Full wwPDB X-ray Structure Validation Report (i)

Sep 26, 2023 – 01:24 AM EDT

PDB ID	:	6BCF
Title	:	I-LtrI G183A bound to cognate substrate (pre-cleavage complex)
Authors	:	Brown, C.; Zhang, K.; McMurrough, T.A.; Gloor, G.B.; Edgell, D.R.; Junop, M.
Deposited on	:	2017-10-20
Resolution	:	2.92 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 (i) 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 (i)) were used in the production of this report:

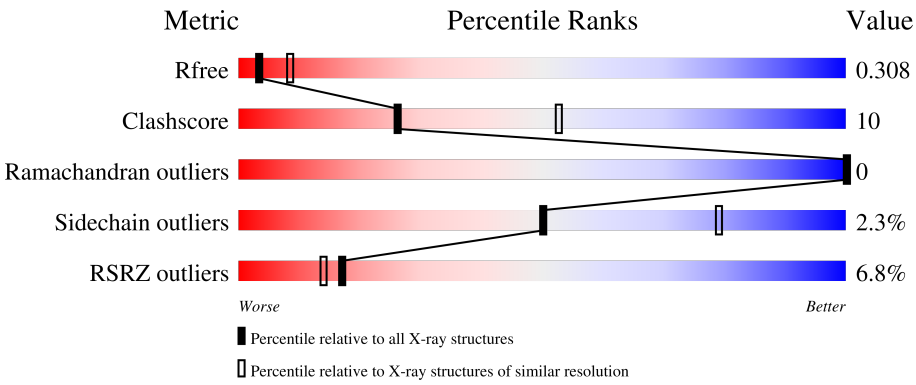
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.92 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	2307 (2.94-2.90)
Clashscore	141614	2531 (2.94-2.90)
Ramachandran outliers	138981	2462 (2.94-2.90)
Sidechain outliers	138945	2464 (2.94-2.90)
RSRZ outliers	127900	2248 (2.94-2.90)

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 <=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	315	7% 68% 22% • 9%
1	D	315	6% 72% 18% • 9%
1	G	315	7% 71% 20% • 9%
1	J	315	9% 67% 23% • 8%
2	B	26	73% 27%

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Mol	Chain	Length	Quality of chain
2	E	26	58% 42%
2	H	26	62% 27% 12%
2	K	26	65% 27% 8%
3	C	26	4% 73% 15% 12%
3	F	26	65% 31%
3	I	26	54% 38% 8%
3	L	26	69% 23% 8%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12768 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 Ribosomal protein 3/homing endonuclease-like fusion protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	288	2126	1364	364	393	5	0	0	0
1	A	288	2130	1369	364	391	6	0	0	0
1	G	288	2120	1363	366	385	6	0	0	0
1	J	289	2131	1366	365	394	6	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	183	ALA	GLY	engineered mutation	UNP C7SWF3
A	183	ALA	GLY	engineered mutation	UNP C7SWF3
G	183	ALA	GLY	engineered mutation	UNP C7SWF3
J	183	ALA	GLY	engineered mutation	UNP C7SWF3

- Molecule 2 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	E	26	534	256	98	155	25	0	0	0
2	B	26	534	256	98	155	25	0	0	0
2	H	26	534	256	98	155	25	0	0	0
2	K	26	534	256	98	155	25	0	0	0

- Molecule 3 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	F	26	529	253	95	155	26	0	0	0
3	C	26	526	253	95	153	25	0	0	0
3	I	26	529	253	95	155	26	0	0	0
3	L	26	529	253	95	155	26	0	0	0

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

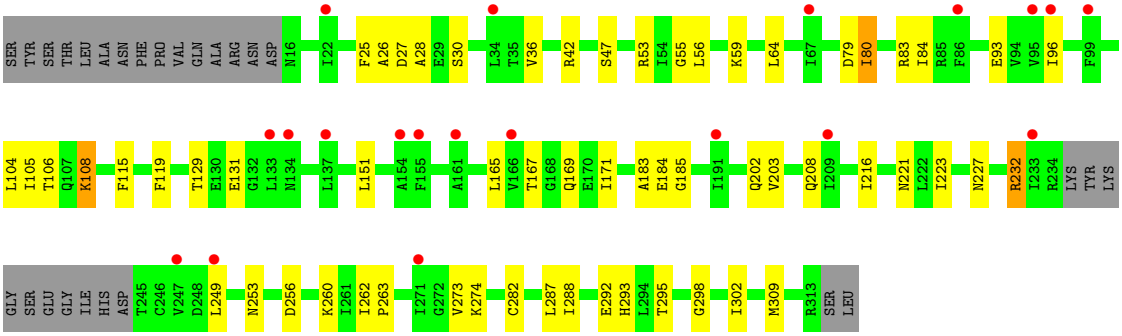
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Ca		
4	D	3	3	3	0	0
4	A	2	2	2	0	0
4	B	1	1	1	0	0
4	G	2	2	2	0	0
4	H	1	1	1	0	0
4	J	2	2	2	0	0
4	L	1	1	1	0	0

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.

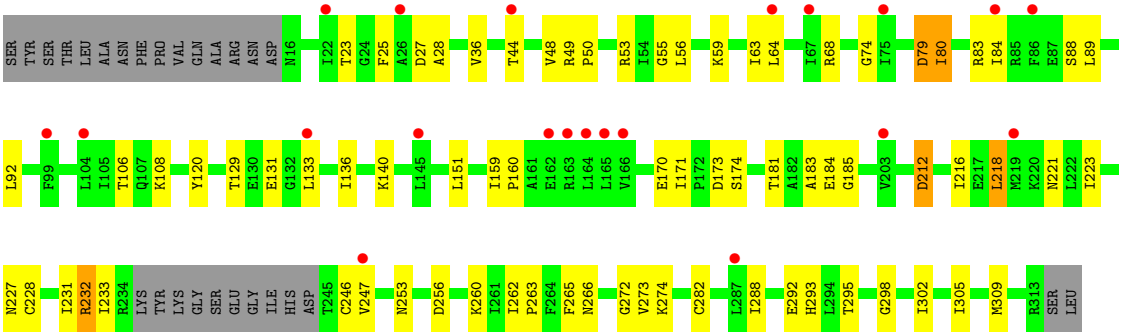
• Molecule 1: Ribosomal protein 3/homing endonuclease-like fusion protein

Chain D:



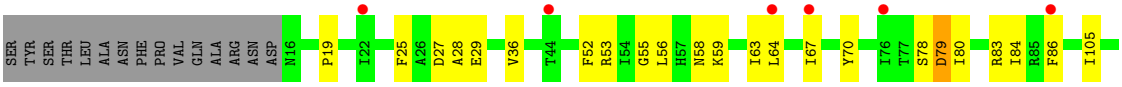
• Molecule 1: Ribosomal protein 3/homing endonuclease-like fusion protein

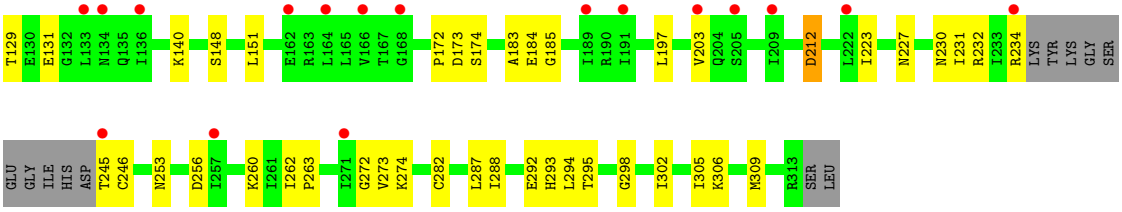
Chain A:



• Molecule 1: Ribosomal protein 3/homing endonuclease-like fusion protein

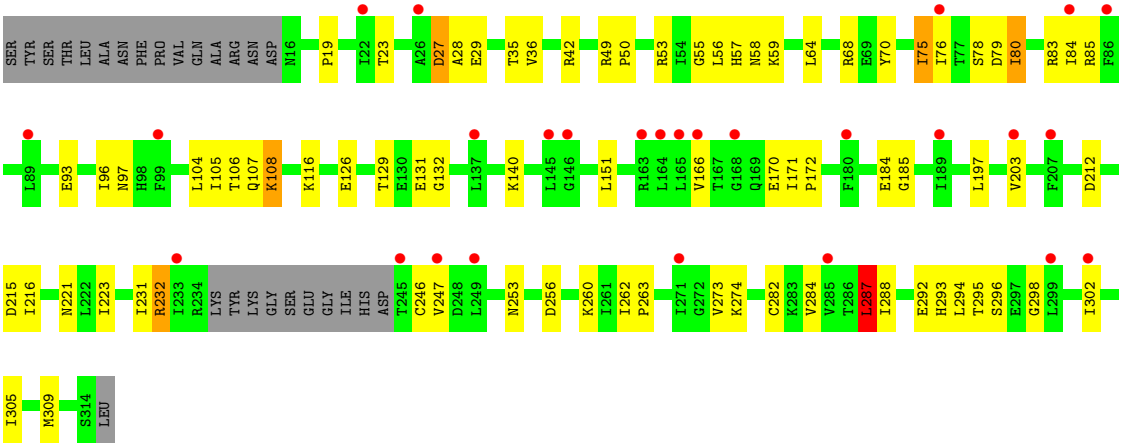
Chain G:





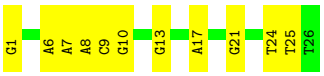
• Molecule 1: Ribosomal protein 3/homing endonuclease-like fusion protein

Chain J:



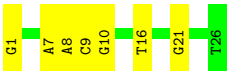
• Molecule 2: DNA (26-MER)

Chain E:



• Molecule 2: DNA (26-MER)

Chain B:

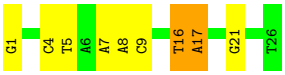


• Molecule 2: DNA (26-MER)

Chain H:

