



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 21, 2021 – 12:04 PM EDT

PDB ID : 7JS3
Title : Structure of the Class II Fructose-1,6-Bisphosphatase from *Francisella tularensis*
Authors : Abad-Zapatero, C.; Wolf, N.M.; Movahedzadeh, F.; Gutka, H.; Selezneva, A.
Deposited on : 2020-08-13
Resolution : 2.40 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.20
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) : 2.20

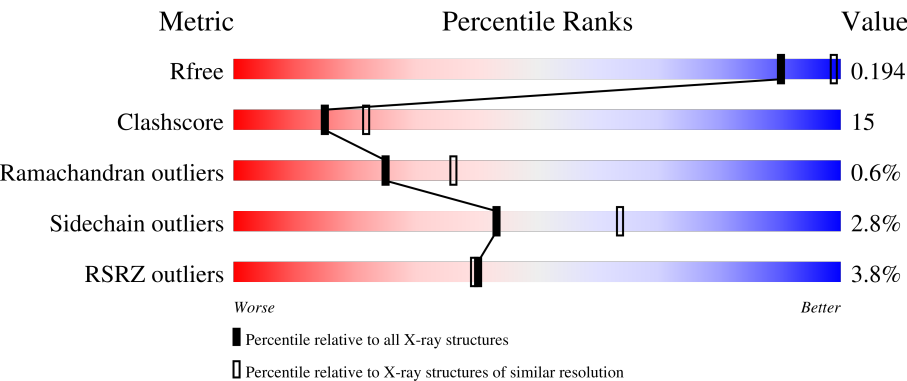
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	348	<div><div>8%</div><div>70%22%7%</div></div>
1	B	348	<div><div>73%19%7%</div></div>
1	C	348	<div><div>8%</div><div>54%38%7%</div></div>
1	D	348	<div><div>4%</div><div>65%27%7%</div></div>

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

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	B	402	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9838 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 Fructose-1,6-bisphosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	1	0
			2409	1494	425	474	16			
1	B	325	Total	C	N	O	S	0	0	0
			2406	1492	425	474	15			
1	C	325	Total	C	N	O	S	0	1	0
			2412	1496	425	476	15			
1	D	325	Total	C	N	O	S	0	0	0
			2406	1492	425	474	15			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A0E2ZJY0
A	-18	GLY	-	expression tag	UNP A0A0E2ZJY0
A	-17	SER	-	expression tag	UNP A0A0E2ZJY0
A	-16	SER	-	expression tag	UNP A0A0E2ZJY0
A	-15	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-14	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-13	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-12	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-11	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-10	HIS	-	expression tag	UNP A0A0E2ZJY0
A	-9	SER	-	expression tag	UNP A0A0E2ZJY0
A	-8	SER	-	expression tag	UNP A0A0E2ZJY0
A	-7	GLY	-	expression tag	UNP A0A0E2ZJY0
A	-6	LEU	-	expression tag	UNP A0A0E2ZJY0
A	-5	VAL	-	expression tag	UNP A0A0E2ZJY0
A	-4	PRO	-	expression tag	UNP A0A0E2ZJY0
A	-3	ARG	-	expression tag	UNP A0A0E2ZJY0
A	-2	GLY	-	expression tag	UNP A0A0E2ZJY0
A	-1	SER	-	expression tag	UNP A0A0E2ZJY0
A	0	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-19	MET	-	initiating methionine	UNP A0A0E2ZJY0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP A0A0E2ZJY0
B	-17	SER	-	expression tag	UNP A0A0E2ZJY0
B	-16	SER	-	expression tag	UNP A0A0E2ZJY0
B	-15	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-14	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-13	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-12	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-11	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-10	HIS	-	expression tag	UNP A0A0E2ZJY0
B	-9	SER	-	expression tag	UNP A0A0E2ZJY0
B	-8	SER	-	expression tag	UNP A0A0E2ZJY0
B	-7	GLY	-	expression tag	UNP A0A0E2ZJY0
B	-6	LEU	-	expression tag	UNP A0A0E2ZJY0
B	-5	VAL	-	expression tag	UNP A0A0E2ZJY0
B	-4	PRO	-	expression tag	UNP A0A0E2ZJY0
B	-3	ARG	-	expression tag	UNP A0A0E2ZJY0
B	-2	GLY	-	expression tag	UNP A0A0E2ZJY0
B	-1	SER	-	expression tag	UNP A0A0E2ZJY0
B	0	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-19	MET	-	initiating methionine	UNP A0A0E2ZJY0
C	-18	GLY	-	expression tag	UNP A0A0E2ZJY0
C	-17	SER	-	expression tag	UNP A0A0E2ZJY0
C	-16	SER	-	expression tag	UNP A0A0E2ZJY0
C	-15	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-14	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-13	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-12	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-11	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-10	HIS	-	expression tag	UNP A0A0E2ZJY0
C	-9	SER	-	expression tag	UNP A0A0E2ZJY0
C	-8	SER	-	expression tag	UNP A0A0E2ZJY0
C	-7	GLY	-	expression tag	UNP A0A0E2ZJY0
C	-6	LEU	-	expression tag	UNP A0A0E2ZJY0
C	-5	VAL	-	expression tag	UNP A0A0E2ZJY0
C	-4	PRO	-	expression tag	UNP A0A0E2ZJY0
C	-3	ARG	-	expression tag	UNP A0A0E2ZJY0
C	-2	GLY	-	expression tag	UNP A0A0E2ZJY0
C	-1	SER	-	expression tag	UNP A0A0E2ZJY0
C	0	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-19	MET	-	initiating methionine	UNP A0A0E2ZJY0
D	-18	GLY	-	expression tag	UNP A0A0E2ZJY0
D	-17	SER	-	expression tag	UNP A0A0E2ZJY0

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP A0A0E2ZJY0
D	-15	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-14	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-13	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-12	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-11	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-10	HIS	-	expression tag	UNP A0A0E2ZJY0
D	-9	SER	-	expression tag	UNP A0A0E2ZJY0
D	-8	SER	-	expression tag	UNP A0A0E2ZJY0
D	-7	GLY	-	expression tag	UNP A0A0E2ZJY0
D	-6	LEU	-	expression tag	UNP A0A0E2ZJY0
D	-5	VAL	-	expression tag	UNP A0A0E2ZJY0
D	-4	PRO	-	expression tag	UNP A0A0E2ZJY0
D	-3	ARG	-	expression tag	UNP A0A0E2ZJY0
D	-2	GLY	-	expression tag	UNP A0A0E2ZJY0
D	-1	SER	-	expression tag	UNP A0A0E2ZJY0
D	0	HIS	-	expression tag	UNP A0A0E2ZJY0

