



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:57 PM GMT

PDB ID : 4GM2
Title : The crystal structure of a peptidase from plasmodium falciparum
Authors : El Bakkouri, M.; Jung, P.; Wernimont, A.K.; Calmettes, C.; Hui, R.; Houry, W.A.; Structural Genomics Consortium (SGC)
Deposited on : 2012-08-15
Resolution : 2.80 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

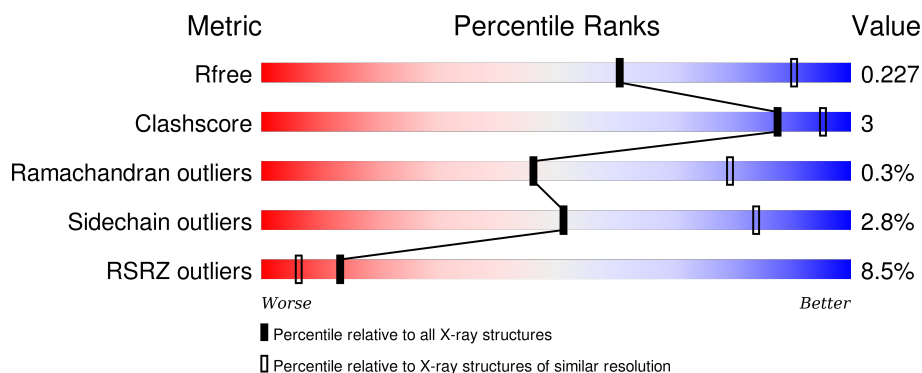
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	205	<div> <div>80%</div> <div>10%</div> <div>11%</div> </div>
1	B	205	<div> <div>2%</div> <div>75%</div> <div>6%</div> <div>20%</div> </div>
1	C	205	<div> <div>6%</div> <div>71%</div> <div>9%</div> <div>19%</div> </div>
1	D	205	<div> <div>15%</div> <div>75%</div> <div>5%</div> <div>20%</div> </div>
1	E	205	<div> <div>7%</div> <div>80%</div> <div>9%</div> <div>11%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	205	<div><div></div><div>12%</div><div>75%</div><div>7%</div><div>17%</div></div>
1	G	205	<div><div></div><div>7%</div><div>73%</div><div>10%</div><div>17%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9611 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1453	931	236	282	4			
1	B	165	Total	C	N	O	S	0	0	0
			1313	845	214	251	3			
1	C	167	Total	C	N	O	S	0	0	0
			1323	849	214	256	4			
1	D	163	Total	C	N	O	S	0	0	0
			1295	832	211	248	4			
1	E	183	Total	C	N	O	S	0	0	0
			1439	926	232	277	4			
1	F	170	Total	C	N	O	S	0	0	0
			1359	875	219	261	4			
1	G	171	Total	C	N	O	S	0	0	0
			1360	872	223	261	4			

There are 147 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
A	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
A	42	SER	-	EXPRESSION TAG	UNP Q8IL98
A	43	SER	-	EXPRESSION TAG	UNP Q8IL98
A	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
A	50	SER	-	EXPRESSION TAG	UNP Q8IL98
A	51	SER	-	EXPRESSION TAG	UNP Q8IL98
A	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
A	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
A	54	GLU	-	EXPRESSION TAG	UNP Q8IL98

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Chain	Residue	Modelled	Actual	Comment	Reference
A	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
A	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
A	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
A	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
A	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
A	60	GLY	-	EXPRESSION TAG	UNP Q8IL98
B	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
B	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
B	42	SER	-	EXPRESSION TAG	UNP Q8IL98
B	43	SER	-	EXPRESSION TAG	UNP Q8IL98
B	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
B	50	SER	-	EXPRESSION TAG	UNP Q8IL98
B	51	SER	-	EXPRESSION TAG	UNP Q8IL98
B	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
B	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
B	54	GLU	-	EXPRESSION TAG	UNP Q8IL98
B	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
B	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
B	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
B	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
B	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
B	60	GLY	-	EXPRESSION TAG	UNP Q8IL98
C	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
C	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
C	42	SER	-	EXPRESSION TAG	UNP Q8IL98
C	43	SER	-	EXPRESSION TAG	UNP Q8IL98
C	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
C	50	SER	-	EXPRESSION TAG	UNP Q8IL98
C	51	SER	-	EXPRESSION TAG	UNP Q8IL98
C	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
C	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
C	54	GLU	-	EXPRESSION TAG	UNP Q8IL98

