



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 02:16 PM BST

PDB ID : 4AAU
EMDB ID: : EMD-2001
Title : ATP-triggered molecular mechanics of the chaperonin GroEL
Authors : Clare, D.K.; Vasishtan, D.; Stagg, S.; Quispe, J.; Farr, G.W.; Topf, M.; Horwich, A.L.; Saibil, H.R.
Deposited on : 2011-12-05
Resolution : 8.50 Å(reported)
Based on PDB ID : 1OEL

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27241

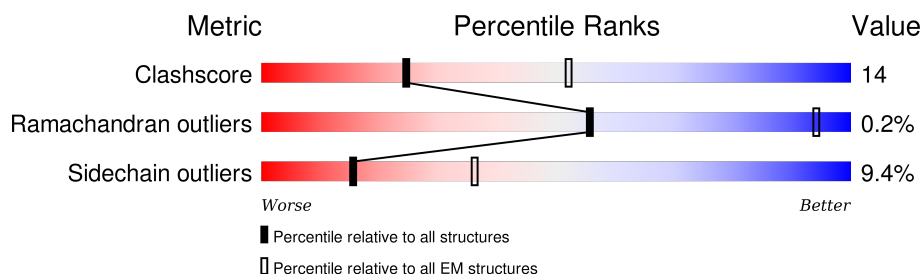
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 8.50 Å.

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







Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	548	60% 25% 9% . .
1	B	548	60% 25% 9% . .
1	C	548	61% 25% 9% . .
1	D	548	60% 25% 9% . .
1	E	548	61% 25% 9% . .
1	F	548	61% 25% 9% . .
1	G	548	61% 25% 9% . .
1	H	548	65% 25% 5% .
1	I	548	65% 25% 5% .

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Mol	Chain	Length	Quality of chain
1	J	548	
1	K	548	
1	L	548	
1	M	548	
1	N	548	

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	PO4	A	1527	-	-	X	-
3	PO4	B	1527	-	-	X	-
3	PO4	C	1527	-	-	X	-
3	PO4	D	1527	-	-	X	-
3	PO4	E	1527	-	-	X	-
3	PO4	F	1527	-	-	X	-
3	PO4	G	1527	-	-	X	-
3	PO4	H	1526	-	-	X	-
3	PO4	I	1526	-	-	X	-
3	PO4	J	1526	-	-	X	-
3	PO4	K	1526	-	-	X	-
3	PO4	L	1526	-	-	X	-
3	PO4	M	1526	-	-	X	-
3	PO4	N	1526	-	-	X	-
4	ATP	M	1527	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 54474 atoms, of which 168 are hydrogens and 0 are deuteriums.

In the tables below, 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 60 KDA CHAPERONIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	B	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	C	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	D	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	E	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	F	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	G	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	H	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	I	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	J	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	K	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	L	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	M	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		
1	N	524	Total	C	N	O	S	0	1
			3846	2391	665	770	20		

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
B	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
D	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
E	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
F	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
G	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
H	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
I	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
J	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
K	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
L	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
M	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5
N	398	ALA	ASP	ENGINEERED MUTATION	UNP P0A6F5

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

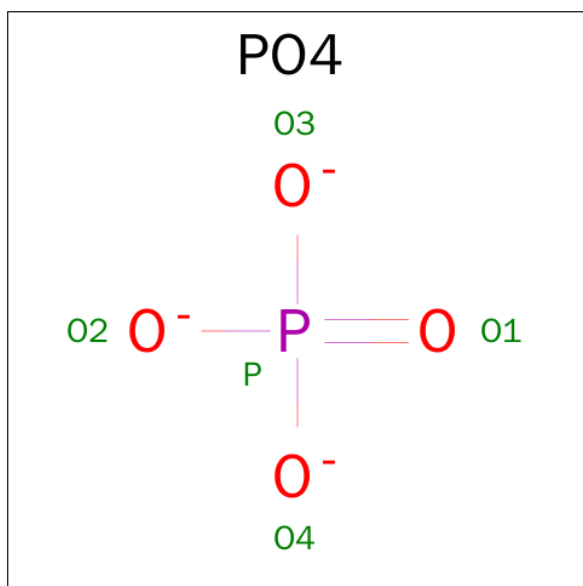
Mol	Chain	Residues	Atoms	AltConf
2	G	1	Total Mg 1 1	0
2	J	1	Total Mg 1 1	0
2	D	1	Total Mg 1 1	0
2	K	1	Total Mg 1 1	0
2	E	1	Total Mg 1 1	0
2	H	1	Total Mg 1 1	0
2	B	1	Total Mg 1 1	0
2	I	1	Total Mg 1 1	0
2	C	1	Total Mg 1 1	0
2	A	1	Total Mg 1 1	0
2	N	1	Total Mg 1 1	0
2	L	1	Total Mg 1 1	0
2	F	1	Total Mg 1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