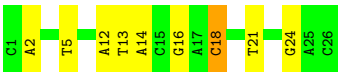
• Molecule 2: DNA (26-MER)

Chain K:



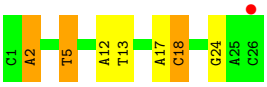
• Molecule 3: DNA (26-MER)

Chain F:



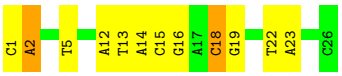
• Molecule 3: DNA (26-MER)

Chain C:



• Molecule 3: DNA (26-MER)

Chain I:



• Molecule 3: DNA (26-MER)

Chain L:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	43.96Å 66.77Å 169.41Å 90.12° 90.08° 90.06°	Depositor
Resolution (Å)	38.99 – 2.92 66.77 – 2.92	Depositor EDS
% Data completeness (in resolution range)	86.9 (38.99-2.92) 80.4 (66.77-2.92)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.24 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.239 , 0.309 0.240 , 0.308	Depositor DCC
R_{free} test set	1867 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	53.3	Xtriage
Anisotropy	1.091	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 41.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.438 for h,-k,-l 0.438 for -h,k,-l 0.437 for -h,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	12768	wwPDB-VP
Average B, all atoms (Å^2)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.72% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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

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	0.50	1/2167 (0.0%)	0.70	2/2958 (0.1%)
1	D	0.48	0/2163	0.70	1/2954 (0.0%)
1	G	0.45	0/2157	0.68	1/2946 (0.0%)
1	J	0.47	0/2168	0.68	2/2961 (0.1%)
2	B	1.05	1/599 (0.2%)	1.22	5/924 (0.5%)
2	E	0.97	0/599	1.19	3/924 (0.3%)
2	H	0.88	0/599	1.19	5/924 (0.5%)
2	K	0.91	0/599	1.22	4/924 (0.4%)
3	C	0.96	0/589	1.24	6/906 (0.7%)
3	F	1.03	1/592 (0.2%)	1.22	7/910 (0.8%)
3	I	0.99	0/592	1.21	4/910 (0.4%)
3	L	0.95	0/592	1.21	6/910 (0.7%)
All	All	0.69	3/13416 (0.0%)	0.93	46/19151 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	232	ARG	CG-CD	-6.50	1.35	1.51
2	B	8	DA	C3'-C2'	-5.32	1.45	1.52
3	F	21	DT	C1'-N1	5.16	1.55	1.49

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	K	16	DT	OP2-P-O3'	11.33	130.13	105.20
1	A	232	ARG	NE-CZ-NH1	-9.48	115.56	120.30
2	E	7	DA	O4'-C1'-N9	9.44	114.61	108.00
2	K	7	DA	O4'-C1'-N9	9.35	114.54	108.00
2	K	17	DA	OP1-P-OP2	-9.21	105.78	119.60
2	B	7	DA	O4'-C1'-N9	9.16	114.41	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	17	DA	OP1-P-OP2	-8.90	106.25	119.60
3	C	5	DT	O4'-C1'-N1	8.85	114.20	108.00
3	C	2	DA	O4'-C1'-N9	8.84	114.19	108.00
3	L	2	DA	O4'-C1'-N9	8.48	113.94	108.00
3	L	5	DT	O4'-C1'-N1	8.38	113.86	108.00
3	F	2	DA	O4'-C1'-N9	7.19	113.03	108.00
1	D	232	ARG	NE-CZ-NH1	-7.01	116.79	120.30
3	C	18	DC	O4'-C1'-N1	6.96	112.87	108.00
3	I	5	DT	O4'-C1'-N1	6.87	112.81	108.00
2	H	7	DA	O4'-C1'-N9	6.80	112.76	108.00
3	I	2	DA	O4'-C1'-N9	6.69	112.69	108.00
2	E	21	DG	O4'-C1'-N9	6.50	112.55	108.00
2	B	21	DG	O4'-C1'-N9	6.47	112.53	108.00
2	B	16	DT	OP1-P-O3'	6.30	119.07	105.20
3	C	24	DG	O4'-C1'-N9	6.28	112.40	108.00
2	H	21	DG	O4'-C1'-N9	6.28	112.39	108.00
1	J	232	ARG	NE-CZ-NH1	-6.23	117.18	120.30
2	E	17	DA	OP1-P-OP2	-6.21	110.28	119.60
3	F	5	DT	O4'-C1'-N1	6.20	112.34	108.00
3	F	18	DC	O4'-C1'-N1	6.17	112.32	108.00
2	H	16	DT	OP2-P-O3'	5.92	118.23	105.20
1	J	287	LEU	CA-CB-CG	5.88	128.83	115.30
3	L	1	DC	O4'-C4'-C3'	-5.79	102.18	104.50
1	G	212	ASP	CB-CG-OD1	5.58	123.32	118.30
3	F	18	DC	C1'-O4'-C4'	-5.56	104.54	110.10
3	F	16	DG	O5'-P-OP2	-5.48	100.77	105.70
2	K	21	DG	O4'-C1'-N9	5.44	111.81	108.00
3	C	18	DC	C1'-O4'-C4'	-5.38	104.72	110.10
2	B	7	DA	C3'-C2'-C1'	-5.25	96.20	102.50
3	F	24	DG	O4'-C1'-N9	5.22	111.65	108.00
3	I	16	DG	O4'-C1'-N9	-5.20	104.36	108.00
3	L	24	DG	O4'-C1'-N9	5.20	111.64	108.00
2	H	7	DA	C3'-C2'-C1'	-5.14	96.33	102.50
3	C	17	DA	O4'-C1'-N9	5.11	111.58	108.00
3	L	16	DG	O4'-C1'-N9	-5.08	104.44	108.00
3	L	5	DT	C3'-C2'-C1'	-5.05	96.43	102.50
3	F	2	DA	OP1-P-O3'	5.05	116.31	105.20
3	I	18	DC	C1'-O4'-C4'	-5.05	105.05	110.10
2	B	8	DA	O4'-C1'-C2'	-5.04	101.87	105.90
1	A	212	ASP	CB-CG-OD1	5.02	122.82	118.30

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	2130	0	2031	52	0
1	D	2126	0	2015	43	0
1	G	2120	0	2019	47	0
1	J	2131	0	2018	58	0
2	B	534	0	296	3	0
2	E	534	0	296	9	0
2	H	534	0	296	13	0
2	K	534	0	296	7	0
3	C	526	0	295	5	0
3	F	529	0	294	6	0
3	I	529	0	294	9	0
3	L	529	0	294	5	0
4	A	2	0	0	0	0
4	B	1	0	0	0	0
4	D	3	0	0	0	0
4	G	2	0	0	0	0
4	H	1	0	0	0	0
4	J	2	0	0	0	0
4	L	1	0	0	0	0
All	All	12768	0	10444	231	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 10.