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

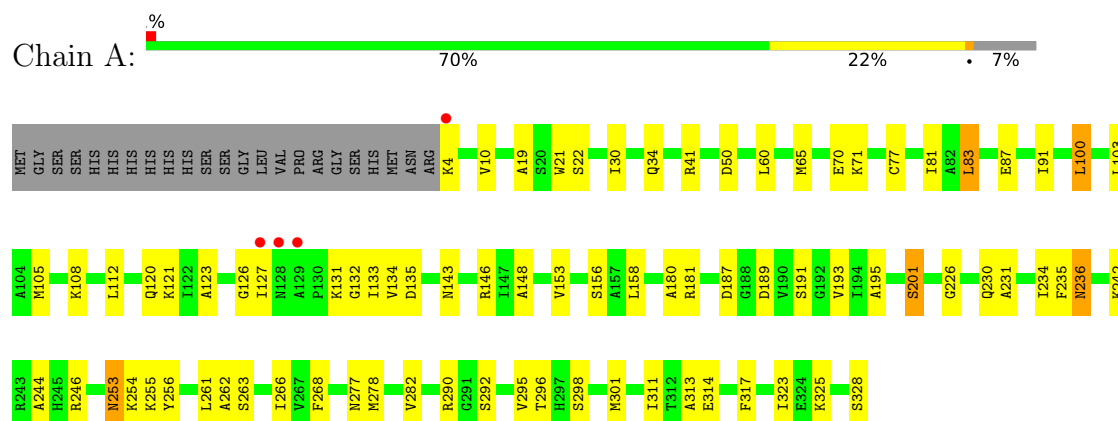
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	59	Total	O	0	0
			59	59		
4	B	55	Total	O	0	0
			55	55		
4	C	21	Total	O	0	0
			21	21		
4	D	30	Total	O	0	0
			30	30		

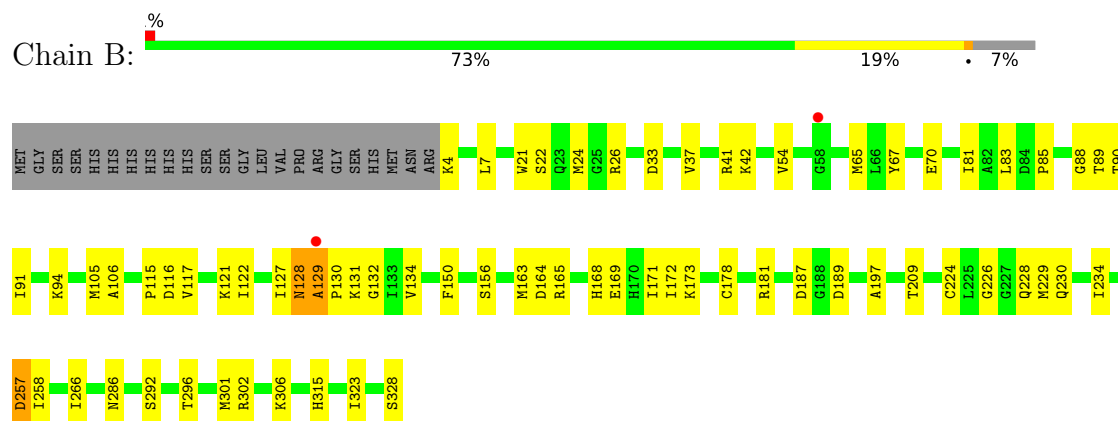
3 Residue-property plots [i](#)

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.

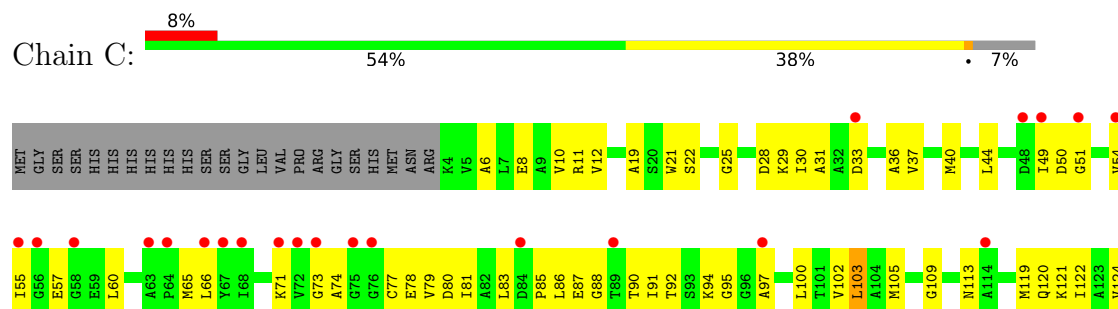
• Molecule 1: Fructose-1,6-bisphosphatase

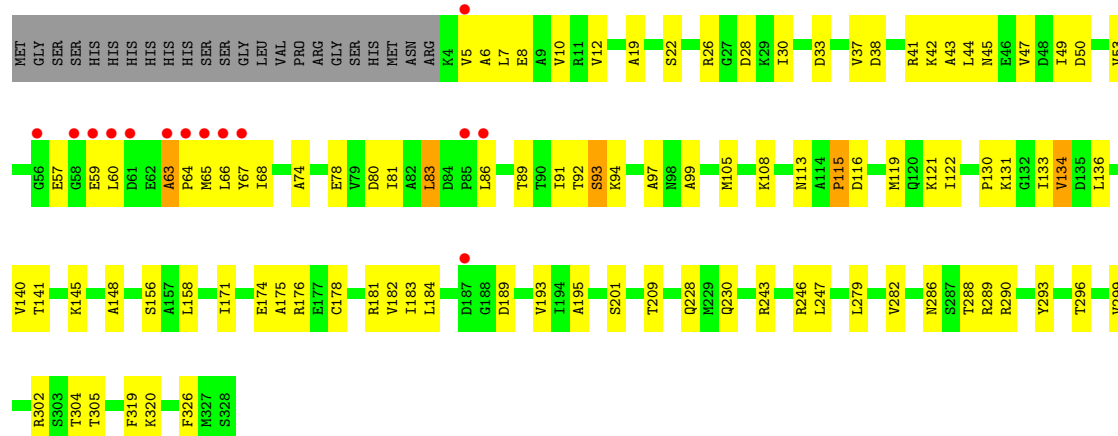


• Molecule 1: Fructose-1,6-bisphosphatase



• Molecule 1: Fructose-1,6-bisphosphatase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.30Å 100.17Å 92.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.11 – 2.40 38.11 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.7 (38.11-2.40) 98.8 (38.11-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0258, PHENIX 1.16-3549	Depositor
R, R_{free}	0.153 , 0.193 0.154 , 0.194	Depositor DCC
R_{free} test set	2732 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	42.8	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.417 for h,-k,-l	Xtriage
Reported twinning fraction	0.568 for H, K, L 0.432 for -h,-k,l	Depositor
Outliers	0 of 53663 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9838	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.80	1/2438 (0.0%)	0.78	1/3290 (0.0%)
1	B	0.74	0/2432	0.74	0/3282
1	C	0.70	0/2441	0.74	0/3294
1	D	0.71	0/2432	0.76	0/3282
All	All	0.74	1/9743 (0.0%)	0.76	1/13148 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	C	0	3
1	D	0	1
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	87	GLU	CD-OE2	-6.07	1.19	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	100	LEU	CB-CG-CD1	-5.08	102.36	111.00

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	236	ASN	Peptide
1	A	83	LEU	Peptide
1	C	109	GLY	Peptide
1	C	186	ASN	Peptide
1	C	83	LEU	Peptide
1	D	83	LEU	Peptide

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2409	0	2430	63	0
1	B	2406	0	2425	67	0
1	C	2412	0	2431	109	0
1	D	2406	0	2425	71	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	6	0	8	1	0
3	B	6	0	8	0	0
3	C	6	0	8	0	0
3	D	18	0	24	2	0
4	A	59	0	0	4	0
4	B	55	0	0	2	0
4	C	21	0	0	1	0
4	D	30	0	0	3	0
All	All	9838	0	9759	294	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 15.

