1:E:230:ILE:HD12	1.88	0.55
1:F:102:ASN:ND2	1:F:299:VAL:O	2.34	0.55
1:C:309:VAL:O	1:C:313:LEU:HG	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:149:THR:HB	1:K:151:SER:H	1.72	0.55
1:M:102:ASN:ND2	1:M:299:VAL:O	2.34	0.55
1:N:61:ARG:HH12	1:N:121:ILE:HD11	1.72	0.55
1:C:71:VAL:HA	1:C:173:ILE:HB	1.88	0.55
1:D:309:VAL:O	1:D:313:LEU:HG	2.07	0.55
1:K:71:VAL:HA	1:K:173:ILE:HB	1.89	0.55
1:B:209:ALA:HA	1:B:220:PHE:HB2	1.88	0.54
1:C:72:HIS:HD2	1:C:162:GLU:HB3	1.73	0.54
1:D:149:THR:HB	1:D:151:SER:H	1.72	0.54
1:K:227:THR:HA	1:K:230:ILE:HD12	1.88	0.54
1:L:240:TRP:CZ3	1:L:265:PRO:HB3	2.41	0.54
1:D:183:ILE:HG13	1:D:185:GLN:H	1.72	0.54
1:H:109:PRO:HB2	1:H:111:VAL:HG12	1.88	0.54
1:I:90:GLN:O	1:I:94:LEU:HG	2.06	0.54
1:M:270:PRO:HG2	1:M:286:PHE:CG	2.43	0.54
1:C:87:ILE:O	1:C:91:ILE:HG12	2.07	0.54
1:J:109:PRO:HB2	1:J:111:VAL:HG12	1.87	0.54
1:M:68:PRO:HB2	1:M:170:TYR:HD1	1.72	0.54
1:I:289:TYR:HA	1:I:298:MET:HB2	1.88	0.54
1:B:153:GLU:O	1:B:159:SER:HB2	2.07	0.54
1:B:240:TRP:CZ3	1:B:265:PRO:HB3	2.42	0.54
1:B:309:VAL:O	1:B:313:LEU:HG	2.08	0.54
1:E:109:PRO:HB2	1:E:111:VAL:HG12	1.89	0.54
1:H:270:PRO:HG2	1:H:286:PHE:CG	2.43	0.54
1:B:68:PRO:HB2	1:B:170:TYR:HD1	1.73	0.54
1:C:227:THR:HA	1:C:230:ILE:HD12	1.88	0.54
1:F:149:THR:HB	1:F:151:SER:H	1.72	0.54
1:H:96:LYS:HA	1:H:101:LEU:HB2	1.89	0.54
1:A:240:TRP:CZ3	1:A:265:PRO:HB3	2.42	0.54
1:B:109:PRO:HB2	1:B:111:VAL:HG12	1.90	0.54
1:C:109:PRO:HB2	1:C:111:VAL:HG12	1.90	0.54
1:F:240:TRP:CZ3	1:F:265:PRO:HB3	2.43	0.54
1:G:227:THR:HA	1:G:230:ILE:HD12	1.88	0.54
1:N:72:HIS:HD2	1:N:162:GLU:HB3	1.72	0.54
1:N:109:PRO:HB2	1:N:111:VAL:HG12	1.89	0.54
1:D:209:ALA:HA	1:D:220:PHE:HB2	1.89	0.54
1:E:183:ILE:HG13	1:E:185:GLN:H	1.72	0.54
1:E:309:VAL:O	1:E:313:LEU:HG	2.07	0.54
1:G:71:VAL:HA	1:G:173:ILE:HB	1.90	0.54
1:G:84:ASP:OD1	1:G:140:ARG:NH2	2.35	0.54
1:G:183:ILE:HG13	1:G:185:GLN:H	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:309:VAL:O	1:J:313:LEU:HG	2.08	0.54
1:L:183:ILE:HG13	1:L:185:GLN:H	1.73	0.54
1:M:309:VAL:O	1:M:313:LEU:HG	2.07	0.54
1:A:149:THR:HB	1:A:151:SER:H	1.73	0.54
1:B:183:ILE:HG13	1:B:185:GLN:H	1.73	0.54
1:E:279:ASN:HA	1:F:279:ASN:H	1.72	0.54
1:J:183:ILE:HG13	1:J:185:GLN:H	1.73	0.54
1:L:149:THR:HB	1:L:151:SER:H	1.73	0.54
1:M:149:THR:HB	1:M:151:SER:H	1.73	0.54
1:I:270:PRO:HG2	1:I:286:PHE:CG	2.43	0.53
1:J:153:GLU:O	1:J:159:SER:HB2	2.07	0.53
1:L:289:TYR:HA	1:L:298:MET:HB2	1.90	0.53
1:N:87:ILE:O	1:N:91:ILE:HG12	2.08	0.53
1:E:68:PRO:HB2	1:E:170:TYR:HD1	1.74	0.53
1:E:270:PRO:HG2	1:E:286:PHE:CG	2.44	0.53
1:N:183:ILE:HG13	1:N:185:GLN:H	1.71	0.53
1:N:240:TRP:CZ3	1:N:265:PRO:HB3	2.43	0.53
1:D:155:LYS:HD3	1:D:200:THR:HA	1.91	0.53
1:G:109:PRO:HB2	1:G:111:VAL:HG12	1.89	0.53
1:G:270:PRO:HG2	1:G:286:PHE:CG	2.44	0.53
1:K:289:TYR:HA	1:K:298:MET:HB2	1.89	0.53
1:D:27:GLU:HB2	1:D:150:THR:HG21	1.91	0.53
1:G:149:THR:HB	1:G:151:SER:H	1.73	0.53
1:C:270:PRO:HG2	1:C:286:PHE:CG	2.43	0.53
1:G:209:ALA:HA	1:G:220:PHE:HB2	1.90	0.53
1:H:72:HIS:HD2	1:H:162:GLU:HB3	1.72	0.53
1:H:129:ASP:HB2	1:H:130:PRO:HD2	1.90	0.53
1:B:71:VAL:HA	1:B:173:ILE:HB	1.89	0.53
1:D:71:VAL:HA	1:D:173:ILE:HB	1.89	0.53
1:M:155:LYS:HD3	1:M:200:THR:HA	1.91	0.53
1:N:149:THR:HB	1:N:151:SER:H	1.72	0.53
1:M:87:ILE:O	1:M:91:ILE:HG12	2.09	0.53
1:A:209:ALA:HA	1:A:220:PHE:HB2	1.91	0.53
1:F:109:PRO:HB2	1:F:111:VAL:HG12	1.90	0.53
1:F:309:VAL:O	1:F:313:LEU:HG	2.08	0.53
1:I:149:THR:HB	1:I:151:SER:H	1.73	0.53
1:I:209:ALA:HA	1:I:220:PHE:HB2	1.91	0.53
1:J:87:ILE:O	1:J:91:ILE:HG12	2.09	0.53
1:K:155:LYS:HD3	1:K:200:THR:HA	1.91	0.53
1:M:201:ASP:OD1	1:M:317:ARG:NH2	2.32	0.53
1:A:183:ILE:HG13	1:A:185:GLN:H	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:84:ASP:HA	1:I:87:ILE:HD12	1.91	0.53
1:N:155:LYS:HD3	1:N:200:THR:HA	1.91	0.53
1:D:72:HIS:HD2	1:D:162:GLU:HB3	1.72	0.53
1:E:129:ASP:HB2	1:E:130:PRO:HD2	1.91	0.53
1:E:149:THR:HB	1:E:151:SER:H	1.74	0.53
1:H:289:TYR:HA	1:H:298:MET:HB2	1.90	0.53
1:I:68:PRO:HB2	1:I:170:TYR:HD1	1.74	0.53
1:J:154:VAL:O	1:J:161:GLY:HA3	2.09	0.53
1:B:149:THR:HB	1:B:151:SER:H	1.73	0.52
1:D:84:ASP:HA	1:D:87:ILE:HD12	1.91	0.52
1:D:87:ILE:O	1:D:91:ILE:HG12	2.09	0.52
1:F:90:GLN:O	1:F:94:LEU:HG	2.09	0.52
1:G:130:PRO:HD3	1:G:165:TRP:CE3	2.44	0.52
1:I:71:VAL:HA	1:I:173:ILE:HB	1.90	0.52
1:J:289:TYR:HA	1:J:298:MET:HB2	1.90	0.52
1:N:309:VAL:O	1:N:313:LEU:HG	2.07	0.52
1:E:130:PRO:HD3	1:E:165:TRP:CE3	2.43	0.52
1:F:90:GLN:NE2	1:F:223:GLY:O	2.42	0.52
1:I:109:PRO:HB2	1:I:111:VAL:HG12	1.89	0.52
1:K:87:ILE:O	1:K:91:ILE:HG12	2.09	0.52
1:M:130:PRO:HD3	1:M:165:TRP:CE3	2.44	0.52
1:N:209:ALA:HA	1:N:220:PHE:HB2	1.91	0.52
1:A:289:TYR:HA	1:A:298:MET:HB2	1.91	0.52
1:B:129:ASP:HB2	1:B:130:PRO:HD2	1.91	0.52
1:G:42:ARG:NH1	1:G:235:ASN:HD21	2.07	0.52
1:K:27:GLU:HB2	1:K:150:THR:HG21	1.91	0.52
1:K:134:GLN:OE1	1:M:180:LYS:NZ	2.43	0.52
1:M:71:VAL:HA	1:M:173:ILE:HB	1.91	0.52
1:M:109:PRO:HB2	1:M:111:VAL:HG12	1.90	0.52
1:N:153:GLU:O	1:N:159:SER:HB2	2.09	0.52
1:F:129:ASP:HB2	1:F:130:PRO:HD2	1.91	0.52
1:H:155:LYS:HD3	1:H:200:THR:HA	1.92	0.52
1:K:42:ARG:NH1	1:K:235:ASN:HD21	2.07	0.52
1:L:155:LYS:HD3	1:L:200:THR:HA	1.92	0.52
1:D:270:PRO:HG2	1:D:286:PHE:CG	2.44	0.52
1:F:61:ARG:HH12	1:F:121:ILE:HD11	1.74	0.52
1:B:72:HIS:HD2	1:B:162:GLU:HB3	1.73	0.52
1:B:130:PRO:HD3	1:B:165:TRP:CE3	2.44	0.52
1:C:155:LYS:HD3	1:C:200:THR:HA	1.91	0.52
1:D:34:ARG:HH22	1:F:323:ARG:C	2.13	0.52
1:E:279:ASN:H	1:F:279:ASN:HA	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:71:VAL:HA	1:F:173:ILE:HB	1.91	0.52
1:F:289:TYR:HA	1:F:298:MET:HB2	1.91	0.52
1:J:270:PRO:HG2	1:J:286:PHE:CG	2.45	0.52
1:A:27:GLU:HB2	1:A:150:THR:HG21	1.91	0.52
1:C:149:THR:HB	1:C:151:SER:H	1.75	0.52
1:K:109:PRO:HB2	1:K:111:VAL:HG12	1.90	0.52
1:L:61:ARG:HH12	1:L:121:ILE:HD11	1.75	0.52
1:L:270:PRO:HG2	1:L:286:PHE:CG	2.44	0.52
1:M:84:ASP:HA	1:M:87:ILE:HD12	1.92	0.52
1:N:27:GLU:HB2	1:N:150:THR:HG21	1.91	0.52
1:B:34:ARG:HH22	1:D:323:ARG:C	2.12	0.52
1:F:130:PRO:HD3	1:F:165:TRP:CE3	2.45	0.52
1:G:192:GLN:OE1	1:G:200:THR:OG1	2.21	0.52
1:G:289:TYR:HA	1:G:298:MET:HB2	1.91	0.52
1:K:130:PRO:HD3	1:K:165:TRP:CE3	2.45	0.52
1:L:109:PRO:HB2	1:L:111:VAL:HG12	1.90	0.52
1:L:209:ALA:HA	1:L:220:PHE:HB2	1.91	0.52
1:M:27:GLU:HB2	1:M:150:THR:HG21	1.92	0.52
1:D:26:MET:HG3	1:D:242:ASP:OD1	2.10	0.52
1:H:84:ASP:HA	1:H:87:ILE:HD12	1.92	0.52
1:H:149:THR:HB	1:H:151:SER:H	1.74	0.52
1:K:192:GLN:OE1	1:K:200:THR:OG1	2.21	0.52
1:A:84:ASP:HA	1:A:87:ILE:HD12	1.92	0.51
1:B:34:ARG:NH1	1:D:322:ALA:O	2.42	0.51
1:C:289:TYR:HA	1:C:298:MET:HB2	1.91	0.51
1:G:96:LYS:HA	1:G:101:LEU:HB2	1.90	0.51
1:G:155:LYS:HD3	1:G:200:THR:HA	1.92	0.51
1:H:201:ASP:OD1	1:H:317:ARG:NH2	2.33	0.51
1:H:309:VAL:O	1:H:313:LEU:HG	2.10	0.51
1:L:99:ARG:NH1	1:L:118:ASP:HB2	2.25	0.51
1:M:26:MET:HG3	1:M:242:ASP:OD1	2.10	0.51
1:N:129:ASP:HB2	1:N:130:PRO:HD2	1.91	0.51
1:A:270:PRO:HG2	1:A:286:PHE:CG	2.45	0.51
1:B:87:ILE:O	1:B:91:ILE:HG12	2.10	0.51
1:G:129:ASP:HB2	1:G:130:PRO:HD2	1.91	0.51
1:A:314:ASP:O	1:A:318:SER:HB3	2.10	0.51
1:B:61:ARG:HH12	1:B:121:ILE:HD11	1.76	0.51
1:E:87:ILE:O	1:E:91:ILE:HG12	2.09	0.51
1:I:61:ARG:HH12	1:I:121:ILE:HD11	1.75	0.51
1:I:129:ASP:HB2	1:I:130:PRO:HD2	1.93	0.51
1:E:155:LYS:HD3	1:E:200:THR:HA	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:27:GLU:HB2	1:J:150:THR:HG21	1.92	0.51
1:N:270:PRO:HG2	1:N:286:PHE:CG	2.45	0.51
1:K:84:ASP:HA	1:K:87:ILE:HD12	1.93	0.51
1:D:289:TYR:HA	1:D:298:MET:HB2	1.92	0.51
1:E:289:TYR:HA	1:E:298:MET:HB2	1.90	0.51
1:F:42:ARG:NH1	1:F:235:ASN:HD21	2.08	0.51
1:G:61:ARG:HH12	1:G:121:ILE:HD11	1.76	0.51
1:G:68:PRO:HB2	1:G:170:TYR:HD1	1.75	0.51
1:K:68:PRO:HB2	1:K:170:TYR:HD1	1.75	0.51
1:K:270:PRO:HG2	1:K:286:PHE:CG	2.46	0.51
1:N:84:ASP:HA	1:N:87:ILE:HD12	1.92	0.51
1:B:155:LYS:HD3	1:B:200:THR:HA	1.93	0.51
1:B:270:PRO:HG2	1:B:286:PHE:CG	2.45	0.51
1:C:240:TRP:CZ3	1:C:265:PRO:HB3	2.46	0.51
1:D:34:ARG:NH1	1:F:322:ALA:O	2.44	0.51
1:D:38:TYR:CE2	1:D:42:ARG:HG3	2.46	0.51
1:D:129:ASP:HB2	1:D:130:PRO:HD2	1.92	0.51
1:J:68:PRO:HB2	1:J:170:TYR:HD1	1.74	0.51
1:J:155:LYS:HD3	1:J:200:THR:HA	1.93	0.51
1:L:90:GLN:NE2	1:L:223:GLY:O	2.44	0.51
1:B:84:ASP:OD1	1:B:140:ARG:NH2	2.39	0.51
1:E:32:LEU:HB3	1:E:38:TYR:CE2	2.46	0.51
1:E:209:ALA:HA	1:E:220:PHE:HB2	1.92	0.51
1:G:153:GLU:O	1:G:159:SER:HB2	2.11	0.51
1:L:87:ILE:O	1:L:91:ILE:HG12	2.11	0.51
1:A:68:PRO:HB2	1:A:170:TYR:HD1	1.76	0.51
1:A:129:ASP:HB2	1:A:130:PRO:HD2	1.93	0.51
1:B:212:THR:HA	1:B:221:ASP:O	2.11	0.51
1:B:289:TYR:HA	1:B:298:MET:HB2	1.91	0.51
1:I:98:PHE:CE1	1:I:227:THR:HG23	2.45	0.51
1:L:129:ASP:HB2	1:L:130:PRO:HD2	1.93	0.51
1:N:289:TYR:HA	1:N:298:MET:HB2	1.91	0.51
1:D:109:PRO:HB2	1:D:111:VAL:HG12	1.92	0.51
1:F:99:ARG:NH1	1:F:118:ASP:HB2	2.26	0.51
1:H:305:GLN:O	1:H:309:VAL:HG23	2.11	0.51
1:I:27:GLU:HB2	1:I:150:THR:HG21	1.93	0.51
1:J:32:LEU:HD13	1:J:38:TYR:CZ	2.46	0.51
1:J:102:ASN:ND2	1:J:299:VAL:O	2.36	0.51
1:J:209:ALA:HA	1:J:220:PHE:HB2	1.93	0.51
1:L:27:GLU:HB2	1:L:150:THR:HG21	1.93	0.51
1:M:289:TYR:HA	1:M:298:MET:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:TYR:CE2	1:B:42:ARG:HG3	2.46	0.50
1:B:84:ASP:HA	1:B:87:ILE:HD12	1.93	0.50
1:D:130:PRO:HD3	1:D:165:TRP:CE3	2.46	0.50
1:H:287:MET:H	1:H:287:MET:CE	2.25	0.50
1:I:87:ILE:O	1:I:91:ILE:HG12	2.11	0.50
1:M:26:MET:HG3	1:M:242:ASP:CG	2.31	0.50
1:M:42:ARG:NH1	1:M:235:ASN:HD21	2.09	0.50
1:N:38:TYR:CE2	1:N:42:ARG:HG3	2.46	0.50
1:A:26:MET:HG3	1:A:242:ASP:CG	2.31	0.50
1:A:26:MET:HG3	1:A:242:ASP:OD1	2.12	0.50
1:B:27:GLU:HB2	1:B:150:THR:HG21	1.92	0.50
1:E:84:ASP:HA	1:E:87:ILE:HD12	1.93	0.50
1:H:209:ALA:HA	1:H:220:PHE:HB2	1.93	0.50
1:J:287:MET:H	1:J:287:MET:CE	2.24	0.50
1:K:201:ASP:OD1	1:K:317:ARG:NH2	2.32	0.50
1:K:209:ALA:HA	1:K:220:PHE:HB2	1.91	0.50
1:L:84:ASP:HA	1:L:87:ILE:HD12	1.93	0.50
1:A:130:PRO:HD3	1:A:165:TRP:CE3	2.46	0.50
1:A:155:LYS:HD3	1:A:200:THR:HA	1.94	0.50
1:E:27:GLU:HB2	1:E:150:THR:HG21	1.92	0.50
1:F:26:MET:HG3	1:F:242:ASP:CG	2.32	0.50
1:G:218:PRO:HA	1:G:220:PHE:N	2.26	0.50
1:H:153:GLU:O	1:H:159:SER:HB2	2.12	0.50
1:A:38:TYR:CE2	1:A:42:ARG:HG3	2.47	0.50
1:D:26:MET:HG3	1:D:242:ASP:CG	2.30	0.50
1:F:68:PRO:HB2	1:F:170:TYR:HD1	1.76	0.50
1:F:209:ALA:HA	1:F:220:PHE:HB2	1.93	0.50
1:G:35:SER:C	1:G:37:SER:H	2.15	0.50
1:L:38:TYR:CE2	1:L:42:ARG:HG3	2.46	0.50
1:M:209:ALA:HA	1:M:220:PHE:HB2	1.92	0.50
1:M:270:PRO:HG2	1:M:286:PHE:CD2	2.47	0.50
1:C:84:ASP:HA	1:C:87:ILE:HD12	1.94	0.50
1:F:84:ASP:HA	1:F:87:ILE:HD12	1.93	0.50
1:F:270:PRO:HG2	1:F:286:PHE:CG	2.47	0.50
1:H:61:ARG:HH12	1:H:121:ILE:HD11	1.75	0.50
1:I:155:LYS:HD3	1:I:200:THR:HA	1.94	0.50
1:I:189:GLY:HA3	1:I:205:ILE:HD13	1.94	0.50
1:J:32:LEU:HB3	1:J:38:TYR:CE2	2.46	0.50
1:J:61:ARG:HH12	1:J:121:ILE:HD11	1.77	0.50
1:K:309:VAL:O	1:K:313:LEU:HG	2.11	0.50
1:B:26:MET:HG3	1:B:242:ASP:CG	2.31	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:ARG:HH12	1:C:121:ILE:HD11	1.77	0.50
1:D:270:PRO:HG2	1:D:286:PHE:CD2	2.46	0.50
1:F:32:LEU:HD13	1:F:38:TYR:CZ	2.47	0.50
1:G:274:HIS:O	1:G:284:ASP:N	2.43	0.50
1:H:130:PRO:HD3	1:H:165:TRP:CE3	2.47	0.50
1:K:26:MET:HG3	1:K:242:ASP:CG	2.32	0.50
1:M:84:ASP:OD1	1:M:140:ARG:NH2	2.38	0.50
1:C:27:GLU:HB2	1:C:150:THR:HG21	1.92	0.50
1:E:32:LEU:HD13	1:E:38:TYR:CZ	2.47	0.50
1:G:75:TRP:CG	1:G:81:ASN:HB2	2.47	0.50
1:I:309:VAL:O	1:I:313:LEU:HG	2.12	0.50
1:J:99:ARG:HB2	1:J:101:LEU:HG	1.94	0.50
1:N:130:PRO:HD3	1:N:165:TRP:CE3	2.46	0.50
1:C:68:PRO:HB2	1:C:170:TYR:HD1	1.76	0.50
1:G:84:ASP:HA	1:G:87:ILE:HD12	1.94	0.50
1:N:26:MET:HG3	1:N:242:ASP:CG	2.32	0.50
1:N:68:PRO:HB2	1:N:170:TYR:HD1	1.77	0.50
1:A:87:ILE:O	1:A:91:ILE:HG12	2.12	0.50
1:C:32:LEU:HB3	1:C:38:TYR:CE2	2.47	0.50
1:E:153:GLU:O	1:E:159:SER:HB2	2.12	0.50
1:N:99:ARG:NH1	1:N:118:ASP:HB2	2.27	0.50
1:A:61:ARG:HH12	1:A:121:ILE:HD11	1.76	0.49
1:E:26:MET:HG3	1:E:242:ASP:CG	2.32	0.49
1:E:71:VAL:HA	1:E:173:ILE:HB	1.93	0.49
1:F:155:LYS:HD3	1:F:200:THR:HA	1.94	0.49
1:H:270:PRO:HG2	1:H:286:PHE:CD2	2.47	0.49
1:J:99:ARG:NH1	1:J:118:ASP:HB2	2.27	0.49
1:J:130:PRO:HD3	1:J:165:TRP:CE3	2.47	0.49
1:N:109:PRO:HD2	1:N:287:MET:SD	2.52	0.49
1:C:209:ALA:HA	1:C:220:PHE:HB2	1.93	0.49
1:E:287:MET:CE	1:E:287:MET:H	2.24	0.49
1:G:87:ILE:O	1:G:91:ILE:HG12	2.12	0.49
1:J:270:PRO:HG2	1:J:286:PHE:CD2	2.47	0.49
1:F:192:GLN:OE1	1:F:200:THR:OG1	2.21	0.49
1:F:309:VAL:HG12	1:F:313:LEU:HD21	1.95	0.49
1:G:26:MET:HG3	1:G:242:ASP:OD1	2.11	0.49
1:I:26:MET:HG3	1:I:242:ASP:CG	2.33	0.49
1:I:270:PRO:HG2	1:I:286:PHE:CD2	2.47	0.49
1:L:98:PHE:CE1	1:L:227:THR:HG23	2.47	0.49
1:M:32:LEU:HD13	1:M:38:TYR:CZ	2.47	0.49
1:A:99:ARG:NH1	1:A:118:ASP:HB2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:129:ASP:HB2	1:C:130:PRO:HD2	1.94	0.49
1:L:42:ARG:NH1	1:L:235:ASN:HD21	2.09	0.49
1:L:218:PRO:HA	1:L:220:PHE:N	2.27	0.49
1:E:274:HIS:O	1:E:284:ASP:N	2.41	0.49
1:H:38:TYR:CE2	1:H:42:ARG:HG3	2.47	0.49
1:K:26:MET:HG3	1:K:242:ASP:OD1	2.12	0.49
1:K:129:ASP:HB2	1:K:130:PRO:HD2	1.93	0.49
1:L:270:PRO:HG2	1:L:286:PHE:CD2	2.48	0.49
1:M:35:SER:C	1:M:37:SER:H	2.16	0.49
1:G:26:MET:HG3	1:G:242:ASP:CG	2.33	0.49
1:I:42:ARG:NH1	1:I:235:ASN:HD21	2.10	0.49
1:L:32:LEU:HD13	1:L:38:TYR:CZ	2.48	0.49
1:M:129:ASP:HB2	1:M:130:PRO:HD2	1.94	0.49
1:M:165:TRP:N	1:M:172:ASN:OD1	2.45	0.49
1:N:26:MET:HG3	1:N:242:ASP:OD1	2.13	0.49
1:H:26:MET:HG3	1:H:242:ASP:CG	2.32	0.49
1:M:287:MET:HE2	1:M:287:MET:H	1.76	0.49
1:N:305:GLN:O	1:N:309:VAL:HG23	2.13	0.49
1:E:94:LEU:O	1:E:98:PHE:HD1	1.96	0.49
1:E:314:ASP:O	1:E:318:SER:HB3	2.12	0.49
1:G:287:MET:CE	1:G:287:MET:H	2.25	0.49
1:H:27:GLU:HB2	1:H:150:THR:HG21	1.95	0.49
1:H:68:PRO:HB2	1:H:170:TYR:HD1	1.78	0.49
1:H:71:VAL:HA	1:H:173:ILE:HB	1.95	0.49
1:H:274:HIS:O	1:H:284:ASP:N	2.42	0.49
1:A:42:ARG:NH1	1:A:235:ASN:HD21	2.11	0.49
1:F:287:MET:H	1:F:287:MET:CE	2.25	0.49
1:G:38:TYR:CE2	1:G:42:ARG:HG3	2.48	0.49
1:H:94:LEU:O	1:H:98:PHE:HD1	1.96	0.49
1:K:99:ARG:NH1	1:K:118:ASP:HB2	2.26	0.49
1:K:99:ARG:HB2	1:K:101:LEU:HG	1.95	0.49
1:M:32:LEU:HB3	1:M:38:TYR:CE2	2.48	0.49
1:C:130:PRO:HD3	1:C:165:TRP:CE3	2.47	0.49
1:E:99:ARG:NH1	1:E:118:ASP:HB2	2.28	0.49
1:E:270:PRO:HG2	1:E:286:PHE:CD2	2.48	0.49
1:G:201:ASP:OD1	1:G:317:ARG:NH2	2.36	0.49
1:H:157:ALA:HA	1:H:161:GLY:O	2.12	0.49
1:I:32:LEU:HB3	1:I:38:TYR:CE2	2.48	0.49
1:L:130:PRO:HD3	1:L:165:TRP:CE3	2.47	0.49
1:L:309:VAL:O	1:L:313:LEU:HG	2.13	0.49
1:B:26:MET:HG3	1:B:242:ASP:OD1	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:LEU:HD13	1:B:38:TYR:CZ	2.48	0.48
1:B:157:ALA:HA	1:B:161:GLY:O	2.13	0.48
1:D:35:SER:C	1:D:37:SER:H	2.15	0.48
1:D:287:MET:H	1:D:287:MET:CE	2.26	0.48
1:F:218:PRO:HA	1:F:220:PHE:N	2.26	0.48
1:I:130:PRO:HD3	1:I:165:TRP:CE3	2.48	0.48
1:I:218:PRO:HA	1:I:220:PHE:N	2.27	0.48
1:K:314:ASP:O	1:K:318:SER:HB3	2.13	0.48
1:L:84:ASP:OD1	1:L:140:ARG:NH2	2.41	0.48
1:M:61:ARG:HH12	1:M:121:ILE:HD11	1.77	0.48
1:M:287:MET:H	1:M:287:MET:CE	2.25	0.48
1:A:309:VAL:O	1:A:313:LEU:HG	2.13	0.48
1:B:218:PRO:HA	1:B:220:PHE:N	2.27	0.48
1:C:26:MET:HG3	1:C:242:ASP:OD1	2.13	0.48
1:C:99:ARG:NH1	1:C:118:ASP:HB2	2.28	0.48
1:C:270:PRO:HG2	1:C:286:PHE:CD2	2.48	0.48
1:D:42:ARG:NH1	1:D:235:ASN:HD21	2.11	0.48
1:D:68:PRO:HB2	1:D:170:TYR:HD1	1.78	0.48
1:D:218:PRO:HA	1:D:220:PHE:N	2.27	0.48
1:E:26:MET:HG3	1:E:242:ASP:OD1	2.13	0.48
1:E:99:ARG:HB2	1:E:101:LEU:HG	1.95	0.48
1:F:87:ILE:O	1:F:91:ILE:HG12	2.13	0.48
1:G:305:GLN:O	1:G:309:VAL:HG23	2.12	0.48
1:I:314:ASP:O	1:I:318:SER:HB3	2.13	0.48
1:J:90:GLN:NE2	1:J:223:GLY:O	2.45	0.48
1:L:26:MET:HG3	1:L:242:ASP:OD1	2.13	0.48
1:F:27:GLU:HB2	1:F:150:THR:HG21	1.95	0.48
1:F:94:LEU:O	1:F:98:PHE:HD1	1.96	0.48