All (231) 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:G:129:THR:HG22	1:G:131:GLU:H	1.38	0.88
1:G:28:ALA:HB1	1:G:184:GLU:HG3	1.57	0.86
1:J:129:THR:HG22	1:J:131:GLU:H	1.39	0.85
2:H:16:DT:H2"	2:H:17:DA:H5"	1.60	0.83
1:J:28:ALA:HB1	1:J:184:GLU:HG3	1.61	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:HG22	1:A:131:GLU:H	1.41	0.83
1:J:53:ARG:HD3	1:J:83:ARG:HD3	1.63	0.81
1:D:295:THR:HG23	1:D:298:GLY:H	1.46	0.80
1:A:28:ALA:HB1	1:A:184:GLU:HG3	1.65	0.79
1:A:295:THR:HG23	1:A:298:GLY:H	1.48	0.79
1:D:129:THR:HG22	1:D:131:GLU:H	1.45	0.78
1:D:28:ALA:HB1	1:D:184:GLU:HG3	1.64	0.77
1:A:53:ARG:HD3	1:A:83:ARG:HD3	1.69	0.75
1:J:295:THR:HG23	1:J:298:GLY:H	1.54	0.73
1:J:185:GLY:O	1:J:274:LYS:NZ	2.20	0.72
1:G:295:THR:HG23	1:G:298:GLY:H	1.54	0.71
1:D:185:GLY:O	1:D:274:LYS:NZ	2.23	0.70
1:D:53:ARG:HD3	1:D:83:ARG:HD3	1.73	0.69
1:G:253:ASN:OD1	1:G:256:ASP:N	2.22	0.69
3:C:12:DA:H2"	3:C:13:DT:H71	1.75	0.68
1:D:253:ASN:OD1	1:D:256:ASP:N	2.23	0.68
2:H:1:DG:HO5'	2:H:1:DG:H8	1.41	0.68
1:J:171:ILE:H	1:J:221:ASN:HD21	1.42	0.66
1:D:171:ILE:H	1:D:221:ASN:HD21	1.43	0.66
1:A:185:GLY:O	1:A:274:LYS:NZ	2.27	0.66
1:J:253:ASN:OD1	1:J:256:ASP:N	2.29	0.66
3:F:12:DA:H2"	3:F:13:DT:H72	1.78	0.65
1:G:53:ARG:HD3	1:G:83:ARG:HD3	1.77	0.65
1:A:262:ILE:O	1:A:266:ASN:ND2	2.24	0.64
1:G:185:GLY:O	1:G:274:LYS:NZ	2.28	0.63
1:G:223:ILE:O	1:G:227:ASN:N	2.32	0.63
1:J:170:GLU:HA	1:J:221:ASN:ND2	2.14	0.63
1:D:64:LEU:HD12	1:D:64:LEU:H	1.62	0.62
1:D:203:VAL:HG11	1:D:302:ILE:HD13	1.82	0.62
1:A:59:LYS:HD2	1:A:273:VAL:HG21	1.80	0.62
1:A:64:LEU:H	1:A:64:LEU:HD12	1.64	0.61
1:D:64:LEU:HD23	1:D:84:ILE:HG22	1.83	0.61
1:G:53:ARG:HD3	1:G:83:ARG:HH11	1.65	0.61
1:J:129:THR:HG22	1:J:131:GLU:N	2.13	0.60
1:G:64:LEU:HD12	1:G:64:LEU:H	1.66	0.60
1:J:64:LEU:HD12	1:J:64:LEU:H	1.66	0.59
1:D:30:SER:O	1:D:108:LYS:NZ	2.35	0.59
1:G:64:LEU:HD23	1:G:84:ILE:HG22	1.84	0.59
1:J:75:ILE:HD11	1:J:85:ARG:NH1	2.16	0.59
1:G:288:ILE:HG23	1:G:293:HIS:CE1	2.36	0.59
1:G:59:LYS:HD2	1:G:273:VAL:HG21	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:93:GLU:O	1:J:96:ILE:HG22	2.04	0.57
1:D:27:ASP:OD1	1:D:106:THR:OG1	2.19	0.56
1:J:64:LEU:HD23	1:J:84:ILE:HG22	1.87	0.56
1:D:104:LEU:HD13	1:D:108:LYS:HG3	1.86	0.56
1:A:173:ASP:OD1	1:A:174:SER:N	2.38	0.56
1:D:288:ILE:HG23	1:D:293:HIS:CE1	2.42	0.55
1:A:129:THR:HG22	1:A:131:GLU:N	2.15	0.55
1:A:262:ILE:HG12	1:A:266:ASN:ND2	2.20	0.55
1:J:284:VAL:HA	1:J:287:LEU:CD1	2.36	0.55
1:D:26:ALA:O	1:D:108:LYS:NZ	2.40	0.55
2:E:24:DT:H2'	2:E:25:DT:H71	1.88	0.55
1:J:262:ILE:HG22	1:J:263:PRO:HD3	1.88	0.55
1:D:93:GLU:O	1:D:96:ILE:HG22	2.06	0.55
1:A:223:ILE:O	1:A:227:ASN:N	2.40	0.55
1:A:23:THR:O	1:A:27:ASP:HB2	2.07	0.54
1:G:129:THR:HG22	1:G:131:GLU:N	2.17	0.54
1:J:212:ASP:HA	1:J:246:CYS:HA	1.89	0.54
1:A:68:ARG:HD3	1:A:74:GLY:O	2.08	0.54
1:J:260:LYS:C	1:J:263:PRO:HD2	2.29	0.53
3:C:12:DA:H2"	3:C:13:DT:C7	2.39	0.53
1:D:129:THR:HG22	1:D:131:GLU:N	2.20	0.52
1:A:292:GLU:O	1:A:298:GLY:HA3	2.08	0.52
1:G:63:ILE:O	1:G:67:ILE:HD12	2.09	0.52
1:G:262:ILE:HG22	1:G:263:PRO:HD3	1.91	0.52
1:D:223:ILE:O	1:D:227:ASN:N	2.43	0.52
1:G:223:ILE:HD11	1:G:231:ILE:HG13	1.92	0.52
1:A:288:ILE:HG23	1:A:293:HIS:CE1	2.45	0.52
3:I:12:DA:H2"	3:I:13:DT:H71	1.91	0.52
1:D:262:ILE:HG22	1:D:263:PRO:HD3	1.92	0.51
3:F:12:DA:H2"	3:F:13:DT:C7	2.39	0.51
1:D:202:GLN:HE22	2:E:6:DA:H62	1.58	0.51
2:H:1:DG:H8	2:H:1:DG:O5'	1.93	0.51
1:D:287:LEU:HD12	1:D:288:ILE:N	2.26	0.51
1:G:260:LYS:C	1:G:263:PRO:HD2	2.30	0.51
1:J:203:VAL:HG11	1:J:302:ILE:HD13	1.92	0.51
1:G:58:ASN:OD1	1:G:78:SER:OG	2.29	0.51
3:L:12:DA:H2"	3:L:13:DT:C7	2.41	0.51
1:D:253:ASN:HB3	1:D:256:ASP:HB2	1.92	0.50
1:J:57:HIS:HB2	2:K:16:DT:OP2	2.11	0.50
2:H:10:DG:H1	3:I:18:DC:N4	2.09	0.50
1:G:29:GLU:HG2	2:H:17:DA:OP2	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:ILE:HG22	1:A:263:PRO:HD3	1.93	0.50
1:A:88:SER:O	1:A:92:LEU:HD13	2.11	0.50
1:J:35:THR:HB	1:J:49:ARG:HG2	1.94	0.50
1:G:292:GLU:O	1:G:298:GLY:HA3	2.11	0.50
1:G:173:ASP:OD1	1:G:174:SER:N	2.45	0.50
1:G:262:ILE:HD11	1:G:282:CYS:SG	2.52	0.50
1:A:36:VAL:HB	1:A:151:LEU:HD11	1.93	0.49
2:E:10:DG:H1	3:F:18:DC:H42	1.61	0.49
1:J:59:LYS:HD2	1:J:273:VAL:HG21	1.94	0.49
1:G:306:LYS:NZ	3:I:19:DG:OP2	2.34	0.49
2:H:16:DT:C2'	2:H:17:DA:H5"	2.37	0.49
3:I:12:DA:H2"	3:I:13:DT:C7	2.43	0.49
1:J:262:ILE:HD11	1:J:282:CYS:SG	2.52	0.49
3:L:12:DA:H2"	3:L:13:DT:H71	1.93	0.49
1:A:253:ASN:OD1	1:A:256:ASP:N	2.41	0.49
1:A:212:ASP:HA	1:A:246:CYS:HA	1.95	0.49
1:G:287:LEU:HD11	1:G:302:ILE:HD11	1.95	0.49
1:G:294:LEU:CD2	2:H:4:DC:H5"	2.42	0.49
1:D:59:LYS:HD2	1:D:273:VAL:HG21	1.95	0.49
1:A:140:LYS:O	1:A:140:LYS:HG3	2.12	0.49
1:J:288:ILE:HG23	1:J:293:HIS:CE1	2.48	0.49
1:D:171:ILE:H	1:D:221:ASN:ND2	2.08	0.48
3:I:1:DC:H2'	3:I:2:DA:C8	2.49	0.48
2:E:10:DG:H1	3:F:18:DC:N4	2.11	0.48
1:D:105:ILE:HG23	1:D:169:GLN:CB	2.44	0.48
1:D:260:LYS:C	1:D:263:PRO:HD2	2.34	0.48
1:J:107:GLN:HA	1:J:166:VAL:HG11	1.95	0.48
1:D:165:LEU:HG	1:D:167:THR:OG1	2.13	0.48
1:D:260:LYS:O	1:D:263:PRO:HD2	2.13	0.48
1:A:53:ARG:HD3	1:A:83:ARG:HH11	1.79	0.48
1:G:287:LEU:HD12	1:G:288:ILE:N	2.28	0.48
1:A:262:ILE:HG12	1:A:266:ASN:HD21	1.78	0.48
1:G:223:ILE:CD1	1:G:231:ILE:HG13	2.44	0.48
1:D:232:ARG:HH22	2:E:9:DC:H6	1.62	0.47
1:D:36:VAL:HB	1:D:151:LEU:HD11	1.96	0.47
1:J:140:LYS:O	1:J:140:LYS:HG3	2.14	0.47
1:J:253:ASN:HB3	1:J:256:ASP:HB2	1.97	0.47
1:D:171:ILE:HG12	1:D:221:ASN:ND2	2.30	0.47
1:G:36:VAL:HB	1:G:151:LEU:HD21	1.96	0.47
1:J:105:ILE:HD11	1:J:172:PRO:HG3	1.96	0.47
2:H:15:DA:H1'	2:H:16:DT:O4'	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:VAL:HG11	1:A:136:ILE:HD13	1.95	0.47
1:J:23:THR:O	1:J:27:ASP:HB2	2.14	0.47
1:D:53:ARG:HD3	1:D:83:ARG:HH11	1.80	0.47
1:A:25:PHE:CE1	1:A:183:ALA:HB2	2.50	0.47
1:A:216:ILE:HG12	1:A:247:VAL:HG11	1.96	0.47
1:A:27:ASP:OD1	1:A:106:THR:OG1	2.19	0.47
1:D:42:ARG:NH1	1:D:47:SER:HA	2.29	0.47
2:H:10:DG:H1	3:I:18:DC:H42	1.63	0.47
1:G:52:PHE:HD2	1:G:86:PHE:CD1	2.33	0.47
1:G:140:LYS:O	1:G:140:LYS:HG3	2.15	0.47
1:A:223:ILE:CD1	1:A:231:ILE:HG13	2.45	0.46
1:G:260:LYS:O	1:G:263:PRO:HD2	2.15	0.46
1:J:42:ARG:HD3	3:L:3:DA:OP2	2.16	0.46
1:J:58:ASN:OD1	1:J:78:SER:OG	2.33	0.46
1:J:107:GLN:HA	1:J:166:VAL:CG1	2.46	0.46
1:D:309:MET:HE2	1:D:309:MET:HB2	1.60	0.46
1:J:126:GLU:O	1:J:132:GLY:HA3	2.16	0.46
1:D:79:ASP:HB2	1:D:80:ILE:H	1.48	0.46
1:G:212:ASP:HA	1:G:246:CYS:HA	1.97	0.46
1:G:63:ILE:HD11	1:G:272:GLY:CA	2.46	0.46
1:G:55:GLY:C	1:G:56:LEU:HD23	2.37	0.45
1:A:64:LEU:HD23	1:A:84:ILE:HG22	1.97	0.45
1:A:181:THR:HG21	1:A:265:PHE:HZ	1.81	0.45
1:G:309:MET:HE2	1:G:309:MET:HB2	1.72	0.45
1:J:294:LEU:CD2	2:K:4:DC:H5"	2.46	0.45
1:D:55:GLY:C	1:D:56:LEU:HD23	2.36	0.45
1:G:79:ASP:HB2	1:G:80:ILE:H	1.46	0.45
1:A:232:ARG:HH12	2:B:9:DC:H5	1.64	0.45
1:G:19:PRO:HB3	1:G:70:TYR:OH	2.16	0.45
1:J:171:ILE:N	1:J:221:ASN:HD21	2.13	0.45
1:A:106:THR:HG22	1:A:218:LEU:HD22	1.98	0.45
1:A:260:LYS:O	1:A:263:PRO:HD2	2.17	0.45
1:J:55:GLY:O	1:J:56:LEU:HD23	2.17	0.45
1:J:292:GLU:O	1:J:295:THR:HG22	2.17	0.45
1:A:302:ILE:O	1:A:305:ILE:N	2.50	0.44
1:A:63:ILE:HD11	1:A:272:GLY:CA	2.47	0.44
1:J:19:PRO:HB3	1:J:70:TYR:OH	2.18	0.44
1:A:309:MET:HE2	1:A:309:MET:HB2	1.65	0.44
1:A:79:ASP:HB2	1:A:80:ILE:H	1.55	0.44
1:J:79:ASP:HB2	1:J:80:ILE:H	1.38	0.43
2:H:13:DG:O6	3:I:14:DA:N6	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:216:ILE:HG12	1:J:247:VAL:HG11	2.00	0.43
1:D:262:ILE:HD11	1:D:282:CYS:SG	2.59	0.43
1:D:292:GLU:O	1:D:295:THR:HG22	2.18	0.43
1:J:151:LEU:HD23	1:J:151:LEU:HA	1.78	0.43
1:J:75:ILE:HD11	1:J:85:ARG:CZ	2.49	0.43
1:J:302:ILE:O	1:J:305:ILE:N	2.51	0.43
2:B:10:DG:H1	3:C:18:DC:N4	2.17	0.43
1:J:309:MET:HB2	1:J:309:MET:HE2	1.72	0.43
2:K:1:DG:HO5'	2:K:1:DG:H8	1.64	0.43
1:A:170:GLU:HA	1:A:221:ASN:ND2	2.34	0.43
3:I:22:DT:H2"	3:I:23:DA:H8	1.83	0.43
1:A:260:LYS:C	1:A:263:PRO:HD2	2.38	0.43
1:G:55:GLY:O	1:G:56:LEU:HD23	2.19	0.43
1:J:53:ARG:HD3	1:J:83:ARG:HH11	1.84	0.43
1:J:36:VAL:HB	1:J:151:LEU:HD11	2.00	0.43
1:A:232:ARG:HG3	1:A:233:ILE:N	2.33	0.42
1:G:148:SER:OG	1:G:151:LEU:HD13	2.20	0.42
1:J:232:ARG:NH1	2:K:8:DA:H2'	2.34	0.42
2:E:13:DG:O6	3:F:14:DA:N6	2.52	0.42
3:L:1:DC:H2'	3:L:2:DA:C8	2.54	0.42
1:A:120:TYR:HD1	1:A:120:TYR:HA	1.65	0.42
1:A:232:ARG:HH11	1:A:232:ARG:HD3	1.56	0.42
1:A:292:GLU:O	1:A:295:THR:HG22	2.19	0.42
1:G:203:VAL:HG11	1:G:302:ILE:HD13	2.00	0.42
1:J:260:LYS:O	1:J:263:PRO:HD2	2.19	0.42
1:D:232:ARG:NH1	2:E:8:DA:H2'	2.34	0.42
1:A:49:ARG:HA	1:A:50:PRO:HD3	1.87	0.42
1:A:223:ILE:HG23	1:A:228:CYS:O	2.19	0.42
1:A:262:ILE:HD11	1:A:282:CYS:SG	2.60	0.42
1:G:105:ILE:HD11	1:G:172:PRO:HG3	2.01	0.42
2:K:4:DC:H2'	2:K:5:DT:C6	2.55	0.42
1:J:75:ILE:O	1:J:75:ILE:HG13	2.20	0.41
3:L:12:DA:H2"	3:L:13:DT:C5	2.56	0.41
1:D:25:PHE:CE1	1:D:183:ALA:HB2	2.55	0.41
1:J:27:ASP:OD2	1:J:105:ILE:N	2.47	0.41
1:J:223:ILE:CD1	1:J:231:ILE:HG13	2.51	0.41
1:D:55:GLY:O	1:D:56:LEU:HD23	2.20	0.41
2:B:1:DG:HO5'	2:B:1:DG:H8	1.65	0.41
1:G:25:PHE:CE1	1:G:183:ALA:HB2	2.55	0.41
2:E:1:DG:C2	3:C:2:DA:C2	3.09	0.41
2:E:13:DG:C6	3:F:14:DA:N6	2.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:GLY:C	1:A:56:LEU:HD23	2.41	0.41
2:H:12:DC:N4	3:I:15:DC:N4	2.69	0.41
1:J:49:ARG:HA	1:J:50:PRO:HD3	1.84	0.41
1:J:68:ARG:HB2	1:J:76:ILE:HD11	2.02	0.41
1:J:104:LEU:HD13	1:J:108:LYS:HG3	2.02	0.41
1:J:106:THR:HA	1:J:215:ASP:OD2	2.21	0.41
1:J:232:ARG:HH12	2:K:9:DC:H6	1.68	0.41
1:G:292:GLU:O	1:G:295:THR:HG22	2.21	0.41
1:J:29:GLU:HG2	2:K:17:DA:OP2	2.20	0.41
1:A:171:ILE:H	1:A:221:ASN:HD21	1.68	0.40
1:G:294:LEU:HD23	2:H:4:DC:H5"	2.03	0.40
1:A:89:LEU:HG	3:C:5:DT:OP2	2.20	0.40
1:A:159:ILE:HA	1:A:160:PRO:HD3	1.92	0.40
1:D:208:GLN:HA	1:D:249:LEU:O	2.21	0.40
1:G:230:ASN:HD21	2:H:7:DA:H3'	1.86	0.40
1:G:305:ILE:HD13	1:G:305:ILE:HA	1.88	0.40
1:J:55:GLY:C	1:J:56:LEU:HD23	2.42	0.40
1:D:115:PHE:O	1:D:119:PHE:N	2.47	0.40
1:A:44:THR:HB	1:A:133:LEU:HD22	2.03	0.40
1:G:234:ARG:HD3	1:G:245:THR:HG22	2.04	0.40
1:J:96:ILE:HD11	1:J:116:LYS:HG3	2.04	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	284/315 (90%)	273 (96%)	11 (4%)	0	100	100
1	D	284/315 (90%)	276 (97%)	8 (3%)	0	100	100
1	G	284/315 (90%)	276 (97%)	8 (3%)	0	100	100
1	J	285/315 (90%)	278 (98%)	7 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1137/1260 (90%)	1103 (97%)	34 (3%)	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	211/282 (75%)	207 (98%)	4 (2%)	57	83
1	D	210/282 (74%)	207 (99%)	3 (1%)	67	88
1	G	208/282 (74%)	204 (98%)	4 (2%)	57	83
1	J	211/282 (75%)	203 (96%)	8 (4%)	33	65
All	All	840/1128 (74%)	821 (98%)	19 (2%)	50	79