All (294) 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:B:83:LEU:HD12	1:B:85:PRO:HD3	1.43	0.98
1:A:127:ILE:HG12	1:A:226:GLY:HA3	1.50	0.93
1:B:83:LEU:CD1	1:B:85:PRO:HD3	1.99	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:121:LYS:NZ	1:C:189:ASP:OD2	2.06	0.89
1:B:22:SER:OG	1:B:323:ILE:HG22	1.73	0.88
1:C:222:LEU:HD22	1:C:227:GLY:HA3	1.55	0.88
1:C:37:VAL:HB	1:C:66:LEU:HD12	1.55	0.87
1:B:127:ILE:HD13	1:B:306:LYS:HE3	1.54	0.87
1:B:127:ILE:CD1	1:B:306:LYS:HE3	2.10	0.81
1:D:289:ARG:NH1	4:D:502:HOH:O	2.11	0.81
1:D:148:ALA:HB2	1:D:158:LEU:HD11	1.62	0.81
1:A:290:ARG:NH1	1:A:314:GLU:OE2	2.11	0.81
1:C:94:LYS:HB2	1:D:184:LEU:HD23	1.64	0.78
1:C:119:MET:HE2	1:C:121:LYS:HE2	1.66	0.78
1:A:131:LYS:HD2	1:A:132:GLY:H	1.51	0.76
1:D:7:LEU:O	1:D:10:VAL:HG12	1.86	0.76
1:C:29:LYS:HD3	1:C:90:THR:HG22	1.67	0.75
1:C:19:ALA:O	1:C:22:SER:OG	2.04	0.75
1:B:197:ALA:HA	1:C:289:ARG:HD3	1.68	0.75
1:A:100:LEU:HD11	1:A:278:MET:CE	2.18	0.73
1:D:130:PRO:HD2	1:D:133:ILE:HD12	1.72	0.72
1:A:121:LYS:NZ	1:A:189:ASP:OD2	2.23	0.72
1:D:33:ASP:HB2	1:D:89:THR:HG21	1.71	0.71
1:D:279:LEU:HD21	1:D:299:VAL:HG23	1.72	0.71
1:A:50:ASP:OD1	1:A:71:LYS:NZ	2.15	0.70
1:C:193:VAL:O	1:C:196:THR:OG1	2.04	0.70
1:C:160:VAL:HB	1:C:182:VAL:HG12	1.74	0.69
1:C:94:LYS:HB2	1:D:184:LEU:CD2	2.23	0.69
1:C:127:ILE:HB	1:C:226:GLY:HA3	1.76	0.68
1:C:290:ARG:HH12	1:C:314:GLU:HB2	1.59	0.67
1:B:37:VAL:HG22	1:B:85:PRO:HG3	1.76	0.67
1:C:25:GLY:HA2	1:C:94:LYS:HA	1.76	0.67
1:C:37:VAL:HB	1:C:66:LEU:CD1	2.23	0.67
1:C:290:ARG:NH1	1:C:314:GLU:HB2	2.10	0.67
1:A:91:ILE:HD11	1:A:187:ASP:HB3	1.78	0.65
1:A:292:SER:OG	1:A:317:PHE:O	2.14	0.65
1:B:130:PRO:O	1:B:228:GLN:NE2	2.30	0.65
1:C:54:VAL:HG23	1:C:55:ILE:HG13	1.79	0.65
1:D:5:VAL:HG23	1:D:8:GLU:HB2	1.76	0.65
1:C:133:ILE:HG13	1:C:134:VAL:N	2.12	0.65
1:A:83:LEU:HD23	1:A:103:LEU:HD13	1.78	0.64
1:A:148:ALA:HB2	1:A:158:LEU:HD11	1.79	0.64
1:C:33:ASP:O	1:C:37:VAL:HG13	1.98	0.64
1:D:243:ARG:HG2	1:D:246:ARG:HH21	1.63	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:CYS:HB2	1:B:301:MET:HE3	1.79	0.63
1:C:124:VAL:HG11	1:C:133:ILE:HD11	1.79	0.63
1:C:239:GLU:O	1:C:242:LYS:HB2	1.99	0.63
1:A:100:LEU:HD11	1:A:278:MET:HE3	1.80	0.63
1:B:128:ASN:HB2	1:B:150:PHE:HE1	1.65	0.62
1:C:37:VAL:HG12	1:C:85:PRO:CG	2.29	0.62
1:C:50:ASP:HB3	1:C:78:GLU:HG2	1.81	0.62
1:C:290:ARG:O	1:C:290:ARG:HG3	1.99	0.62
1:B:21:TRP:HD1	1:B:24:MET:HE3	1.63	0.62
1:A:100:LEU:HD11	1:A:278:MET:HE1	1.82	0.61
1:A:234:ILE:HG22	1:A:236:ASN:ND2	2.14	0.61
1:A:234:ILE:CG2	1:A:236:ASN:ND2	2.64	0.60
1:A:41:ARG:NH2	1:A:70:GLU:OE2	2.33	0.60
1:D:53:VAL:HG13	1:D:83:LEU:HD13	1.81	0.60
1:B:121:LYS:NZ	1:B:189:ASP:OD2	2.34	0.60
1:D:122:ILE:HG12	1:D:134:VAL:HG21	1.84	0.60
1:D:91:ILE:HA	1:D:94:LYS:HB3	1.83	0.59
1:D:47:VAL:HG12	1:D:49:ILE:HG12	1.84	0.59
1:D:63:ALA:HB3	1:D:64:PRO:HD3	1.83	0.59
1:C:113:ASN:O	1:C:247:LEU:HD21	2.02	0.59
1:B:169:GLU:O	1:B:173:LYS:HG2	2.03	0.59
1:C:153:VAL:HG23	1:D:319:PHE:HB3	1.85	0.59
1:D:59:GLU:O	1:D:67:TYR:HB2	2.03	0.59
1:A:235:PHE:CE1	1:A:244:ALA:HB2	2.37	0.58
1:B:164:ASP:O	1:B:165:ARG:HD3	2.03	0.58
1:A:180:ALA:O	1:B:26:ARG:NH1	2.35	0.58
1:D:41:ARG:NH2	1:D:65:MET:O	2.37	0.58
1:C:79:VAL:HG23	1:C:81:ILE:HD11	1.84	0.58
1:A:133:ILE:HG23	1:A:134:VAL:HG13	1.86	0.58
1:C:266:ILE:HB	1:C:303:SER:HB3	1.85	0.58
1:B:22:SER:OG	1:B:323:ILE:O	2.22	0.57
1:D:326:PHE:O	4:D:503:HOH:O	2.17	0.57
1:D:57:GLU:O	1:D:64:PRO:HG3	2.05	0.57
1:C:49:ILE:HD11	1:C:77:CYS:HB2	1.87	0.57
1:D:53:VAL:O	1:D:68:ILE:HA	2.04	0.57
1:B:128:ASN:HB2	1:B:150:PHE:CE1	2.39	0.57
1:A:126:GLY:HA2	1:D:289:ARG:HG2	1.86	0.57
1:B:115:PRO:O	1:B:117:VAL:HG12	2.05	0.57
1:C:144:LEU:CB	1:C:155:MET:HE3	2.34	0.57
1:C:144:LEU:HB2	1:C:155:MET:HE3	1.86	0.56
1:B:7:LEU:HD12	1:C:11:ARG:HE	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:SER:HA	1:B:323:ILE:CG2	2.36	0.56
1:C:54:VAL:HG11	1:C:80:ASP:HB3	1.87	0.56
1:B:229:MET:HB3	1:B:258:ILE:HG13	1.88	0.56
1:C:164:ASP:N	1:C:185:ILE:O	2.35	0.56
1:A:127:ILE:O	1:A:127:ILE:HG22	2.06	0.55
1:C:95:GLY:HA3	1:D:183:ILE:HB	1.87	0.55
1:C:235:PHE:CE2	1:C:252:LEU:HG	2.41	0.55
1:A:131:LYS:N	4:A:506:HOH:O	2.37	0.55
1:D:81:ILE:HD12	1:D:105:MET:HG2	1.89	0.54
1:B:302:ARG:O	1:B:306:LYS:HA	2.06	0.54
1:C:165:ARG:HG3	1:C:167:ARG:CZ	2.38	0.54
1:B:33:ASP:O	1:B:37:VAL:HG23	2.08	0.54
1:B:106:ALA:HB2	1:B:266:ILE:HD13	1.89	0.54
