



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:13 PM UTC

PDB ID : 8FMM / pdb_00008fmm
Title : Complex structure of wild type Troponin complex
Authors : Wang, P.; Ahmed, M.; Sadek, H.
Deposited on : 2022-12-23
Resolution : 3.11 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

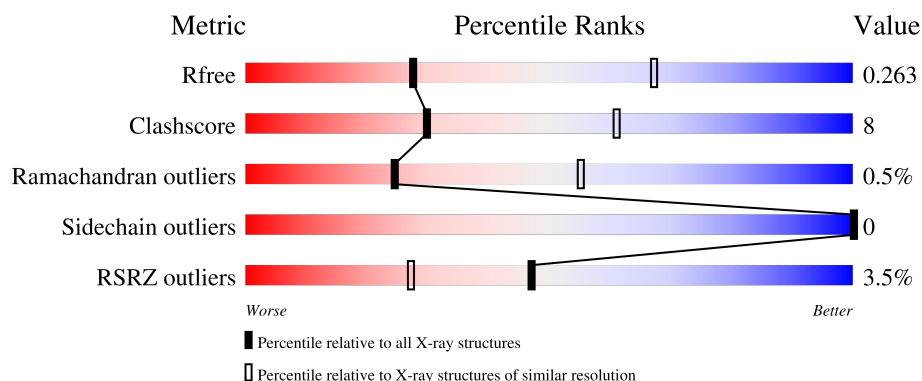
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.11 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1816 (3.14-3.10)
Clashscore	190562	1906 (3.14-3.10)
Ramachandran outliers	187476	1802 (3.14-3.10)
Sidechain outliers	187428	1802 (3.14-3.10)
RSRZ outliers	180081	1816 (3.14-3.10)

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	164	2% 77% 18% 5%
1	D	164	4% 73% 21% 5%
2	B	109	3% 52% 13% 35%
2	E	109	% 50% 12% 39%
3	C	135	% 60% 21% 19%

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Mol	Chain	Length	Quality of chain			
			5%	59%	19%	21%
3	F	135				

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5394 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 Troponin C, slow skeletal and cardiac muscles.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	0	0	0
			1238	769	188	270	11			
1	D	155	Total	C	N	O	S	0	0	0
			1234	766	188	270	10			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLN	-	expression tag	UNP P63316
A	-1	GLY	-	expression tag	UNP P63316
A	0	SER	-	expression tag	UNP P63316
A	35	SER	CYS	conflict	UNP P63316
A	84	SER	CYS	conflict	UNP P63316
A	115	GLU	ASP	conflict	UNP P63316
D	-2	GLN	-	expression tag	UNP P63316
D	-1	GLY	-	expression tag	UNP P63316
D	0	SER	-	expression tag	UNP P63316
D	35	SER	CYS	conflict	UNP P63316
D	84	SER	CYS	conflict	UNP P63316
D	115	GLU	ASP	conflict	UNP P63316

- Molecule 2 is a protein called Troponin T, cardiac muscle.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	71	Total	C	N	O	0	0	0
			619	387	116	116			
2	E	67	Total	C	N	O	0	0	0
			583	367	108	108			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	180	GLN	-	expression tag	UNP P45379
B	181	GLY	-	expression tag	UNP P45379
B	182	SER	-	expression tag	UNP P45379
E	180	GLN	-	expression tag	UNP P45379
E	181	GLY	-	expression tag	UNP P45379
E	182	SER	-	expression tag	UNP P45379

- Molecule 3 is a protein called Troponin I, cardiac muscle.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	110	Total	C	N	O	S	0	0	0
			869	540	160	167	2			
3	F	106	Total	C	N	O	S	0	0	0
			845	525	156	162	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	80	ALA	CYS	conflict	UNP P19429
C	97	ALA	CYS	conflict	UNP P19429
F	80	ALA	CYS	conflict	UNP P19429
F	97	ALA	CYS	conflict	UNP P19429

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Ca	0	0
			3	3		
4	D	3	Total	Ca	0	0
			3	3		

LYS
ALA
LYS
VAL
THR
GLY
ARG
TRP
LYS

● Molecule 3: Troponin I, cardiac muscle



GLU
PRO
HIS
ALA
LYS
LYS
LYS
LYS
SER
K40
I41
S44
R45
K46
L47
K50
L53
L54
E71
R79
G89
L93
L96
A97
R98
Q99
R103
K106
E110
D113
I114
E115
A116
K117
V118
T119
D127
L128
K131
R136
GLY
LYS
PHE
LYS
ARG
PRO

THR
LEU
ARG
ARG
VAL
ARG
I149
S150
A151
D152
A153
M154
M155
G160
A161
ARG
ALA
LYS
GLU
SER

● Molecule 3: Troponin I, cardiac muscle



GLU
PRO
HIS
ALA
LYS
LYS
SER
K40
I41
S42
R45
K50
Q81
E84
L85
A86
G87
L88
G89
F90
A91
E92
L93
Q94
D95
L96
A97
R98
Q99
L100
H101
A102
R103
V104
R111
I114
V118
E124
I125
A126
D127
L128
T129
Q130
K131
I132
G137
LYS
PHE
LYS