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Mol	Chain	Residues	Atoms		AltConf
2	M	1	Total	Mg	0
			1	1	

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



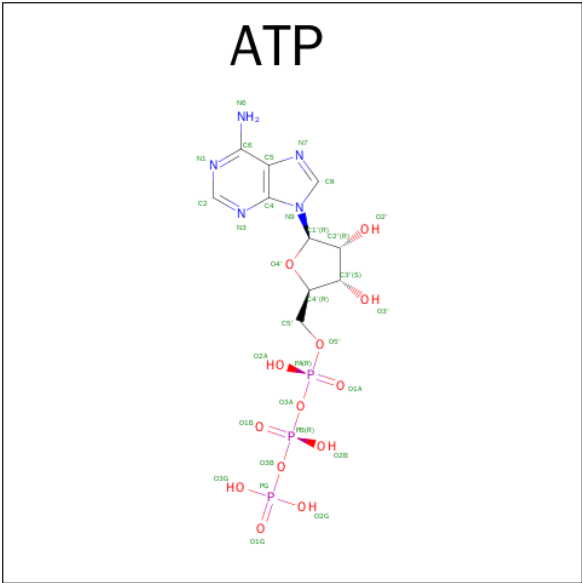
Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	P	0
			1	1	
3	B	1	Total	P	0
			1	1	
3	C	1	Total	P	0
			1	1	
3	D	1	Total	P	0
			1	1	
3	E	1	Total	P	0
			1	1	
3	F	1	Total	P	0
			1	1	
3	G	1	Total	P	0
			1	1	
3	H	1	Total	P	0
			1	1	
3	I	1	Total	P	0
			1	1	
3	J	1	Total	P	0
			1	1	

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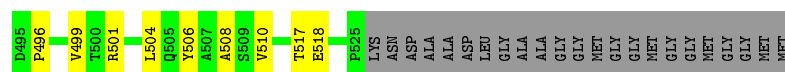
Mol	Chain	Residues	Atoms		AltConf
3	K	1	Total	P	0
			1	1	
3	L	1	Total	P	0
			1	1	
3	M	1	Total	P	0
			1	1	
3	N	1	Total	P	0
			1	1	

- Molecule 4 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: C₁₀H₁₆N₅O₁₃P₃).



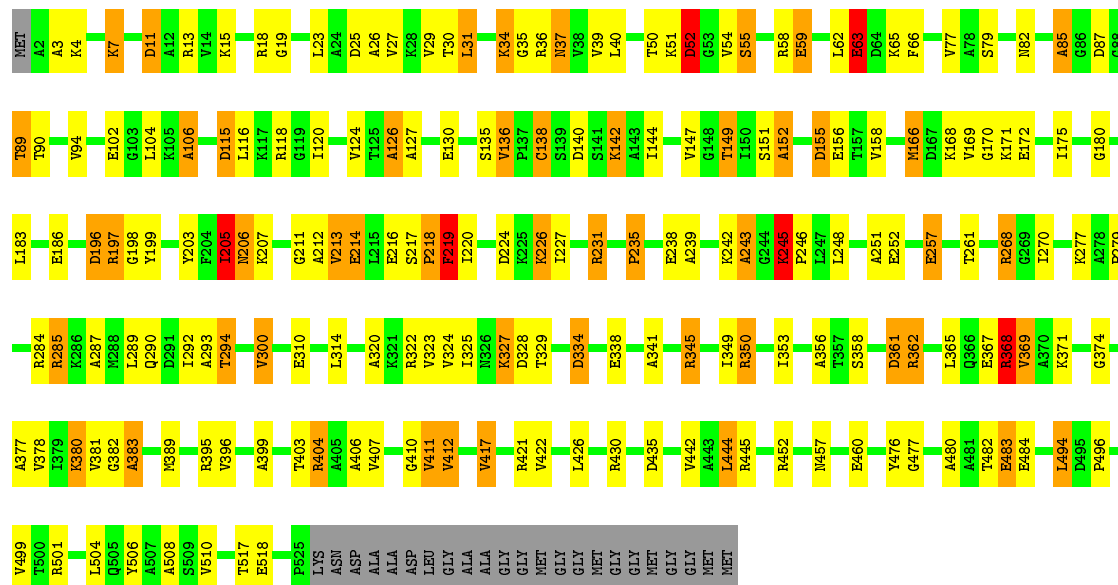
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Mol	Chain	Residues	Atoms						AltConf
4	H	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	I	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	J	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	K	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	L	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	M	1	Total 43	C 10	H 12	N 5	O 13	P 3	0
4	N	1	Total 43	C 10	H 12	N 5	O 13	P 3	0