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Chain	Residue	Modelled	Actual	Comment	Reference
C	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
C	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
C	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
C	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
C	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
C	60	GLY	-	EXPRESSION TAG	UNP Q8IL98
D	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
D	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
D	42	SER	-	EXPRESSION TAG	UNP Q8IL98
D	43	SER	-	EXPRESSION TAG	UNP Q8IL98
D	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
D	50	SER	-	EXPRESSION TAG	UNP Q8IL98
D	51	SER	-	EXPRESSION TAG	UNP Q8IL98
D	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
D	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
D	54	GLU	-	EXPRESSION TAG	UNP Q8IL98
D	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
D	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
D	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
D	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
D	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
D	60	GLY	-	EXPRESSION TAG	UNP Q8IL98
E	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
E	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
E	42	SER	-	EXPRESSION TAG	UNP Q8IL98
E	43	SER	-	EXPRESSION TAG	UNP Q8IL98
E	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
E	50	SER	-	EXPRESSION TAG	UNP Q8IL98
E	51	SER	-	EXPRESSION TAG	UNP Q8IL98
E	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
E	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
E	54	GLU	-	EXPRESSION TAG	UNP Q8IL98

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Chain	Residue	Modelled	Actual	Comment	Reference
E	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
E	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
E	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
E	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
E	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
E	60	GLY	-	INITIATING METHIONINE	UNP Q8IL98
F	40	MET	-	EXPRESSION TAG	UNP Q8IL98
F	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
F	42	SER	-	EXPRESSION TAG	UNP Q8IL98
F	43	SER	-	EXPRESSION TAG	UNP Q8IL98
F	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
F	50	SER	-	EXPRESSION TAG	UNP Q8IL98
F	51	SER	-	EXPRESSION TAG	UNP Q8IL98
F	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
F	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
F	54	GLU	-	EXPRESSION TAG	UNP Q8IL98
F	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
F	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
F	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
F	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
F	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
F	60	GLY	-	EXPRESSION TAG	UNP Q8IL98
G	40	MET	-	INITIATING METHIONINE	UNP Q8IL98
G	41	GLY	-	EXPRESSION TAG	UNP Q8IL98
G	42	SER	-	EXPRESSION TAG	UNP Q8IL98
G	43	SER	-	EXPRESSION TAG	UNP Q8IL98
G	44	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	45	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	46	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	47	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	48	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	49	HIS	-	EXPRESSION TAG	UNP Q8IL98
G	50	SER	-	EXPRESSION TAG	UNP Q8IL98
G	51	SER	-	EXPRESSION TAG	UNP Q8IL98
G	52	GLY	-	EXPRESSION TAG	UNP Q8IL98
G	53	ARG	-	EXPRESSION TAG	UNP Q8IL98
G	54	GLU	-	EXPRESSION TAG	UNP Q8IL98

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Chain	Residue	Modelled	Actual	Comment	Reference
G	55	ASN	-	EXPRESSION TAG	UNP Q8IL98
G	56	LEU	-	EXPRESSION TAG	UNP Q8IL98
G	57	TYR	-	EXPRESSION TAG	UNP Q8IL98
G	58	PHE	-	EXPRESSION TAG	UNP Q8IL98
G	59	GLN	-	EXPRESSION TAG	UNP Q8IL98
G	60	GLY	-	EXPRESSION TAG	UNP Q8IL98