ARG
PRO
THR
LEU
ARG
ARG
VAL
ARG
ARG
ILE
SER
ALA
D152
A153
M154
L159
GLY
ALA
ARG
ALA
LYS
GLU
SER

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.07Å 168.22Å 69.68Å 90.00° 101.40° 90.00°	Depositor
Resolution (Å)	41.24 – 3.11 41.24 – 3.11	Depositor EDS
% Data completeness (in resolution range)	87.0 (41.24-3.11) 87.0 (41.24-3.11)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.96 (at 3.12Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.230 , 0.264 0.228 , 0.263	Depositor DCC
R_{free} test set	731 reflections (4.28%)	wwPDB-VP
Wilson B-factor (Å ²)	36.2	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	5394	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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: CA

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	0.16	0/1250	0.36	0/1670
1	D	0.16	0/1246	0.44	1/1665 (0.1%)
2	B	0.14	0/625	0.37	0/831
2	E	0.16	0/589	0.42	0/783
3	C	0.15	0/872	0.37	0/1164
3	F	0.15	0/848	0.39	0/1131
All	All	0.15	0/5430	0.39	1/7244 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	54	PRO	N-CA-C	5.12	123.01	112.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1238	0	1171	22	0
1	D	1234	0	1163	26	0
2	B	619	0	634	9	0
2	E	583	0	600	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	869	0	910	21	0
3	F	845	0	884	21	0
4	A	3	0	0	0	0
4	D	3	0	0	0	0
All	All	5394	0	5362	91	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 8.