1:A:131:LYS:HD2	1:A:132:GLY:N	2.21	0.54
1:C:119:MET:HE3	1:C:121:LYS:HG2	1.90	0.54
1:D:45:ASN:O	1:D:74:ALA:HB2	2.08	0.54
1:A:290:ARG:NH2	4:A:510:HOH:O	2.41	0.54
1:B:90:THR:O	1:B:94:LYS:HG3	2.08	0.54
1:B:257:ASP:OD1	1:B:257:ASP:N	2.41	0.54
1:B:54:VAL:HG12	1:B:81:ILE:O	2.07	0.54
1:D:176:ARG:HE	1:D:182:VAL:HB	1.73	0.54
1:B:21:TRP:HD1	1:B:24:MET:CE	2.20	0.53
1:C:290:ARG:NH1	1:C:314:GLU:CB	2.70	0.53
1:A:41:ARG:HD3	1:A:65:MET:CE	2.39	0.53
1:C:221:ALA:HB2	1:C:268:PHE:CE2	2.43	0.53
1:D:304:THR:OG1	1:D:305:THR:N	2.40	0.53
1:B:83:LEU:HD12	1:B:85:PRO:CD	2.27	0.53
1:A:21:TRP:CZ3	1:A:323:ILE:HD11	2.43	0.53
1:B:171:ILE:HG13	1:B:172:ILE:N	2.25	0.52
1:C:222:LEU:HD22	1:C:227:GLY:CA	2.34	0.52
1:B:21:TRP:CD1	1:B:24:MET:HE3	2.43	0.52
1:A:135:ASP:OD1	1:A:255:LYS:NZ	2.42	0.52
1:B:127:ILE:CD1	1:B:306:LYS:CE	2.87	0.52
1:C:37:VAL:HG12	1:C:85:PRO:HG2	1.90	0.52
1:A:195:ALA:HB1	1:A:201:SER:OG	2.09	0.52
1:C:147:ILE:HD11	1:C:205:VAL:HG21	1.91	0.52
1:A:242:LYS:O	1:A:246:ARG:HG2	2.10	0.52
1:B:42:LYS:HD3	1:B:328:SER:HB3	1.92	0.52
1:C:12:VAL:HG11	1:C:103:LEU:HD13	1.92	0.52
1:C:80:ASP:O	1:C:105:MET:HA	2.08	0.52
1:D:108:LYS:HD3	1:D:108:LYS:C	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ILE:CG1	1:A:226:GLY:HA3	2.32	0.51
1:C:176:ARG:HD3	1:D:26:ARG:HA	1.92	0.51
1:D:286:ASN:HB2	4:D:505:HOH:O	2.09	0.51
1:A:41:ARG:NH1	1:A:65:MET:HA	2.24	0.51
1:C:30:ILE:HG13	1:C:31:ALA:N	2.26	0.51
1:C:21:TRP:O	1:D:181:ARG:NH2	2.43	0.51
1:C:190:VAL:HG22	1:C:214:GLU:HB3	1.92	0.51
1:C:173:LYS:O	1:C:177:GLU:HG3	2.11	0.51
1:D:30:ILE:HD12	1:D:30:ILE:H	1.76	0.51
1:C:254:LYS:HD3	1:C:256:TYR:CE2	2.45	0.51
1:B:156:SER:O	1:B:181:ARG:HD2	2.11	0.51
1:A:81:ILE:HD13	1:A:105:MET:HG2	1.93	0.50
1:D:156:SER:O	1:D:181:ARG:HD2	2.12	0.50
1:D:279:LEU:HD21	1:D:299:VAL:CG2	2.40	0.50
1:C:44:LEU:O	1:C:73:GLY:HA2	2.11	0.50
1:D:86:LEU:HD21	1:D:92:THR:OG1	2.10	0.50
1:D:86:LEU:HD23	1:D:89:THR:HA	1.92	0.50
1:C:143:ASN:O	1:C:147:ILE:HG23	2.11	0.50
1:D:171:ILE:HG21	3:D:403:GOL:O3	2.12	0.50
1:B:122:ILE:HG12	1:B:134:VAL:CG2	2.42	0.50
1:D:41:ARG:HE	1:D:66:LEU:HD12	1.77	0.50
1:D:175:ALA:O	1:D:178:CYS:HB2	2.12	0.50
1:A:30:ILE:HD12	4:A:531:HOH:O	2.11	0.49
1:B:37:VAL:HG22	1:B:85:PRO:CG	2.40	0.49
1:A:143:ASN:OD1	1:A:146:ARG:NH2	2.46	0.49
1:B:88:GLY:HA2	4:B:547:HOH:O	2.12	0.49
1:B:127:ILE:HD11	1:B:306:LYS:CE	2.42	0.49
1:D:53:VAL:HG13	1:D:83:LEU:CD1	2.42	0.49
1:C:184:LEU:HD13	4:C:510:HOH:O	2.12	0.49
1:D:140:VAL:HB	1:D:174:GLU:OE2	2.12	0.49
1:B:127:ILE:HD11	1:B:306:LYS:HZ1	1.78	0.48
1:D:37:VAL:HG21	1:D:65:MET:HB2	1.95	0.48
1:A:234:ILE:HG22	1:A:236:ASN:HD22	1.79	0.48
1:A:253:ASN:N	1:A:253:ASN:OD1	2.46	0.48
1:B:129:ALA:N	1:B:130:PRO:CD	2.76	0.48
1:C:57:GLU:HB2	1:C:66:LEU:HB3	1.94	0.48
1:D:80:ASP:CG	1:D:108:LYS:HA	2.33	0.48
1:B:91:ILE:HD11	1:B:187:ASP:HB3	1.95	0.48
1:C:245:HIS:HA	1:C:249:ILE:O	2.14	0.48
1:C:266:ILE:O	1:C:302:ARG:HA	2.13	0.48
1:B:127:ILE:HD12	1:C:290:ARG:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:ARG:HB3	1:C:167:ARG:HG2	1.96	0.48
1:D:91:ILE:HG13	1:D:99:ALA:HB2	1.95	0.48
1:A:19:ALA:O	1:A:22:SER:HB2	2.14	0.48
1:C:159:VAL:CG2	1:C:181:ARG:NH2	2.77	0.48
1:A:181:ARG:HG2	1:B:24:MET:O	2.14	0.48
1:B:41:ARG:NH1	1:B:65:MET:HG2	2.28	0.48
1:B:226:GLY:O	4:B:503:HOH:O	2.20	0.48
1:B:22:SER:HA	1:B:323:ILE:HG21	1.94	0.48
1:C:238:GLU:HG3	1:C:241:ILE:HB	1.96	0.48
1:A:234:ILE:CG2	1:A:236:ASN:HD22	2.27	0.47
1:B:163:MET:HB2	1:B:168:HIS:CE1	2.50	0.47
1:D:41:ARG:HB2	1:D:66:LEU:HD13	1.97	0.47
1:B:4:LYS:HZ2	1:B:4:LYS:N	2.11	0.47
1:D:130:PRO:O	1:D:228:GLN:HG3	2.14	0.47
1:B:131:LYS:HG2	1:B:132:GLY:H	1.80	0.47
1:C:295:VAL:HA	1:C:313:ALA:O	2.14	0.47
1:A:230:GLN:OE1	1:A:255:LYS:HD2	2.15	0.47
1:B:209:THR:HG21	1:B:234:ILE:HD11	1.97	0.47
1:C:234:ILE:HG22	1:C:235:PHE:H	1.79	0.47
1:C:122:ILE:HD13	1:C:134:VAL:HB	1.97	0.47
1:C:211:GLY:HA3	1:C:214:GLU:OE1	2.14	0.47
1:B:229:MET:HB3	1:B:258:ILE:CG1	2.45	0.46
1:C:120:GLN:OE1	1:C:136:LEU:HB3	2.15	0.46
1:D:290:ARG:HD3	1:D:293:TYR:CE1	2.51	0.46
1:B:163:MET:HB2	1:B:168:HIS:HE1	1.81	0.46
1:C:137:ASP:OD2	1:C:255:LYS:NZ	2.30	0.46
1:D:131:LYS:HE3	1:D:131:LYS:HB2	1.81	0.46
1:C:325:LYS:HG2	1:C:325:LYS:O	2.16	0.45
1:A:41:ARG:CZ	1:A:70:GLU:OE2	2.64	0.45
1:B:22:SER:HA	1:B:323:ILE:HG22	1.99	0.45
1:B:286:ASN:HA	1:B:292:SER:HA	1.98	0.45
1:C:139:SER:OG	1:C:142:ASN:HB2	2.16	0.45
1:A:254:LYS:HA	3:A:403:GOL:H31	1.98	0.45