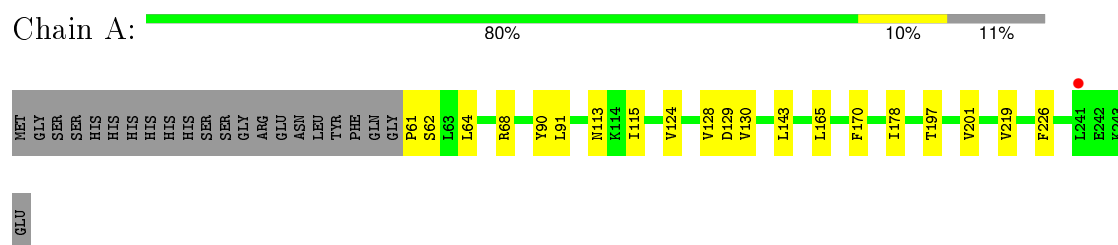
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	17	Total O 17 17	0	0
2	B	10	Total O 10 10	0	0
2	C	11	Total O 11 11	0	0
2	D	5	Total O 5 5	0	0
2	E	5	Total O 5 5	0	0
2	F	9	Total O 9 9	0	0
2	G	12	Total O 12 12	0	0

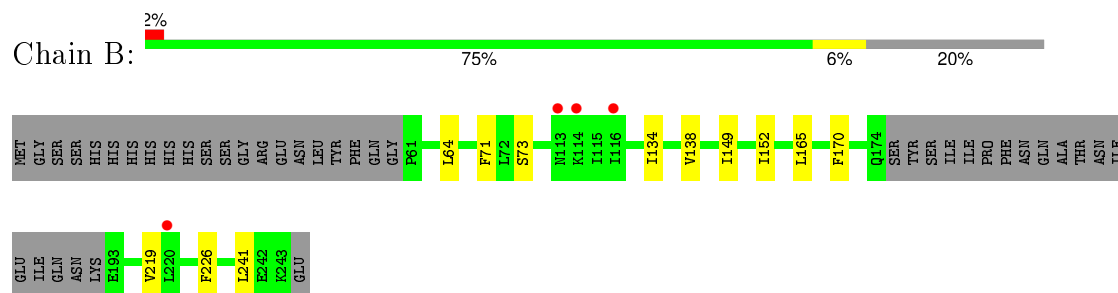
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

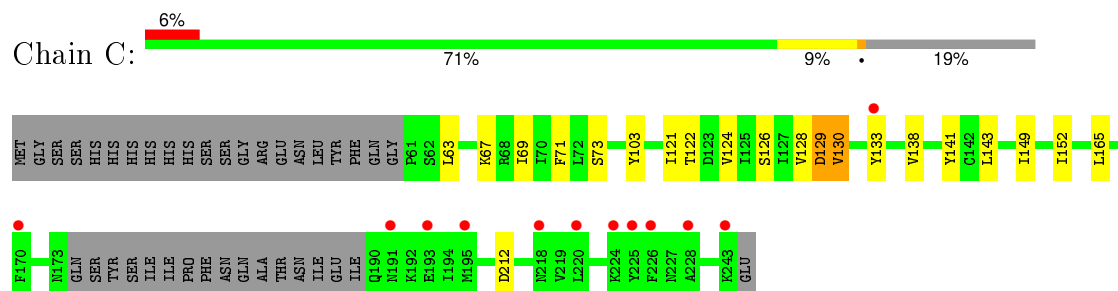
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



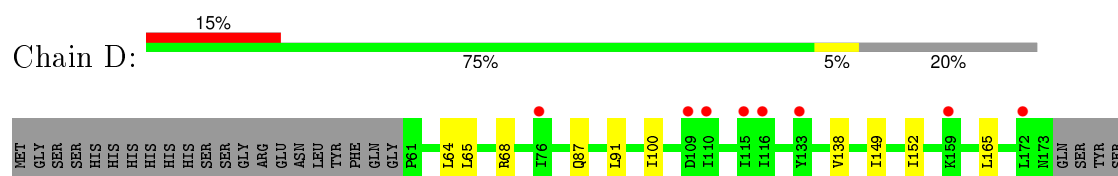
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