All (91) 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:D:52:PRO:HB2	1:D:57:LEU:HD11	1.61	0.81
3:F:88:LEU:HD12	3:F:92:GLU:HB3	1.64	0.79
1:A:46:ARG:NH1	1:A:51:ASN:OD1	2.25	0.70
1:A:121:LEU:HB3	1:A:128:ILE:HG12	1.76	0.67
1:D:129:THR:HG22	1:D:131:ASP:H	1.66	0.61
1:D:99:ASP:OD1	1:D:102:ARG:NH2	2.34	0.60
2:B:214:GLU:OE2	3:C:98:ARG:NH2	2.23	0.59
2:E:217:LYS:HD3	3:F:102:ALA:HB2	1.82	0.59
1:D:158:LYS:HD2	1:D:158:LYS:O	2.03	0.59
1:A:156:PHE:CE1	3:C:50:LYS:HE2	2.38	0.59
2:E:265:ARG:NH2	3:F:124:GLU:OE1	2.38	0.57
1:D:60:MET:HA	1:D:63:GLU:HG2	1.86	0.56
3:F:92:GLU:HA	3:F:95:ASP:HB3	1.88	0.56
1:D:9:VAL:HG21	1:D:79:VAL:HG12	1.87	0.56
2:B:204:GLU:HA	2:B:207:LYS:HB3	1.88	0.56
2:E:221:ILE:HD12	3:F:97:ALA:HB1	1.87	0.55
1:D:53:THR:HG22	1:D:54:PRO:HD2	1.87	0.55
1:A:117:LEU:HD12	1:A:137:MET:HE2	1.87	0.55
1:A:104:PHE:HZ	3:C:53:LEU:HB3	1.72	0.54
1:D:37:SER:HB3	1:D:71:THR:HG22	1.89	0.53
1:D:62:ASP:HA	1:D:65:ASP:HB3	1.91	0.52
3:C:115:GLU:O	3:C:119:THR:OG1	2.26	0.52
1:A:11:GLN:OE1	3:C:44:SER:OG	2.22	0.52
3:C:150:SER:HB3	3:C:153:ALA:HB3	1.91	0.52
1:D:27:PHE:CZ	1:D:44:VAL:HG11	2.45	0.51
2:B:264:LEU:HD13	3:C:128:LEU:HB2	1.93	0.51
1:D:56:GLU:O	1:D:60:MET:HG2	2.11	0.51
3:F:92:GLU:O	3:F:96:LEU:N	2.27	0.51
1:A:124:THR:HG23	1:A:126:GLU:H	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:ASN:HD21	1:A:145:ASP:CG	2.20	0.50
3:F:88:LEU:HG	3:F:93:LEU:HG	1.93	0.50
2:E:222:ASP:HA	2:E:224:LEU:HD23	1.94	0.49
2:E:221:ILE:O	2:E:224:LEU:HB3	2.12	0.49
1:A:135:GLU:HA	1:A:138:LYS:HB2	1.94	0.49
1:D:115:GLU:O	1:D:119:ILE:HG12	2.13	0.48
1:D:38:THR:HG22	1:D:57:LEU:HB3	1.95	0.48
2:E:240:ILE:HG13	3:F:104:VAL:HG22	1.94	0.48
1:D:79:VAL:O	1:D:83:ARG:HG3	2.14	0.47
1:A:4:ILE:HG13	3:C:47:LEU:HD13	1.95	0.47
3:C:127:ASP:O	3:C:131:LYS:HG2	2.15	0.47
1:D:27:PHE:CE1	3:F:154:MET:HE2	2.50	0.47
1:D:29:LEU:HD12	1:D:29:LEU:HA	1.77	0.47
3:C:89:GLY:O	3:C:93:LEU:HB2	2.14	0.47
3:C:96:LEU:HA	3:C:99:GLN:HG2	1.96	0.47
1:D:2:ASP:O	1:D:4:ILE:N	2.48	0.47
2:E:206:GLU:HB3	2:E:209:LYS:NZ	2.30	0.47
1:A:81:MET:HE1	3:C:154:MET:HE3	1.96	0.46
2:B:218:VAL:O	2:B:222:ASP:HB2	2.14	0.46
2:B:226:GLU:O	2:B:230:ARG:HG3	2.14	0.46
1:D:157:MET:HE1	3:F:50:LYS:HB3	1.97	0.46
2:B:236:LEU:O	2:B:240:ILE:HG13	2.15	0.46
1:A:48:LEU:HD12	1:A:48:LEU:O	2.16	0.45
1:A:27:PHE:CE1	1:A:44:VAL:HG21	2.51	0.45
1:D:150:TYR:HB3	2:E:267:ARG:HD2	1.99	0.45
3:F:127:ASP:O	3:F:131:LYS:HG3	2.16	0.45
1:D:14:GLU:H	1:D:14:GLU:CD	2.25	0.45
1:D:118:LYS:HG2	1:D:133:ILE:HG21	1.98	0.45
1:D:54:PRO:HA	1:D:57:LEU:HB2	1.99	0.45
3:F:128:LEU:O	3:F:132:ILE:HG22	2.17	0.45
2:B:225:ASN:O	2:B:229:LEU:HB3	2.17	0.44
1:A:127:THR:O	1:A:127:THR:OG1	2.34	0.44
1:A:11:GLN:HE21	1:A:11:GLN:HB3	1.58	0.44
1:A:157:MET:HE3	3:C:54:LEU:HD21	2.00	0.44
3:C:151:ALA:O	3:C:155:MET:HG2	2.18	0.44
1:A:22:ALA:O	1:A:26:ILE:HG13	2.17	0.44
3:C:41:ILE:HG21	3:C:45:ARG:HG2	2.00	0.44
3:F:42:SER:HB2	3:F:45:ARG:HB2	2.00	0.44
2:E:224:LEU:HD22	2:E:229:LEU:HD11	2.00	0.43
3:C:113:ASP:O	3:C:117:LYS:HG3	2.19	0.43
1:D:114:LEU:HG	1:D:118:LYS:HE3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:114:ILE:O	3:F:118:VAL:HG23	2.19	0.43
3:F:100:LEU:O	3:F:104:VAL:HG23	2.19	0.42
2:B:215:ARG:HA	2:B:218:VAL:HG22	2.00	0.42
1:D:64:VAL:HG21	1:D:80:MET:HE3	2.01	0.42
2:E:244:GLU:OE1	2:E:247:LYS:NZ	2.36	0.42
3:C:106:LYS:O	3:C:110:GLU:HG3	2.18	0.42
2:B:248:PHE:HZ	3:C:71:GLU:HG2	1.84	0.42
1:A:104:PHE:CZ	3:C:53:LEU:HB3	2.54	0.42
3:C:79:ARG:O	3:C:103:ARG:NH2	2.42	0.42
3:F:125:ILE:O	3:F:129:THR:OG1	2.29	0.42
1:A:27:PHE:HZ	3:C:154:MET:SD	2.43	0.41
2:E:247:LYS:HG3	3:F:111:ARG:HA	2.02	0.41
3:F:93:LEU:HD23	3:F:93:LEU:HA	1.92	0.41
3:F:95:ASP:OD1	3:F:99:GLN:NE2	2.53	0.41
3:F:81:GLN:H	3:F:81:GLN:CD	2.28	0.41
1:A:31:ALA:HB2	1:A:40:GLU:CD	2.45	0.41
2:E:264:LEU:HD13	3:F:128:LEU:HB2	2.02	0.41
1:A:118:LYS:HG2	1:A:133:ILE:HG21	2.03	0.41
1:A:120:MET:HE3	1:A:120:MET:HB3	1.89	0.41
1:D:2:ASP:C	1:D:4:ILE:H	2.29	0.40
1:D:53:THR:CG2	1:D:54:PRO:HD2	2.50	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	151/164 (92%)	140 (93%)	11 (7%)	0	100	100
1	D	151/164 (92%)	138 (91%)	11 (7%)	2 (1%)	9	33
2	B	69/109 (63%)	66 (96%)	3 (4%)	0	100	100
2	E	65/109 (60%)	61 (94%)	4 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	106/135 (78%)	98 (92%)	8 (8%)	0	100	100
3	F	102/135 (76%)	97 (95%)	4 (4%)	1 (1%)	12	39
All	All	644/816 (79%)	600 (93%)	41 (6%)	3 (0%)	24	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	3	ASP
1	D	54	PRO
3	F	87	GLY