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	80	ILE
1	D	108	LYS
1	D	216	ILE
1	A	79	ASP
1	A	80	ILE
1	A	108	LYS
1	A	218	LEU
1	G	27	ASP
1	G	79	ASP
1	G	197	LEU
1	G	232	ARG
1	J	27	ASP
1	J	75	ILE
1	J	80	ILE
1	J	97	ASN
1	J	108	LYS
1	J	197	LEU

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Mol	Chain	Res	Type
1	J	287	LEU
1	J	296	SER

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

Mol	Chain	Res	Type
1	D	221	ASN
1	A	58	ASN
1	A	175	ASN
1	A	221	ASN
1	G	211	GLN
1	J	221	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 12 ligands modelled in this entry, 12 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	288/315 (91%)	0.31	21 (7%)	15	12	46, 74, 99, 126	0
1	D	288/315 (91%)	0.24	20 (6%)	16	13	46, 73, 94, 112	0
1	G	288/315 (91%)	0.29	23 (7%)	12	10	49, 74, 96, 123	0
1	J	289/315 (91%)	0.41	27 (9%)	8	6	51, 75, 97, 135	0
2	B	26/26 (100%)	-0.60	0	100	100	54, 69, 87, 90	0
2	E	26/26 (100%)	-0.51	0	100	100	56, 69, 92, 95	0
2	H	26/26 (100%)	-0.54	0	100	100	60, 71, 90, 96	0
2	K	26/26 (100%)	-0.53	0	100	100	60, 76, 101, 106	0
3	C	26/26 (100%)	-0.39	1 (3%)	40	37	54, 74, 91, 95	0
3	F	26/26 (100%)	-0.37	0	100	100	53, 77, 85, 94	0
3	I	26/26 (100%)	-0.42	0	100	100	54, 79, 86, 96	0
3	L	26/26 (100%)	-0.43	0	100	100	58, 82, 97, 104	0
All	All	1361/1468 (92%)	0.19	92 (6%)	17	14	46, 74, 97, 135	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	67	ILE	6.6
1	J	166	VAL	5.4
1	G	67	ILE	5.2
1	J	84	ILE	4.9
1	D	133	LEU	4.9
1	J	164	LEU	4.8
1	A	166	VAL	4.8
1	G	162	GLU	4.2
1	J	163	ARG	4.2
1	J	168	GLY	4.1
1	A	163	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	D	166	VAL	3.9
1	D	209	ILE	3.7
1	A	84	ILE	3.7
1	G	166	VAL	3.7
1	J	165	LEU	3.6
1	A	164	LEU	3.6
1	G	209	ILE	3.5
1	J	233	ILE	3.5
1	G	245	THR	3.5
1	D	191	ILE	3.5
1	G	133	LEU	3.4
1	J	86	PHE	3.3
1	J	247	VAL	3.3
1	A	99	PHE	3.2
1	J	302	ILE	3.2
1	G	189	ILE	3.1
1	D	95	VAL	3.1
1	D	137	LEU	3.1
1	D	155	PHE	3.1
1	G	22	ILE	3.1
1	A	64	LEU	3.0
1	J	26	ALA	2.9
1	J	203	VAL	2.9
1	A	203	VAL	2.9
1	D	134	ASN	2.9
1	J	99	PHE	2.9
1	J	285	VAL	2.8
1	G	222	LEU	2.7
1	J	89	LEU	2.7
1	D	22	ILE	2.7
1	A	219	MET	2.7
1	G	136	ILE	2.6
1	G	86	PHE	2.6
1	D	99	PHE	2.6
1	D	67	ILE	2.6
1	A	247	VAL	2.6
1	G	44	THR	2.5
1	G	76	ILE	2.5
1	J	76	ILE	2.5
1	J	189	ILE	2.5
1	A	287	LEU	2.5
1	A	86	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	J	245	THR	2.5
1	A	75	ILE	2.5
1	J	271	ILE	2.4
1	G	205	SER	2.4
1	D	233	ILE	2.4
1	G	134	ASN	2.4
1	G	64	LEU	2.4
1	J	145	LEU	2.4
1	G	234	ARG	2.4
1	G	168	GLY	2.4
1	A	165	LEU	2.3
1	A	145	LEU	2.3
1	A	162	GLU	2.3
1	A	133	LEU	2.3
1	D	96	ILE	2.2
1	J	207	PHE	2.2
1	J	180	PHE	2.2
3	C	26	DC	2.2
1	G	191	ILE	2.2
1	G	271	ILE	2.2
1	G	257	ILE	2.2
1	D	154	ALA	2.2
1	A	44	THR	2.2
1	D	34	LEU	2.2
1	D	249	LEU	2.2
1	D	247	VAL	2.2
1	G	203	VAL	2.2
1	A	26	ALA	2.1
1	D	161	ALA	2.1
1	J	249	LEU	2.1
1	J	299	LEU	2.1
1	J	22	ILE	2.1
1	A	104	LEU	2.0
1	G	164	LEU	2.0
1	D	271	ILE	2.0
1	D	86	PHE	2.0
1	J	137	LEU	2.0
1	A	22	ILE	2.0
1	J	146	GLY	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)