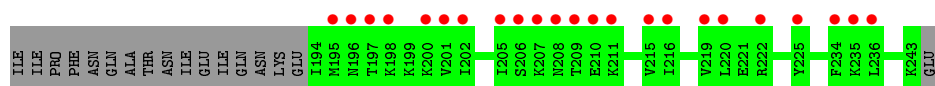


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

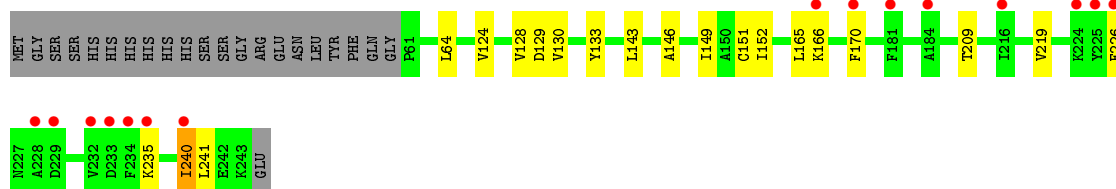
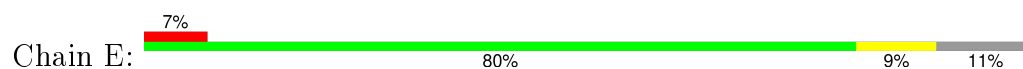


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

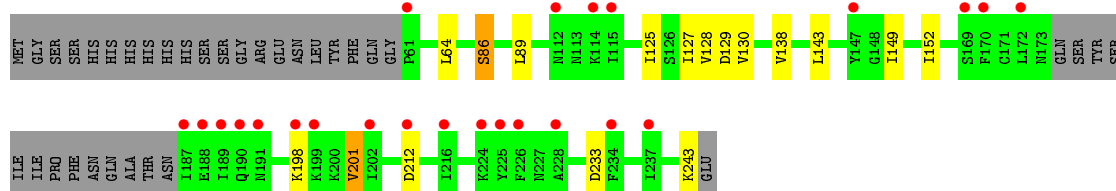
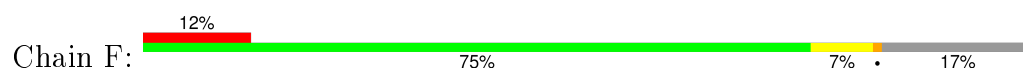




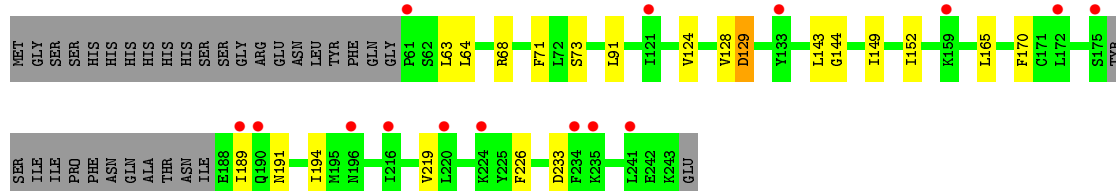
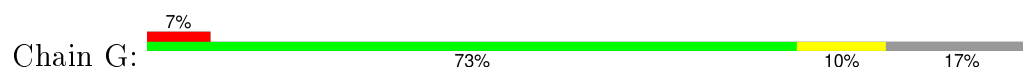
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	182.90 Å 102.16 Å 96.41 Å 90.00° 114.53° 90.00°	Depositor
Resolution (Å)	48.79 – 2.80 48.74 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.79-2.80) 99.6 (48.74-2.80)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 2.81 Å)	Xtriage
Refinement program	REFMAC 5.7.0027	Depositor
R, R_{free}	0.199 , 0.229 0.198 , 0.227	Depositor DCC
R_{free} test set	2006 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	84.3	Xtriage
Anisotropy	0.280	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 61.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 39735 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9611	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.32	0/1475	0.47	0/1995
1	B	0.33	0/1333	0.46	0/1799
1	C	0.34	0/1343	0.49	0/1813
1	D	0.32	0/1315	0.45	0/1774
1	E	0.32	0/1463	0.47	0/1979
1	F	0.32	0/1379	0.47	0/1859
1	G	0.33	0/1380	0.47	0/1859
All	All	0.33	0/9688	0.47	0/13078