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	137/144 (95%)	137 (100%)	0	100	100
1	D	136/144 (94%)	136 (100%)	0	100	100
2	B	67/97 (69%)	67 (100%)	0	100	100
2	E	63/97 (65%)	63 (100%)	0	100	100
3	C	90/112 (80%)	90 (100%)	0	100	100
3	F	88/112 (79%)	88 (100%)	0	100	100
All	All	581/706 (82%)	581 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
1	A	143	ASN
3	C	48	GLN
1	D	18	ASN

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Mol	Chain	Res	Type
3	F	48	GLN
3	F	101	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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 [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	155/164 (94%)	0.35	3 (1%) 66 45	12, 36, 68, 86	0
1	D	155/164 (94%)	0.69	7 (4%) 38 21	23, 53, 86, 107	0
2	B	71/109 (65%)	0.28	3 (4%) 40 22	14, 29, 54, 66	0
2	E	67/109 (61%)	0.20	1 (1%) 72 52	7, 32, 55, 80	0
3	C	110/135 (81%)	0.43	2 (1%) 67 46	15, 39, 62, 84	0
3	F	106/135 (78%)	0.55	7 (6%) 24 12	7, 37, 77, 96	0
All	All	664/816 (81%)	0.45	23 (3%) 47 27	7, 39, 77, 107	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	86	ALA	3.6
3	F	91	ALA	3.5
1	A	161	GLU	2.9
1	D	57	LEU	2.8
1	D	95	GLU	2.6
3	C	161	ALA	2.6
2	B	272	GLN	2.6
1	D	127	THR	2.5
2	B	204	GLU	2.5
1	D	134	GLU	2.4
3	F	90	PHE	2.4
3	F	154	MET	2.4
1	D	124	THR	2.4
1	D	155	GLU	2.2
1	A	127	THR	2.2
1	A	129	THR	2.2
3	F	132	ILE	2.2
3	C	160	GLY	2.2
3	F	159	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
3	F	84	GLU	2.1
2	B	202	GLN	2.1
2	E	226	GLU	2.1
1	D	158	LYS	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

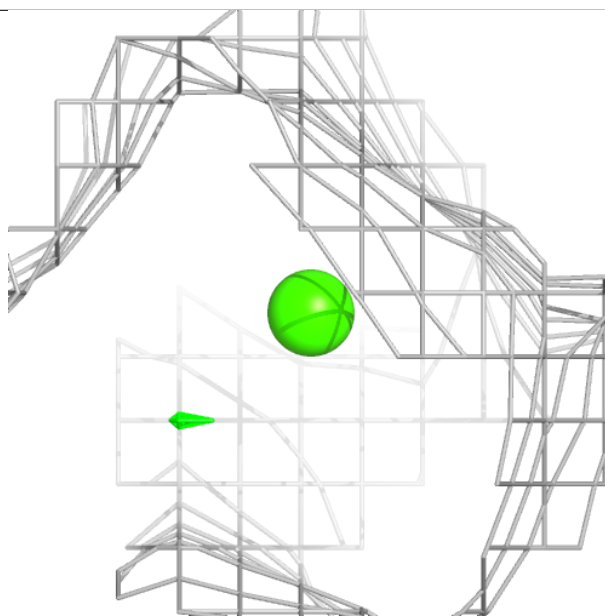
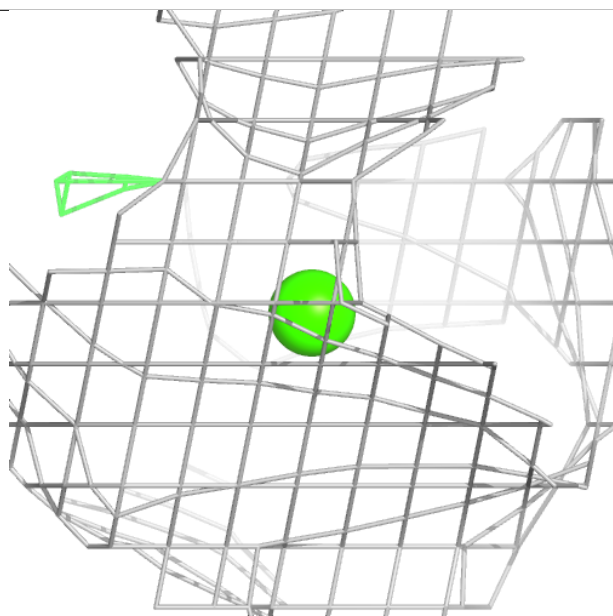
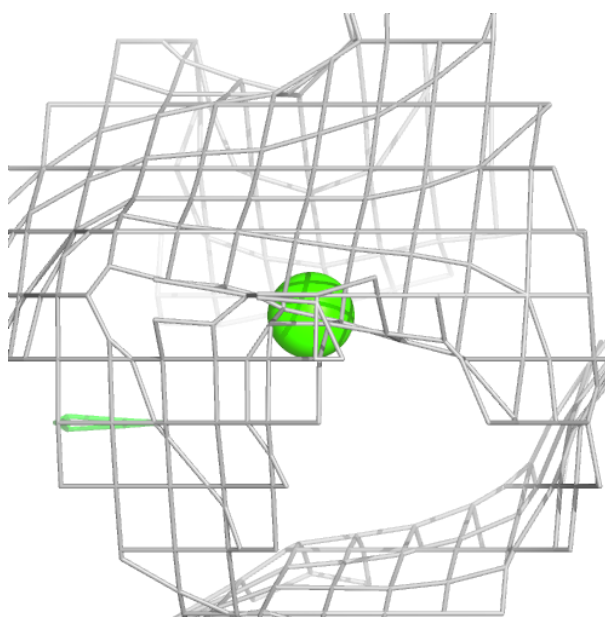
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CA	A	203	1/1	0.95	0.06	47,47,47,47	0
4	CA	D	201	1/1	0.97	0.07	70,70,70,70	0
4	CA	D	202	1/1	0.97	0.04	30,30,30,30	0
4	CA	A	202	1/1	0.99	0.02	22,22,22,22	0
4	CA	A	201	1/1	0.99	0.12	29,29,29,29	0
4	CA	D	203	1/1	0.99	0.03	34,34,34,34	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.