1:H:87:ILE:O	1:H:91:ILE:HG12	2.13	0.48
1:H:99:ARG:HB2	1:H:101:LEU:HG	1.95	0.48
1:I:35:SER:C	1:I:37:SER:H	2.17	0.48
1:I:99:ARG:HB2	1:I:101:LEU:HG	1.96	0.48
1:K:35:SER:C	1:K:37:SER:H	2.16	0.48
1:C:94:LEU:O	1:C:98:PHE:HD1	1.97	0.48
1:E:109:PRO:HD2	1:E:287:MET:SD	2.53	0.48
1:K:287:MET:H	1:K:287:MET:CE	2.26	0.48
1:N:35:SER:C	1:N:37:SER:H	2.17	0.48
1:A:287:MET:H	1:A:287:MET:CE	2.26	0.48
1:E:35:SER:C	1:E:37:SER:H	2.16	0.48
1:H:314:ASP:O	1:H:318:SER:HB3	2.13	0.48
1:I:75:TRP:CG	1:I:81:ASN:HB2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:218:PRO:HA	1:J:220:PHE:N	2.28	0.48
1:A:218:PRO:HA	1:A:220:PHE:N	2.28	0.48
1:B:42:ARG:NH1	1:B:235:ASN:HD21	2.10	0.48
1:B:314:ASP:O	1:B:318:SER:HB3	2.14	0.48
1:C:75:TRP:CG	1:C:81:ASN:HB2	2.48	0.48
1:D:99:ARG:HB2	1:D:101:LEU:HG	1.95	0.48
1:I:26:MET:HG3	1:I:242:ASP:OD1	2.13	0.48
1:I:75:TRP:HB2	1:I:77:GLU:O	2.13	0.48
1:I:99:ARG:NH1	1:I:118:ASP:HB2	2.28	0.48
1:N:32:LEU:HD13	1:N:38:TYR:CZ	2.49	0.48
1:B:75:TRP:CG	1:B:81:ASN:HB2	2.48	0.48
1:F:32:LEU:HB3	1:F:38:TYR:CE2	2.48	0.48
1:G:38:TYR:HD1	1:G:196:GLY:HA2	1.78	0.48
1:I:32:LEU:HD13	1:I:38:TYR:CZ	2.49	0.48
1:I:155:LYS:HA	1:I:174:TRP:NE1	2.29	0.48
1:J:35:SER:C	1:J:37:SER:H	2.16	0.48
1:A:32:LEU:HD13	1:A:38:TYR:CZ	2.48	0.48
1:B:35:SER:C	1:B:37:SER:H	2.17	0.48
1:H:32:LEU:HD13	1:H:38:TYR:CZ	2.49	0.48
1:J:153:GLU:CD	1:J:153:GLU:H	2.17	0.48
1:K:32:LEU:HD13	1:K:38:TYR:CZ	2.49	0.48
1:K:38:TYR:CE2	1:K:42:ARG:HG3	2.48	0.48
1:B:99:ARG:HB2	1:B:101:LEU:HG	1.96	0.48
1:C:35:SER:C	1:C:37:SER:H	2.17	0.48
1:D:61:ARG:HH12	1:D:121:ILE:HD11	1.79	0.48
1:E:61:ARG:HH12	1:E:121:ILE:HD11	1.78	0.48
1:E:155:LYS:HA	1:E:174:TRP:NE1	2.28	0.48
1:H:75:TRP:CG	1:H:81:ASN:HB2	2.49	0.48
1:J:129:ASP:HB2	1:J:130:PRO:HD2	1.96	0.48
1:A:75:TRP:HB2	1:A:77:GLU:O	2.14	0.48
1:C:26:MET:HG3	1:C:242:ASP:CG	2.35	0.48
1:F:38:TYR:CE2	1:F:42:ARG:HG3	2.49	0.48
1:I:38:TYR:CE2	1:I:42:ARG:HG3	2.49	0.48
1:J:75:TRP:CG	1:J:81:ASN:HB2	2.49	0.48
1:N:270:PRO:HG2	1:N:286:PHE:CD2	2.48	0.48
1:C:32:LEU:HD13	1:C:38:TYR:CZ	2.48	0.47
1:D:32:LEU:HD13	1:D:38:TYR:CZ	2.49	0.47
1:E:42:ARG:NH1	1:E:235:ASN:HD21	2.12	0.47
1:H:189:GLY:HA3	1:H:205:ILE:HD13	1.95	0.47
1:N:90:GLN:NE2	1:N:223:GLY:O	2.45	0.47
1:G:270:PRO:HG2	1:G:286:PHE:CD2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:90:GLN:NE2	1:H:223:GLY:O	2.46	0.47
1:J:192:GLN:OE1	1:J:200:THR:OG1	2.21	0.47
1:K:61:ARG:HH12	1:K:121:ILE:HD11	1.78	0.47
1:M:99:ARG:NH1	1:M:118:ASP:HB2	2.30	0.47
1:B:270:PRO:HG2	1:B:286:PHE:CD2	2.50	0.47
1:C:314:ASP:O	1:C:318:SER:HB3	2.15	0.47
1:D:90:GLN:NE2	1:D:223:GLY:O	2.46	0.47
1:G:27:GLU:HB2	1:G:150:THR:HG21	1.96	0.47
1:G:99:ARG:HB2	1:G:101:LEU:HG	1.97	0.47
1:G:99:ARG:NH1	1:G:118:ASP:HB2	2.28	0.47
1:H:218:PRO:HA	1:H:220:PHE:N	2.27	0.47
1:J:314:ASP:O	1:J:318:SER:HB3	2.14	0.47
1:K:305:GLN:O	1:K:309:VAL:HG23	2.15	0.47
1:N:94:LEU:O	1:N:98:PHE:HD1	1.97	0.47
1:N:218:PRO:HA	1:N:220:PHE:N	2.29	0.47
1:N:238:HIS:O	1:N:254:ASP:HB3	2.14	0.47
1:C:38:TYR:CE2	1:C:42:ARG:HG3	2.49	0.47
1:L:194:PRO:HD2	1:L:235:ASN:ND2	2.30	0.47
1:L:287:MET:CE	1:L:287:MET:H	2.26	0.47
1:M:38:TYR:CE2	1:M:42:ARG:HG3	2.49	0.47
1:N:42:ARG:NH1	1:N:235:ASN:HD21	2.11	0.47
1:N:157:ALA:HA	1:N:161:GLY:O	2.14	0.47
1:A:201:ASP:OD1	1:A:317:ARG:NH2	2.33	0.47
1:C:99:ARG:HB2	1:C:101:LEU:HG	1.96	0.47
1:H:26:MET:HG3	1:H:242:ASP:OD1	2.14	0.47
1:H:152:ASP:OD2	1:H:155:LYS:NZ	2.33	0.47
1:K:98:PHE:CE1	1:K:227:THR:HG23	2.50	0.47
1:B:189:GLY:HA3	1:B:205:ILE:HD13	1.96	0.47
1:C:98:PHE:CE1	1:C:227:THR:HG23	2.50	0.47
1:H:35:SER:C	1:H:37:SER:H	2.18	0.47
1:I:152:ASP:OD2	1:I:155:LYS:NZ	2.39	0.47
1:J:84:ASP:HA	1:J:87:ILE:HD12	1.96	0.47
1:M:94:LEU:O	1:M:98:PHE:HD1	1.97	0.47
1:A:189:GLY:HA3	1:A:205:ILE:HD13	1.97	0.47
1:A:270:PRO:HG2	1:A:286:PHE:CD2	2.50	0.47
1:B:99:ARG:NH1	1:B:118:ASP:HB2	2.30	0.47
1:B:192:GLN:OE1	1:B:200:THR:OG1	2.21	0.47
1:B:287:MET:CE	1:B:287:MET:H	2.27	0.47
1:C:201:ASP:OD1	1:C:317:ARG:NH2	2.35	0.47
1:D:75:TRP:CG	1:D:81:ASN:HB2	2.50	0.47
1:F:26:MET:HG3	1:F:242:ASP:OD1	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:65:VAL:HG21	1:F:321:LEU:HD22	1.97	0.47
1:F:75:TRP:HB2	1:F:77:GLU:O	2.14	0.47
1:G:32:LEU:HB3	1:G:38:TYR:CE2	2.49	0.47
1:I:192:GLN:OE1	1:I:200:THR:OG1	2.20	0.47
1:L:26:MET:HG3	1:L:242:ASP:CG	2.34	0.47
1:L:75:TRP:HB2	1:L:77:GLU:O	2.15	0.47
1:L:153:GLU:CD	1:L:153:GLU:H	2.18	0.47
1:M:189:GLY:HA3	1:M:205:ILE:HD13	1.97	0.47
1:N:194:PRO:HD2	1:N:235:ASN:ND2	2.30	0.47
1:A:32:LEU:HB3	1:A:38:TYR:CE2	2.50	0.47
1:A:212:THR:HA	1:A:221:ASP:O	2.14	0.47
1:A:278:SER:HB2	1:B:279:ASN:C	2.35	0.47
1:C:218:PRO:HA	1:C:220:PHE:N	2.29	0.47
1:L:238:HIS:O	1:L:254:ASP:HB3	2.15	0.47
1:M:309:VAL:HG12	1:M:313:LEU:HD21	1.96	0.47
1:N:314:ASP:O	1:N:318:SER:HB3	2.14	0.47
1:B:194:PRO:HD2	1:B:235:ASN:ND2	2.29	0.47
1:B:249:PRO:O	1:B:252:ARG:HG3	2.15	0.47
1:B:274:HIS:O	1:B:284:ASP:N	2.46	0.47
1:E:75:TRP:CG	1:E:81:ASN:HB2	2.50	0.47
1:E:192:GLN:OE1	1:E:200:THR:OG1	2.24	0.47
1:I:84:ASP:OD1	1:I:140:ARG:NH2	2.38	0.47
1:I:305:GLN:O	1:I:309:VAL:HG23	2.15	0.47
1:L:274:HIS:O	1:L:284:ASP:N	2.46	0.47
1:C:84:ASP:OD1	1:C:140:ARG:NH2	2.42	0.47
1:D:94:LEU:O	1:D:98:PHE:HD1	1.98	0.47
1:E:84:ASP:OD1	1:E:140:ARG:NH2	2.38	0.47
1:G:94:LEU:O	1:G:98:PHE:HD1	1.98	0.47
1:G:314:ASP:O	1:G:318:SER:HB3	2.14	0.47
1:K:84:ASP:OD1	1:K:140:ARG:NH2	2.41	0.47
1:L:68:PRO:HB2	1:L:170:TYR:HD1	1.80	0.47
1:L:75:TRP:CG	1:L:81:ASN:HB2	2.50	0.47
1:L:94:LEU:O	1:L:98:PHE:HD1	1.97	0.47
1:A:134:GLN:OE1	1:C:180:LYS:NZ	2.49	0.46
1:B:179:LEU:HD23	1:B:179:LEU:HA	1.78	0.46
1:C:42:ARG:NH1	1:C:235:ASN:HD21	2.12	0.46
1:D:309:VAL:HG12	1:D:313:LEU:HD21	1.98	0.46
1:D:314:ASP:O	1:D:318:SER:HB3	2.14	0.46
1:G:32:LEU:HD13	1:G:38:TYR:CZ	2.50	0.46
1:G:179:LEU:HD23	1:G:179:LEU:HA	1.80	0.46
1:I:220:PHE:HB3	1:I:225:THR:OG1	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:21:ARG:NH1	1:J:225:THR:HG23	2.30	0.46
1:J:98:PHE:CE1	1:J:227:THR:HG23	2.50	0.46
1:L:35:SER:C	1:L:37:SER:H	2.17	0.46
1:M:99:ARG:HB2	1:M:101:LEU:HG	1.97	0.46
1:N:155:LYS:HA	1:N:174:TRP:NE1	2.30	0.46
1:C:212:THR:HA	1:C:221:ASP:O	2.15	0.46
1:C:274:HIS:O	1:C:284:ASP:N	2.44	0.46
1:F:314:ASP:O	1:F:318:SER:HB3	2.15	0.46
1:I:212:THR:HA	1:I:221:ASP:O	2.15	0.46
1:K:32:LEU:HB3	1:K:38:TYR:CE2	2.50	0.46
1:K:94:LEU:O	1:K:98:PHE:HD1	1.99	0.46
1:M:98:PHE:CE1	1:M:227:THR:HG23	2.50	0.46
1:M:192:GLN:OE1	1:M:200:THR:OG1	2.24	0.46
1:D:99:ARG:NH1	1:D:118:ASP:HB2	2.30	0.46
1:E:38:TYR:CE2	1:E:42:ARG:HG3	2.50	0.46
1:E:218:PRO:HA	1:E:220:PHE:N	2.30	0.46
1:E:305:GLN:O	1:E:309:VAL:HG23	2.14	0.46
1:G:194:PRO:HD2	1:G:235:ASN:ND2	2.30	0.46
1:L:32:LEU:HB3	1:L:38:TYR:CE2	2.51	0.46
1:N:75:TRP:CG	1:N:81:ASN:HB2	2.50	0.46
1:B:238:HIS:O	1:B:254:ASP:HB3	2.15	0.46
1:C:109:PRO:HD2	1:C:287:MET:SD	2.55	0.46
1:L:189:GLY:HA3	1:L:205:ILE:HD13	1.97	0.46
1:M:274:HIS:O	1:M:284:ASP:N	2.44	0.46
1:N:230:ILE:O	1:N:233:TRP:N	2.46	0.46
1:N:274:HIS:O	1:N:284:ASP:N	2.46	0.46
1:C:21:ARG:NH1	1:C:225:THR:HG23	2.31	0.46
1:D:38:TYR:HD1	1:D:196:GLY:HA2	1.81	0.46
1:F:98:PHE:CE1	1:F:227:THR:HG23	2.51	0.46
1:F:274:HIS:O	1:F:284:ASP:N	2.47	0.46
1:G:212:THR:HA	1:G:221:ASP:O	2.16	0.46
1:I:179:LEU:HD23	1:I:179:LEU:HA	1.76	0.46
1:J:94:LEU:O	1:J:98:PHE:HD1	1.99	0.46
1:J:212:THR:HA	1:J:221:ASP:O	2.15	0.46
1:K:218:PRO:HA	1:K:220:PHE:N	2.28	0.46
1:M:157:ALA:HA	1:M:161:GLY:O	2.16	0.46
1:M:220:PHE:HB3	1:M:225:THR:OG1	2.15	0.46
1:C:134:GLN:OE1	1:E:180:LYS:NZ	2.48	0.46
1:E:21:ARG:NH1	1:E:225:THR:HG23	2.31	0.46
1:E:90:GLN:NE2	1:E:223:GLY:O	2.48	0.46
1:H:42:ARG:NH1	1:H:235:ASN:HD21	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:157:ALA:HA	1:I:161:GLY:O	2.16	0.46
1:J:26:MET:HG3	1:J:242:ASP:CG	2.35	0.46
1:J:109:PRO:HD2	1:J:287:MET:SD	2.56	0.46
1:N:65:VAL:HG21	1:N:321:LEU:HD22	1.98	0.46
1:B:90:GLN:NE2	1:B:223:GLY:O	2.46	0.46
1:D:212:THR:HA	1:D:221:ASP:O	2.15	0.46
1:D:238:HIS:O	1:D:254:ASP:HB3	2.16	0.46
1:J:179:LEU:HD23	1:J:179:LEU:HA	1.82	0.46
1:M:155:LYS:HA	1:M:174:TRP:NE1	2.31	0.46
1:M:314:ASP:O	1:M:318:SER:HB3	2.16	0.46
1:B:68:PRO:HB2	1:B:170:TYR:CD1	2.51	0.46
1:E:36:ALA:HA	1:E:39:VAL:HB	1.98	0.46
1:J:26:MET:HG3	1:J:242:ASP:OD1	2.16	0.46
1:J:305:GLN:O	1:J:309:VAL:HG23	2.14	0.46
1:L:155:LYS:HA	1:L:174:TRP:NE1	2.31	0.46
1:M:212:THR:HA	1:M:221:ASP:O	2.15	0.46
1:M:305:GLN:O	1:M:309:VAL:HG23	2.16	0.46
1:N:287:MET:CE	1:N:287:MET:H	2.28	0.46
1:A:194:PRO:HD2	1:A:235:ASN:ND2	2.31	0.46
1:A:287:MET:H	1:A:287:MET:HE2	1.79	0.46
1:A:305:GLN:O	1:A:309:VAL:HG23	2.15	0.46
1:C:287:MET:H	1:C:287:MET:CE	2.29	0.46
1:D:194:PRO:HD2	1:D:235:ASN:ND2	2.31	0.46
1:F:165:TRP:N	1:F:172:ASN:OD1	2.47	0.46
1:G:155:LYS:HA	1:G:174:TRP:NE1	2.31	0.46
1:H:75:TRP:HB2	1:H:77:GLU:O	2.16	0.46
1:K:270:PRO:HG2	1:K:286:PHE:CD2	2.50	0.46
1:K:274:HIS:O	1:K:284:ASP:N	2.47	0.46
1:L:99:ARG:HB2	1:L:101:LEU:HG	1.97	0.46
1:L:212:THR:HA	1:L:221:ASP:O	2.16	0.46
1:N:75:TRP:HB2	1:N:77:GLU:O	2.16	0.46
1:N:99:ARG:HB2	1:N:101:LEU:HG	1.97	0.46
1:C:75:TRP:HB2	1:C:77:GLU:O	2.15	0.46
1:C:192:GLN:OE1	1:C:200:THR:OG1	2.21	0.46
1:D:32:LEU:HB3	1:D:38:TYR:CE2	2.51	0.46
1:D:153:GLU:CD	1:D:153:GLU:H	2.19	0.46
1:E:249:PRO:O	1:E:252:ARG:HG3	2.16	0.46
1:F:35:SER:C	1:F:37:SER:H	2.18	0.46
1:F:75:TRP:CG	1:F:81:ASN:HB2	2.50	0.46
1:G:90:GLN:NE2	1:G:223:GLY:O	2.45	0.46
1:H:99:ARG:NH1	1:H:118:ASP:HB2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:230:ILE:O	1:J:233:TRP:N	2.45	0.46
1:M:249:PRO:O	1:M:252:ARG:HG3	2.16	0.46
1:N:32:LEU:HB3	1:N:38:TYR:CE2	2.51	0.46
1:A:35:SER:C	1:A:37:SER:H	2.19	0.45
1:B:32:LEU:HB3	1:B:38:TYR:CE2	2.51	0.45
1:B:305:GLN:O	1:B:309:VAL:HG23	2.15	0.45
1:C:189:GLY:HA3	1:C:205:ILE:HD13	1.97	0.45
1:C:322:ALA:O	1:E:34:ARG:NH1	2.50	0.45
1:H:32:LEU:HB3	1:H:38:TYR:CE2	2.50	0.45
1:H:212:THR:HA	1:H:221:ASP:O	2.16	0.45
1:I:287:MET:H	1:I:287:MET:CE	2.29	0.45
1:J:300:MET:HE2	1:J:300:MET:HB2	1.87	0.45
1:L:249:PRO:O	1:L:252:ARG:HG3	2.16	0.45
1:M:75:TRP:CG	1:M:81:ASN:HB2	2.52	0.45
1:A:38:TYR:HD1	1:A:196:GLY:HA2	1.81	0.45
1:B:38:TYR:HD1	1:B:196:GLY:HA2	1.81	0.45
1:D:157:ALA:HA	1:D:161:GLY:O	2.17	0.45
1:F:109:PRO:HD2	1:F:287:MET:SD	2.56	0.45
1:F:212:THR:HA	1:F:221:ASP:O	2.17	0.45
1:G:309:VAL:O	1:G:313:LEU:HG	2.16	0.45
1:H:98:PHE:CE1	1:H:227:THR:HG23	2.52	0.45
1:I:274:HIS:O	1:I:284:ASP:N	2.45	0.45
1:N:165:TRP:N	1:N:172:ASN:OD1	2.46	0.45
1:A:90:GLN:NE2	1:A:223:GLY:O	2.50	0.45
1:A:99:ARG:HB2	1:A:101:LEU:HG	1.97	0.45
1:A:155:LYS:HA	1:A:174:TRP:NE1	2.31	0.45
1:A:238:HIS:O	1:A:254:ASP:HB3	2.17	0.45
1:B:309:VAL:HG12	1:B:313:LEU:HD21	1.98	0.45
1:F:99:ARG:HB2	1:F:101:LEU:HG	1.98	0.45
1:G:165:TRP:N	1:G:172:ASN:OD1	2.45	0.45
1:H:34:ARG:NH1	1:J:322:ALA:O	2.49	0.45
1:K:21:ARG:NH1	1:K:225:THR:HG23	2.31	0.45
1:K:75:TRP:HB2	1:K:77:GLU:O	2.16	0.45
1:L:86:GLN:O	1:L:89:SER:OG	2.33	0.45
1:M:218:PRO:HA	1:M:220:PHE:N	2.29	0.45
1:B:94:LEU:O	1:B:98:PHE:HD1	1.99	0.45
1:D:249:PRO:O	1:D:252:ARG:HG3	2.17	0.45
1:G:86:GLN:O	1:G:89:SER:OG	2.32	0.45
1:K:38:TYR:HD1	1:K:196:GLY:HA2	1.82	0.45
1:K:153:GLU:CD	1:K:153:GLU:H	2.18	0.45
1:M:21:ARG:NH1	1:M:225:THR:HG23	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:153:GLU:CD	1:M:153:GLU:H	2.19	0.45
1:A:153:GLU:CD	1:A:153:GLU:H	2.19	0.45
1:A:220:PHE:HB3	1:A:225:THR:OG1	2.16	0.45
1:D:201:ASP:OD1	1:D:317:ARG:NH2	2.34	0.45
1:A:94:LEU:O	1:A:98:PHE:HD1	1.99	0.45
1:B:287:MET:H	1:B:287:MET:HE2	1.81	0.45
1:G:157:ALA:HA	1:G:161:GLY:O	2.17	0.45
1:H:34:ARG:HH22	1:J:323:ARG:C	2.20	0.45
1:H:270:PRO:HG2	1:H:286:PHE:CB	2.47	0.45
1:K:194:PRO:HD2	1:K:235:ASN:ND2	2.32	0.45
1:N:287:MET:H	1:N:287:MET:HE2	1.81	0.45
1:F:305:GLN:O	1:F:309:VAL:HG23	2.17	0.45
1:G:194:PRO:HA	1:G:317:ARG:NH2	2.32	0.45
1:I:68:PRO:HB2	1:I:170:TYR:CD1	2.52	0.45
1:J:309:VAL:HG12	1:J:313:LEU:HD21	1.99	0.45
1:L:109:PRO:HD2	1:L:287:MET:SD	2.56	0.45
1:M:68:PRO:HB2	1:M:170:TYR:CD1	2.51	0.45
1:B:21:ARG:NH1	1:B:225:THR:HG23	2.31	0.45
1:B:98:PHE:CE1	1:B:227:THR:HG23	2.52	0.45
1:B:220:PHE:HB3	1:B:225:THR:OG1	2.16	0.45
1:C:153:GLU:CD	1:C:153:GLU:H	2.19	0.45
1:C:181:SER:OG	1:C:183:ILE:HG12	2.16	0.45
1:E:75:TRP:HB2	1:E:77:GLU:O	2.16	0.45
1:E:322:ALA:O	1:G:34:ARG:NH1	2.50	0.45
1:I:94:LEU:O	1:I:98:PHE:HD1	2.00	0.45
1:J:189:GLY:HA3	1:J:205:ILE:HD13	1.99	0.45
1:J:220:PHE:HB3	1:J:225:THR:OG1	2.16	0.45
1:K:109:PRO:HD2	1:K:287:MET:SD	2.57	0.45
1:M:38:TYR:HD1	1:M:196:GLY:HA2	1.82	0.45
1:N:21:ARG:NH1	1:N:225:THR:HG23	2.32	0.45
1:N:38:TYR:HD1	1:N:196:GLY:HA2	1.80	0.45
1:B:199:GLU:H	1:B:199:GLU:HG3	1.57	0.45
1:C:194:PRO:HD2	1:C:235:ASN:ND2	2.32	0.45
1:C:220:PHE:HB3	1:C:225:THR:OG1	2.16	0.45
1:F:220:PHE:HB3	1:F:225:THR:OG1	2.17	0.45
1:G:220:PHE:HB3	1:G:225:THR:OG1	2.17	0.45
1:H:65:VAL:HG21	1:H:321:LEU:HD22	1.97	0.45
1:H:199:GLU:H	1:H:199:GLU:HG3	1.55	0.45
1:H:220:PHE:HB3	1:H:225:THR:OG1	2.16	0.45
1:H:238:HIS:O	1:H:254:ASP:HB3	2.17	0.45
1:K:189:GLY:HA3	1:K:205:ILE:HD13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:230:ILE:O	1:K:233:TRP:N	2.48	0.45
1:L:152:ASP:OD2	1:L:155:LYS:NZ	2.44	0.45
1:A:165:TRP:N	1:A:172:ASN:OD1	2.47	0.45
1:A:279:ASN:C	1:B:278:SER:HB2	2.37	0.45
1:B:230:ILE:O	1:B:233:TRP:N	2.48	0.45
1:E:323:ARG:C	1:G:34:ARG:HH22	2.19	0.45
1:F:36:ALA:HA	1:F:39:VAL:HB	1.99	0.45
1:G:189:GLY:HA3	1:G:205:ILE:HD13	1.99	0.45
1:G:199:GLU:H	1:G:199:GLU:HG3	1.56	0.45
1:H:194:PRO:HA	1:H:317:ARG:NH2	2.32	0.45
1:J:75:TRP:HB2	1:J:77:GLU:O	2.17	0.45
1:J:155:LYS:HA	1:J:174:TRP:NE1	2.31	0.45
1:K:75:TRP:CG	1:K:81:ASN:HB2	2.52	0.45
1:K:157:ALA:HA	1:K:161:GLY:O	2.17	0.45
1:L:230:ILE:O	1:L:233:TRP:N	2.48	0.45
1:C:90:GLN:NE2	1:C:223:GLY:O	2.50	0.44
1:F:189:GLY:HA3	1:F:205:ILE:HD13	1.99	0.44
1:G:109:PRO:HD2	1:G:287:MET:SD	2.57	0.44
1:J:42:ARG:NH1	1:J:235:ASN:HD21	2.15	0.44
1:N:249:PRO:O	1:N:252:ARG:HG3	2.16	0.44
1:F:21:ARG:HG3	1:F:188:LEU:HD12	1.99	0.44
1:F:270:PRO:HG2	1:F:286:PHE:CD2	2.53	0.44
1:I:153:GLU:CD	1:I:153:GLU:H	2.20	0.44
1:J:68:PRO:HB2	1:J:170:TYR:CD1	2.52	0.44
1:K:212:THR:HA	1:K:221:ASP:O	2.18	0.44
1:L:157:ALA:HA	1:L:161:GLY:O	2.16	0.44
1:L:194:PRO:HA	1:L:317:ARG:NH2	2.32	0.44
1:M:90:GLN:NE2	1:M:223:GLY:O	2.50	0.44
1:N:98:PHE:CE1	1:N:227:THR:HG23	2.51	0.44
1:C:38:TYR:HD1	1:C:196:GLY:HA2	1.82	0.44
1:C:129:ASP:HB3	1:C:135:THR:HB	2.00	0.44
1:C:165:TRP:N	1:C:172:ASN:OD1	2.50	0.44
1:F:194:PRO:HD2	1:F:235:ASN:ND2	2.33	0.44
1:G:98:PHE:CE1	1:G:227:THR:HG23	2.53	0.44
1:H:300:MET:HE2	1:H:300:MET:HB2	1.92	0.44
1:J:274:HIS:O	1:J:284:ASP:N	2.46	0.44
1:L:36:ALA:HA	1:L:39:VAL:HB	1.99	0.44
1:M:199:GLU:H	1:M:199:GLU:HG3	1.58	0.44
1:N:212:THR:HA	1:N:221:ASP:O	2.17	0.44
1:D:305:GLN:O	1:D:309:VAL:HG23	2.16	0.44
1:J:38:TYR:CE2	1:J:42:ARG:HG3	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:36:ALA:HA	1:N:39:VAL:HB	2.00	0.44
1:D:179:LEU:HD23	1:D:179:LEU:HA	1.80	0.44
1:D:269:CYS:HA	1:D:270:PRO:HD3	1.83	0.44
1:E:38:TYR:HD1	1:E:196:GLY:HA2	1.82	0.44
1:E:153:GLU:CD	1:E:153:GLU:H	2.19	0.44
1:H:155:LYS:HA	1:H:174:TRP:NE1	2.33	0.44
1:K:249:PRO:O	1:K:252:ARG:HG3	2.18	0.44
1:M:270:PRO:HG2	1:M:286:PHE:CB	2.48	0.44
1:C:323:ARG:C	1:E:34:ARG:HH22	2.21	0.44
1:D:181:SER:OG	1:D:183:ILE:HG12	2.18	0.44
1:D:220:PHE:HB3	1:D:225:THR:OG1	2.17	0.44
1:F:181:SER:OG	1:F:183:ILE:HG12	2.17	0.44
1:I:165:TRP:O	1:I:172:ASN:ND2	2.46	0.44
1:I:249:PRO:O	1:I:252:ARG:HG3	2.17	0.44
1:J:157:ALA:HA	1:J:161:GLY:O	2.17	0.44
1:L:201:ASP:OD1	1:L:317:ARG:NH2	2.35	0.44
1:A:21:ARG:NH1	1:A:225:THR:HG23	2.32	0.44
1:A:157:ALA:HA	1:A:161:GLY:O	2.17	0.44
1:B:65:VAL:HG21	1:B:321:LEU:HD22	2.00	0.44
1:C:309:VAL:HG12	1:C:313:LEU:HD21	1.98	0.44