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1453	0	1472	9	0
1	B	1313	0	1343	6	0
1	C	1323	0	1340	11	0
1	D	1295	0	1320	5	0
1	E	1439	0	1450	12	0
1	F	1359	0	1399	8	0
1	G	1360	0	1392	9	0
2	A	17	0	0	0	0
2	B	10	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	11	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	9	0	0	0	0
2	G	12	0	0	0	0
All	All	9611	0	9716	50	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 3.

All (50) 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:A:124:VAL:O	1:A:128:VAL:HG12	1.86	0.76
1:E:240:ILE:HG13	1:E:241:LEU:H	1.59	0.67
1:A:129:ASP:HB3	1:G:165:LEU:HD13	1.77	0.65
1:G:124:VAL:O	1:G:128:VAL:HG12	1.97	0.64
1:B:165:LEU:HD13	1:E:129:ASP:HB3	1.80	0.63
1:G:170:PHE:HB2	1:G:226:PHE:HB2	1.80	0.62
1:E:170:PHE:HB2	1:E:226:PHE:HB2	1.81	0.62
1:B:170:PHE:HB2	1:B:226:PHE:HB2	1.82	0.61
1:C:165:LEU:HD13	1:F:129:ASP:HB3	1.84	0.59
1:C:124:VAL:O	1:C:128:VAL:HG12	2.03	0.58
1:E:166:LYS:HA	1:E:240:ILE:HD11	1.90	0.54
1:C:126:SER:O	1:C:130:VAL:HG12	2.09	0.53
1:D:68:ARG:HG2	1:D:91:LEU:HD22	1.90	0.53
1:B:241:LEU:HD13	1:E:133:TYR:CE1	2.44	0.53
1:E:149:ILE:HA	1:E:152:ILE:HD12	1.92	0.52
1:A:165:LEU:HD13	1:C:129:ASP:HB3	1.91	0.52
1:C:165:LEU:HD11	1:F:130:VAL:HG12	1.93	0.51
1:D:165:LEU:HD13	1:G:129:ASP:HB3	1.93	0.51
1:G:149:ILE:HA	1:G:152:ILE:HD12	1.93	0.51
1:D:149:ILE:HA	1:D:152:ILE:HD12	1.93	0.50
1:C:67:LYS:HB2	1:C:69:ILE:HD12	1.93	0.50
1:B:149:ILE:HA	1:B:152:ILE:HD12	1.92	0.49
1:E:124:VAL:O	1:E:128:VAL:HG13	2.14	0.48
1:E:143:LEU:HA	1:E:165:LEU:HD12	1.95	0.48
1:C:71:PHE:HE2	1:F:86:SER:HA	1.79	0.48
1:E:240:ILE:HG13	1:E:241:LEU:N	2.28	0.47
1:F:149:ILE:HA	1:F:152:ILE:HD12	1.98	0.46
1:G:71:PHE:CZ	1:G:73:SER:HB3	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ARG:HG2	1:A:91:LEU:HD22	1.98	0.45
1:C:71:PHE:CZ	1:C:73:SER:HB3	2.52	0.44
1:A:115:ILE:HD13	1:A:178:ILE:HG12	1.99	0.44
1:B:134:ILE:O	1:F:243:LYS:HE3	2.17	0.44
1:C:149:ILE:HA	1:C:152:ILE:HD12	2.00	0.43
1:F:198:LYS:O	1:F:201:VAL:HG12	2.19	0.43
1:F:125:ILE:O	1:F:128:VAL:HG12	2.19	0.43
1:G:191:ASN:HA	1:G:194:ILE:HD12	2.01	0.42
1:B:71:PHE:CZ	1:B:73:SER:HB3	2.55	0.42
1:D:100:ILE:HB	1:D:138:VAL:HG12	2.01	0.42
1:A:197:THR:O	1:A:201:VAL:HG12	2.20	0.42
1:F:127:ILE:O	1:F:130:VAL:HG22	2.19	0.42
1:C:103:TYR:CD2	1:C:141:TYR:HB2	2.56	0.41
1:A:170:PHE:HB2	1:A:226:PHE:HB2	2.02	0.41
1:E:146:ALA:O	1:E:151:CYS:HB2	2.20	0.41
1:E:209:THR:HA	1:E:235:LYS:HB3	2.02	0.41
1:E:130:VAL:HA	1:E:133:TYR:HB3	2.03	0.41
1:D:65:LEU:HD11	1:D:87:GLN:HG2	2.02	0.41
1:A:90:TYR:CE1	1:G:63:LEU:HG	2.56	0.41
1:G:68:ARG:HG2	1:G:91:LEU:HD22	2.03	0.40
1:C:121:ILE:O	1:C:124:VAL:HG12	2.22	0.40
1:A:61:PRO:HB2	1:A:62:SER:H	1.65	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	181/205 (88%)	173 (96%)	7 (4%)	1 (1%)	30	65
1	B	161/205 (78%)	155 (96%)	6 (4%)	0	100	100
1	C	163/205 (80%)	157 (96%)	6 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	159/205 (78%)	152 (96%)	7 (4%)	0	100	100
1	E	181/205 (88%)	170 (94%)	10 (6%)	1 (1%)	30	65
1	F	166/205 (81%)	161 (97%)	5 (3%)	0	100	100
1	G	167/205 (82%)	161 (96%)	4 (2%)	2 (1%)	16	47
All	All	1178/1435 (82%)	1129 (96%)	45 (4%)	4 (0%)	46	79