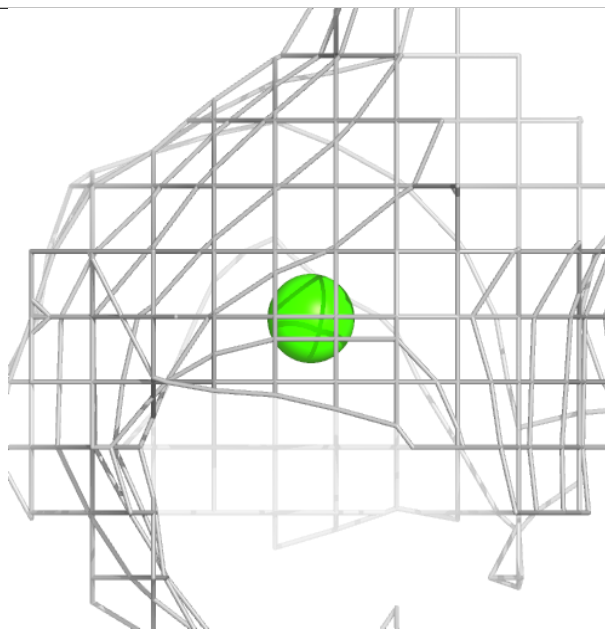
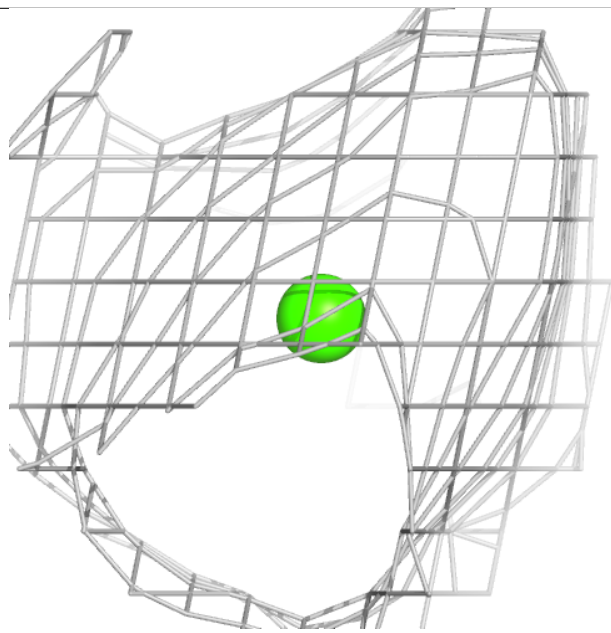
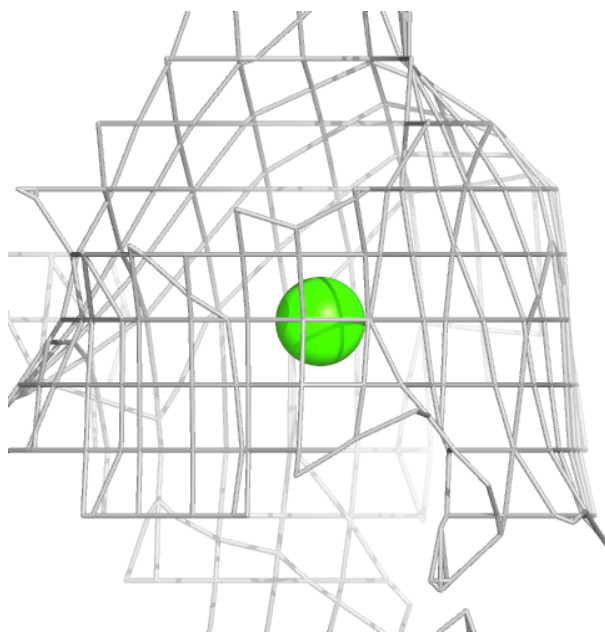
Electron density around CA A 203:

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



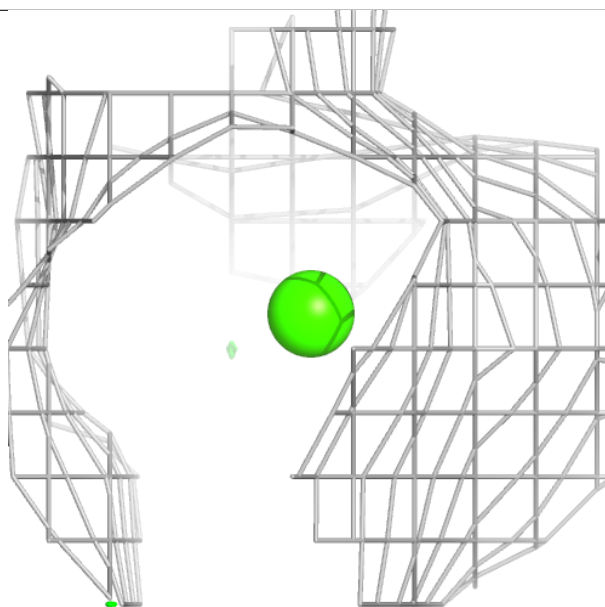
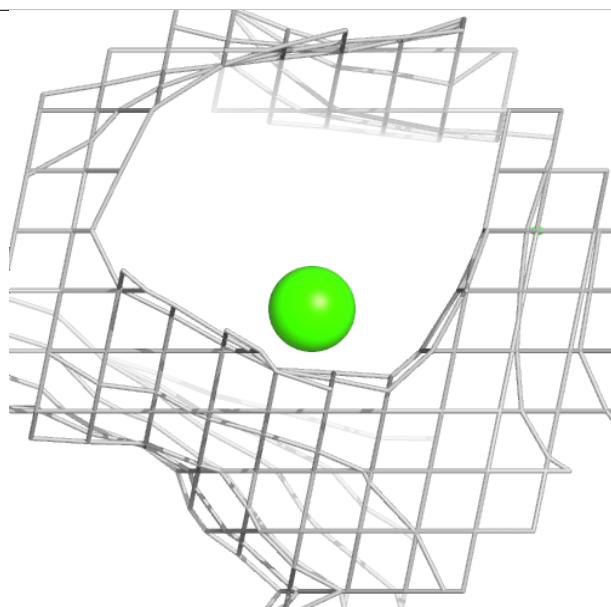
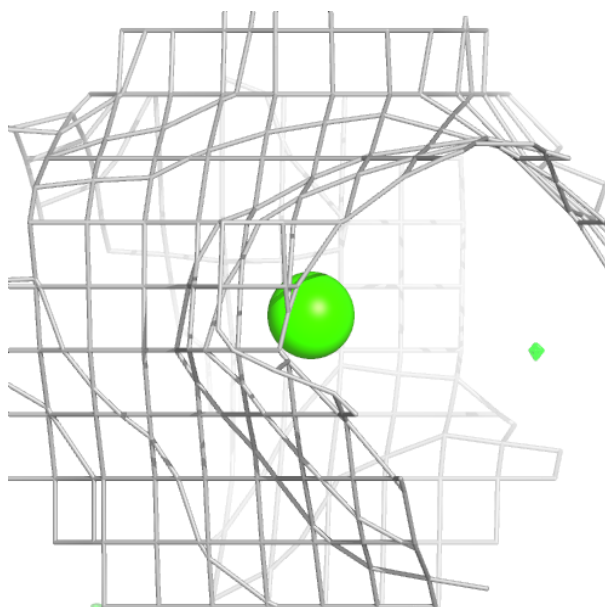
Electron density around CA D 201:

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



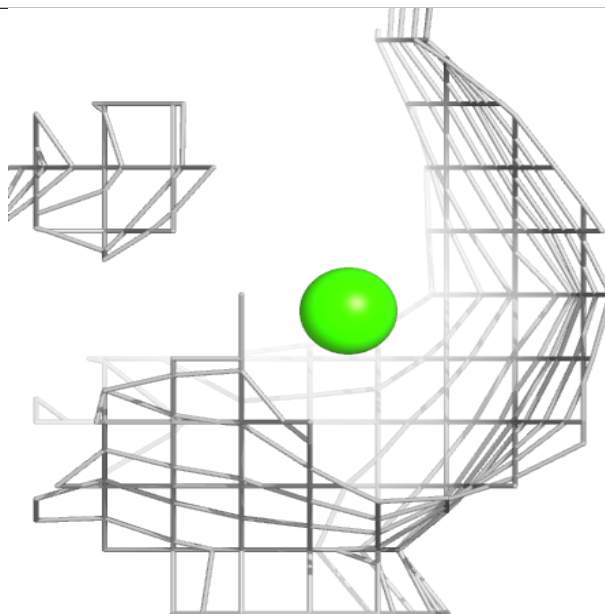
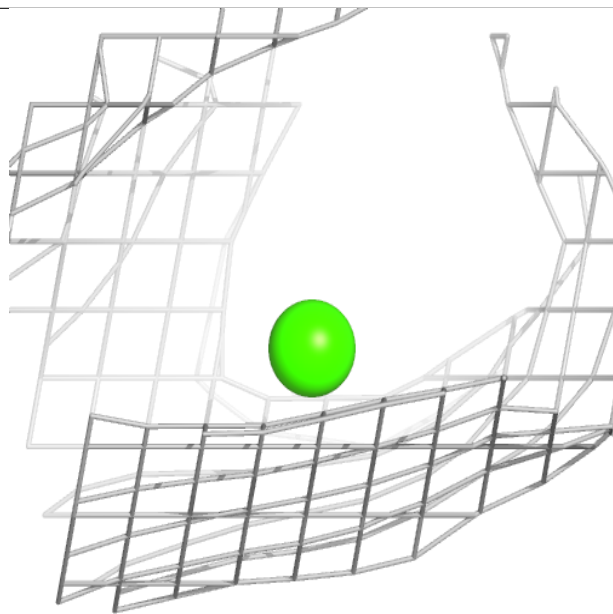
Electron density around CA D 202:

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



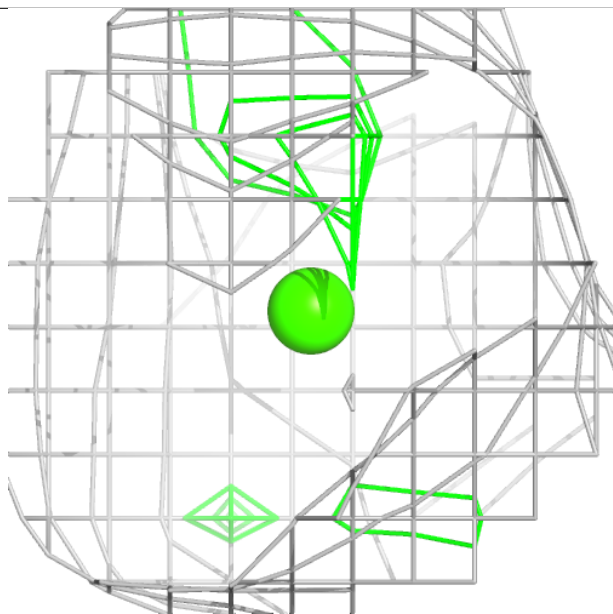
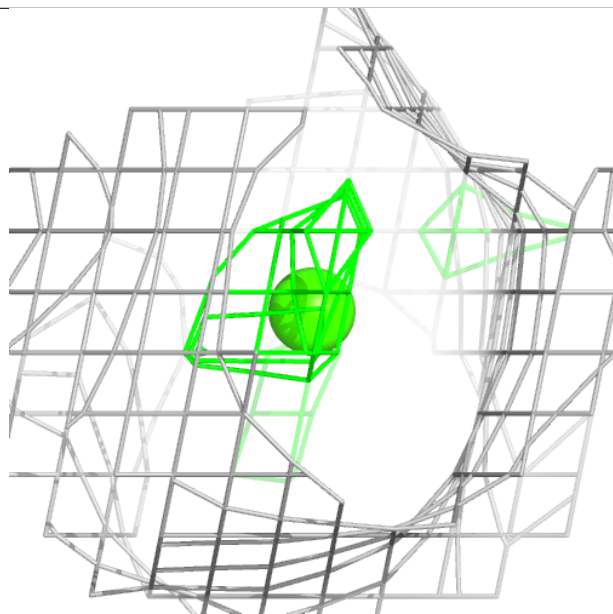
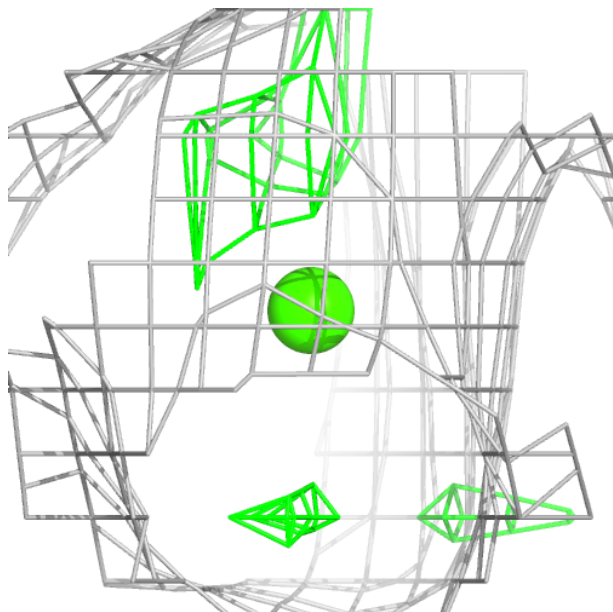
Electron density around CA A 202:

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and green (positive)



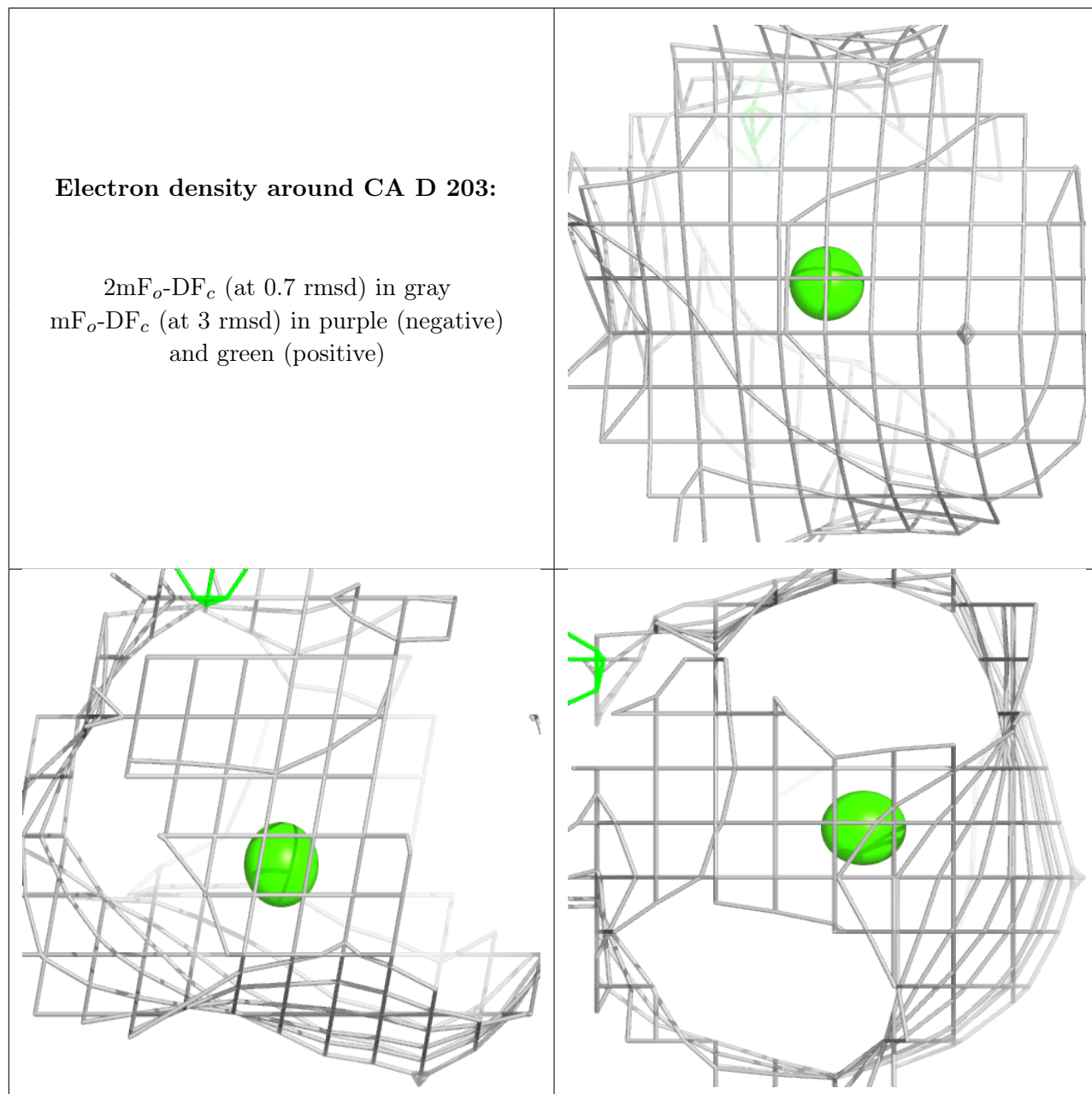
Electron density around CA A 201:

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



Electron density around CA D 203:

$2mF_o-DF_c$ (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.