1:D:98:PHE:CE1	1:D:227:THR:HG23	2.52	0.44
1:F:157:ALA:HA	1:F:161:GLY:O	2.17	0.44
1:H:102:ASN:OD1	1:H:102:ASN:N	2.51	0.44
1:K:90:GLN:NE2	1:K:223:GLY:O	2.50	0.44
1:L:192:GLN:OE1	1:L:200:THR:OG1	2.21	0.44
1:N:220:PHE:HB3	1:N:225:THR:OG1	2.17	0.44
1:A:98:PHE:CE1	1:A:227:THR:HG23	2.53	0.44
1:D:65:VAL:HG21	1:D:321:LEU:HD22	1.99	0.44
1:D:75:TRP:HB2	1:D:77:GLU:O	2.17	0.44
1:E:102:ASN:OD1	1:E:102:ASN:N	2.51	0.44
1:F:153:GLU:CD	1:F:153:GLU:H	2.21	0.44
1:K:220:PHE:HB3	1:K:225:THR:OG1	2.17	0.44
1:M:36:ALA:HA	1:M:39:VAL:HB	1.99	0.44
1:M:238:HIS:O	1:M:254:ASP:HB3	2.18	0.44
1:N:189:GLY:HA3	1:N:205:ILE:HD13	1.99	0.44
1:E:238:HIS:O	1:E:254:ASP:HB3	2.17	0.44
1:F:38:TYR:HD1	1:F:196:GLY:HA2	1.82	0.44
1:F:206:VAL:HG13	1:F:208:ALA:H	1.83	0.44
1:G:273:PRO:HB2	1:G:283:GLY:HA3	1.99	0.44
1:E:157:ALA:HA	1:E:161:GLY:O	2.17	0.43
1:E:212:THR:HA	1:E:221:ASP:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:21:ARG:NH1	1:F:225:THR:HG23	2.33	0.43
1:C:249:PRO:O	1:C:252:ARG:HG3	2.18	0.43
1:D:84:ASP:OD1	1:D:140:ARG:NH2	2.40	0.43
1:D:189:GLY:HA3	1:D:205:ILE:HD13	2.00	0.43
1:E:220:PHE:HB3	1:E:225:THR:OG1	2.18	0.43
1:E:309:VAL:HG12	1:E:313:LEU:HD21	1.99	0.43
1:I:38:TYR:HD1	1:I:196:GLY:HA2	1.83	0.43
1:J:206:VAL:HG13	1:J:208:ALA:H	1.83	0.43
1:L:21:ARG:NH1	1:L:225:THR:HG23	2.33	0.43
1:A:84:ASP:OD1	1:A:140:ARG:NH2	2.39	0.43
1:H:309:VAL:HG12	1:H:313:LEU:HD21	2.00	0.43
1:K:272:TYR:HA	1:K:273:PRO:HA	1.87	0.43
1:M:65:VAL:HG21	1:M:321:LEU:HD22	2.00	0.43
1:N:84:ASP:OD1	1:N:140:ARG:NH2	2.42	0.43
1:E:98:PHE:CE1	1:E:227:THR:HG23	2.53	0.43
1:I:65:VAL:HG21	1:I:321:LEU:HD22	2.00	0.43
1:I:269:CYS:HA	1:I:270:PRO:HD3	1.83	0.43
1:L:309:VAL:HG12	1:L:313:LEU:HD21	1.99	0.43
1:N:181:SER:OG	1:N:183:ILE:HG12	2.19	0.43
1:A:279:ASN:O	1:B:278:SER:HB2	2.19	0.43
1:B:270:PRO:HG2	1:B:286:PHE:CB	2.49	0.43
1:C:270:PRO:HG2	1:C:286:PHE:CB	2.48	0.43
1:E:108:VAL:CG2	1:E:300:MET:HG2	2.48	0.43
1:I:86:GLN:O	1:I:89:SER:OG	2.36	0.43
1:K:68:PRO:HB2	1:K:170:TYR:CD1	2.53	0.43
1:A:46:GLU:HB3	1:B:50:LEU:HD21	2.00	0.43
1:B:155:LYS:HA	1:B:174:TRP:NE1	2.33	0.43
1:D:109:PRO:HD2	1:D:287:MET:SD	2.58	0.43
1:F:102:ASN:N	1:F:102:ASN:OD1	2.51	0.43
1:G:129:ASP:HB3	1:G:135:THR:HB	2.01	0.43
1:G:249:PRO:O	1:G:252:ARG:HG3	2.18	0.43
1:I:109:PRO:HD2	1:I:287:MET:SD	2.59	0.43
1:I:249:PRO:HB2	1:I:263:ALA:CB	2.46	0.43
1:J:108:VAL:CG2	1:J:300:MET:HG2	2.49	0.43
1:J:165:TRP:N	1:J:172:ASN:OD1	2.45	0.43
1:K:224:ARG:HD2	1:K:289:TYR:CE1	2.40	0.43
1:A:75:TRP:CG	1:A:81:ASN:HB2	2.53	0.43
1:C:305:GLN:O	1:C:309:VAL:HG23	2.17	0.43
1:D:21:ARG:HG3	1:D:188:LEU:HD12	2.01	0.43
1:D:270:PRO:HG2	1:D:286:PHE:CB	2.49	0.43
1:F:269:CYS:HA	1:F:270:PRO:HD3	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:179:LEU:HD23	1:H:179:LEU:HA	1.78	0.43
1:I:90:GLN:NE2	1:I:223:GLY:O	2.48	0.43
1:K:155:LYS:HA	1:K:174:TRP:NE1	2.33	0.43
1:L:165:TRP:N	1:L:172:ASN:OD1	2.49	0.43
1:M:75:TRP:HB2	1:M:77:GLU:O	2.19	0.43
1:M:300:MET:HE2	1:M:300:MET:HB2	1.94	0.43
1:A:192:GLN:OE1	1:A:200:THR:OG1	2.22	0.43
1:C:65:VAL:HG21	1:C:321:LEU:HD22	2.01	0.43
1:G:65:VAL:HG21	1:G:321:LEU:HD22	2.00	0.43
1:G:68:PRO:HB2	1:G:170:TYR:CD1	2.53	0.43
1:G:75:TRP:HB2	1:G:77:GLU:O	2.19	0.43
1:I:309:VAL:HG12	1:I:313:LEU:HD21	1.99	0.43
1:J:180:LYS:HB3	1:J:180:LYS:HE2	1.76	0.43
1:K:36:ALA:HA	1:K:39:VAL:HB	2.01	0.43
1:M:194:PRO:HD2	1:M:235:ASN:ND2	2.34	0.43
1:N:102:ASN:N	1:N:102:ASN:OD1	2.52	0.43
1:N:199:GLU:H	1:N:199:GLU:HG3	1.55	0.43
1:A:199:GLU:H	1:A:199:GLU:HG3	1.59	0.43
1:B:36:ALA:HA	1:B:39:VAL:HB	2.01	0.43
1:C:287:MET:H	1:C:287:MET:HE2	1.83	0.43
1:E:116:ILE:HG12	1:E:300:MET:HE3	2.00	0.43
1:F:194:PRO:HA	1:F:317:ARG:NH2	2.34	0.43
1:H:21:ARG:NH1	1:H:225:THR:HG23	2.34	0.43
1:H:192:GLN:NE2	1:H:196:GLY:HA3	2.34	0.43
1:L:314:ASP:O	1:L:318:SER:HB3	2.18	0.43
1:N:68:PRO:HB2	1:N:170:TYR:CD1	2.54	0.43
1:B:165:TRP:N	1:B:172:ASN:OD1	2.49	0.43
1:D:102:ASN:OD1	1:D:102:ASN:N	2.52	0.43
1:E:270:PRO:HG2	1:E:286:PHE:CB	2.49	0.43
1:H:38:TYR:HD1	1:H:196:GLY:HA2	1.82	0.43
1:K:129:ASP:HB3	1:K:135:THR:HB	2.01	0.43
1:L:180:LYS:HB3	1:L:180:LYS:HE2	1.75	0.43
1:N:153:GLU:CD	1:N:153:GLU:H	2.21	0.43
1:N:270:PRO:HG2	1:N:286:PHE:CB	2.49	0.43
1:A:309:VAL:HG12	1:A:313:LEU:HD21	2.00	0.42
1:B:75:TRP:HB2	1:B:77:GLU:O	2.19	0.42
1:B:165:TRP:O	1:B:172:ASN:ND2	2.47	0.42
1:D:155:LYS:HA	1:D:174:TRP:NE1	2.34	0.42
1:D:274:HIS:O	1:D:284:ASP:N	2.47	0.42
1:D:180:LYS:HE2	1:D:180:LYS:HB3	1.75	0.42
1:E:194:PRO:HD2	1:E:235:ASN:ND2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:155:LYS:HA	1:F:174:TRP:NE1	2.33	0.42
1:H:272:TYR:HA	1:H:273:PRO:HA	1.88	0.42
1:I:194:PRO:HD2	1:I:235:ASN:ND2	2.35	0.42
1:K:95:ASN:O	1:K:101:LEU:HD12	2.19	0.42
1:M:108:VAL:CG2	1:M:300:MET:HG2	2.49	0.42
1:A:109:PRO:HD2	1:A:287:MET:SD	2.59	0.42
1:B:224:ARG:HD2	1:B:289:TYR:CE1	2.42	0.42
1:C:238:HIS:O	1:C:254:ASP:HB3	2.19	0.42
1:D:165:TRP:N	1:D:172:ASN:OD1	2.50	0.42
1:E:65:VAL:HG21	1:E:321:LEU:HD22	2.01	0.42
1:G:153:GLU:CD	1:G:153:GLU:H	2.21	0.42
1:I:270:PRO:HG2	1:I:286:PHE:CB	2.49	0.42
1:D:108:VAL:CG2	1:D:300:MET:HG2	2.49	0.42
1:F:249:PRO:O	1:F:252:ARG:HG3	2.19	0.42
1:H:84:ASP:OD1	1:H:140:ARG:NH2	2.42	0.42
1:H:153:GLU:CD	1:H:153:GLU:H	2.21	0.42
1:J:238:HIS:O	1:J:254:ASP:HB3	2.20	0.42
1:L:38:TYR:HD1	1:L:196:GLY:HA2	1.83	0.42
1:L:102:ASN:OD1	1:L:102:ASN:N	2.52	0.42
1:A:36:ALA:HA	1:A:39:VAL:HB	2.00	0.42
1:A:152:ASP:OD2	1:A:155:LYS:NZ	2.41	0.42
1:A:194:PRO:HA	1:A:317:ARG:NH2	2.34	0.42
1:C:245:ARG:HD3	1:C:245:ARG:N	2.33	0.42
1:D:36:ALA:HA	1:D:39:VAL:HB	2.00	0.42
1:E:165:TRP:N	1:E:172:ASN:OD1	2.46	0.42
1:F:68:PRO:HB2	1:F:170:TYR:CD1	2.54	0.42
1:G:21:ARG:NH1	1:G:225:THR:HG23	2.35	0.42
1:I:238:HIS:O	1:I:254:ASP:HB3	2.19	0.42
1:N:180:LYS:HB3	1:N:180:LYS:HE2	1.75	0.42
1:B:153:GLU:CD	1:B:153:GLU:H	2.21	0.42
1:B:269:CYS:HA	1:B:270:PRO:HD3	1.83	0.42
1:C:157:ALA:HA	1:C:161:GLY:O	2.19	0.42
1:E:68:PRO:HB2	1:E:170:TYR:CD1	2.53	0.42
1:F:29:HIS:O	1:F:33:LEU:HG	2.19	0.42
1:F:222:LYS:HB3	1:F:224:ARG:HG2	2.01	0.42
1:J:201:ASP:OD1	1:J:317:ARG:NH2	2.33	0.42
1:K:165:TRP:O	1:K:172:ASN:ND2	2.47	0.42
1:L:270:PRO:HG2	1:L:286:PHE:CB	2.50	0.42
1:M:181:SER:OG	1:M:183:ILE:HG12	2.19	0.42
1:A:230:ILE:O	1:A:233:TRP:N	2.47	0.42
1:C:68:PRO:HB2	1:C:170:TYR:CD1	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:287:MET:H	1:D:287:MET:HE2	1.85	0.42
1:G:238:HIS:O	1:G:254:ASP:HB3	2.20	0.42
1:I:29:HIS:O	1:I:33:LEU:HG	2.20	0.42
1:I:206:VAL:HG13	1:I:208:ALA:H	1.84	0.42
1:J:249:PRO:O	1:J:252:ARG:HG3	2.20	0.42
1:K:108:VAL:CG2	1:K:300:MET:HG2	2.50	0.42
1:L:220:PHE:HB3	1:L:225:THR:OG1	2.19	0.42
1:L:305:GLN:O	1:L:309:VAL:HG23	2.19	0.42
1:A:68:PRO:HB2	1:A:170:TYR:CD1	2.54	0.42
1:A:165:TRP:O	1:A:172:ASN:ND2	2.47	0.42
1:C:102:ASN:OD1	1:C:102:ASN:N	2.52	0.42
1:E:70:VAL:HG21	1:E:165:TRP:CE2	2.55	0.42
1:F:224:ARG:HD2	1:F:289:TYR:CE1	2.43	0.42
1:I:36:ALA:HA	1:I:39:VAL:HB	2.01	0.42
1:K:238:HIS:O	1:K:254:ASP:HB3	2.19	0.42
1:M:129:ASP:HB3	1:M:135:THR:HB	2.00	0.42
1:N:201:ASP:OD1	1:N:317:ARG:NH2	2.37	0.42
1:A:245:ARG:N	1:A:245:ARG:HD3	2.35	0.42
1:B:129:ASP:HB3	1:B:135:THR:HB	2.02	0.42
1:F:179:LEU:HD23	1:F:179:LEU:HA	1.76	0.42
1:I:21:ARG:NH1	1:I:225:THR:HG23	2.35	0.42
1:J:36:ALA:HA	1:J:39:VAL:HB	2.01	0.42
1:J:70:VAL:HG21	1:J:165:TRP:CE2	2.55	0.42
1:J:129:ASP:HB3	1:J:135:THR:HB	2.01	0.42
1:K:102:ASN:OD1	1:K:102:ASN:N	2.53	0.42
1:K:194:PRO:HA	1:K:317:ARG:NH2	2.35	0.42
1:N:129:ASP:HB3	1:N:135:THR:HB	2.02	0.42
1:N:222:LYS:HB3	1:N:224:ARG:HG2	2.02	0.42
1:N:300:MET:HE2	1:N:300:MET:HB2	1.93	0.42
1:A:274:HIS:O	1:A:284:ASP:N	2.46	0.42
1:C:36:ALA:HA	1:C:39:VAL:HB	2.02	0.42
1:G:180:LYS:HE2	1:G:180:LYS:HB3	1.76	0.42
1:G:270:PRO:HG2	1:G:286:PHE:CB	2.50	0.42
1:J:194:PRO:HA	1:J:317:ARG:NH2	2.35	0.42
1:J:270:PRO:HG2	1:J:286:PHE:CB	2.50	0.42
1:K:309:VAL:HG12	1:K:313:LEU:HD21	2.01	0.42
1:A:222:LYS:HB3	1:A:224:ARG:HG2	2.02	0.41
1:B:164:ALA:HB2	1:B:174:TRP:HZ2	1.84	0.41
1:D:21:ARG:NH1	1:D:225:THR:HG23	2.35	0.41
1:E:32:LEU:HD22	1:E:38:TYR:CE1	2.55	0.41
1:I:194:PRO:HA	1:I:317:ARG:NH2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:102:ASN:OD1	1:J:102:ASN:N	2.53	0.41
1:J:181:SER:OG	1:J:183:ILE:HG12	2.20	0.41
1:N:164:ALA:HB2	1:N:174:TRP:HZ2	1.84	0.41
1:A:108:VAL:HG12	1:A:109:PRO:O	2.21	0.41
1:A:129:ASP:HB3	1:A:135:THR:HB	2.02	0.41
1:A:300:MET:HE2	1:A:300:MET:HB2	1.95	0.41
1:B:109:PRO:HD2	1:B:287:MET:SD	2.60	0.41
1:C:155:LYS:HA	1:C:174:TRP:NE1	2.36	0.41
1:C:269:CYS:HA	1:C:270:PRO:HD3	1.84	0.41
1:F:90:GLN:HE22	1:F:223:GLY:C	2.23	0.41
1:F:230:ILE:O	1:F:233:TRP:N	2.51	0.41
1:G:30:HIS:HB2	1:G:244:LEU:CD1	2.50	0.41
1:G:108:VAL:HG12	1:G:109:PRO:O	2.20	0.41
1:J:32:LEU:HB3	1:J:38:TYR:CD2	2.56	0.41
1:L:245:ARG:HD3	1:L:245:ARG:N	2.36	0.41
1:A:50:LEU:HD11	1:A:54:GLN:OE1	2.21	0.41
1:A:206:VAL:HG13	1:A:208:ALA:H	1.86	0.41
1:I:129:ASP:HB3	1:I:135:THR:HB	2.03	0.41
1:I:165:TRP:N	1:I:172:ASN:OD1	2.49	0.41
1:L:129:ASP:HB3	1:L:135:THR:HB	2.01	0.41
1:L:273:PRO:HB2	1:L:283:GLY:HA3	2.03	0.41
1:B:108:VAL:HG12	1:B:109:PRO:O	2.20	0.41
1:C:70:VAL:HG21	1:C:165:TRP:CE2	2.56	0.41
1:H:36:ALA:HA	1:H:39:VAL:HB	2.02	0.41
1:H:109:PRO:HD2	1:H:287:MET:SD	2.60	0.41
1:I:300:MET:HE2	1:I:300:MET:HB2	1.99	0.41
1:K:70:VAL:HG21	1:K:165:TRP:CE2	2.56	0.41
1:L:65:VAL:HG21	1:L:321:LEU:HD22	2.02	0.41
1:N:108:VAL:CG2	1:N:300:MET:HG2	2.50	0.41
1:B:108:VAL:CG2	1:B:300:MET:HG2	2.51	0.41
1:B:249:PRO:HA	1:B:252:ARG:NE	2.35	0.41
1:G:50:LEU:HD21	1:H:46:GLU:HB3	2.02	0.41
1:H:108:VAL:CG2	1:H:300:MET:HG2	2.50	0.41
1:H:222:LYS:HB3	1:H:224:ARG:HG2	2.01	0.41
1:I:249:PRO:HA	1:I:252:ARG:NE	2.36	0.41
1:J:29:HIS:O	1:J:33:LEU:HG	2.21	0.41
1:J:165:TRP:O	1:J:172:ASN:ND2	2.44	0.41
1:K:65:VAL:HG21	1:K:321:LEU:HD22	2.02	0.41
1:D:273:PRO:HB2	1:D:283:GLY:HA3	2.02	0.41
1:H:129:ASP:HB3	1:H:135:THR:HB	2.03	0.41
1:H:194:PRO:HD2	1:H:235:ASN:ND2	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:38:TYR:HD1	1:J:196:GLY:HA2	1.83	0.41
1:A:270:PRO:HG2	1:A:286:PHE:CB	2.50	0.41
1:B:102:ASN:OD1	1:B:102:ASN:N	2.53	0.41
1:B:273:PRO:HB2	1:B:283:GLY:HA3	2.02	0.41
1:F:32:LEU:HB3	1:F:38:TYR:CD2	2.56	0.41
1:F:129:ASP:HB3	1:F:135:THR:HB	2.03	0.41
1:I:70:VAL:HG21	1:I:165:TRP:CE2	2.56	0.41
1:I:201:ASP:OD1	1:I:317:ARG:NH2	2.35	0.41
1:K:165:TRP:N	1:K:172:ASN:OD1	2.50	0.41
1:L:249:PRO:HA	1:L:252:ARG:NE	2.36	0.41
1:M:179:LEU:HA	1:M:179:LEU:HD23	1.79	0.41
1:N:249:PRO:HA	1:N:252:ARG:NE	2.36	0.41
1:A:181:SER:OG	1:A:183:ILE:HG12	2.20	0.41
1:D:129:ASP:HB3	1:D:135:THR:HB	2.02	0.41
1:D:313:LEU:HA	1:D:317:ARG:HB2	2.03	0.41
1:H:68:PRO:HB2	1:H:170:TYR:CD1	2.55	0.41
1:H:206:VAL:HG13	1:H:208:ALA:H	1.85	0.41
1:L:181:SER:OG	1:L:183:ILE:HG12	2.20	0.41
1:M:269:CYS:HA	1:M:270:PRO:HD3	1.84	0.41
1:N:273:PRO:HB2	1:N:283:GLY:HA3	2.03	0.41
1:A:65:VAL:HG21	1:A:321:LEU:HD22	2.02	0.41
1:A:86:GLN:O	1:A:89:SER:OG	2.39	0.41
1:A:278:SER:HB2	1:B:279:ASN:O	2.21	0.41
1:B:180:LYS:HB3	1:B:180:LYS:HE2	1.73	0.41
1:C:95:ASN:O	1:C:101:LEU:HD12	2.21	0.41
1:C:194:PRO:HG2	1:C:235:ASN:HB2	2.03	0.41
1:C:206:VAL:HG13	1:C:208:ALA:H	1.86	0.41
1:C:272:TYR:HA	1:C:273:PRO:HA	1.89	0.41
1:E:189:GLY:HA3	1:E:205:ILE:HD13	2.02	0.41
1:E:249:PRO:HB2	1:E:263:ALA:CB	2.48	0.41
1:F:86:GLN:O	1:F:89:SER:OG	2.38	0.41
1:F:180:LYS:HB3	1:F:180:LYS:HE2	1.76	0.41
1:F:227:THR:O	1:F:230:ILE:HB	2.21	0.41
1:F:270:PRO:HG2	1:F:286:PHE:CB	2.51	0.41
1:F:272:TYR:HA	1:F:286:PHE:HZ	1.86	0.41
1:G:36:ALA:HA	1:G:39:VAL:HB	2.02	0.41
1:G:102:ASN:N	1:G:102:ASN:OD1	2.53	0.41
1:G:165:TRP:O	1:G:172:ASN:ND2	2.49	0.41
1:H:180:LYS:HB3	1:H:180:LYS:HE2	1.78	0.41
1:H:273:PRO:HB2	1:H:283:GLY:HA3	2.03	0.41
1:I:313:LEU:HA	1:I:317:ARG:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:222:LYS:HB3	1:L:224:ARG:HG2	2.03	0.41
1:L:287:MET:H	1:L:287:MET:HE2	1.86	0.41
1:M:29:HIS:O	1:M:33:LEU:HG	2.21	0.41
1:N:152:ASP:OD2	1:N:155:LYS:NZ	2.42	0.41
1:N:245:ARG:HD3	1:N:245:ARG:N	2.36	0.41
1:A:257:ASP:HB3	1:B:251:SER:HA	2.02	0.41
1:D:68:PRO:HB3	1:D:126:ALA:HB2	2.03	0.41
1:D:90:GLN:HE22	1:D:223:GLY:C	2.25	0.41
1:E:129:ASP:HB3	1:E:135:THR:HB	2.03	0.41
1:F:249:PRO:HB2	1:F:263:ALA:CB	2.46	0.41
1:I:180:LYS:HE2	1:I:180:LYS:HB3	1.74	0.41
1:I:222:LYS:HB3	1:I:224:ARG:HG2	2.03	0.41
1:M:50:LEU:HD21	1:N:46:GLU:HB3	2.01	0.41
1:A:95:ASN:O	1:A:101:LEU:HD12	2.21	0.40
1:C:29:HIS:O	1:C:33:LEU:HG	2.21	0.40
1:D:61:ARG:O	1:D:61:ARG:HG2	2.21	0.40
1:D:224:ARG:HD2	1:D:289:TYR:CE1	2.44	0.40
1:F:300:MET:HE2	1:F:300:MET:HB2	1.96	0.40
1:H:29:HIS:O	1:H:33:LEU:HG	2.20	0.40
1:H:76:ASN:HB3	1:H:80:GLU:OE2	2.21	0.40
1:H:230:ILE:O	1:H:233:TRP:N	2.51	0.40
1:J:32:LEU:HD22	1:J:38:TYR:CE1	2.56	0.40
1:K:29:HIS:O	1:K:33:LEU:HG	2.22	0.40
1:L:194:PRO:HD2	1:L:235:ASN:HD22	1.87	0.40
1:M:165:TRP:O	1:M:172:ASN:ND2	2.48	0.40
1:N:194:PRO:HA	1:N:317:ARG:NH2	2.36	0.40
1:B:181:SER:OG	1:B:183:ILE:HG12	2.22	0.40
1:E:30:HIS:HB2	1:E:244:LEU:CD1	2.51	0.40
1:F:108:VAL:CG2	1:F:300:MET:HG2	2.50	0.40
1:G:269:CYS:HA	1:G:270:PRO:HD3	1.84	0.40
1:K:181:SER:OG	1:K:183:ILE:HG12	2.20	0.40
1:K:313:LEU:HA	1:K:317:ARG:HB2	2.03	0.40
1:M:21:ARG:HG3	1:M:188:LEU:HD12	2.02	0.40
1:M:102:ASN:N	1:M:102:ASN:OD1	2.52	0.40
1:B:246:PHE:HB2	1:D:64:ILE:HG22	2.03	0.40
1:B:272:TYR:HA	1:B:273:PRO:HA	1.88	0.40
1:C:32:LEU:HB3	1:C:38:TYR:CD2	2.56	0.40
1:C:222:LYS:HB3	1:C:224:ARG:HG2	2.03	0.40
1:J:249:PRO:HB2	1:J:263:ALA:CB	2.48	0.40
1:K:64:ILE:HG22	1:M:246:PHE:HB2	2.03	0.40
1:K:270:PRO:HG2	1:K:286:PHE:CB	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:109:PRO:HD2	1:M:287:MET:SD	2.62	0.40
1:N:164:ALA:HB2	1:N:174:TRP:CZ2	2.57	0.40
1:A:70:VAL:HG21	1:A:165:TRP:CE2	2.57	0.40
1:A:249:PRO:O	1:A:252:ARG:HG3	2.21	0.40
1:A:322:ALA:O	1:C:34:ARG:NH1	2.54	0.40
1:C:180:LYS:HE2	1:C:180:LYS:HB3	1.77	0.40
1:E:95:ASN:O	1:E:101:LEU:HD12	2.21	0.40
1:E:287:MET:H	1:E:287:MET:HE2	1.85	0.40
1:F:313:LEU:HA	1:F:317:ARG:HB2	2.04	0.40
1:H:313:LEU:HA	1:H:317:ARG:HB2	2.04	0.40
1:I:32:LEU:HB3	1:I:38:TYR:CD2	2.57	0.40
1:J:164:ALA:HB2	1:J:174:TRP:HZ2	1.86	0.40
1:J:245:ARG:N	1:J:245:ARG:HD3	2.37	0.40
1:M:249:PRO:HA	1:M:252:ARG:NE	2.35	0.40
1:E:32:LEU:HB3	1:E:38:TYR:CD2	2.56	0.40
1:E:165:TRP:O	1:E:172:ASN:ND2	2.49	0.40
1:F:245:ARG:HD3	1:F:245:ARG:N	2.36	0.40
1:F:273:PRO:HB2	1:F:283:GLY:HA3	2.03	0.40
1:G:127:THR:HG22	1:G:134:GLN:HG3	2.04	0.40
1:K:179:LEU:HD23	1:K:179:LEU:HA	1.80	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:SER:OG	1:B:113:SER:OG[4_566]	2.18	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	304/360 (84%)	286 (94%)	18 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	304/360 (84%)	287 (94%)	17 (6%)	0	100	100
1	C	304/360 (84%)	289 (95%)	15 (5%)	0	100	100
1	D	304/360 (84%)	289 (95%)	15 (5%)	0	100	100
1	E	304/360 (84%)	286 (94%)	18 (6%)	0	100	100
1	F	304/360 (84%)	287 (94%)	17 (6%)	0	100	100
1	G	304/360 (84%)	288 (95%)	16 (5%)	0	100	100
1	H	304/360 (84%)	287 (94%)	17 (6%)	0	100	100
1	I	304/360 (84%)	287 (94%)	17 (6%)	0	100	100
1	J	304/360 (84%)	289 (95%)	15 (5%)	0	100	100
1	K	304/360 (84%)	288 (95%)	16 (5%)	0	100	100
1	L	304/360 (84%)	288 (95%)	16 (5%)	0	100	100
1	M	304/360 (84%)	288 (95%)	16 (5%)	0	100	100
1	N	304/360 (84%)	287 (94%)	17 (6%)	0	100	100
All	All	4256/5040 (84%)	4026 (95%)	230 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	262/310 (84%)	233 (89%)	29 (11%)	6	25
1	B	262/310 (84%)	232 (88%)	30 (12%)	5	23
1	C	262/310 (84%)	234 (89%)	28 (11%)	6	26
1	D	262/310 (84%)	233 (89%)	29 (11%)	6	25
1	E	262/310 (84%)	233 (89%)	29 (11%)	6	25
1	F	262/310 (84%)	231 (88%)	31 (12%)	5	23
1	G	262/310 (84%)	232 (88%)	30 (12%)	5	23
1	H	262/310 (84%)	234 (89%)	28 (11%)	6	26