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	189	ILE
1	E	240	ILE
1	A	113	ASN
1	G	144	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	167/190 (88%)	163 (98%)	4 (2%)	57	87
1	B	151/190 (80%)	148 (98%)	3 (2%)	63	90
1	C	152/190 (80%)	144 (95%)	8 (5%)	28	61
1	D	149/190 (78%)	148 (99%)	1 (1%)	88	97
1	E	163/190 (86%)	161 (99%)	2 (1%)	78	95
1	F	158/190 (83%)	150 (95%)	8 (5%)	29	63
1	G	157/190 (83%)	152 (97%)	5 (3%)	46	80
All	All	1097/1330 (82%)	1066 (97%)	31 (3%)	51	84

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	LEU
1	A	130	VAL

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Mol	Chain	Res	Type
1	A	143	LEU
1	A	219	VAL
1	B	64	LEU
1	B	138	VAL
1	B	219	VAL
1	C	63	LEU
1	C	122	THR
1	C	129	ASP
1	C	130	VAL
1	C	133	TYR
1	C	138	VAL
1	C	143	LEU
1	C	212	ASP
1	D	64	LEU
1	E	64	LEU
1	E	219	VAL
1	F	64	LEU
1	F	86	SER
1	F	89	LEU
1	F	138	VAL
1	F	143	LEU
1	F	201	VAL
1	F	212	ASP
1	F	233	ASP
1	G	64	LEU
1	G	129	ASP
1	G	143	LEU
1	G	219	VAL
1	G	233	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	174	GLN
1	C	190	GLN
1	F	190	GLN