1:A:195:ALA:HB2	1:A:277:ASN:HB3	1.97	0.45
1:C:155:MET:HE1	1:C:178:CYS:SG	2.57	0.45
1:A:254:LYS:HD3	1:A:256:TYR:CE2	2.52	0.45
1:C:159:VAL:CG2	1:C:181:ARG:HH21	2.30	0.45
1:A:4:LYS:HD3	1:D:8:GLU:HG3	1.98	0.45
1:A:34:GLN:HG3	4:A:558:HOH:O	2.16	0.45
1:C:102:VAL:HG12	1:C:270:ALA:HB2	1.97	0.45
1:B:127:ILE:HG13	1:C:290:ARG:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:MET:O	1:C:209:THR:HA	2.16	0.44
1:C:139:SER:O	1:C:143:ASN:ND2	2.45	0.44
1:C:317:PHE:CE2	1:C:319:PHE:HA	2.53	0.44
1:B:105:MET:HB2	1:B:105:MET:HE2	1.62	0.44
1:B:115:PRO:HG2	1:B:117:VAL:HG11	1.99	0.44
1:D:289:ARG:HE	1:D:289:ARG:HB3	1.57	0.44
1:A:123:ALA:CB	1:A:193:VAL:HG13	2.48	0.44
1:B:67:TYR:CE1	1:B:70:GLU:HB2	2.53	0.44
1:B:296:THR:HG23	1:B:315:HIS:CD2	2.53	0.44
1:C:49:ILE:O	1:C:73:GLY:HA3	2.18	0.44
1:A:41:ARG:HD3	1:A:65:MET:HE3	2.00	0.44
1:C:129:ALA:HA	1:C:131:LYS:HE2	1.99	0.44
1:C:124:VAL:HG11	1:C:133:ILE:CD1	2.47	0.44
1:B:131:LYS:HG2	1:B:132:GLY:N	2.32	0.44
1:C:155:MET:CE	1:C:178:CYS:SG	3.06	0.43
1:D:189:ASP:O	1:D:193:VAL:HG23	2.18	0.43
1:A:189:ASP:OD1	1:A:189:ASP:N	2.51	0.43
1:A:262:ALA:HB1	1:A:266:ILE:HD13	1.99	0.43
1:D:141:THR:O	1:D:145:LYS:HG3	2.18	0.43
1:A:156:SER:O	1:A:181:ARG:HD2	2.18	0.43
1:C:317:PHE:HE2	1:C:319:PHE:HA	1.84	0.43
1:B:127:ILE:HG12	1:B:226:GLY:HA3	2.00	0.43
1:C:57:GLU:HG3	1:C:66:LEU:HD13	1.99	0.43
1:D:44:LEU:HD23	1:D:44:LEU:HA	1.79	0.43
1:C:159:VAL:HG22	1:C:181:ARG:NH2	2.34	0.43
1:A:246:ARG:HA	1:A:246:ARG:NE	2.33	0.43
1:C:293:TYR:CE2	1:C:314:GLU:HG2	2.53	0.43
1:D:50:ASP:HB3	1:D:78:GLU:HA	2.01	0.43
1:D:91:ILE:HD12	1:D:97:ALA:O	2.17	0.43
1:C:37:VAL:HG12	1:C:85:PRO:HG3	1.99	0.43
1:C:235:PHE:CZ	1:C:252:LEU:HG	2.53	0.43
1:D:28:ASP:N	1:D:93:SER:OG	2.51	0.43
1:C:284:ARG:HG3	1:C:319:PHE:CE2	2.53	0.43
1:D:86:LEU:HD23	1:D:89:THR:CA	2.49	0.43
1:C:10:VAL:HG11	1:C:311:ILE:HD12	2.01	0.43
1:C:222:LEU:O	1:C:226:GLY:N	2.47	0.43
1:D:136:LEU:HD12	1:D:230:GLN:HG2	2.01	0.43
1:A:317:PHE:CZ	1:A:325:LYS:HB2	2.54	0.42
1:C:147:ILE:HD11	1:C:205:VAL:CG2	2.49	0.42
1:B:33:ASP:CG	1:B:89:THR:HG21	2.40	0.42
1:C:28:ASP:HB3	1:C:31:ALA:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:282:VAL:HG22	1:D:296:THR:HG22	2.01	0.42
1:C:120:GLN:HG2	1:C:136:LEU:HD13	2.01	0.42
1:C:129:ALA:N	1:C:130:PRO:HD3	2.34	0.42
1:A:112:LEU:HD22	1:A:261:LEU:HD23	2.02	0.42
1:A:4:LYS:HG2	1:D:8:GLU:HB3	2.02	0.42
1:A:10:VAL:HG21	1:A:311:ILE:HD12	2.02	0.42
1:D:121:LYS:NZ	1:D:189:ASP:OD2	2.53	0.42
1:C:133:ILE:HG13	1:C:134:VAL:H	1.81	0.42
1:D:12:VAL:HG13	1:D:43:ALA:HB3	2.01	0.42
1:A:120:GLN:O	1:A:231:ALA:HA	2.20	0.42
1:A:296:THR:OG1	1:A:313:ALA:HB3	2.20	0.41
1:C:268:PHE:HB3	1:C:301:MET:HG3	2.01	0.41
1:D:122:ILE:HG12	1:D:134:VAL:CG2	2.49	0.41
1:C:170:HIS:CG	1:C:171:ILE:N	2.89	0.41
1:A:268:PHE:HB3	1:A:301:MET:HG3	2.02	0.41
1:C:49:ILE:O	1:C:74:ALA:N	2.45	0.41
1:A:189:ASP:O	1:A:193:VAL:HG23	2.20	0.41
1:A:191:SER:HA	1:A:278:MET:HE2	2.03	0.41
1:C:154:HIS:HB2	1:D:320:LYS:O	2.21	0.41
1:B:127:ILE:HD11	1:B:306:LYS:NZ	2.36	0.41
1:A:41:ARG:HH11	1:A:65:MET:HA	1.86	0.41
1:A:328:SER:O	1:A:328:SER:OG	2.38	0.41
1:B:129:ALA:H	1:B:130:PRO:HD3	1.85	0.41
1:C:100:LEU:HA	1:C:273:VAL:HG23	2.02	0.41
1:D:115:PRO:O	1:D:116:ASP:C	2.59	0.41
1:B:224:CYS:HB2	1:B:301:MET:CE	2.47	0.41
1:C:8:GLU:OE1	1:C:11:ARG:NH2	2.53	0.41
1:C:36:ALA:HB1	1:C:85:PRO:O	2.21	0.41
1:C:81:ILE:HD12	1:C:105:MET:HE1	2.03	0.41
1:C:87[A]:GLU:HG2	1:C:214:GLU:HG2	2.03	0.41
1:D:6:ALA:HB2	1:D:302:ARG:NH1	2.36	0.41
1:A:282:VAL:HA	1:A:295:VAL:O	2.21	0.41
1:B:115:PRO:HG2	1:B:117:VAL:CG1	2.51	0.41
1:C:86:LEU:HD21	1:C:92:THR:OG1	2.20	0.41
1:C:88:GLY:CA	1:C:91:ILE:HG12	2.51	0.40
1:C:91:ILE:HD12	1:C:97:ALA:O	2.20	0.40
1:D:136:LEU:HA	3:D:403:GOL:H12	2.03	0.40
1:C:51:GLY:O	1:C:71:LYS:HA	2.20	0.40
1:D:119:MET:O	1:D:209:THR:HA	2.21	0.40
1:C:65:MET:HG2	1:C:65:MET:O	2.21	0.40
1:C:120:GLN:HB3	1:C:234:ILE:HD11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:19:ALA:O	1:D:22:SER:HB2	2.21	0.40
1:D:195:ALA:O	1:D:201:SER:OG	2.23	0.40
1:B:94:LYS:HE2	1:B:94:LYS:HB3	1.93	0.40
1:D:290:ARG:HD3	1:D:293:TYR:CD1	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	324/348 (93%)	307 (95%)	17 (5%)	0	100	100
1	B	323/348 (93%)	298 (92%)	23 (7%)	2 (1%)	25	36
1	C	324/348 (93%)	287 (89%)	34 (10%)	3 (1%)	17	25
1	D	323/348 (93%)	296 (92%)	24 (7%)	3 (1%)	17	25
All	All	1294/1392 (93%)	1188 (92%)	98 (8%)	8 (1%)	25	36