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	I	262/310 (84%)	233 (89%)	29 (11%)	6	25
1	J	262/310 (84%)	234 (89%)	28 (11%)	6	26
1	K	262/310 (84%)	232 (88%)	30 (12%)	5	23
1	L	262/310 (84%)	233 (89%)	29 (11%)	6	25
1	M	262/310 (84%)	230 (88%)	32 (12%)	5	22
1	N	262/310 (84%)	232 (88%)	30 (12%)	5	23
All	All	3668/4340 (84%)	3256 (89%)	412 (11%)	6	24

All (412) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	HIS
1	A	37	SER
1	A	54	GLN
1	A	59	ILE
1	A	61	ARG
1	A	63	GLU
1	A	64	ILE
1	A	69	VAL
1	A	72	HIS
1	A	84	ASP
1	A	93	ILE
1	A	105	VAL
1	A	123	PHE
1	A	125	LEU
1	A	128	LYS
1	A	141	THR
1	A	145	VAL
1	A	149	THR
1	A	150	THR
1	A	153	GLU
1	A	156	PHE
1	A	192	GLN
1	A	220	PHE
1	A	239	ILE
1	A	247	GLU
1	A	248	ASP
1	A	267	PHE
1	A	286	PHE
1	A	287	MET

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Mol	Chain	Res	Type
1	B	30	HIS
1	B	37	SER
1	B	54	GLN
1	B	59	ILE
1	B	61	ARG
1	B	63	GLU
1	B	64	ILE
1	B	69	VAL
1	B	72	HIS
1	B	78	GLU
1	B	84	ASP
1	B	93	ILE
1	B	105	VAL
1	B	123	PHE
1	B	125	LEU
1	B	128	LYS
1	B	141	THR
1	B	145	VAL
1	B	149	THR
1	B	150	THR
1	B	153	GLU
1	B	156	PHE
1	B	192	GLN
1	B	220	PHE
1	B	239	ILE
1	B	247	GLU
1	B	248	ASP
1	B	267	PHE
1	B	286	PHE
1	B	287	MET
1	C	30	HIS
1	C	37	SER
1	C	54	GLN
1	C	59	ILE
1	C	61	ARG
1	C	63	GLU
1	C	64	ILE
1	C	72	HIS
1	C	84	ASP
1	C	93	ILE
1	C	105	VAL
1	C	123	PHE

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Mol	Chain	Res	Type
1	C	125	LEU
1	C	128	LYS
1	C	141	THR
1	C	145	VAL
1	C	149	THR
1	C	150	THR
1	C	153	GLU
1	C	156	PHE
1	C	192	GLN
1	C	220	PHE
1	C	239	ILE
1	C	247	GLU
1	C	248	ASP
1	C	267	PHE
1	C	286	PHE
1	C	287	MET
1	D	30	HIS
1	D	37	SER
1	D	54	GLN
1	D	59	ILE
1	D	61	ARG
1	D	63	GLU
1	D	64	ILE
1	D	69	VAL
1	D	72	HIS
1	D	84	ASP
1	D	93	ILE
1	D	105	VAL
1	D	123	PHE
1	D	125	LEU
1	D	128	LYS
1	D	141	THR
1	D	145	VAL
1	D	149	THR
1	D	150	THR
1	D	153	GLU
1	D	156	PHE
1	D	192	GLN
1	D	220	PHE
1	D	239	ILE
1	D	247	GLU
1	D	248	ASP

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Mol	Chain	Res	Type
1	D	267	PHE
1	D	286	PHE
1	D	287	MET
1	E	30	HIS
1	E	37	SER
1	E	54	GLN
1	E	59	ILE
1	E	61	ARG
1	E	63	GLU
1	E	64	ILE
1	E	69	VAL
1	E	72	HIS
1	E	84	ASP
1	E	93	ILE
1	E	105	VAL
1	E	123	PHE
1	E	125	LEU
1	E	128	LYS
1	E	141	THR
1	E	145	VAL
1	E	149	THR
1	E	150	THR
1	E	153	GLU
1	E	156	PHE
1	E	192	GLN
1	E	220	PHE
1	E	239	ILE
1	E	247	GLU
1	E	248	ASP
1	E	267	PHE
1	E	286	PHE
1	E	287	MET
1	F	30	HIS
1	F	37	SER
1	F	54	GLN
1	F	59	ILE
1	F	61	ARG
1	F	63	GLU
1	F	64	ILE
1	F	69	VAL
1	F	72	HIS
1	F	84	ASP

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Mol	Chain	Res	Type
1	F	93	ILE
1	F	105	VAL
1	F	106	SER
1	F	123	PHE
1	F	125	LEU
1	F	128	LYS
1	F	141	THR
1	F	145	VAL
1	F	149	THR
1	F	150	THR
1	F	153	GLU
1	F	156	PHE
1	F	169	ARG
1	F	192	GLN
1	F	220	PHE
1	F	239	ILE
1	F	247	GLU
1	F	248	ASP
1	F	267	PHE
1	F	286	PHE
1	F	287	MET
1	G	30	HIS
1	G	37	SER
1	G	54	GLN
1	G	59	ILE
1	G	61	ARG
1	G	63	GLU
1	G	64	ILE
1	G	69	VAL
1	G	72	HIS
1	G	84	ASP
1	G	93	ILE
1	G	105	VAL
1	G	106	SER
1	G	123	PHE
1	G	125	LEU
1	G	128	LYS
1	G	141	THR
1	G	145	VAL
1	G	149	THR
1	G	150	THR
1	G	153	GLU

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Mol	Chain	Res	Type
1	G	156	PHE
1	G	192	GLN
1	G	220	PHE
1	G	239	ILE
1	G	247	GLU
1	G	248	ASP
1	G	267	PHE
1	G	286	PHE
1	G	287	MET
1	H	30	HIS
1	H	37	SER
1	H	54	GLN
1	H	59	ILE
1	H	61	ARG
1	H	63	GLU
1	H	64	ILE
1	H	72	HIS
1	H	84	ASP
1	H	93	ILE
1	H	105	VAL
1	H	123	PHE
1	H	125	LEU
1	H	128	LYS
1	H	141	THR
1	H	145	VAL
1	H	149	THR
1	H	150	THR
1	H	153	GLU
1	H	156	PHE
1	H	192	GLN
1	H	220	PHE
1	H	239	ILE
1	H	247	GLU
1	H	248	ASP
1	H	267	PHE
1	H	286	PHE
1	H	287	MET
1	I	30	HIS
1	I	37	SER
1	I	54	GLN
1	I	59	ILE
1	I	61	ARG

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Mol	Chain	Res	Type
1	I	63	GLU
1	I	64	ILE
1	I	69	VAL
1	I	72	HIS
1	I	84	ASP
1	I	93	ILE
1	I	105	VAL
1	I	123	PHE
1	I	125	LEU
1	I	128	LYS
1	I	141	THR
1	I	145	VAL
1	I	149	THR
1	I	150	THR
1	I	153	GLU
1	I	156	PHE
1	I	192	GLN
1	I	220	PHE
1	I	239	ILE
1	I	247	GLU
1	I	248	ASP
1	I	267	PHE
1	I	286	PHE
1	I	287	MET
1	J	30	HIS
1	J	37	SER
1	J	54	GLN
1	J	59	ILE
1	J	61	ARG
1	J	63	GLU
1	J	64	ILE
1	J	72	HIS
1	J	84	ASP
1	J	93	ILE
1	J	105	VAL
1	J	123	PHE
1	J	125	LEU
1	J	128	LYS
1	J	141	THR
1	J	145	VAL
1	J	149	THR
1	J	150	THR

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Mol	Chain	Res	Type
1	J	153	GLU
1	J	156	PHE
1	J	192	GLN
1	J	220	PHE
1	J	239	ILE
1	J	247	GLU
1	J	248	ASP
1	J	267	PHE
1	J	286	PHE
1	J	287	MET
1	K	30	HIS
1	K	37	SER
1	K	54	GLN
1	K	59	ILE
1	K	61	ARG
1	K	63	GLU
1	K	64	ILE
1	K	69	VAL
1	K	72	HIS
1	K	84	ASP
1	K	93	ILE
1	K	105	VAL
1	K	106	SER
1	K	123	PHE
1	K	125	LEU
1	K	128	LYS
1	K	141	THR
1	K	145	VAL
1	K	149	THR
1	K	150	THR
1	K	153	GLU
1	K	156	PHE
1	K	192	GLN
1	K	220	PHE
1	K	239	ILE
1	K	247	GLU
1	K	248	ASP
1	K	267	PHE
1	K	286	PHE
1	K	287	MET
1	L	30	HIS
1	L	37	SER

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Mol	Chain	Res	Type
1	L	54	GLN
1	L	59	ILE
1	L	61	ARG
1	L	63	GLU
1	L	64	ILE
1	L	69	VAL
1	L	72	HIS
1	L	84	ASP
1	L	93	ILE
1	L	105	VAL
1	L	123	PHE
1	L	125	LEU
1	L	128	LYS
1	L	141	THR
1	L	145	VAL
1	L	149	THR
1	L	150	THR
1	L	153	GLU
1	L	156	PHE
1	L	192	GLN
1	L	220	PHE
1	L	239	ILE
1	L	247	GLU
1	L	248	ASP
1	L	267	PHE
1	L	286	PHE
1	L	287	MET
1	M	30	HIS
1	M	37	SER
1	M	54	GLN
1	M	59	ILE
1	M	61	ARG
1	M	63	GLU
1	M	64	ILE
1	M	69	VAL
1	M	72	HIS
1	M	84	ASP
1	M	93	ILE
1	M	105	VAL
1	M	123	PHE
1	M	125	LEU
1	M	128	LYS

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Mol	Chain	Res	Type
1	M	141	THR
1	M	145	VAL
1	M	149	THR
1	M	150	THR
1	M	153	GLU
1	M	156	PHE
1	M	169	ARG
1	M	192	GLN
1	M	220	PHE
1	M	239	ILE
1	M	247	GLU
1	M	248	ASP
1	M	252	ARG
1	M	267	PHE
1	M	270	PRO
1	M	286	PHE
1	M	287	MET
1	N	30	HIS
1	N	37	SER
1	N	54	GLN
1	N	59	ILE
1	N	61	ARG
1	N	63	GLU
1	N	64	ILE
1	N	69	VAL
1	N	72	HIS
1	N	84	ASP
1	N	93	ILE
1	N	105	VAL
1	N	123	PHE
1	N	125	LEU
1	N	128	LYS
1	N	141	THR
1	N	145	VAL
1	N	149	THR
1	N	150	THR
1	N	153	GLU
1	N	156	PHE
1	N	192	GLN
1	N	220	PHE
1	N	239	ILE
1	N	247	GLU

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Mol	Chain	Res	Type
1	N	248	ASP
1	N	259	THR
1	N	267	PHE
1	N	286	PHE
1	N	287	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	HIS
1	A	72	HIS
1	A	235	ASN
1	B	29	HIS
1	B	72	HIS
1	B	134	GLN
1	B	235	ASN
1	C	29	HIS
1	C	72	HIS
1	C	235	ASN
1	D	29	HIS
1	D	72	HIS
1	D	235	ASN
1	E	29	HIS
1	E	72	HIS
1	E	235	ASN
1	F	29	HIS
1	F	72	HIS
1	F	235	ASN
1	G	29	HIS
1	G	72	HIS
1	G	235	ASN
1	H	29	HIS
1	H	72	HIS
1	H	235	ASN
1	I	29	HIS
1	I	72	HIS
1	I	235	ASN
1	J	29	HIS
1	J	72	HIS
1	J	235	ASN
1	K	29	HIS
1	K	72	HIS

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Mol	Chain	Res	Type
1	K	235	ASN
1	L	29	HIS
1	L	72	HIS
1	L	235	ASN
1	M	29	HIS
1	M	72	HIS
1	M	134	GLN
1	M	235	ASN
1	N	29	HIS
1	N	72	HIS
1	N	235	ASN

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 28 ligands modelled in this entry, 28 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	306/360 (85%)	0.28	9 (2%) 51 41	164, 186, 210, 221	0
1	B	306/360 (85%)	0.24	6 (1%) 65 56	151, 170, 187, 204	0
1	C	306/360 (85%)	0.46	15 (4%) 29 25	159, 184, 207, 220	0
1	D	306/360 (85%)	0.28	10 (3%) 46 37	147, 169, 191, 214	0
1	E	306/360 (85%)	0.53	18 (5%) 22 19	145, 166, 188, 221	0
1	F	306/360 (85%)	0.31	13 (4%) 36 30	143, 166, 187, 209	0
1	G	306/360 (85%)	0.47	16 (5%) 27 24	143, 160, 180, 200	0
1	H	306/360 (85%)	0.28	10 (3%) 46 37	120, 149, 164, 190	0
1	I	306/360 (85%)	0.51	19 (6%) 20 17	142, 172, 199, 219	0
1	J	306/360 (85%)	0.26	6 (1%) 65 56	137, 160, 184, 198	0
1	K	306/360 (85%)	0.54	20 (6%) 18 15	159, 186, 211, 228	0
1	L	306/360 (85%)	0.26	7 (2%) 60 51	144, 168, 193, 208	0
1	M	306/360 (85%)	0.43	20 (6%) 18 15	160, 180, 200, 220	0
1	N	306/360 (85%)	0.26	7 (2%) 60 51	141, 163, 182, 193	0
All	All	4284/5040 (85%)	0.36	176 (4%) 37 30	120, 170, 200, 228	0

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	210	PHE	5.9
1	I	210	PHE	4.6
1	F	189	GLY	4.6
1	C	237	TYR	4.5
1	F	188	LEU	4.4
1	G	21	ARG	4.1
1	I	237	TYR	4.1
1	K	222	LYS	4.1

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Mol	Chain	Res	Type	RSRZ
1	C	210	PHE	4.0
1	B	21	ARG	4.0
1	I	21	ARG	4.0
1	M	222	LYS	3.9
1	C	74	VAL	3.6
1	N	285	MET	3.6
1	F	210	PHE	3.6
1	A	210	PHE	3.5
1	I	188	LEU	3.5
1	K	122	GLU	3.4
1	I	296	LYS	3.4
1	E	74	VAL	3.4
1	E	175	VAL	3.4
1	C	138	ILE	3.3
1	G	210	PHE	3.3
1	E	210	PHE	3.3
1	L	210	PHE	3.3
1	E	21	ARG	3.3
1	M	188	LEU	3.3
1	M	210	PHE	3.3
1	F	138	ILE	3.2
1	G	285	MET	3.2
1	E	191	ALA	3.2
1	J	210	PHE	3.2
1	E	212	THR	3.2
1	G	69	VAL	3.2
1	B	210	PHE	3.1
1	M	264	ASP	3.1
1	K	74	VAL	3.0
1	D	210	PHE	3.0
1	M	21	ARG	3.0
1	A	69	VAL	3.0
1	G	175	VAL	3.0
1	K	64	ILE	3.0
1	F	137	GLY	3.0
1	C	165	TRP	3.0
1	K	61	ARG	3.0
1	D	245	ARG	2.9
1	N	112	TRP	2.9
1	I	74	VAL	2.9
1	K	296	LYS	2.9
1	I	189	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	H	175	VAL	2.8
1	C	21	ARG	2.8
1	K	21	ARG	2.8
1	M	237	TYR	2.8
1	F	237	TYR	2.8
1	H	21	ARG	2.8
1	K	237	TYR	2.8
1	A	235	ASN	2.8
1	D	56	PHE	2.8
1	M	291	ASP	2.8
1	M	296	LYS	2.8
1	I	255	GLU	2.7
1	L	137	GLY	2.7
1	H	210	PHE	2.7
1	G	66	LYS	2.7
1	E	70	VAL	2.7
1	A	188	LEU	2.7
1	B	245	ARG	2.7
1	J	21	ARG	2.7
1	G	82	ILE	2.7
1	F	190	TYR	2.7
1	D	175	VAL	2.7
1	G	188	LEU	2.7
1	C	137	GLY	2.6
1	E	308	ARG	2.6
1	K	245	ARG	2.6
1	M	245	ARG	2.6
1	G	301	PHE	2.6
1	D	74	VAL	2.6
1	C	69	VAL	2.6
1	C	175	VAL	2.6
1	K	211	GLY	2.6
1	I	165	TRP	2.6
1	N	188	LEU	2.6
1	M	175	VAL	2.5
1	J	61	ARG	2.5
1	H	170	TYR	2.5
1	J	237	TYR	2.5
1	K	212	THR	2.5
1	M	301	PHE	2.5
1	E	69	VAL	2.5
1	E	66	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	54	GLN	2.4
1	F	57	ARG	2.4
1	K	190	TYR	2.4
1	J	74	VAL	2.4
1	E	222	LYS	2.4
1	D	57	ARG	2.4
1	A	74	VAL	2.4
1	I	245	ARG	2.4
1	M	288	ASN	2.4
1	A	323	ARG	2.4
1	C	61	ARG	2.4
1	M	112	TRP	2.4
1	A	189	GLY	2.4
1	E	138	ILE	2.3
1	L	162	GLU	2.3
1	C	188	LEU	2.3
1	H	74	VAL	2.3
1	D	188	LEU	2.3
1	F	69	VAL	2.3
1	N	210	PHE	2.3
1	E	82	ILE	2.3
1	A	285	MET	2.3
1	H	291	ASP	2.3
1	E	188	LEU	2.3
1	D	54	GLN	2.3
1	G	291	ASP	2.3
1	M	212	THR	2.3
1	I	57	ARG	2.3
1	A	21	ARG	2.3
1	L	301	PHE	2.3
1	B	285	MET	2.2
1	J	24	GLY	2.2
1	G	74	VAL	2.2
1	C	255	GLU	2.2
1	B	188	LEU	2.2
1	F	285	MET	2.2
1	G	308	ARG	2.2
1	I	66	LYS	2.2
1	F	175	VAL	2.2
1	K	191	ALA	2.2
1	L	138	ILE	2.2
1	G	286	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	301	PHE	2.2
1	M	282	ASN	2.2
1	D	82	ILE	2.2
1	C	195	GLY	2.2
1	L	237	TYR	2.2
1	K	18	PRO	2.2
1	E	152	ASP	2.2
1	G	237	TYR	2.2
1	H	237	TYR	2.2
1	K	175	VAL	2.2
1	E	71	VAL	2.2
1	N	190	TYR	2.2
1	I	323	ARG	2.1
1	M	170	TYR	2.1
1	D	272	TYR	2.1
1	I	190	TYR	2.1
1	M	239	ILE	2.1
1	B	301	PHE	2.1
1	E	148	PHE	2.1
1	N	21	ARG	2.1
1	I	162	GLU	2.1
1	C	71	VAL	2.1
1	H	176	CYS	2.1
1	M	299	VAL	2.1
1	M	262	GLN	2.1
1	C	301	PHE	2.1
1	F	289	TYR	2.1
1	I	169	ARG	2.1
1	I	56	PHE	2.1
1	K	58	ALA	2.1
1	K	188	LEU	2.1
1	K	189	GLY	2.1
1	M	260	PRO	2.1
1	H	190	TYR	2.1
1	H	78	GLU	2.1
1	I	308	ARG	2.0
1	L	70	VAL	2.0
1	I	69	VAL	2.0
1	G	18	PRO	2.0
1	N	114	ASN	2.0
1	K	119	LEU	2.0
1	G	282	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 monosaccharides 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.