5.3.3 RNA ⓘ

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	183/205 (89%)	0.30	1 (0%) 91 88	58, 79, 116, 127	0
1	B	165/205 (80%)	0.27	4 (2%) 62 50	64, 99, 162, 190	0
1	C	167/205 (81%)	0.35	12 (7%) 18 10	61, 84, 147, 204	0
1	D	163/205 (79%)	0.77	31 (19%) 2 1	77, 115, 180, 198	0
1	E	183/205 (89%)	0.42	15 (8%) 14 7	70, 104, 142, 171	0
1	F	170/205 (82%)	0.75	24 (14%) 4 2	61, 97, 166, 223	0
1	G	171/205 (83%)	0.51	15 (8%) 12 6	62, 88, 139, 213	0
All	All	1202/1435 (83%)	0.48	102 (8%) 13 6	58, 95, 159, 223	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	195	MET	8.3
1	F	187	ILE	7.0
1	G	189	ILE	6.0
1	F	190	GLN	5.9
1	F	189	ILE	5.2
1	F	114	LYS	5.2
1	D	225	TYR	5.1
1	F	170	PHE	4.8
1	F	191	ASN	4.8
1	D	220	LEU	4.5
1	D	216	ILE	4.4
1	E	226	PHE	4.2
1	F	226	PHE	4.1
1	D	196	ASN	4.1
1	C	224	LYS	4.1
1	D	219	VAL	4.1
1	D	234	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
1	F	172	LEU	3.9
1	D	208	ASN	3.8
1	D	209	THR	3.8
1	C	195	MET	3.7
1	D	159	LYS	3.6
1	G	172	LEU	3.5
1	F	147	TYR	3.4
1	E	225	TYR	3.3
1	D	116	ILE	3.3
1	C	170	PHE	3.3
1	G	61	PRO	3.3
1	G	224	LYS	3.3
1	G	159	LYS	3.2
1	E	233	ASP	3.2
1	E	234	PHE	3.1
1	E	166	LYS	3.1
1	D	133	TYR	3.0
1	D	76	ILE	3.0
1	D	115	ILE	3.0
1	F	224	LYS	3.0
1	F	199	LYS	3.0
1	G	241	LEU	2.9
1	C	193	GLU	2.9
1	D	236	LEU	2.9
1	G	196	ASN	2.9
1	E	170	PHE	2.9
1	F	115	ILE	2.9
1	F	225	TYR	2.9
1	B	220	LEU	2.8
1	G	220	LEU	2.8
1	G	235	LYS	2.8
1	E	184	ALA	2.8
1	F	169	SER	2.7
1	E	224	LYS	2.7
1	G	190	GLN	2.7
1	G	133	TYR	2.7
1	F	216	ILE	2.7
1	C	225	TYR	2.7
1	B	114	LYS	2.6
1	C	226	PHE	2.6
1	B	116	ILE	2.6
1	F	112	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	222	ARG	2.6
1	D	235	LYS	2.6
1	G	234	PHE	2.6
1	C	191	ASN	2.6
1	D	109	ASP	2.5
1	B	113	ASN	2.5
1	F	234	PHE	2.5
1	F	228	ALA	2.5
1	F	237	ILE	2.4
1	E	235	LYS	2.4
1	D	201	VAL	2.4
1	E	232	VAL	2.4
1	F	198	LYS	2.4
1	D	207	LYS	2.3
1	F	61	PRO	2.3
1	D	211	LYS	2.3
1	C	133	TYR	2.3
1	C	218	ASN	2.3
1	F	188	GLU	2.3
1	D	110	ILE	2.3
1	E	181	PHE	2.3
1	E	240	ILE	2.2
1	E	228	ALA	2.2
1	F	212	ASP	2.2
1	D	202	ILE	2.2
1	D	200	LYS	2.2
1	G	216	ILE	2.2
1	D	210	GLU	2.2
1	A	241	LEU	2.2
1	E	229	ASP	2.2
1	D	215	VAL	2.2
1	D	206	SER	2.1
1	F	202	ILE	2.1
1	D	172	LEU	2.1
1	G	175	SER	2.1
1	D	197	THR	2.1
1	G	121	ILE	2.1
1	C	243	LYS	2.1
1	D	198	LYS	2.1
1	D	205	ILE	2.1
1	C	220	LEU	2.1
1	E	216	ILE	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	228	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.