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	116	ASP
1	C	60	LEU
1	D	63	ALA
1	C	6	ALA
1	C	172	ILE
1	D	115	PRO
1	B	129	ALA
1	D	60	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	252/271 (93%)	243 (96%)	9 (4%)	35	54
1	B	251/271 (93%)	247 (98%)	4 (2%)	62	79
1	C	252/271 (93%)	243 (96%)	9 (4%)	35	54
1	D	251/271 (93%)	244 (97%)	7 (3%)	43	63
All	All	1006/1084 (93%)	977 (97%)	29 (3%)	43	62

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LEU
1	A	77[A]	CYS
1	A	77[B]	CYS
1	A	108	LYS
1	A	153	VAL
1	A	201	SER
1	A	253	ASN
1	A	263	SER
1	A	298	SER
1	B	128	ASN
1	B	178	CYS
1	B	230	GLN
1	B	257	ASP
1	C	40	MET
1	C	103	LEU
1	C	178	CYS
1	C	184	LEU
1	C	223	LYS
1	C	250	THR
1	C	263	SER
1	C	302	ARG
1	C	312	THR
1	D	38	ASP
1	D	42	LYS

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Mol	Chain	Res	Type
1	D	93	SER
1	D	113	ASN
1	D	134	VAL
1	D	247	LEU
1	D	288	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	200	ASN
1	A	236	ASN
1	C	45	ASN
1	D	113	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 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	D	401	-	5,5,5	0.11	0	5,5,5	0.27	0
3	GOL	B	402	-	5,5,5	0.16	0	5,5,5	0.33	0
3	GOL	A	403	-	5,5,5	0.22	0	5,5,5	0.84	0
3	GOL	D	403	-	5,5,5	0.10	0	5,5,5	0.31	0
3	GOL	D	402	-	5,5,5	0.19	0	5,5,5	0.49	0
3	GOL	C	402	-	5,5,5	0.08	0	5,5,5	0.31	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	D	401	-	-	2/4/4/4	-
3	GOL	B	402	-	-	2/4/4/4	-
3	GOL	A	403	-	-	0/4/4/4	-
3	GOL	D	403	-	-	1/4/4/4	-
3	GOL	D	402	-	-	2/4/4/4	-
3	GOL	C	402	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	401	GOL	O1-C1-C2-C3
3	D	402	GOL	C1-C2-C3-O3
3	B	402	GOL	O1-C1-C2-C3
3	D	401	GOL	O1-C1-C2-O2
3	D	402	GOL	O2-C2-C3-O3
3	B	402	GOL	O1-C1-C2-O2
3	D	403	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	GOL	1	0
3	D	403	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	325/348 (93%)	-0.38	4 (1%) 79 77	19, 33, 72, 94	0
1	B	325/348 (93%)	-0.27	2 (0%) 89 88	23, 41, 71, 109	0
1	C	325/348 (93%)	0.40	29 (8%) 9 9	33, 65, 118, 149	0
1	D	325/348 (93%)	0.01	14 (4%) 35 33	26, 52, 87, 169	0
All	All	1300/1392 (93%)	-0.06	49 (3%) 40 39	19, 47, 98, 169	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	59	GLU	9.3
1	D	58	GLY	8.2
1	C	55	ILE	8.0
1	C	56	GLY	5.8
1	C	66	LEU	5.5
1	C	75	GLY	5.2
1	C	67	TYR	5.1
1	D	67	TYR	4.9
1	C	54	VAL	4.8
1	D	66	LEU	4.5
1	B	129	ALA	4.4
1	D	61	ASP	4.4
1	C	72	VAL	4.2
1	C	234	ILE	4.1
1	A	4	LYS	3.5
1	D	85	PRO	3.5
1	A	127	ILE	3.4
1	C	68	ILE	3.1
1	C	245	HIS	3.1
1	D	60	LEU	3.0
1	D	63	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	253	ASN	2.9
1	D	64	PRO	2.9
1	B	58	GLY	2.9
1	C	48	ASP	2.9
1	C	97	ALA	2.9
1	D	86	LEU	2.7
1	C	114	ALA	2.7
1	C	236	ASN	2.7
1	A	128	ASN	2.6
1	C	51	GLY	2.6
1	D	5	VAL	2.6
1	C	84	ASP	2.5
1	C	58	GLY	2.5
1	C	89	THR	2.5
1	D	56	GLY	2.4
1	C	73	GLY	2.4
1	A	129	ALA	2.4
1	C	71	LYS	2.3
1	C	132	GLY	2.3
1	C	170	HIS	2.3
1	C	76	GLY	2.2
1	C	64	PRO	2.2
1	D	187	ASP	2.2
1	C	129	ALA	2.2
1	C	33	ASP	2.1
1	D	65	MET	2.1
1	C	49	ILE	2.0
1	C	63	ALA	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,