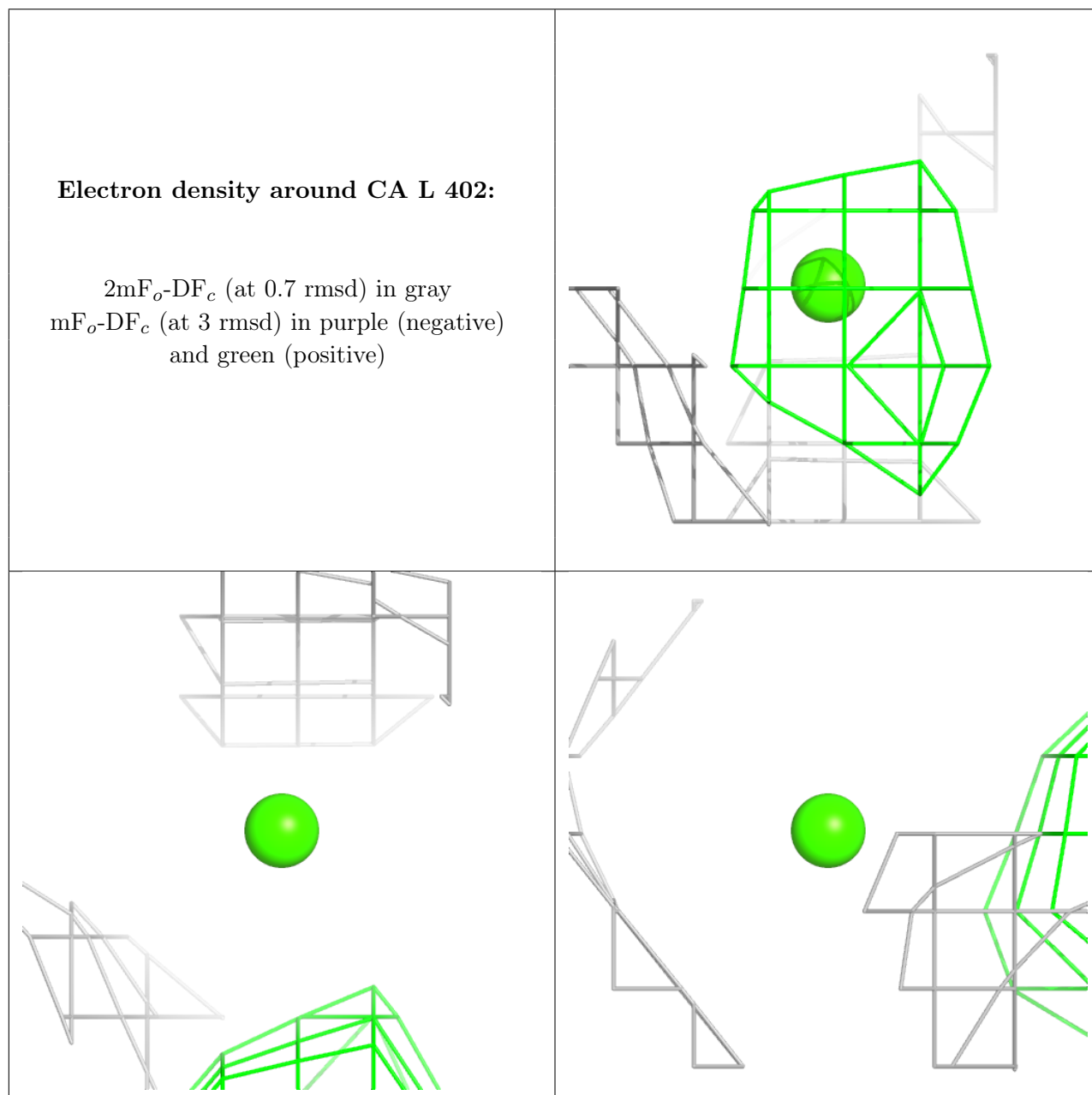
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	CA	L	402	1/1	0.71	0.14	162,162,162,162	0
3	CA	G	402	1/1	0.78	0.19	162,162,162,162	0
2	ZN	N	401	1/1	0.81	0.16	159,159,159,159	0
2	ZN	C	401	1/1	0.86	0.15	180,180,180,180	0
3	CA	E	402	1/1	0.86	0.14	177,177,177,177	0
3	CA	B	402	1/1	0.87	0.24	183,183,183,183	0
2	ZN	M	401	1/1	0.87	0.12	177,177,177,177	0
3	CA	J	402	1/1	0.88	0.17	163,163,163,163	0
2	ZN	G	401	1/1	0.88	0.18	159,159,159,159	0
3	CA	D	402	1/1	0.89	0.19	168,168,168,168	0
3	CA	C	402	1/1	0.90	0.21	188,188,188,188	0
2	ZN	A	401	1/1	0.91	0.13	179,179,179,179	0
2	ZN	D	401	1/1	0.92	0.13	156,156,156,156	0
3	CA	M	402	1/1	0.92	0.18	190,190,190,190	0
3	CA	I	402	1/1	0.93	0.11	171,171,171,171	0
2	ZN	B	401	1/1	0.93	0.16	149,149,149,149	0
3	CA	F	402	1/1	0.94	0.20	172,172,172,172	0
2	ZN	H	401	1/1	0.94	0.17	136,136,136,136	0
2	ZN	L	401	1/1	0.94	0.23	145,145,145,145	0
2	ZN	F	401	1/1	0.95	0.17	140,140,140,140	0
2	ZN	E	401	1/1	0.95	0.14	152,152,152,152	0
3	CA	A	402	1/1	0.96	0.20	193,193,193,193	0
3	CA	K	402	1/1	0.96	0.12	191,191,191,191	0
3	CA	H	402	1/1	0.97	0.19	159,159,159,159	0
2	ZN	I	401	1/1	0.97	0.09	142,142,142,142	0
2	ZN	K	401	1/1	0.97	0.13	174,174,174,174	0
2	ZN	J	401	1/1	0.99	0.19	142,142,142,142	0

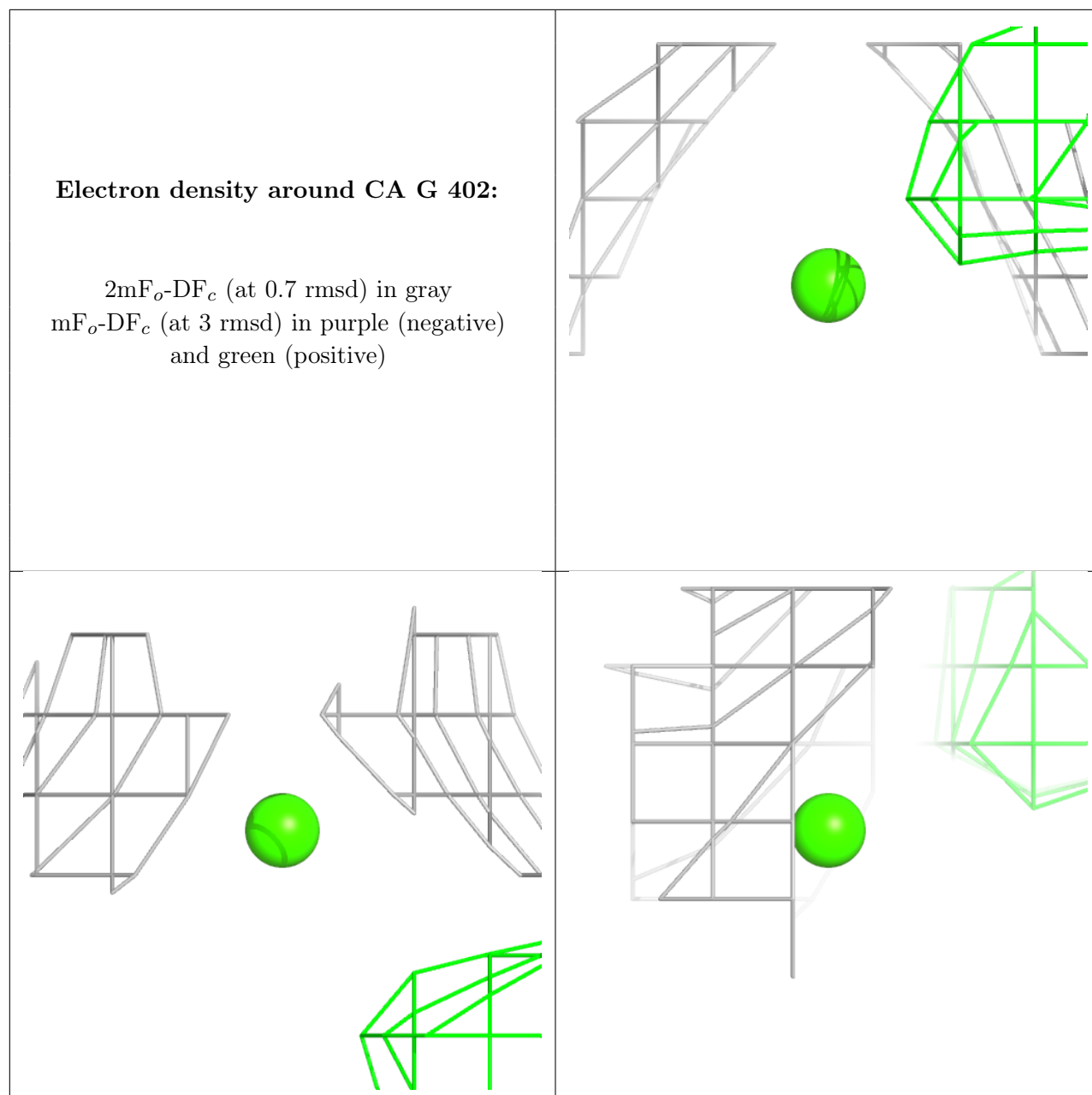
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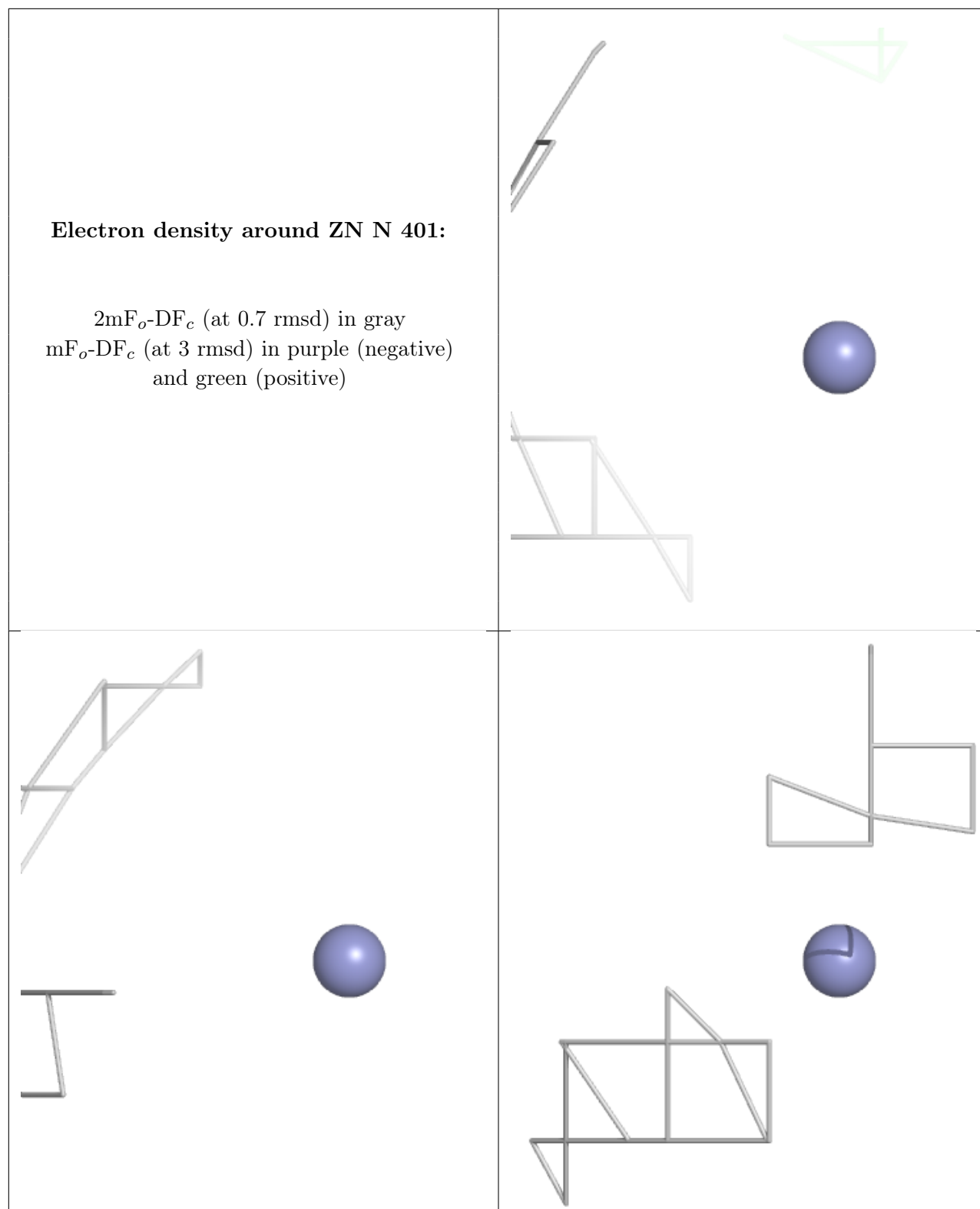
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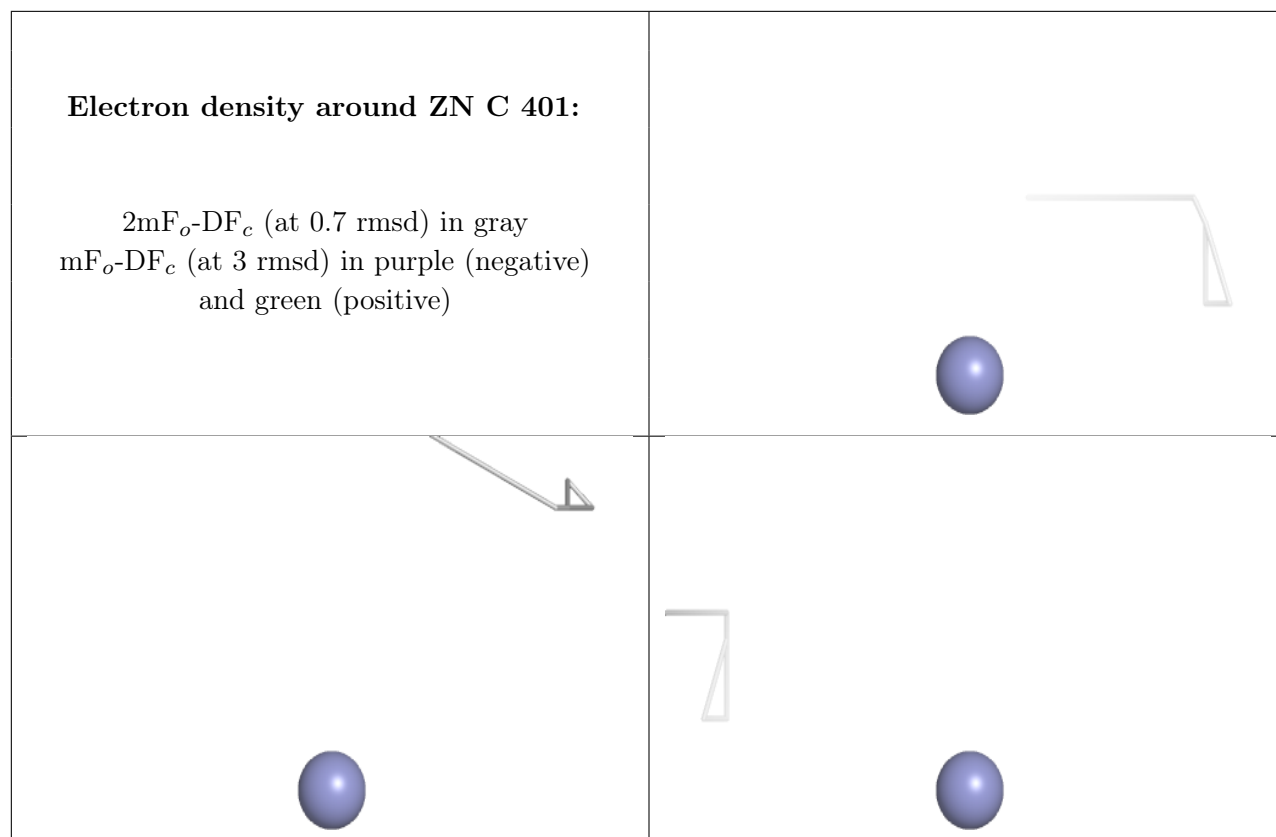
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	CA	N	402	1/1	0.99	0.19	168,168,168,168	0

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.



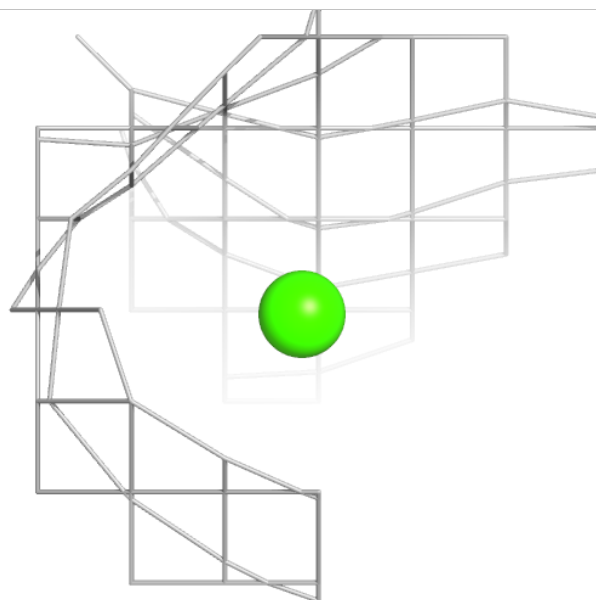
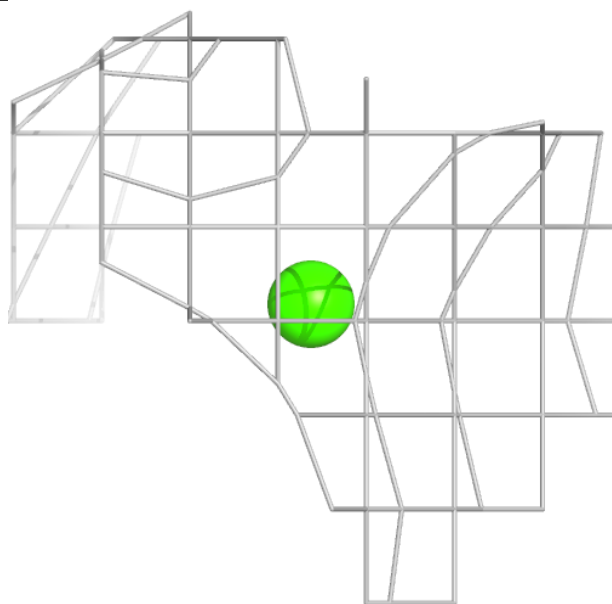
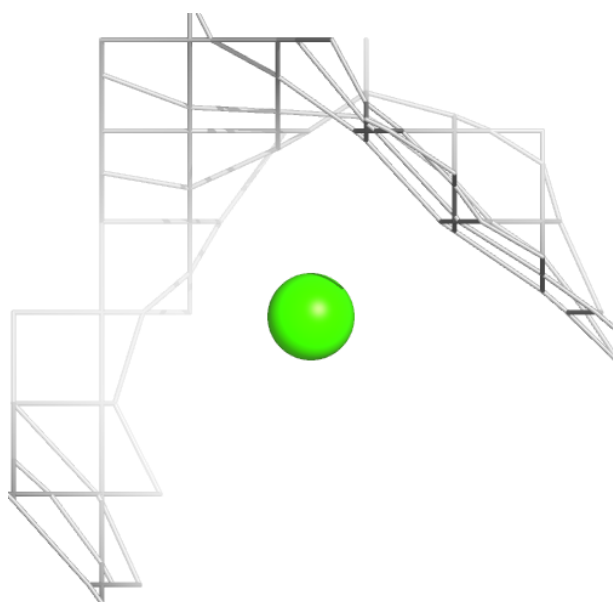






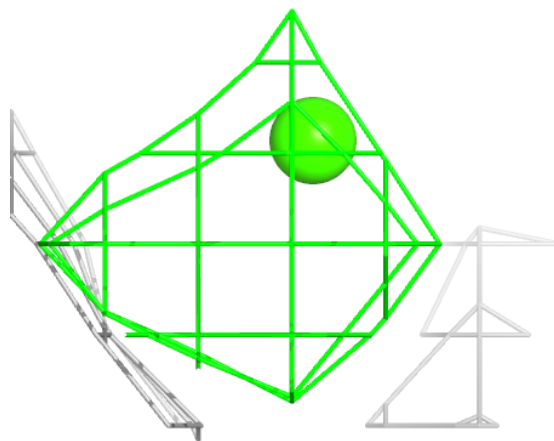
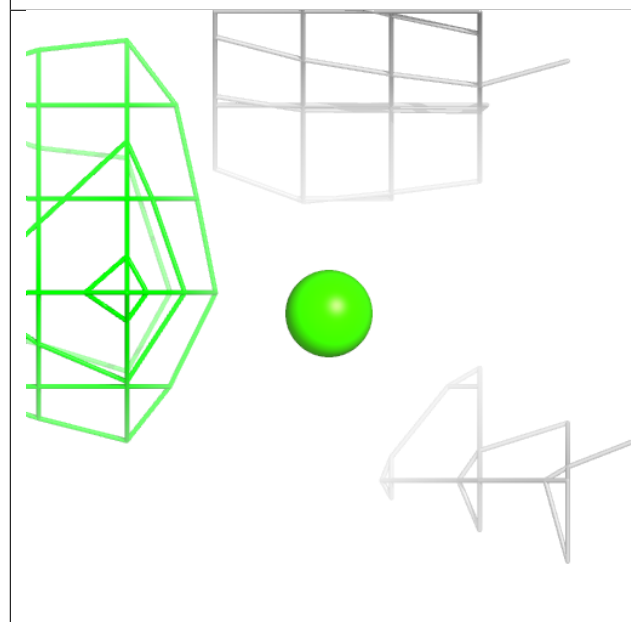
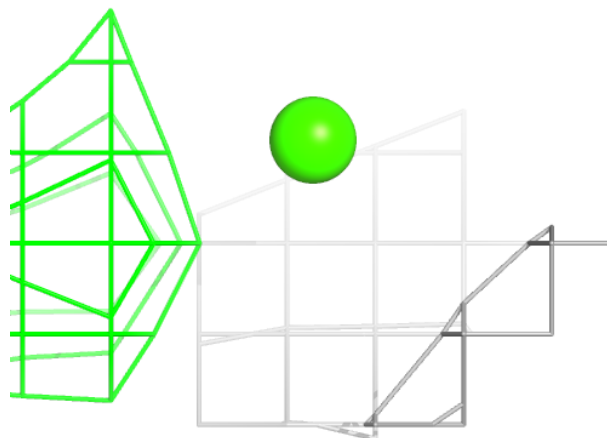
Electron density around CA E 402:

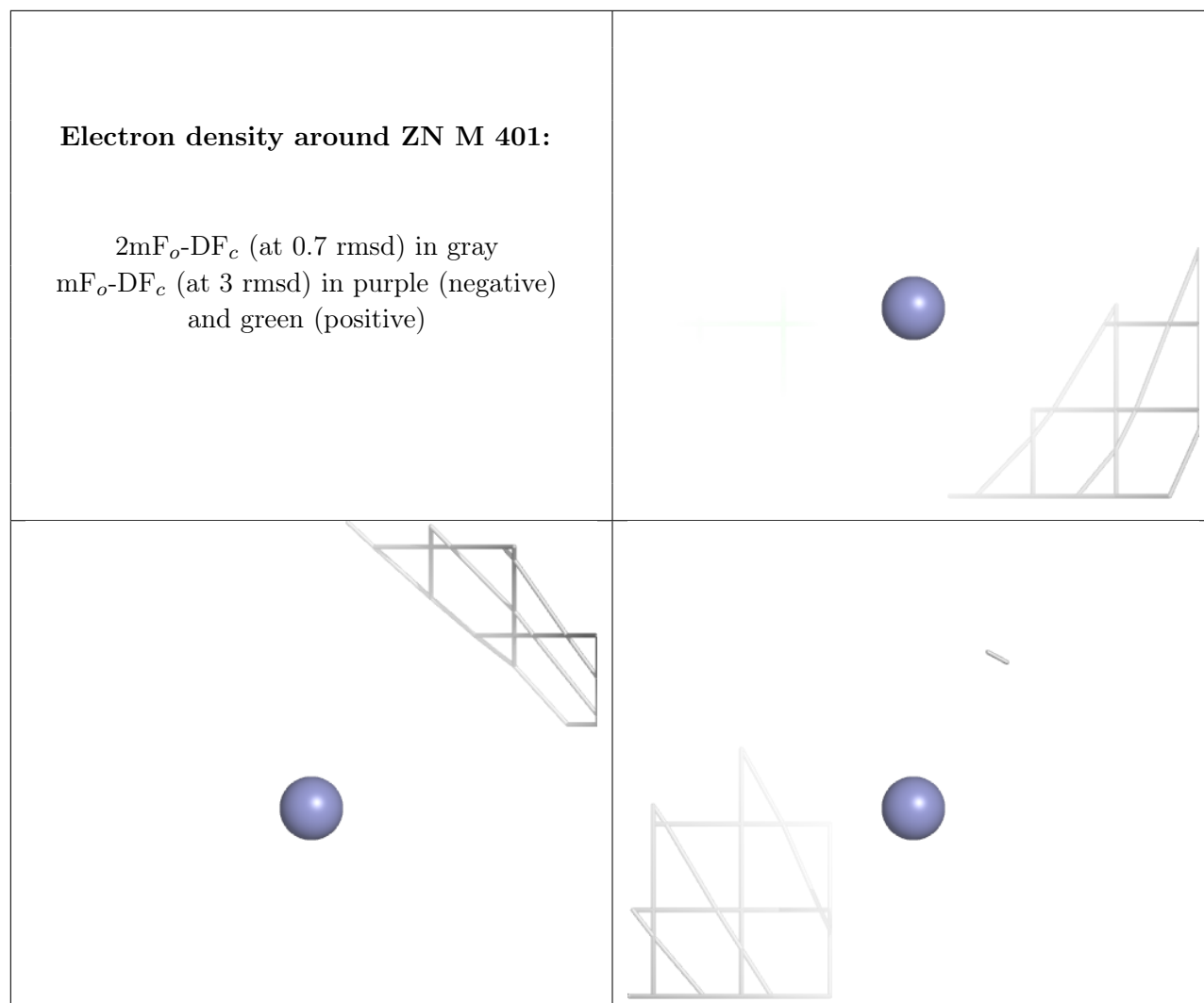
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

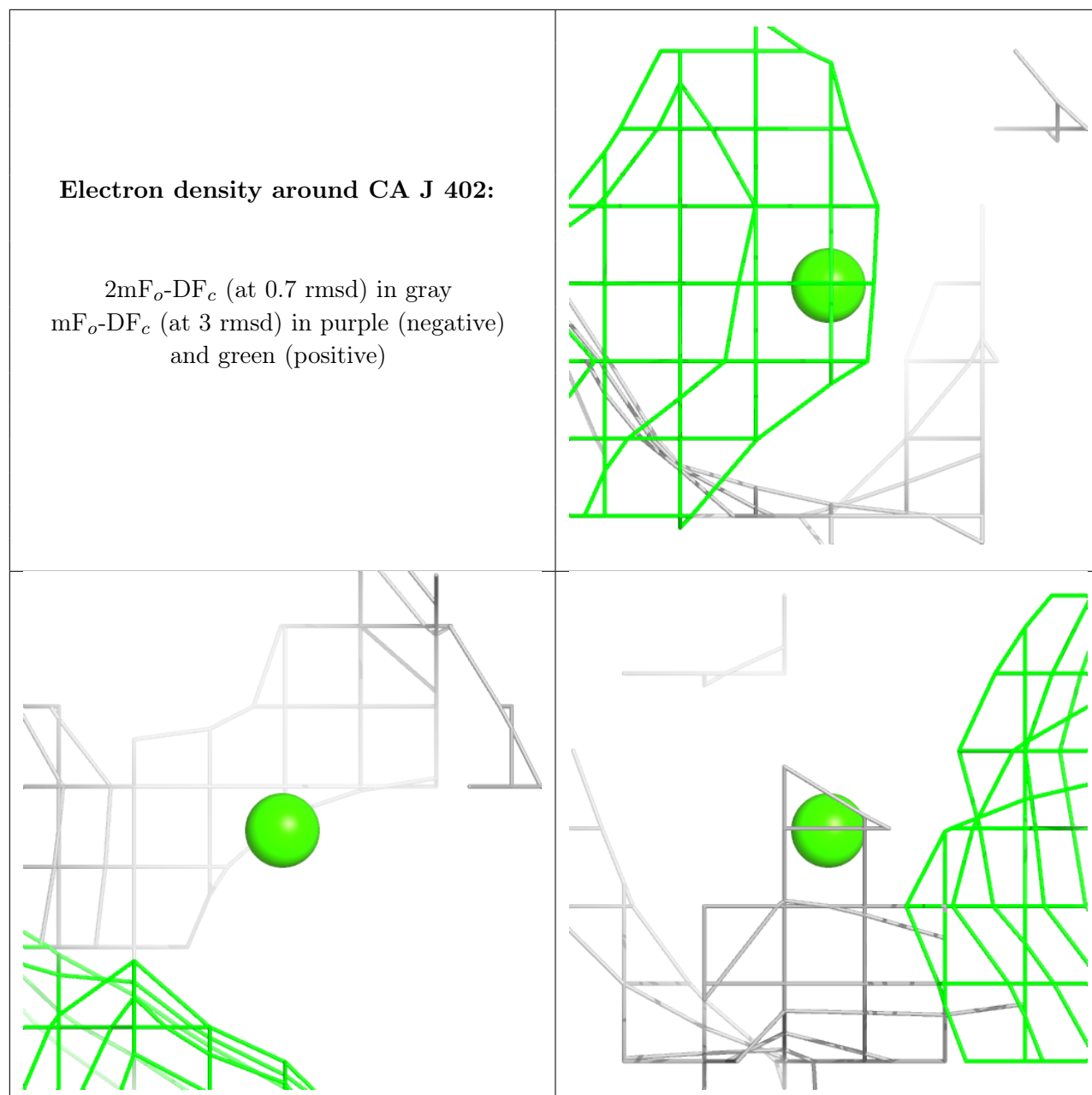


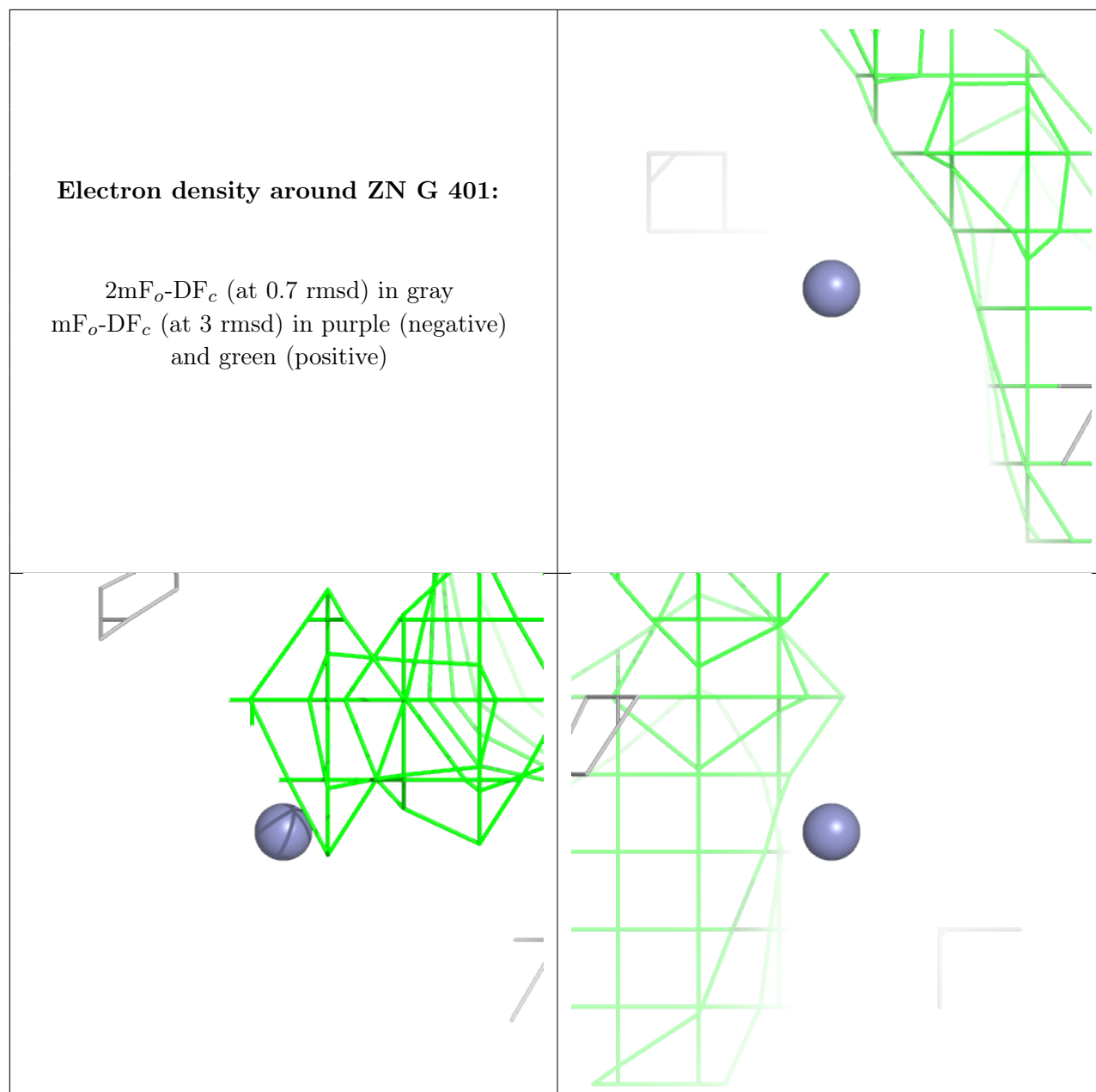
Electron density around CA B 402:

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and green (positive)



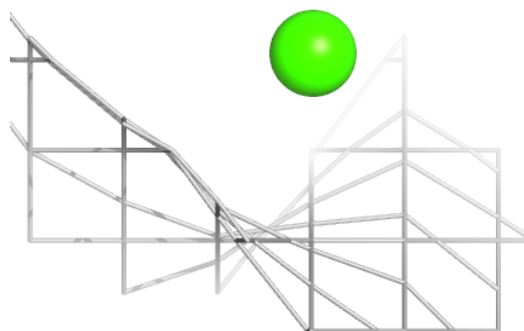
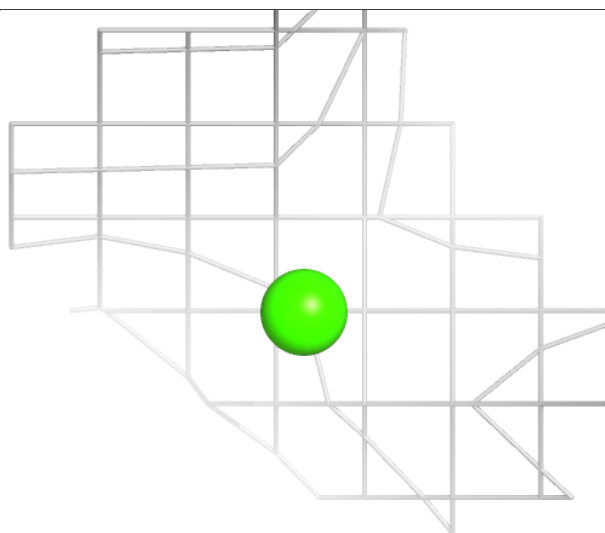
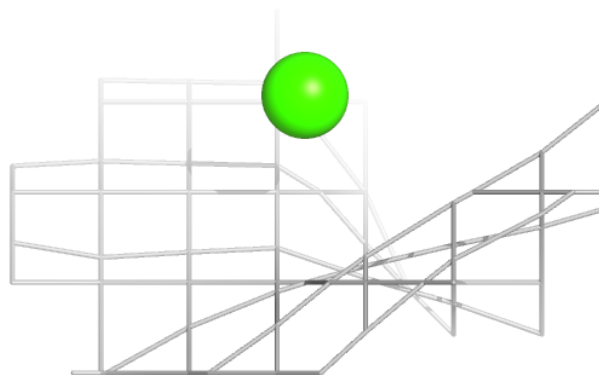






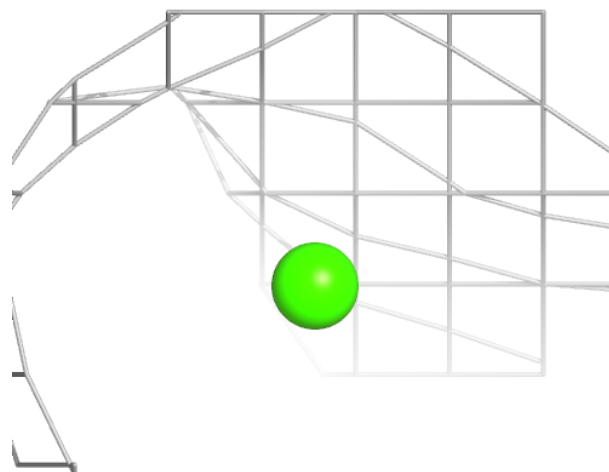
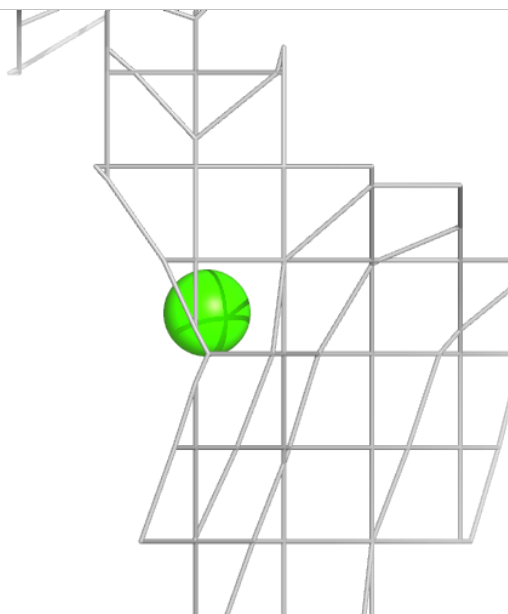
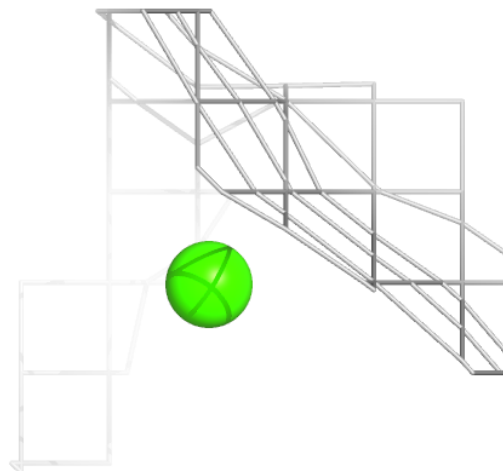
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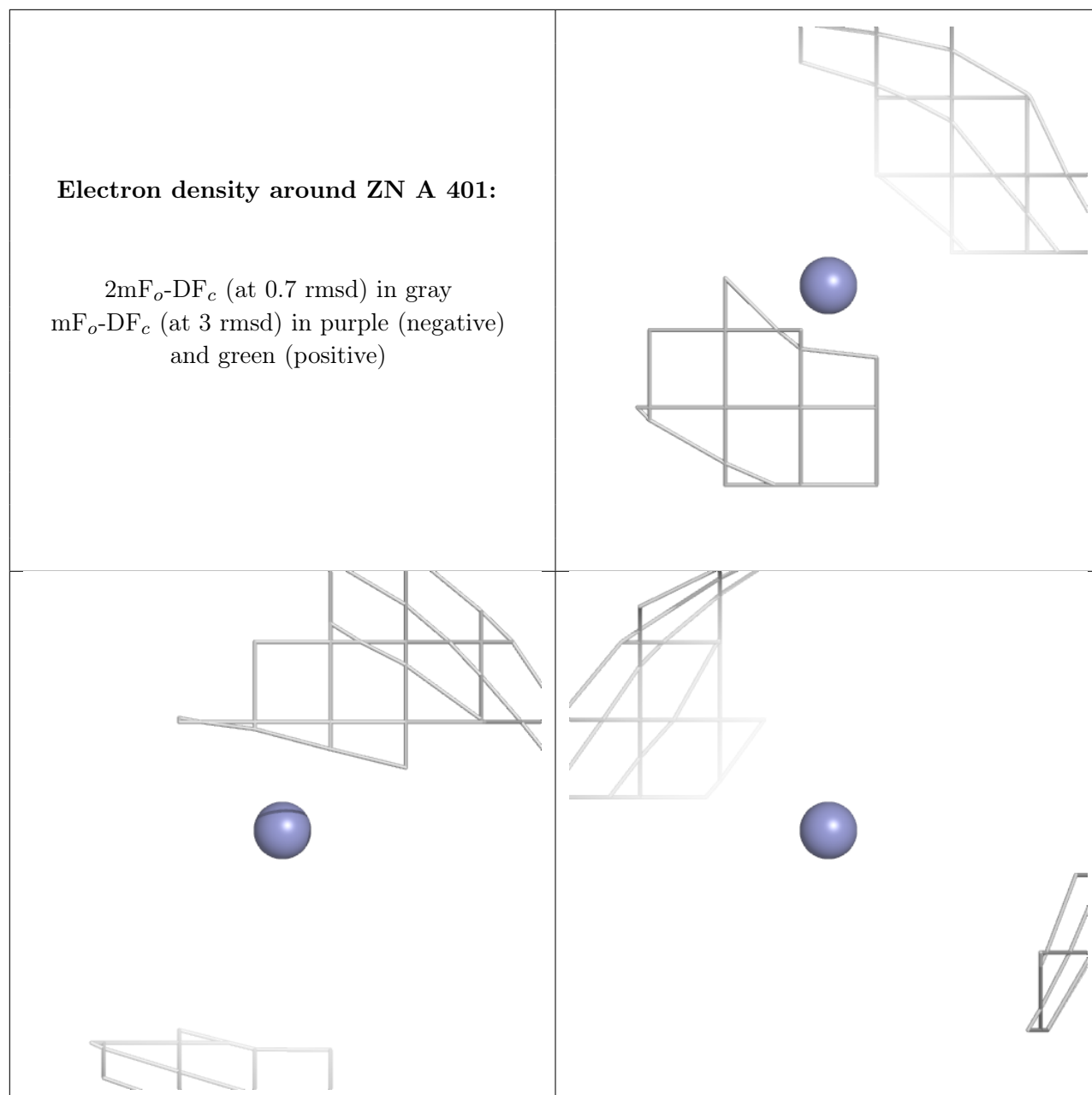
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and green (positive)