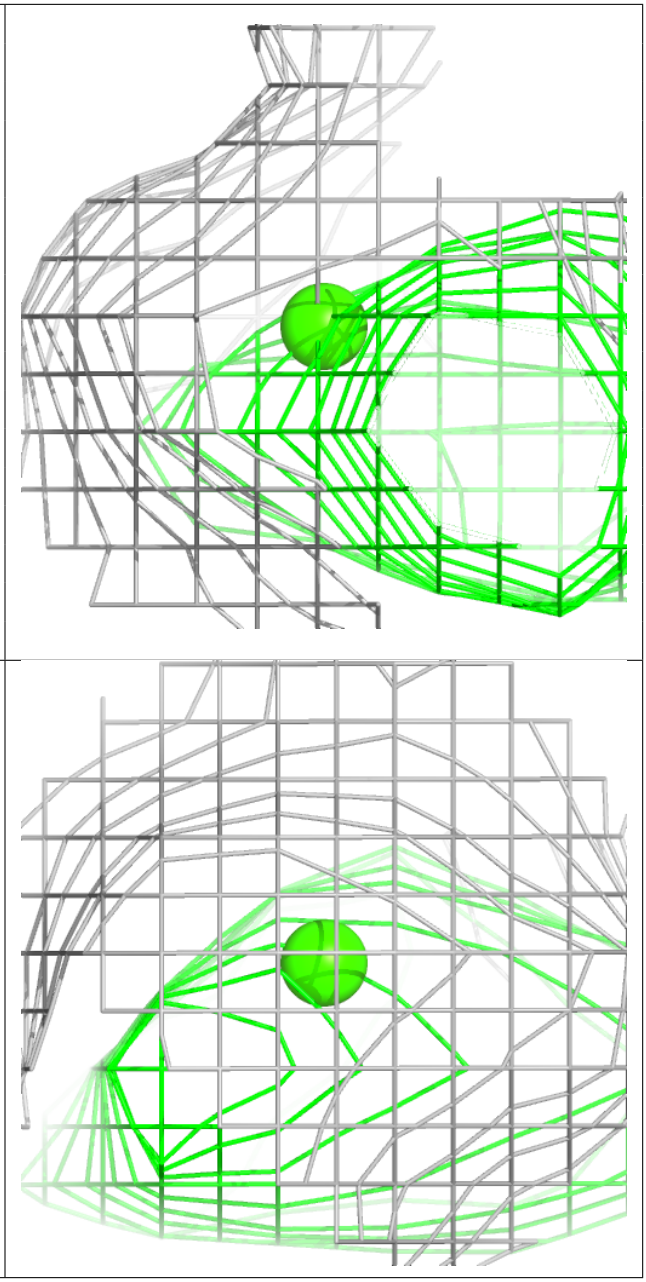
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	CA	A	401	1/1	0.89	0.28	73,73,73,73	0
4	CA	J	402	1/1	0.93	0.26	80,80,80,80	0
4	CA	G	402	1/1	0.94	0.24	63,63,63,63	0
4	CA	A	402	1/1	0.94	0.28	56,56,56,56	0
4	CA	H	101	1/1	0.95	0.09	30,30,30,30	0
4	CA	D	402	1/1	0.95	0.30	65,65,65,65	0
4	CA	G	401	1/1	0.96	0.37	90,90,90,90	0
4	CA	J	401	1/1	0.96	0.31	72,72,72,72	0
4	CA	B	101	1/1	0.96	0.15	30,30,30,30	0
4	CA	D	401	1/1	0.97	0.31	77,77,77,77	0
4	CA	D	403	1/1	0.97	0.05	30,30,30,30	0
4	CA	L	101	1/1	0.99	0.09	30,30,30,30	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.

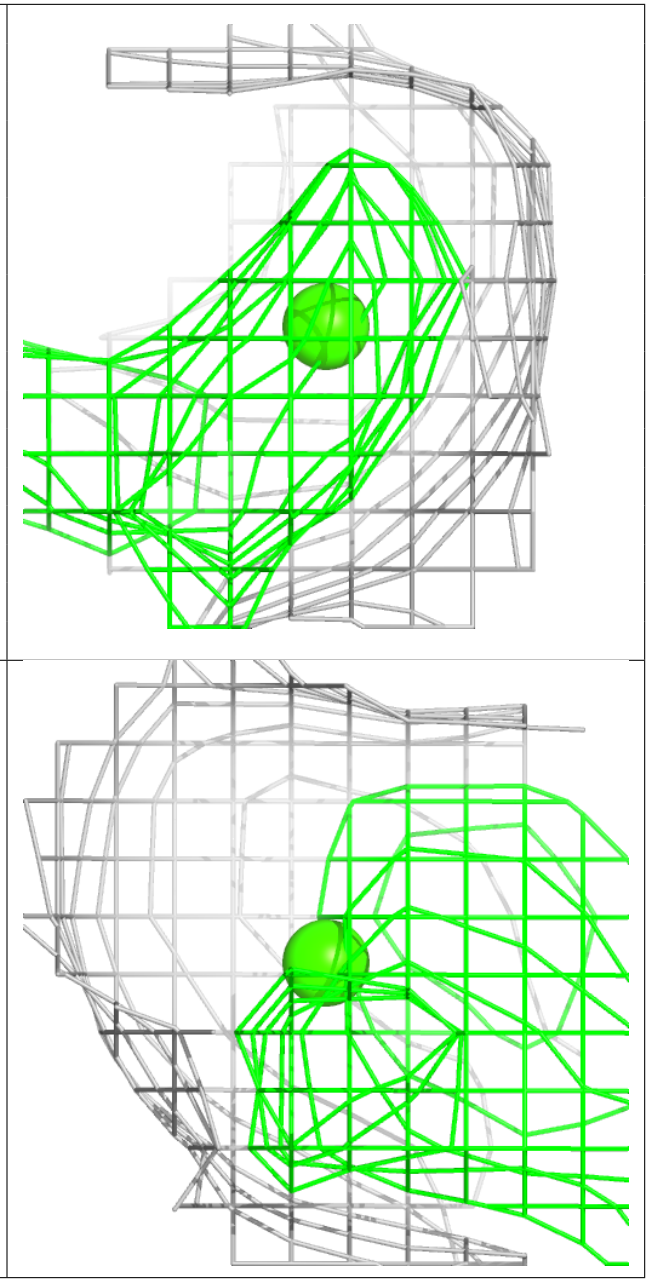
Electron density around CA A 401:

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



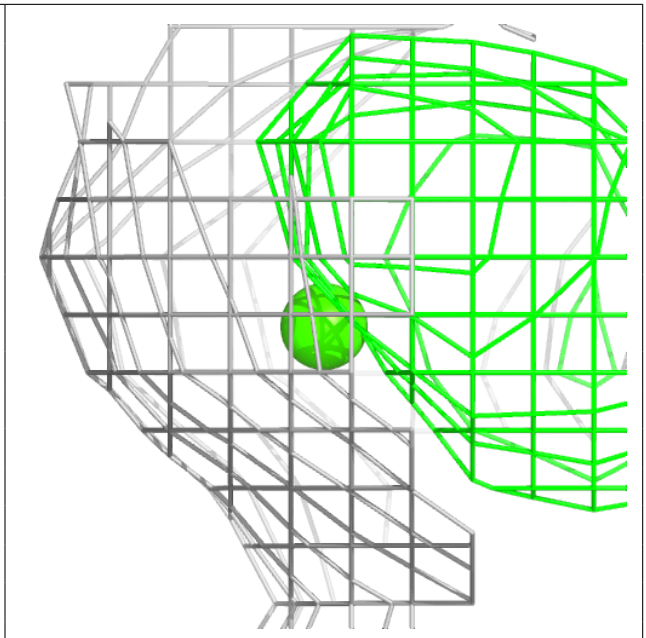
Electron density around CA J 402:

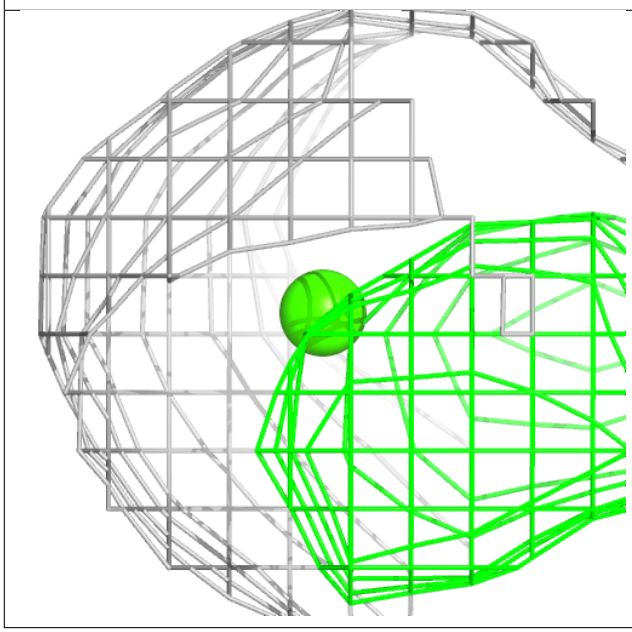
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

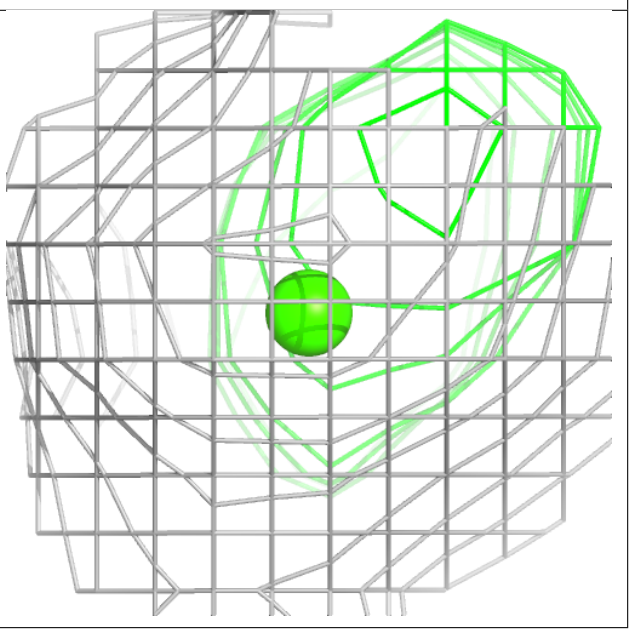


Electron density around CA G 402:

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

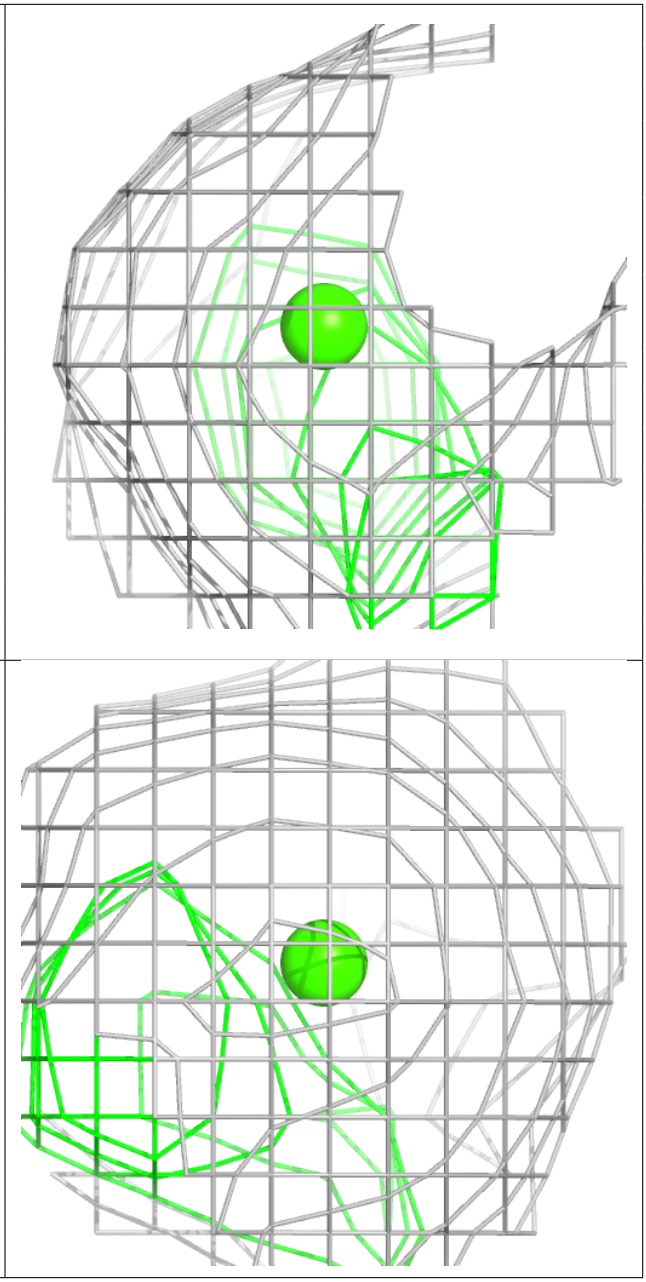






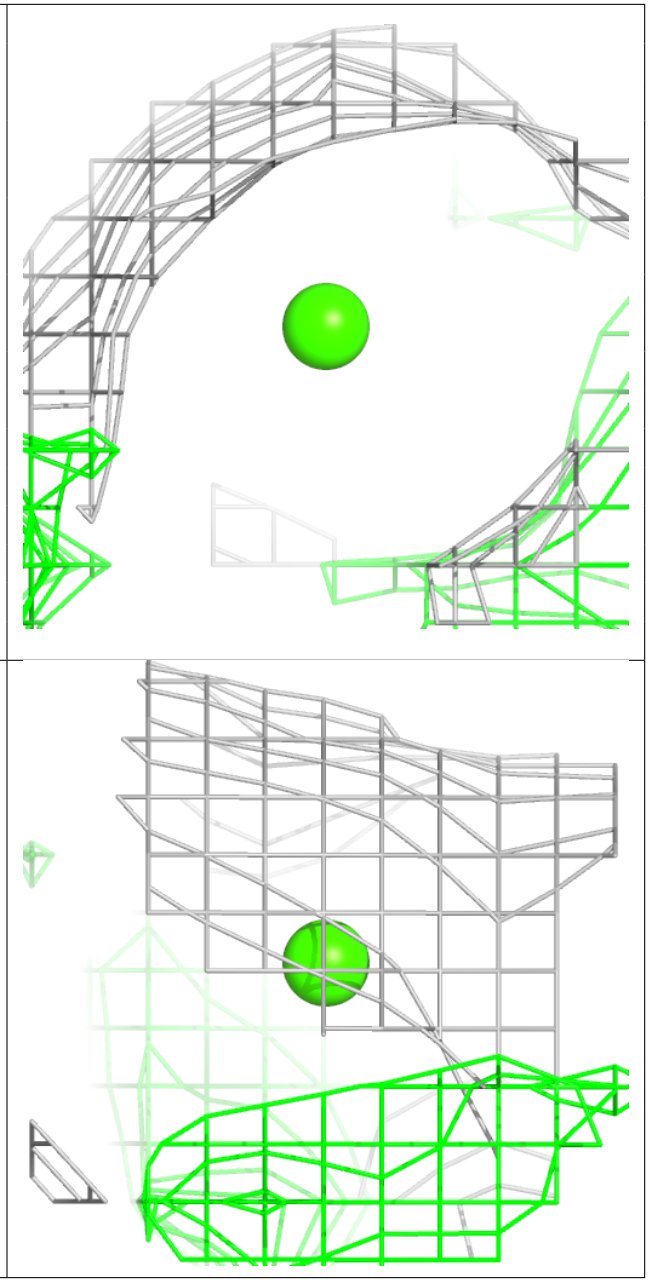
Electron density around CA A 402:

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



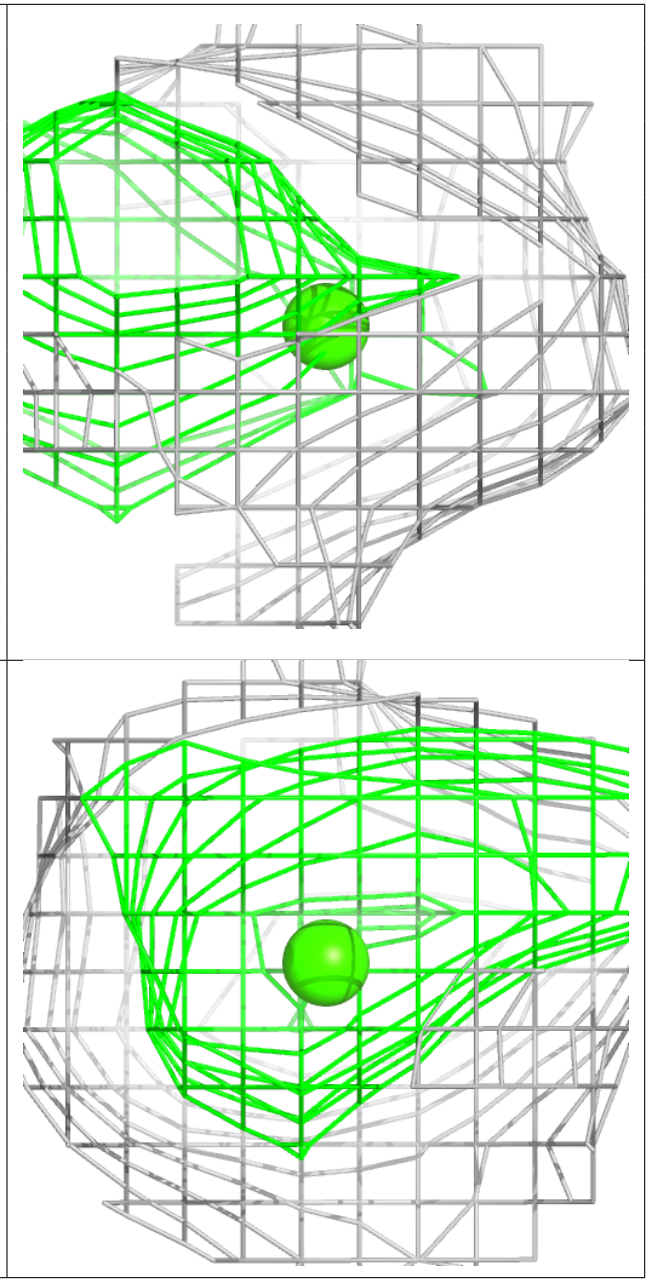
Electron density around CA H 101:

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



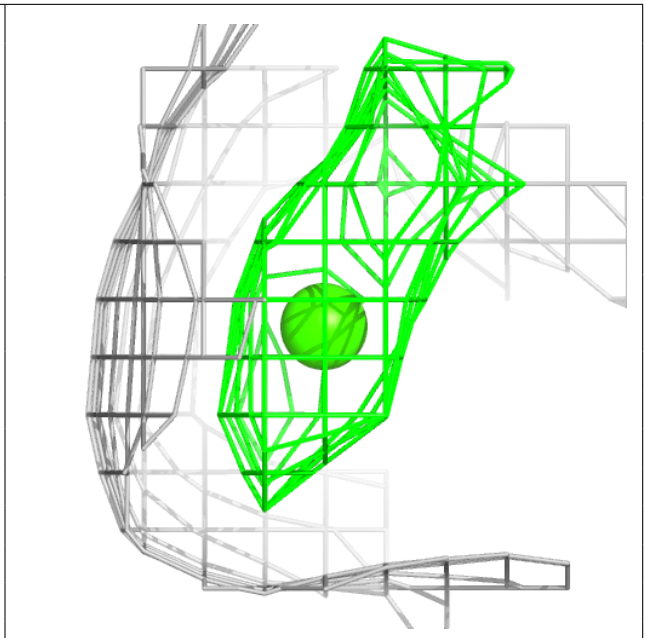
Electron density around CA D 402:

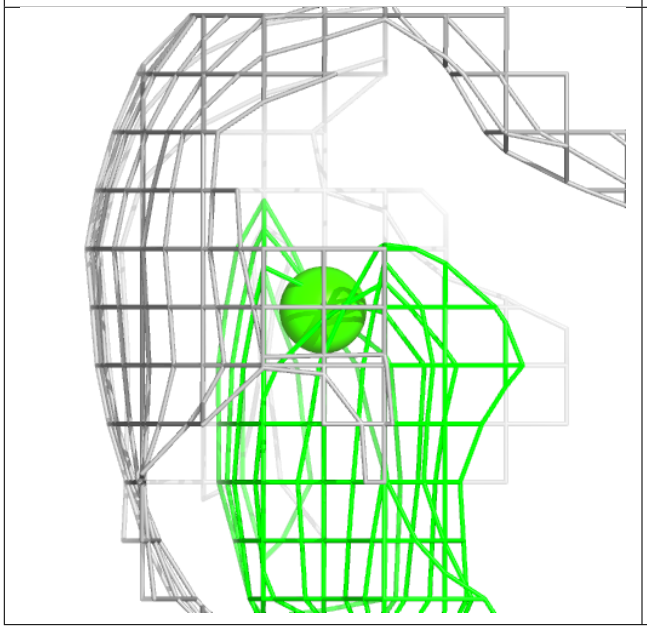
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

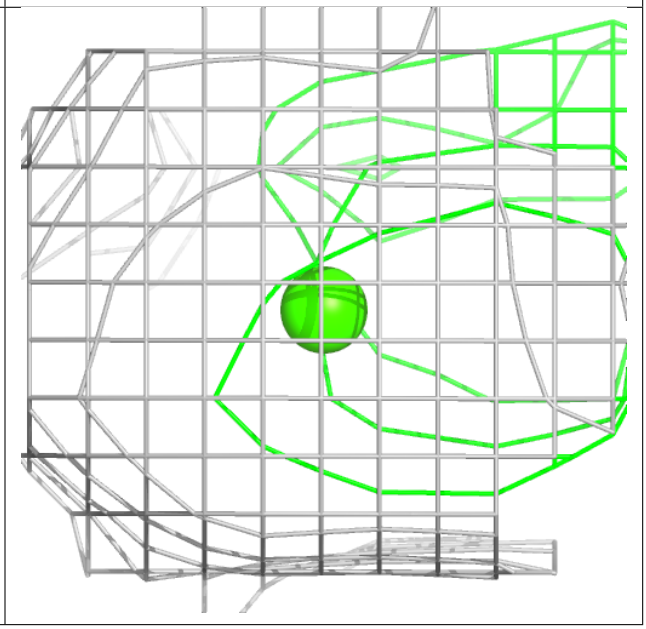


Electron density around CA G 401:

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

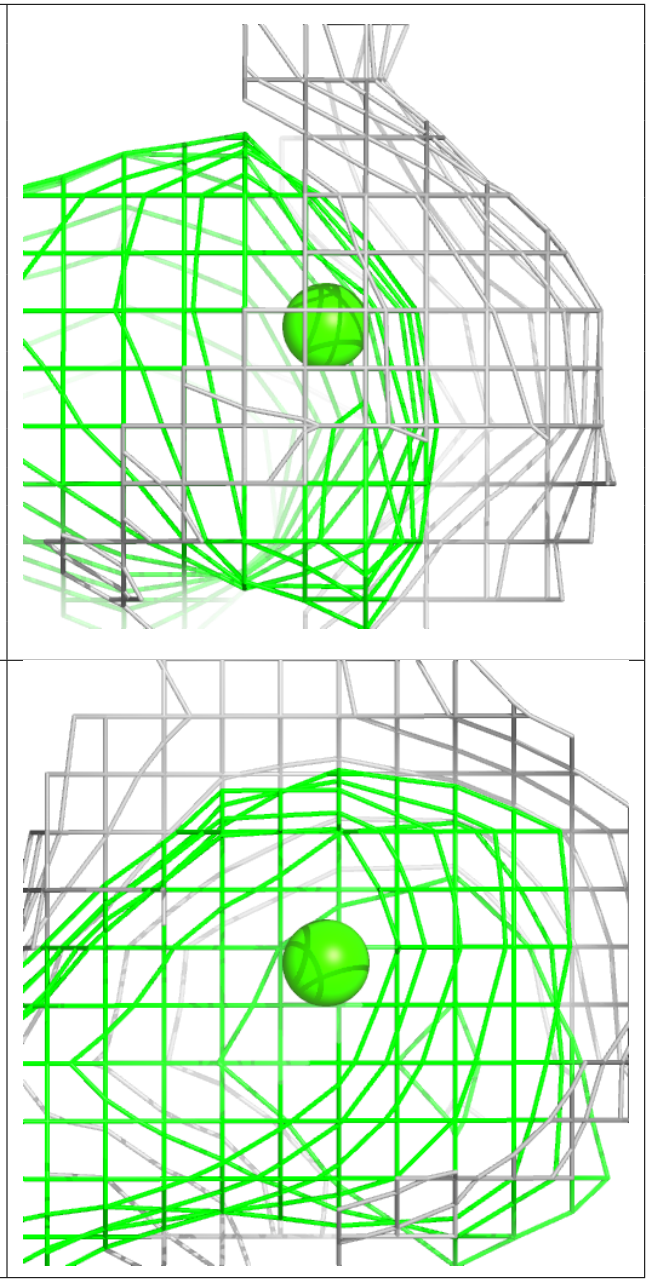






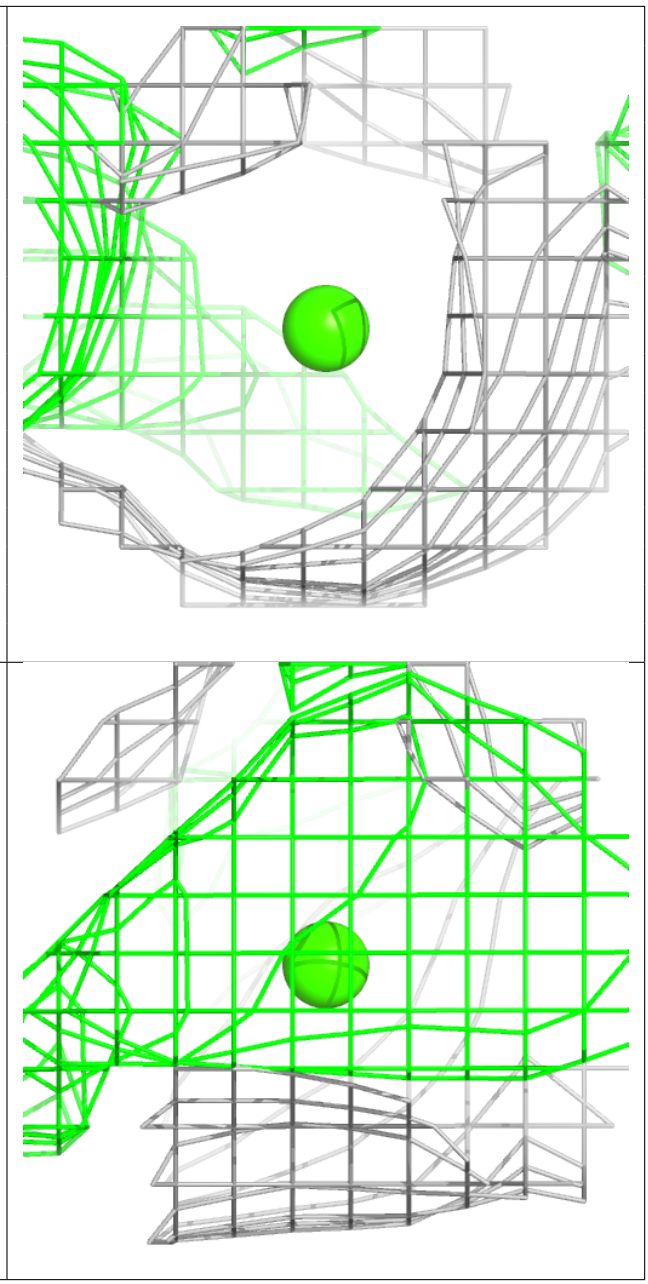
Electron density around CA J 401:

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



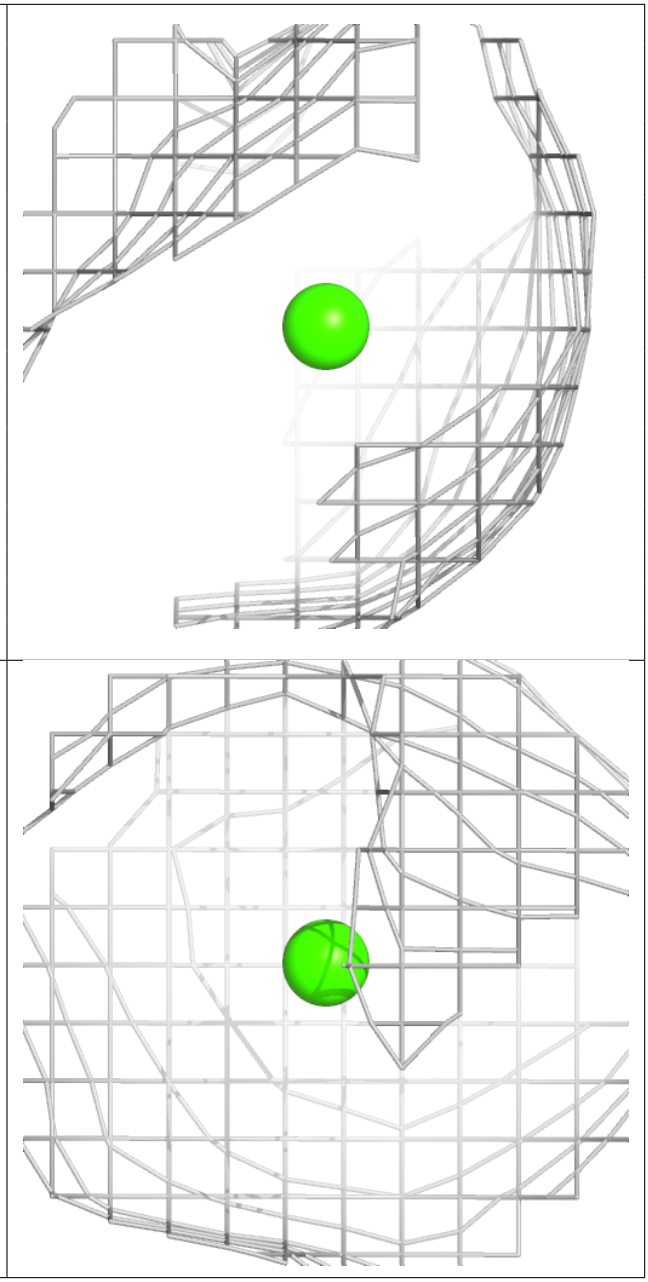
Electron density around CA B 101:

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



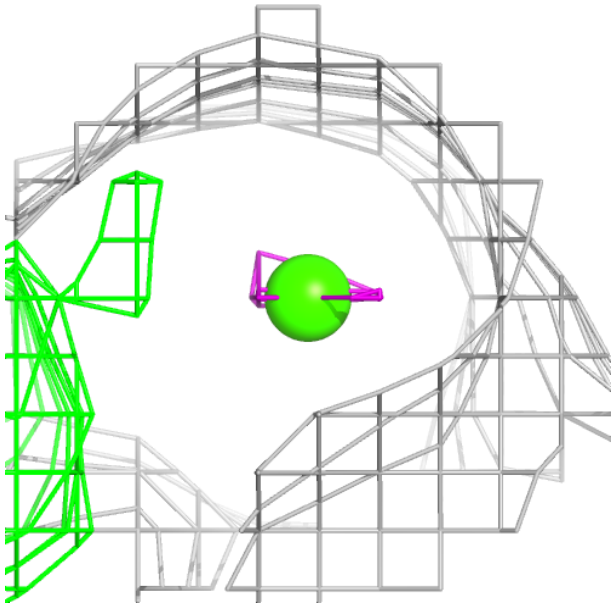
Electron density around CA D 401:

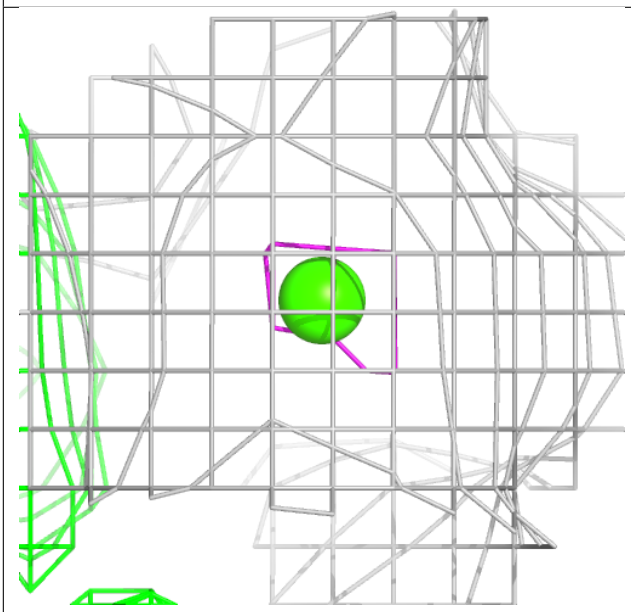
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

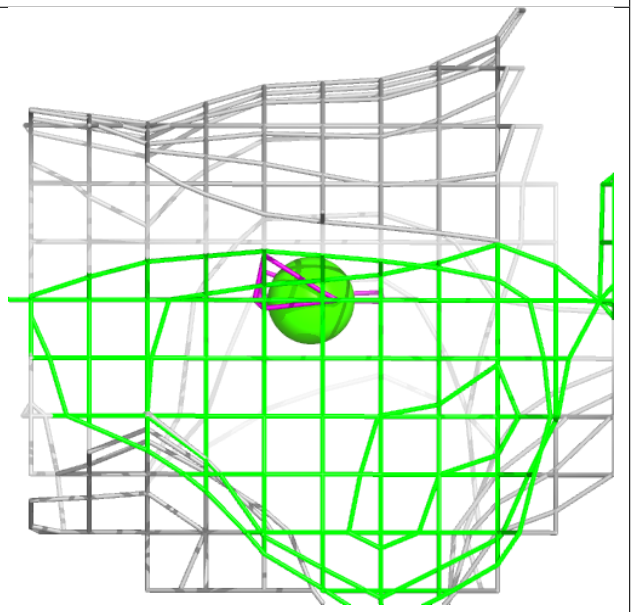


Electron density around CA D 403:

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

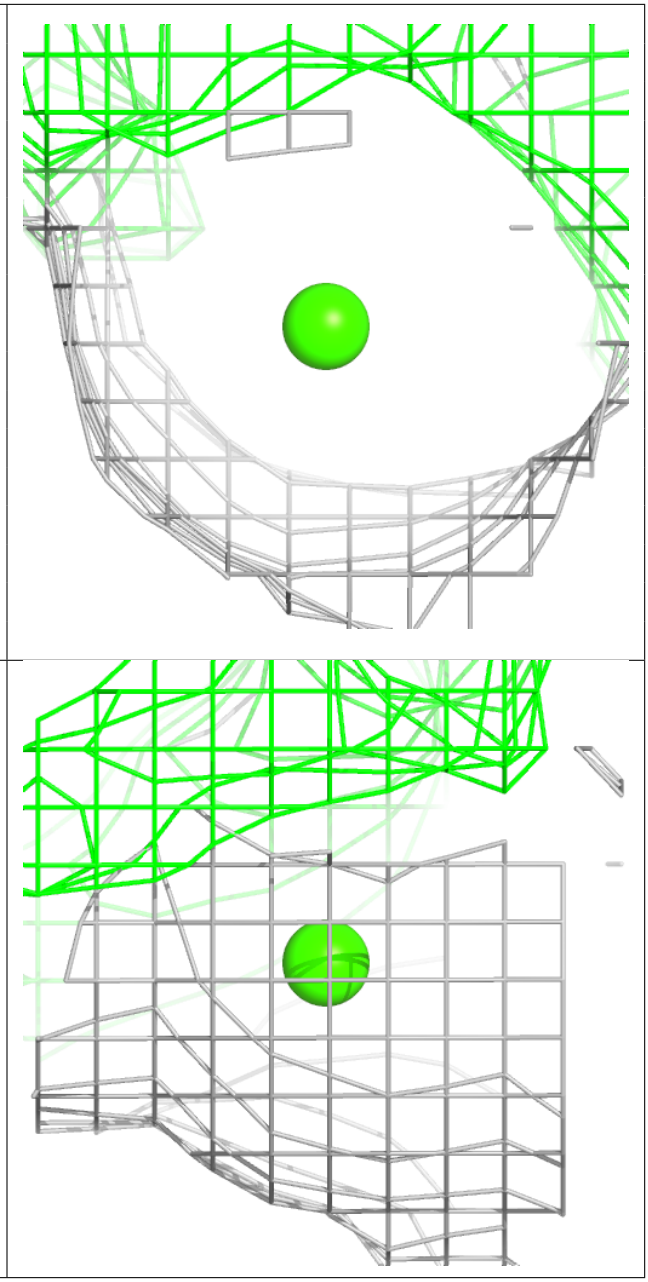






Electron density around CA L 101:

$2mF_o$-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 (i)

There are no such residues in this entry.