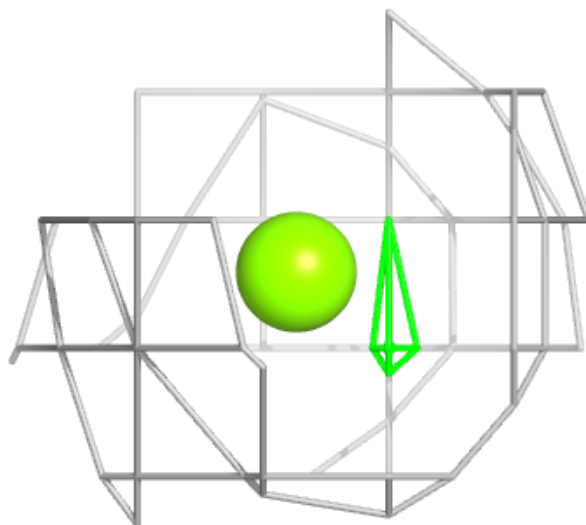
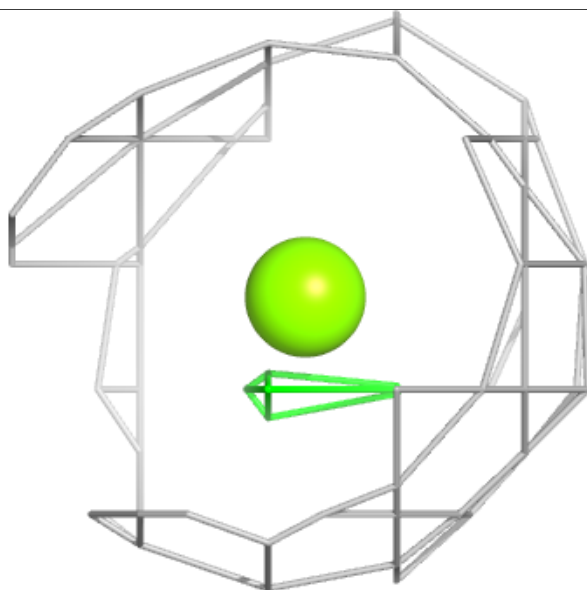
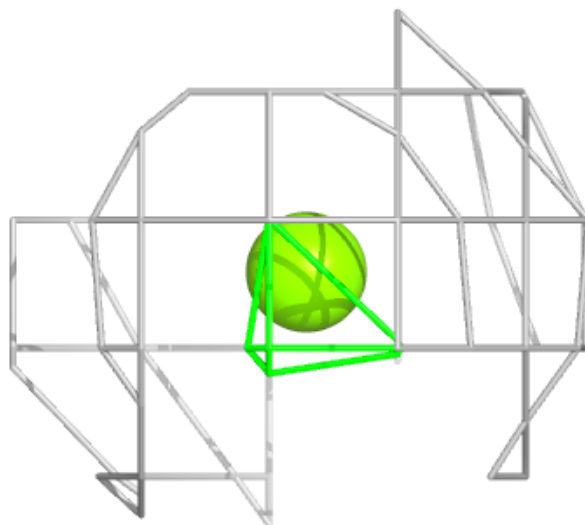
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	402	6/6	0.53	0.44	102,109,109,112	0
2	MG	C	401	1/1	0.79	0.18	68,68,68,68	0
3	GOL	C	402	6/6	0.80	0.13	65,73,76,78	0
3	GOL	D	401	6/6	0.80	0.21	74,84,85,87	0
3	GOL	D	402	6/6	0.82	0.38	60,72,81,88	0
3	GOL	D	403	6/6	0.89	0.21	63,65,67,71	0
2	MG	A	401	1/1	0.92	0.13	31,31,31,31	0
2	MG	B	401	1/1	0.92	0.11	39,39,39,39	0
3	GOL	A	403	6/6	0.96	0.12	38,43,45,47	0
2	MG	A	402	1/1	0.96	0.07	27,27,27,27	0

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

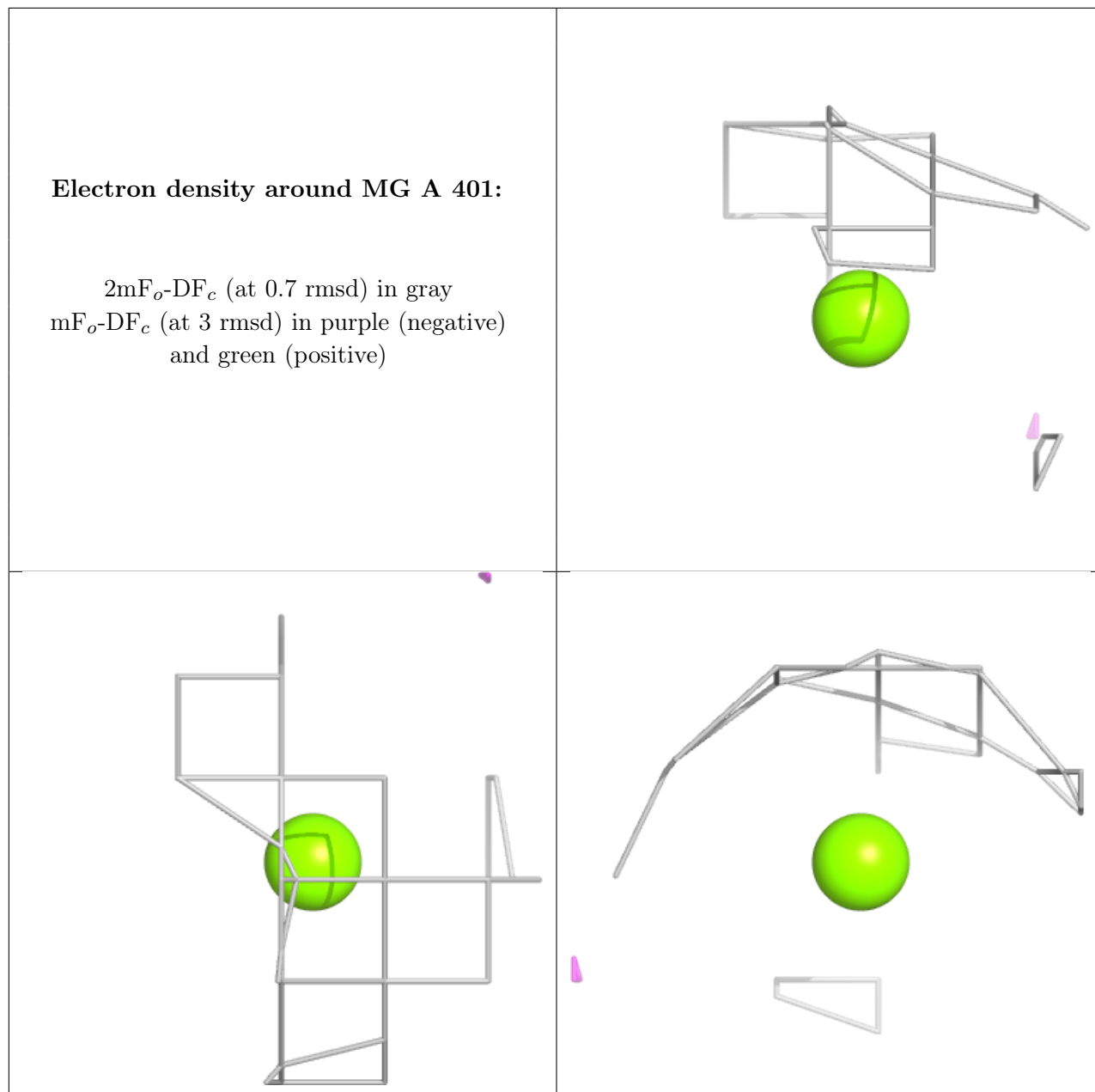
Electron density around MG C 401:

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



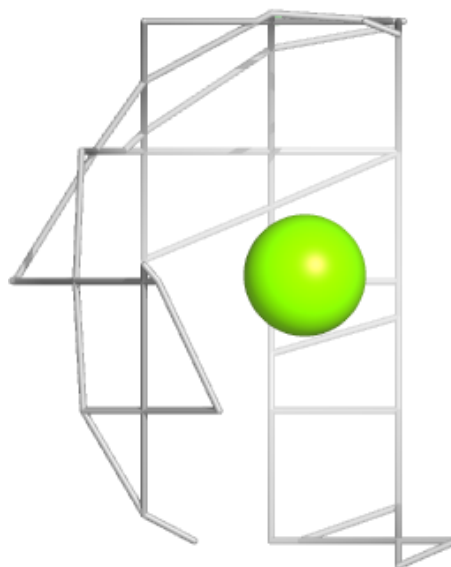
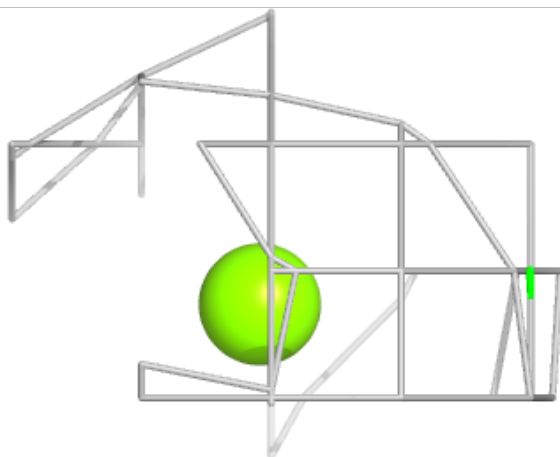
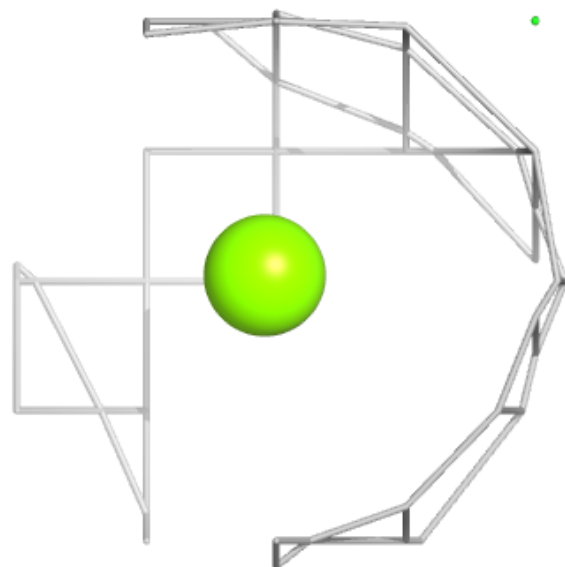
Electron density around MG A 401:

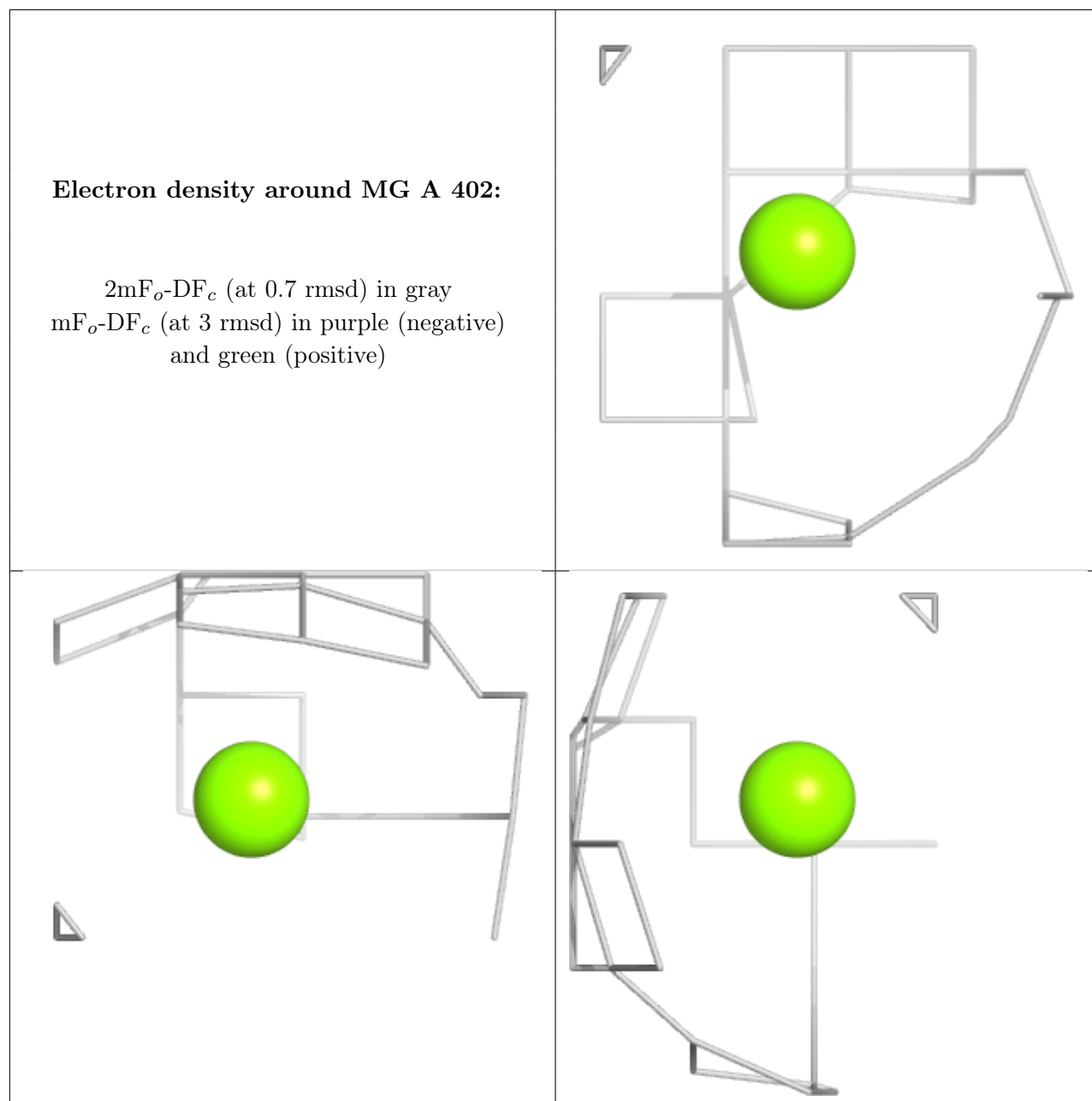
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around MG B 401:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.