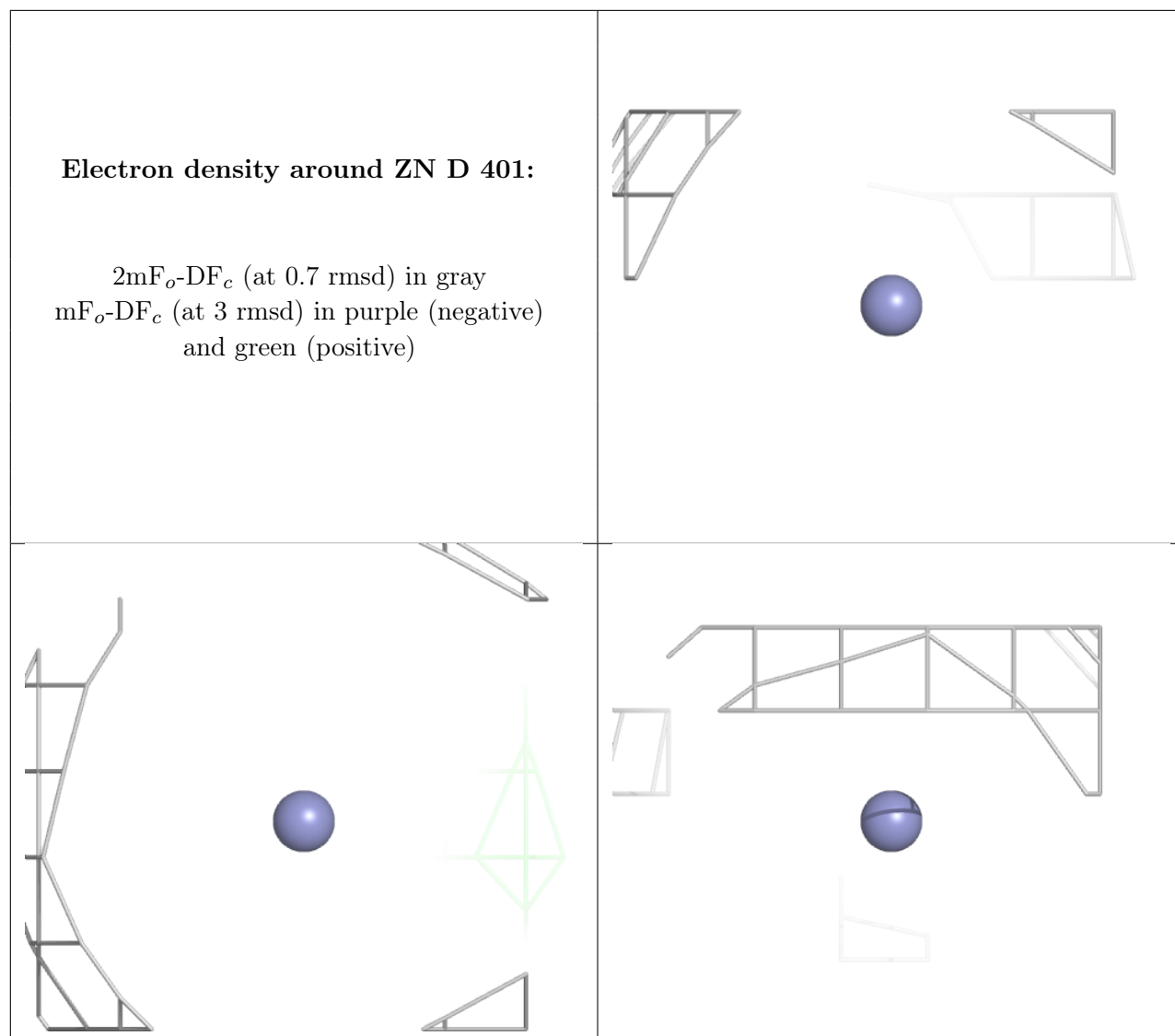


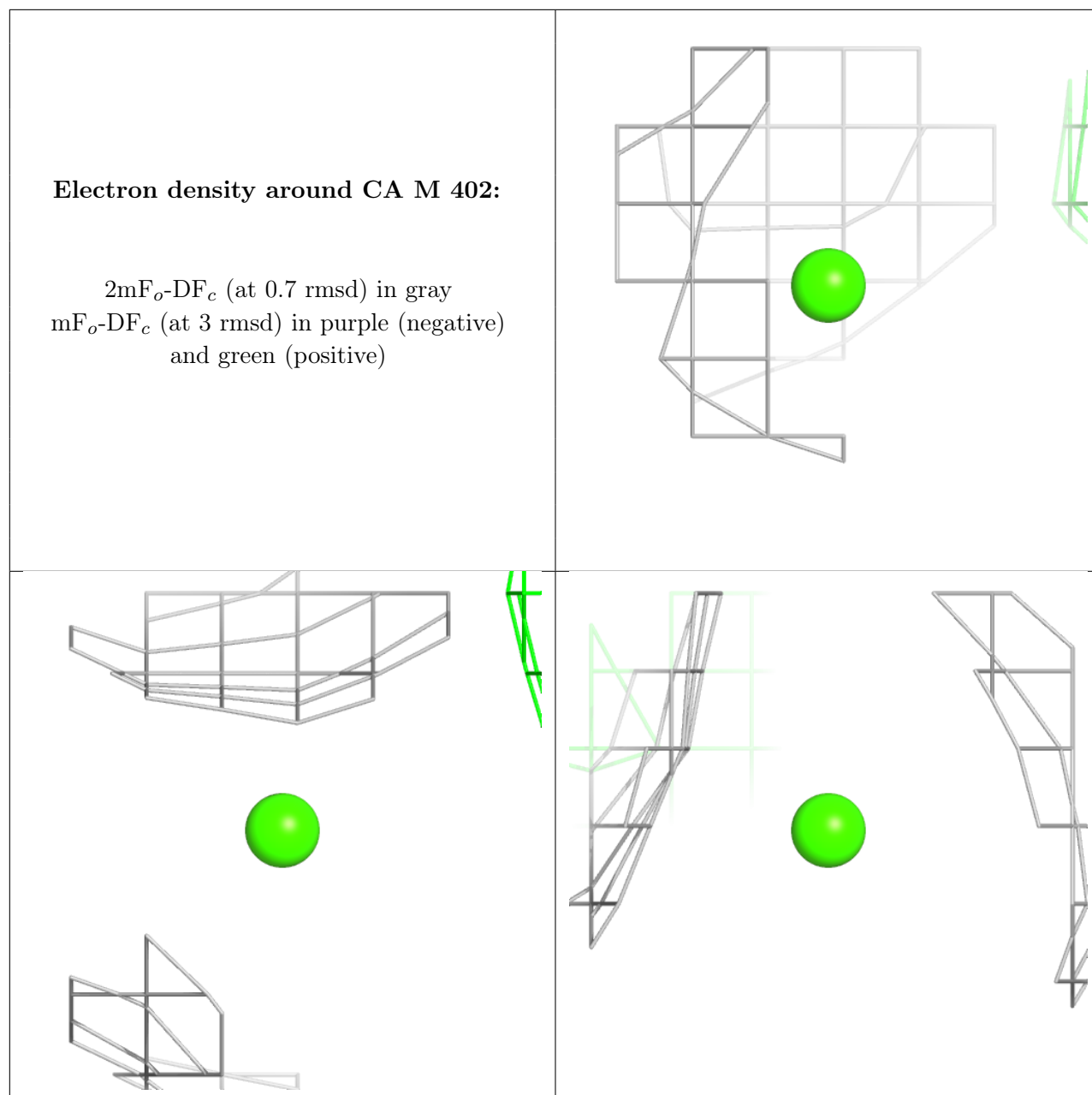
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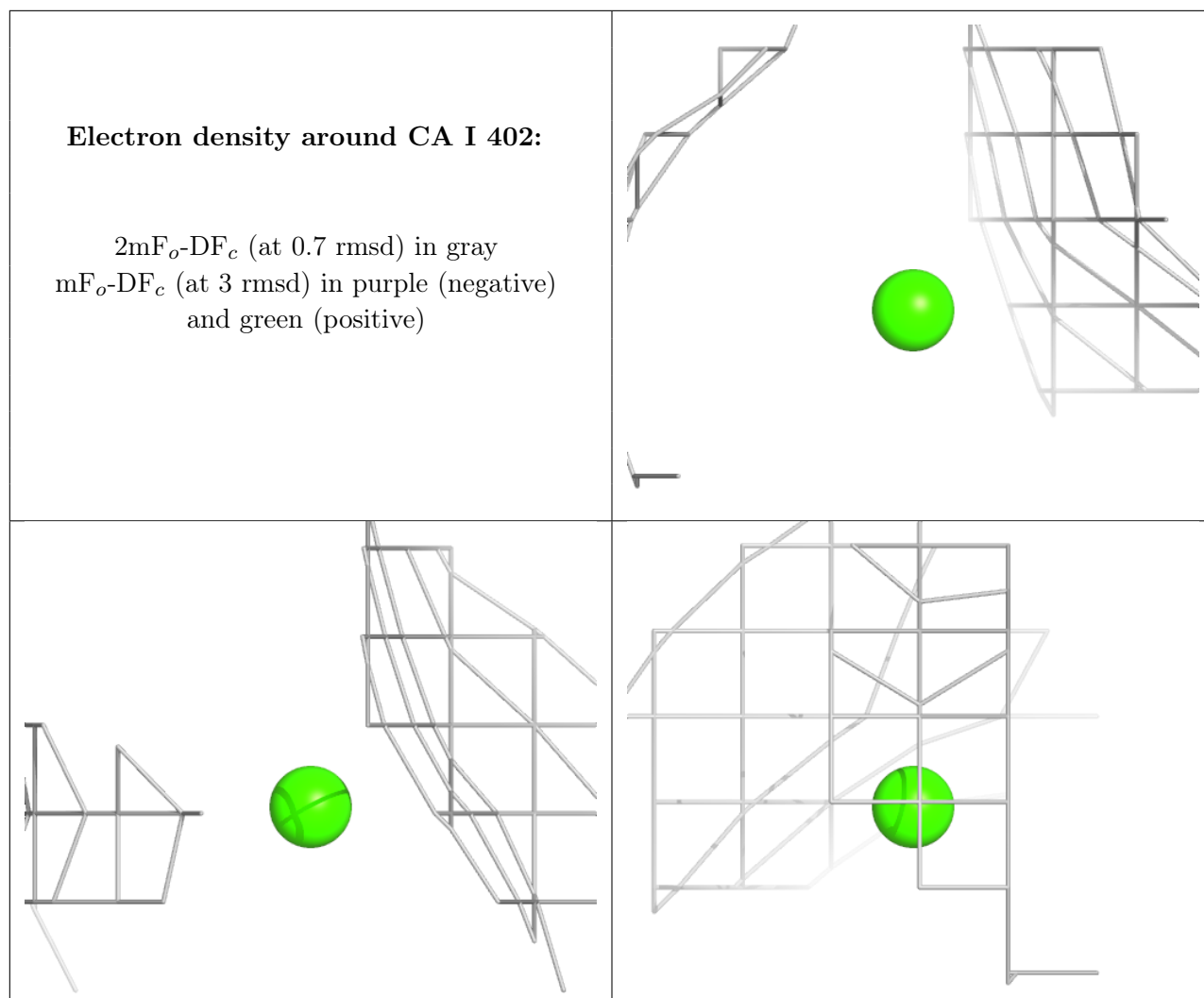
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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

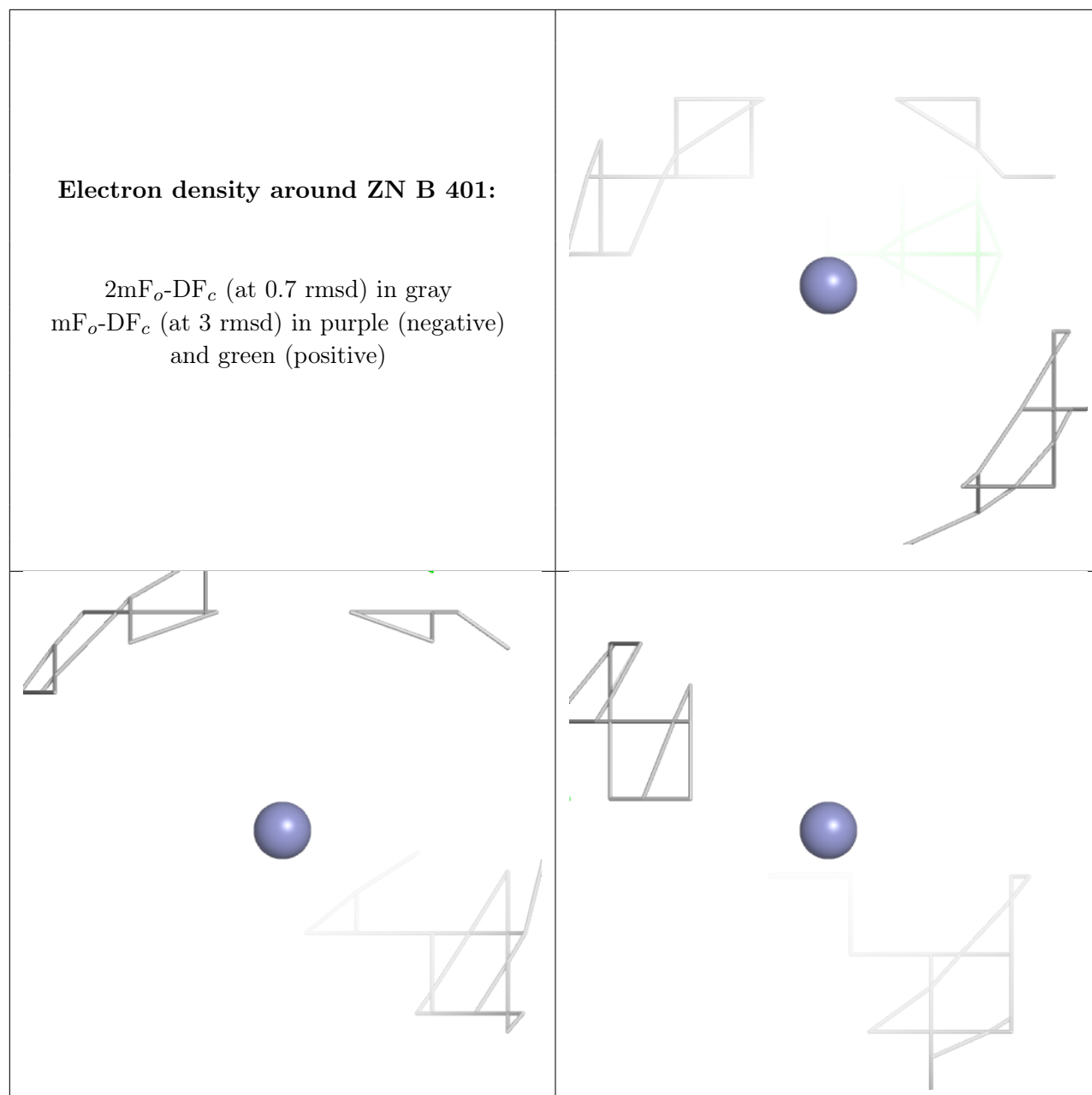






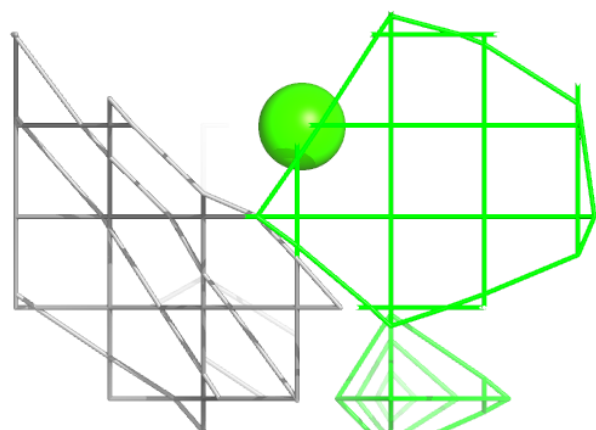
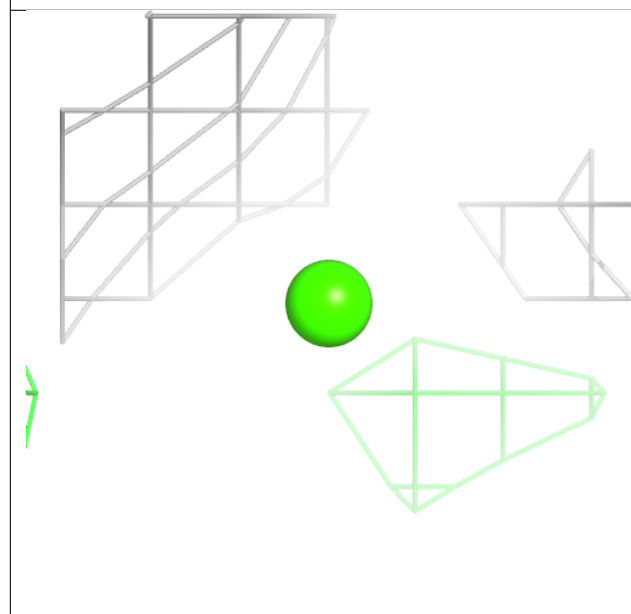
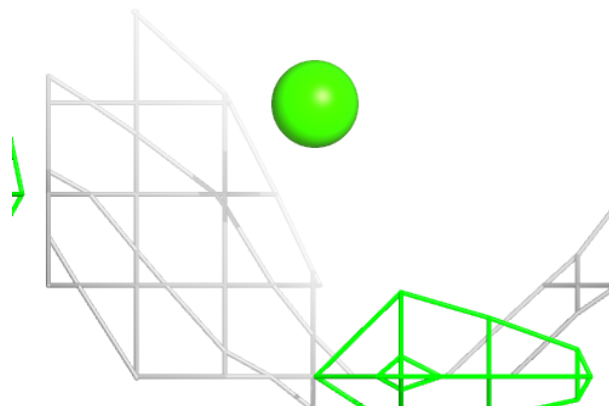


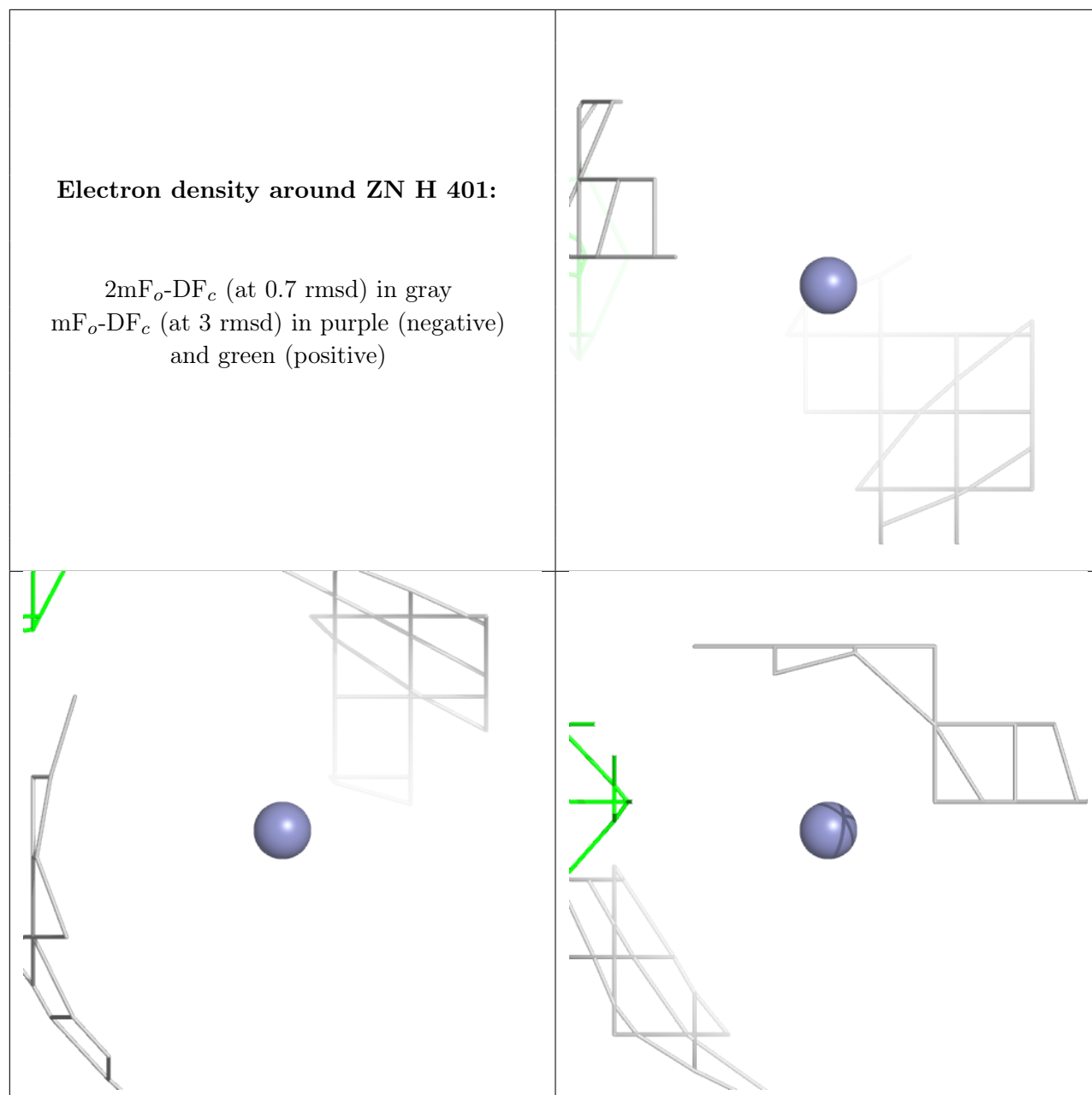


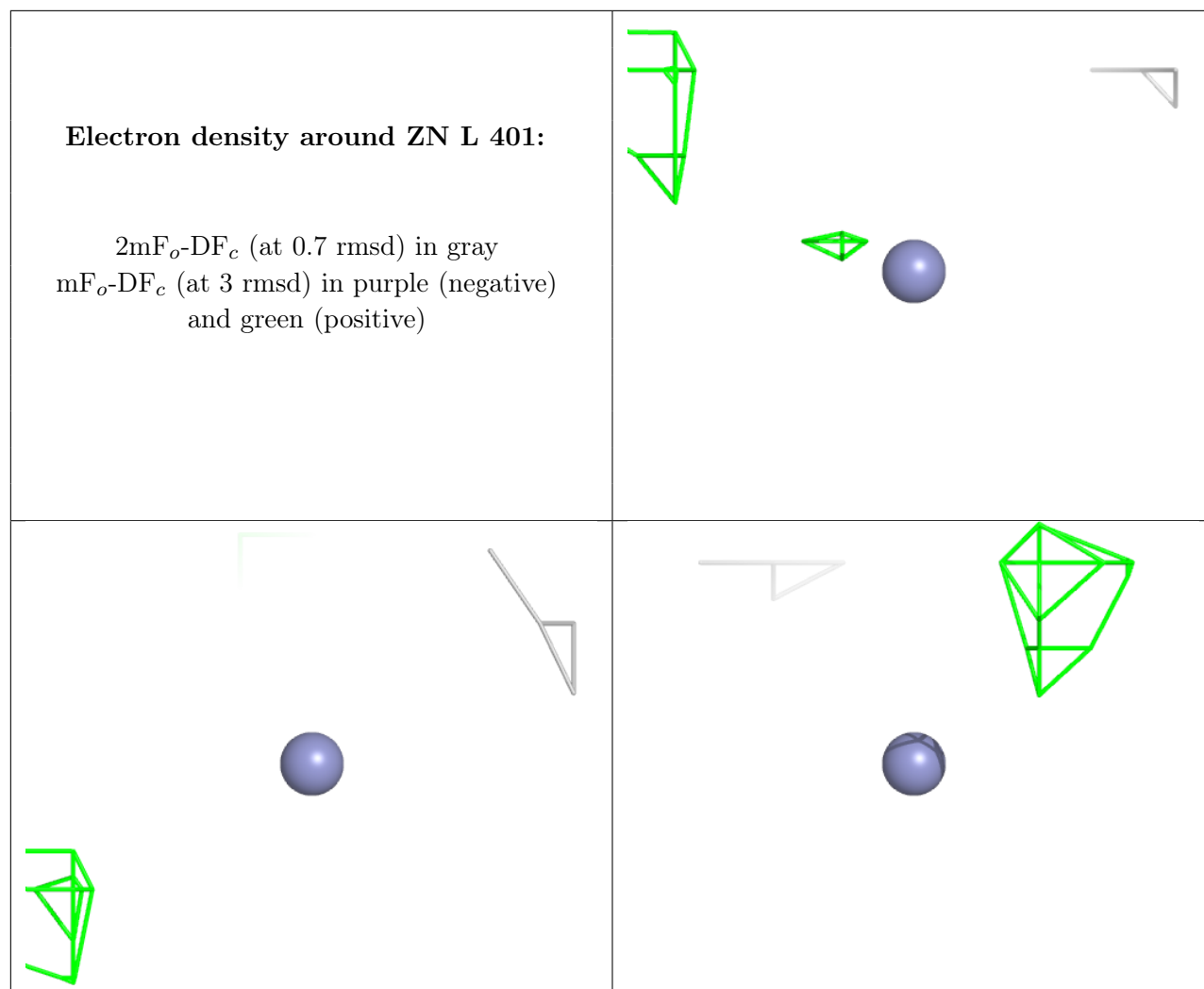


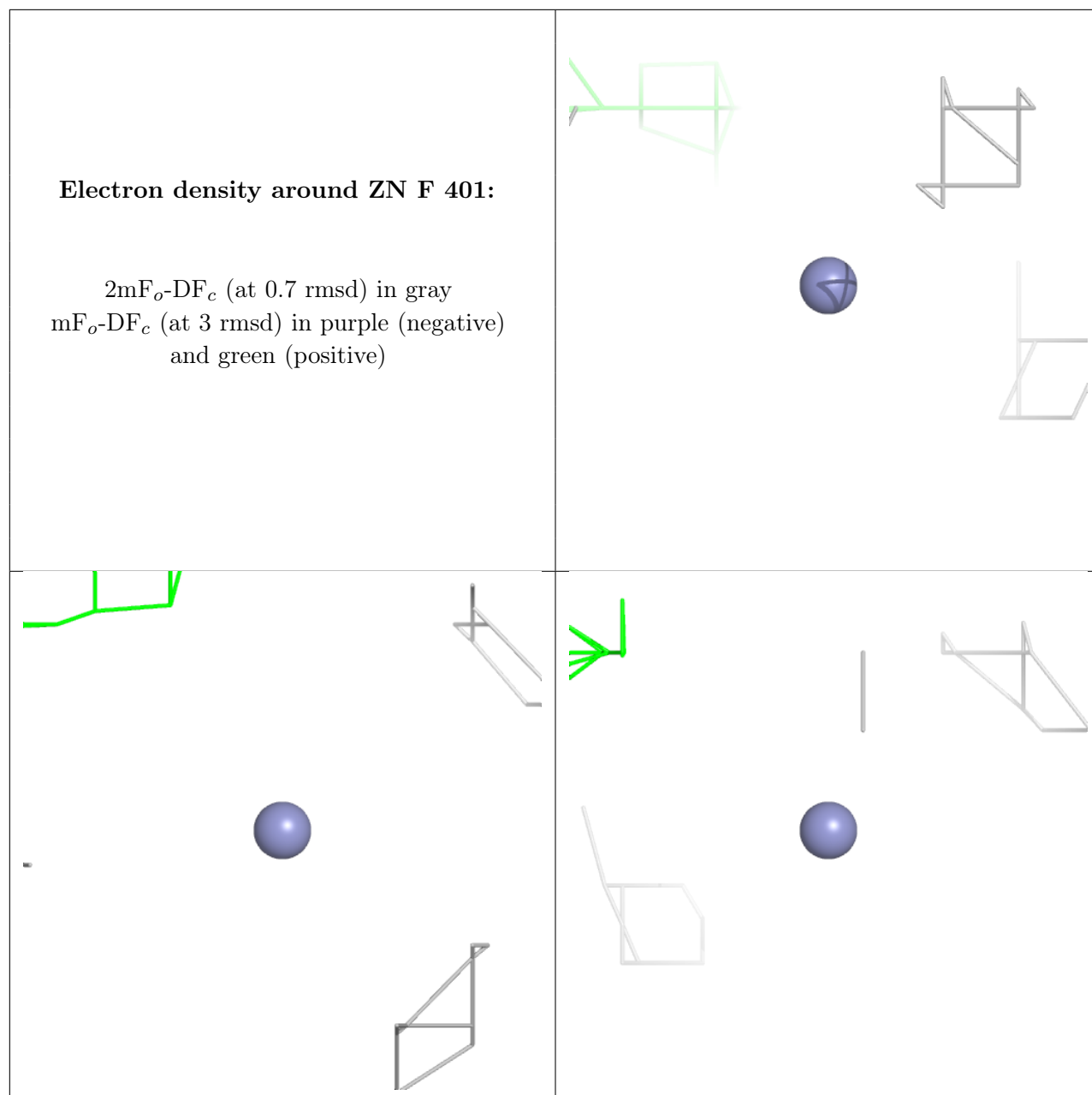
Electron density around CA F 402:

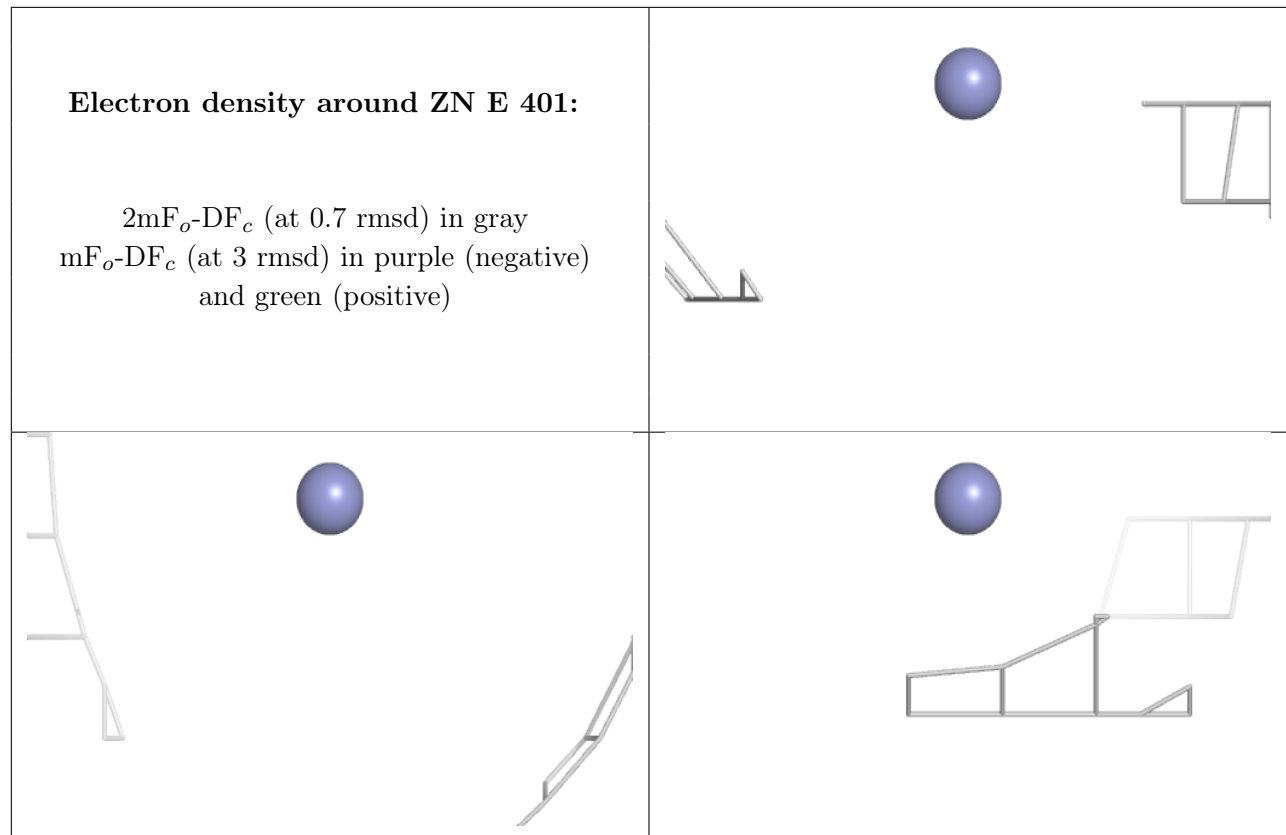
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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





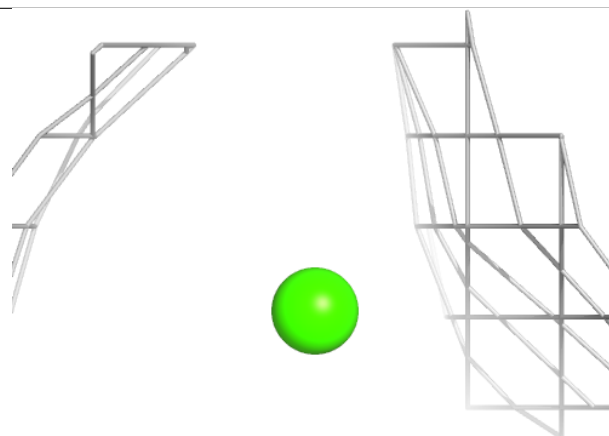
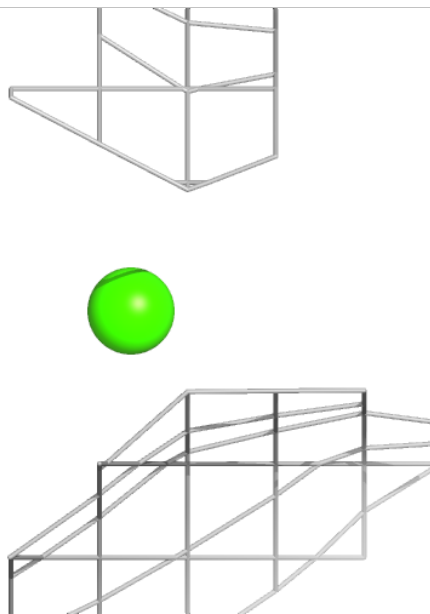
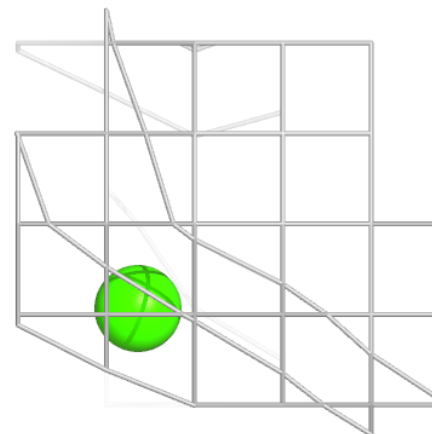


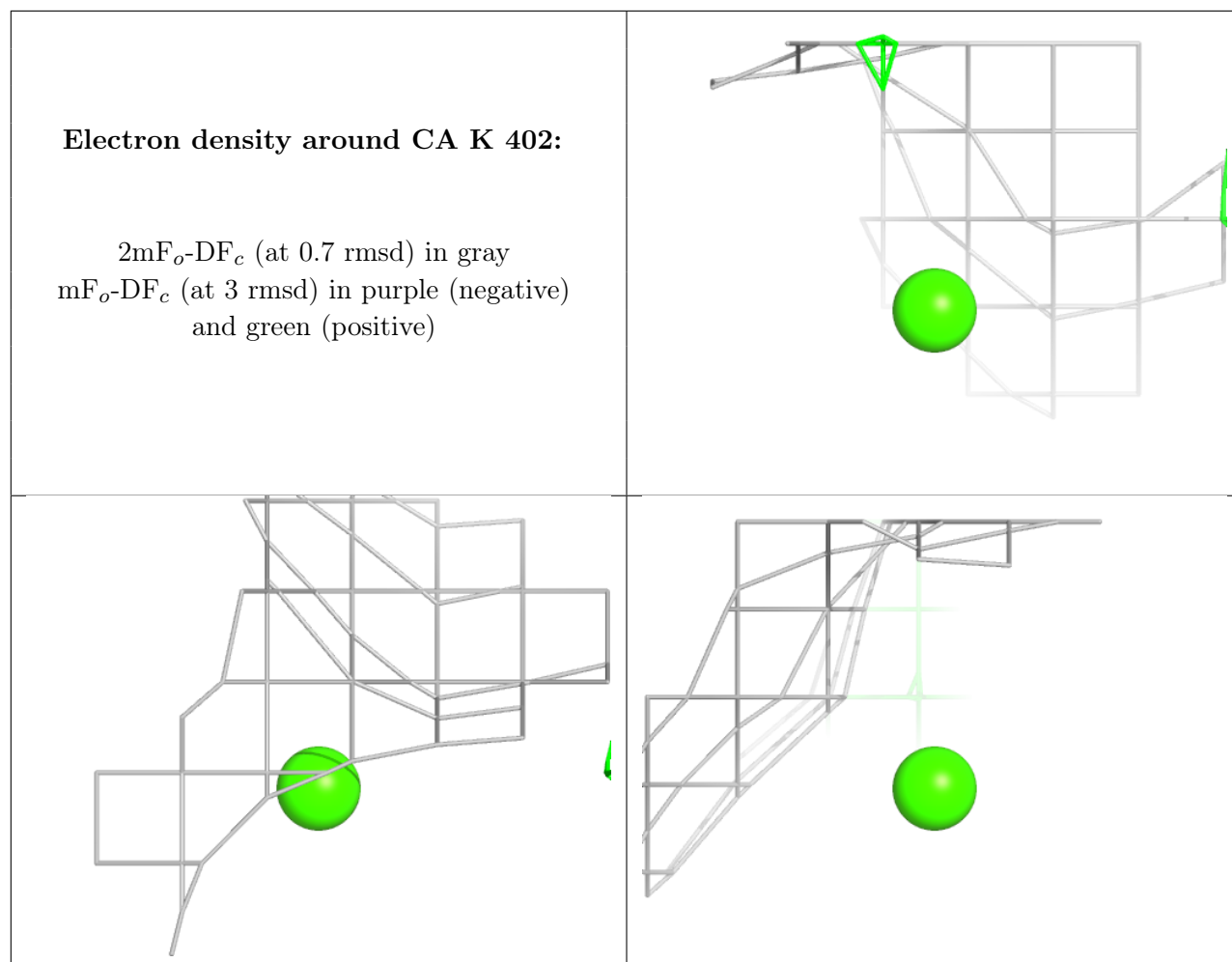


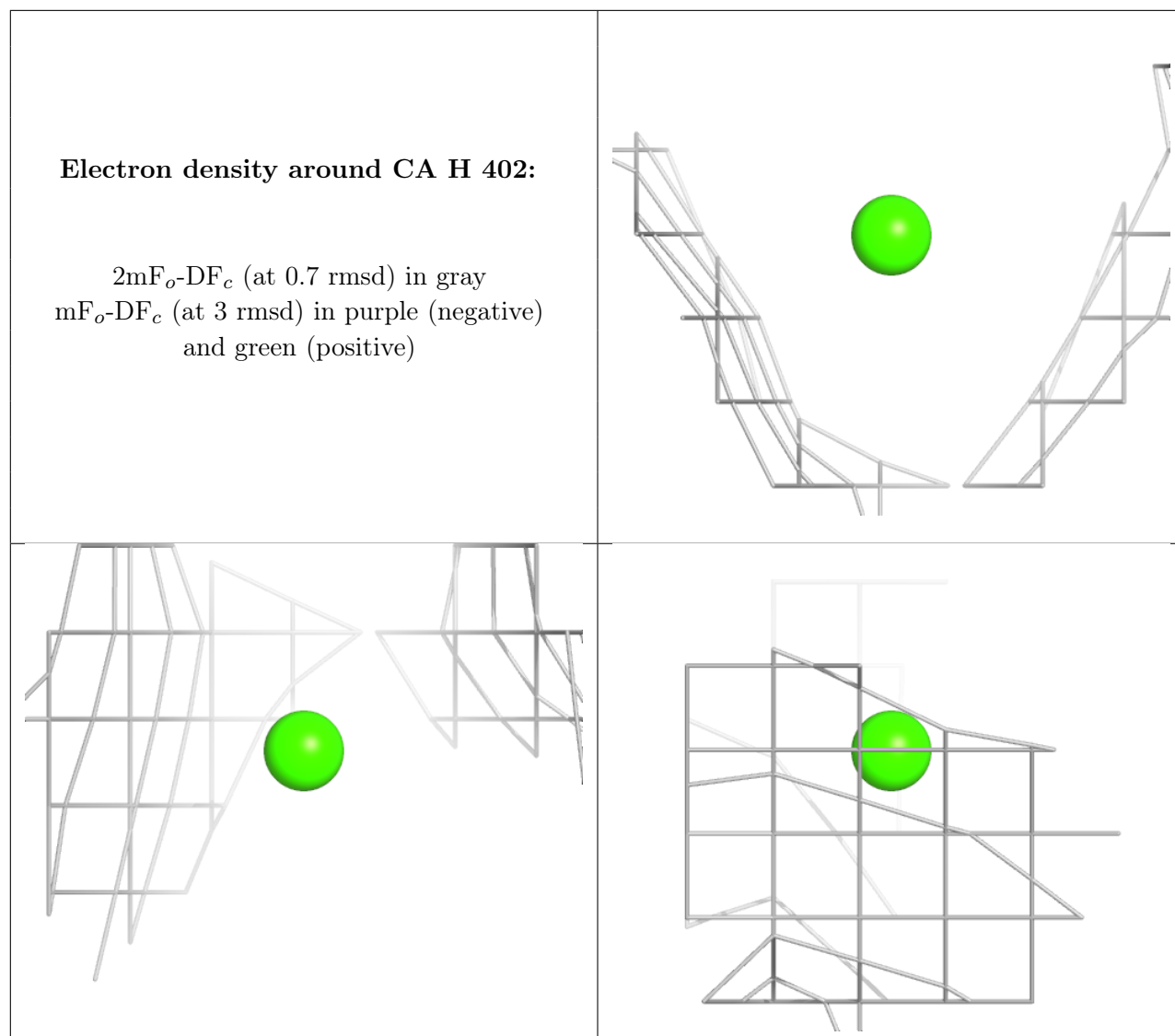


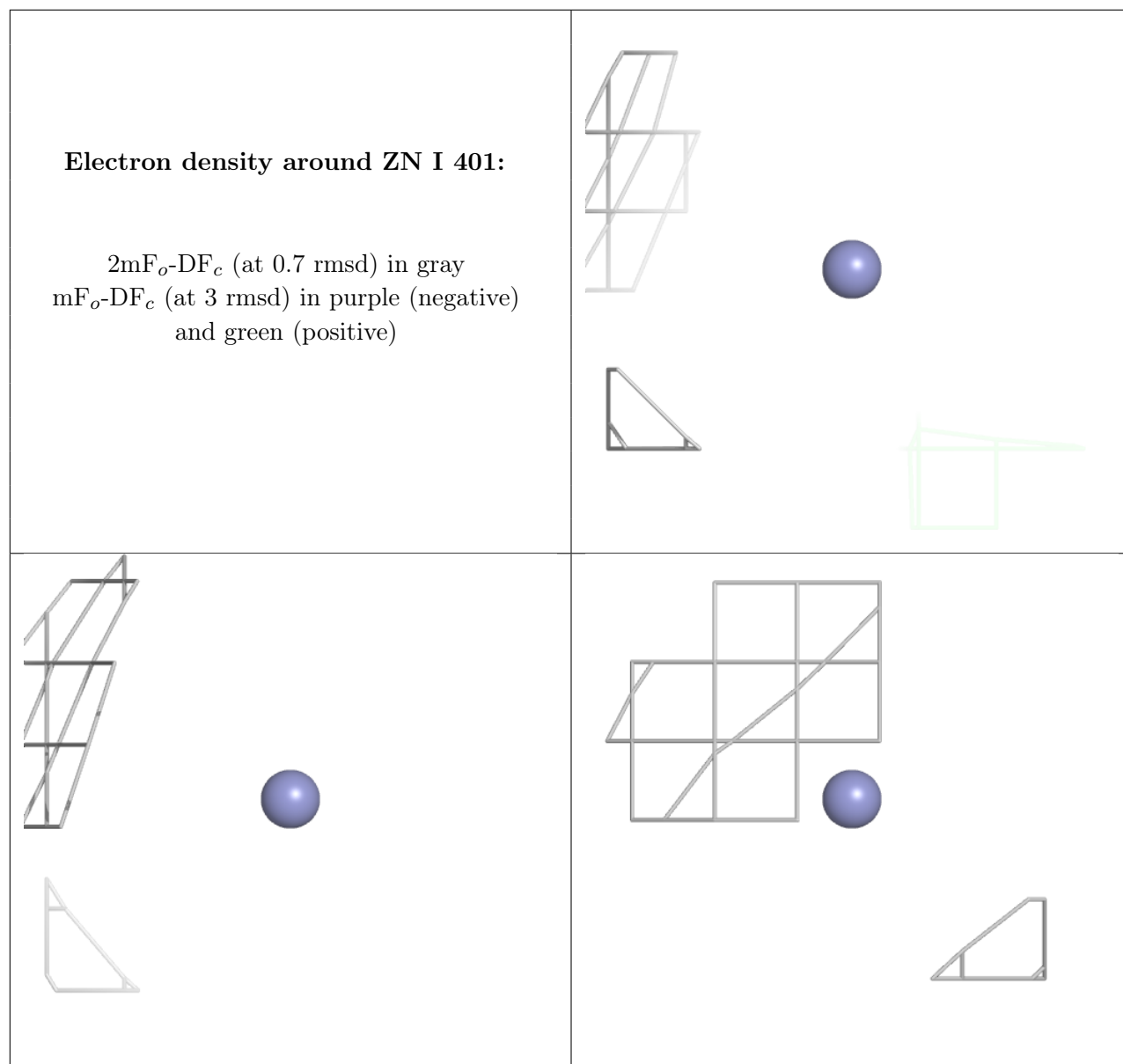
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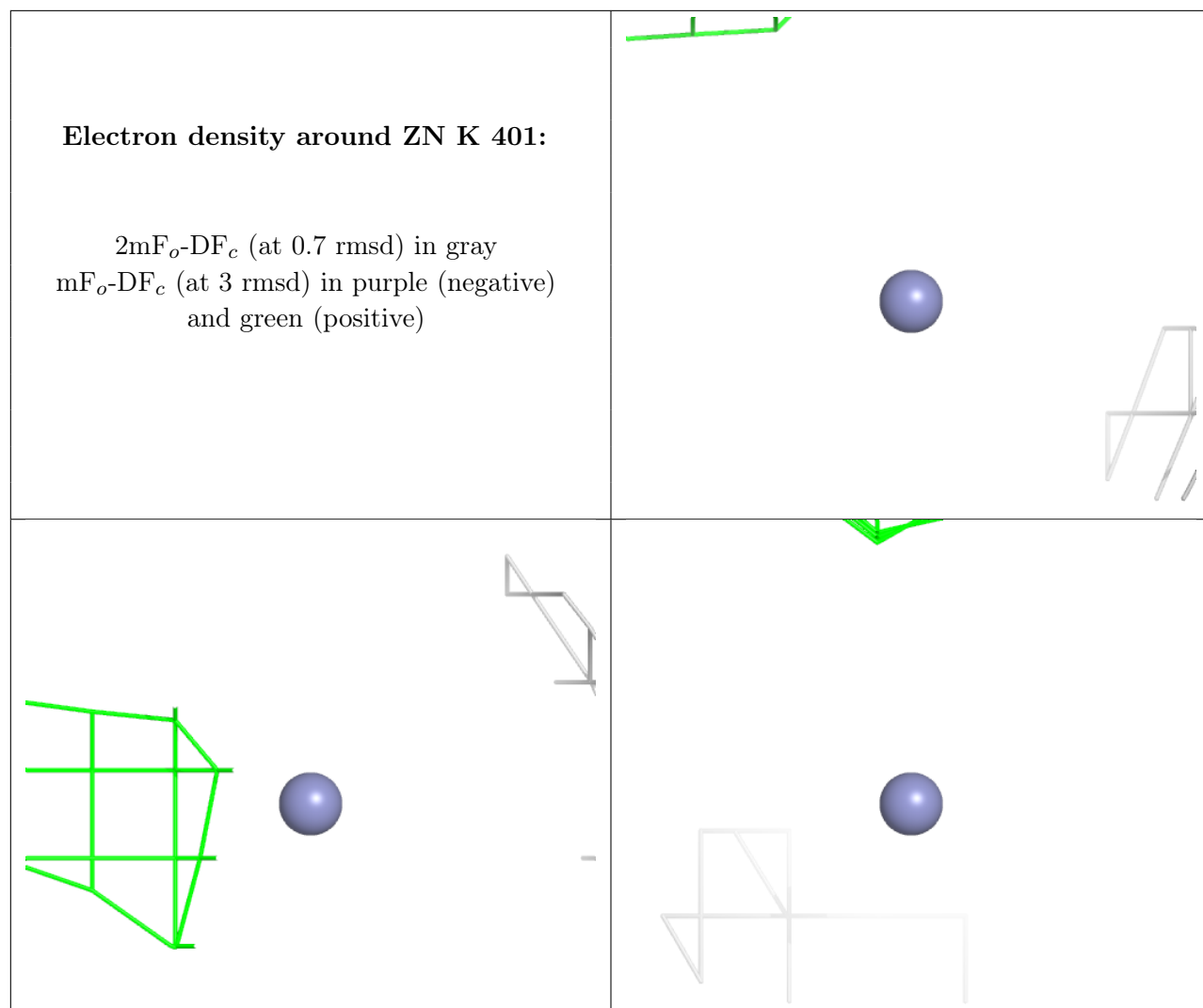
$2mF_o-DF_c$ (at 0.7 rnsd) in gray
 mF_o-DF_c (at 3 rnsd) in purple (negative)
and green (positive)

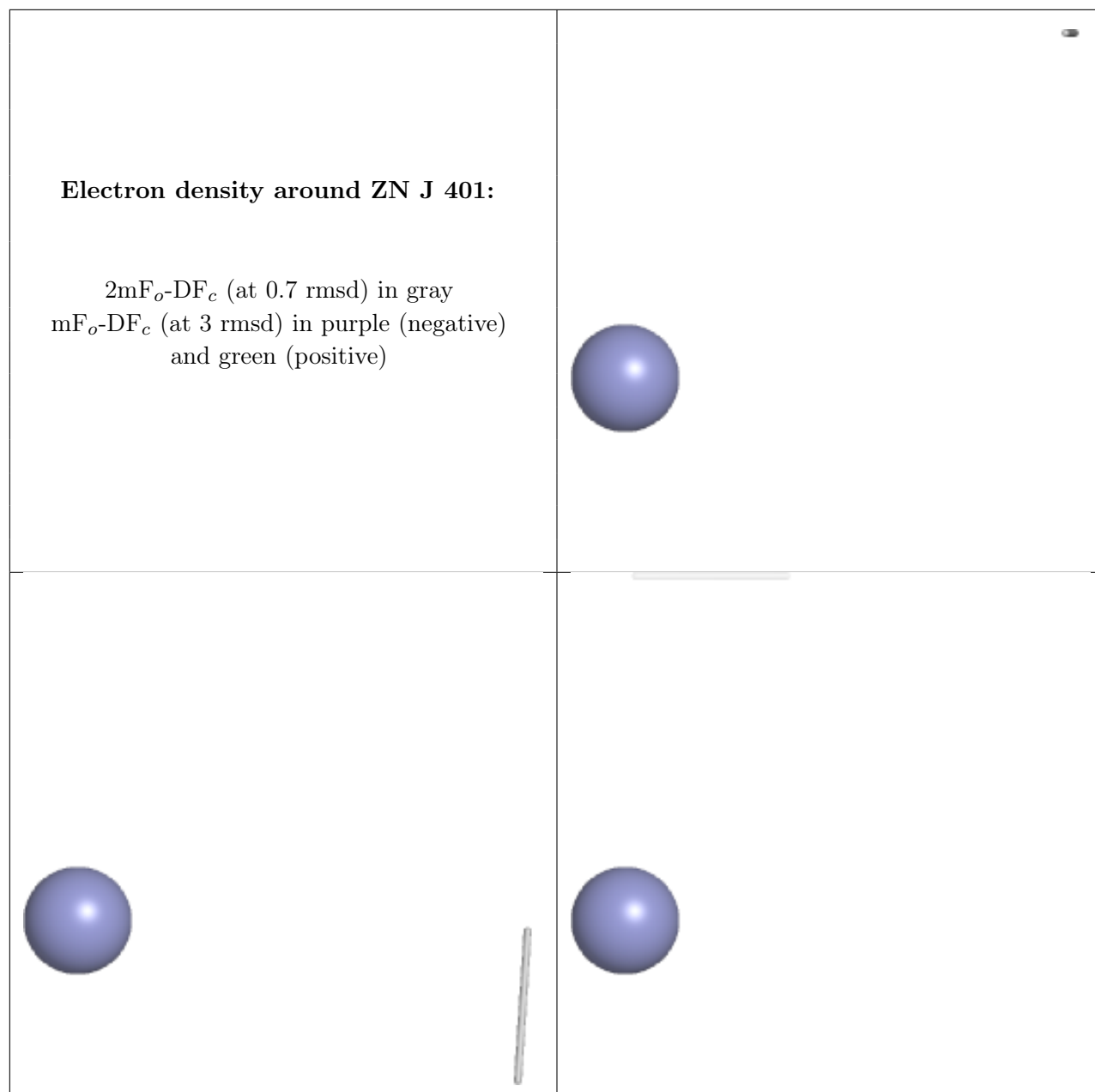


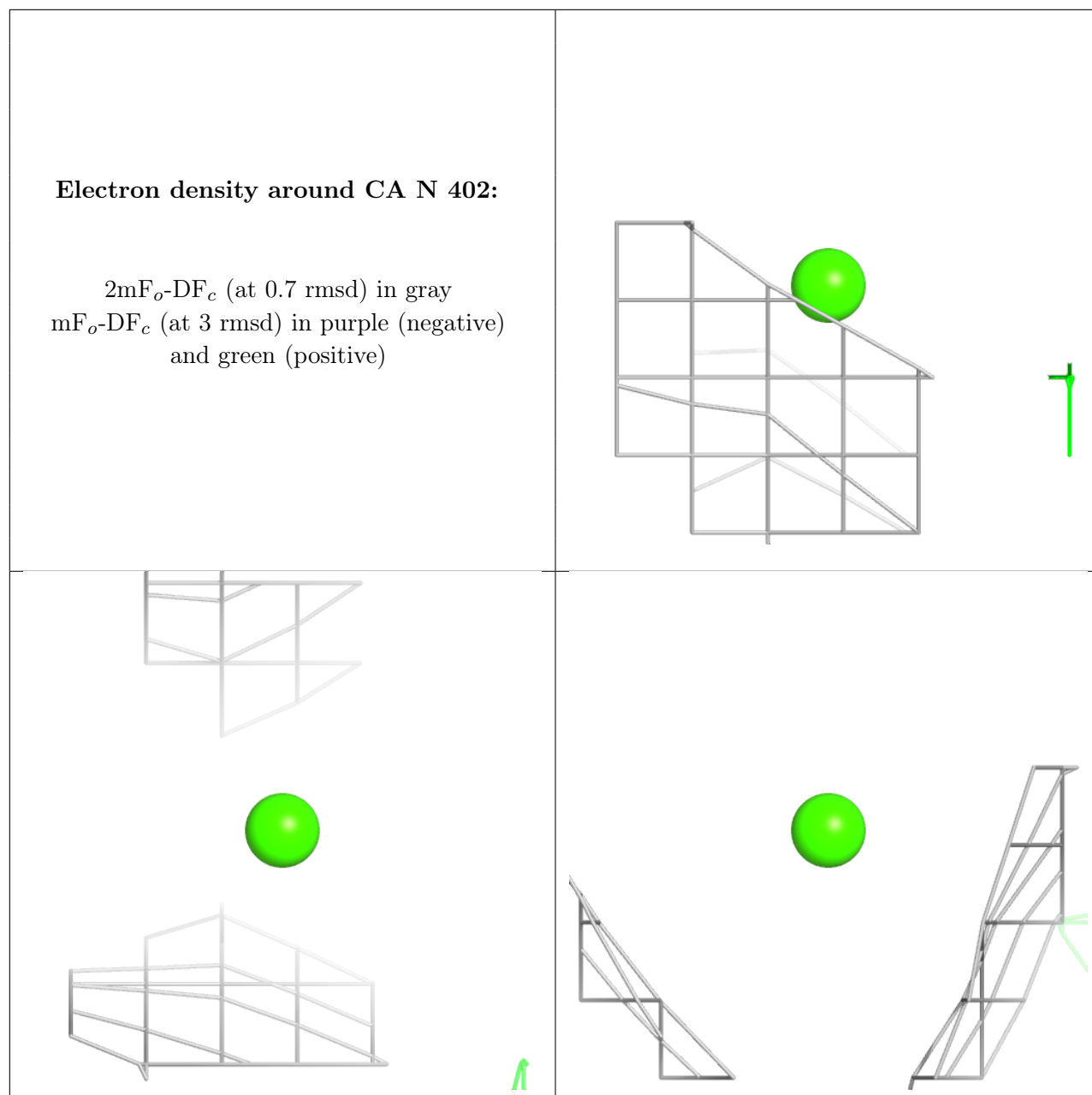












6.5 Other polymers [i](#)

There are no such residues in this entry.