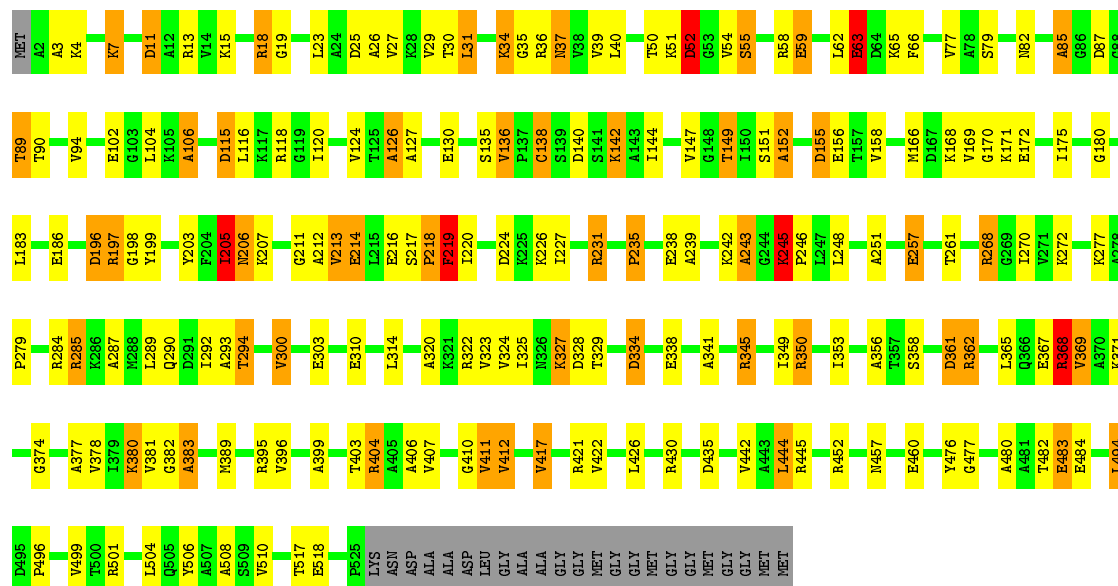
• Molecule 1: 60 KDA CHAPERONIN

Chain C: 61% 25% 9%



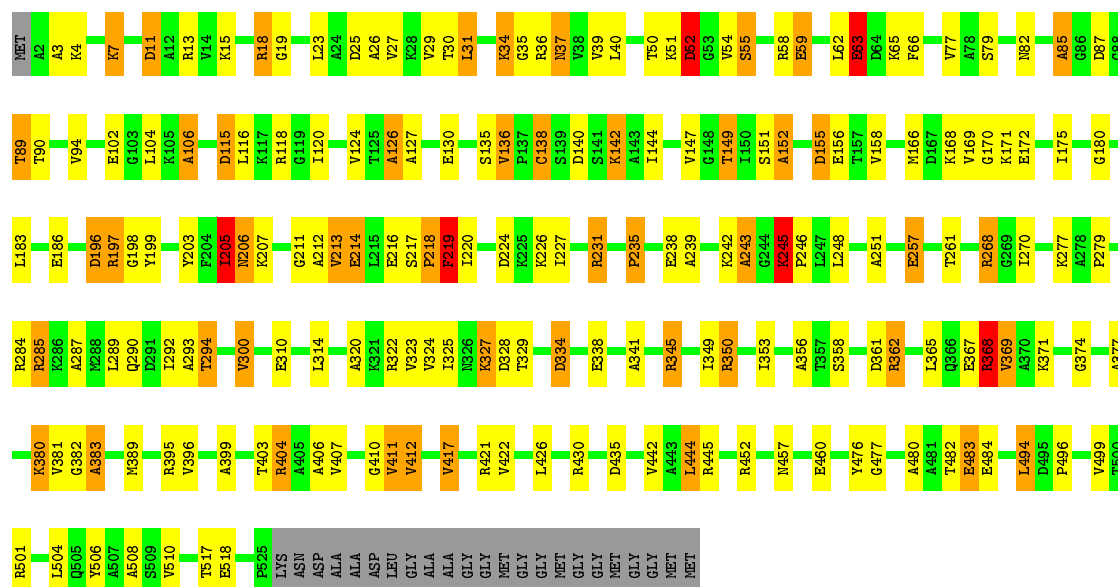
• Molecule 1: 60 KDA CHAPERONIN

Chain D: 60% 25% 9%

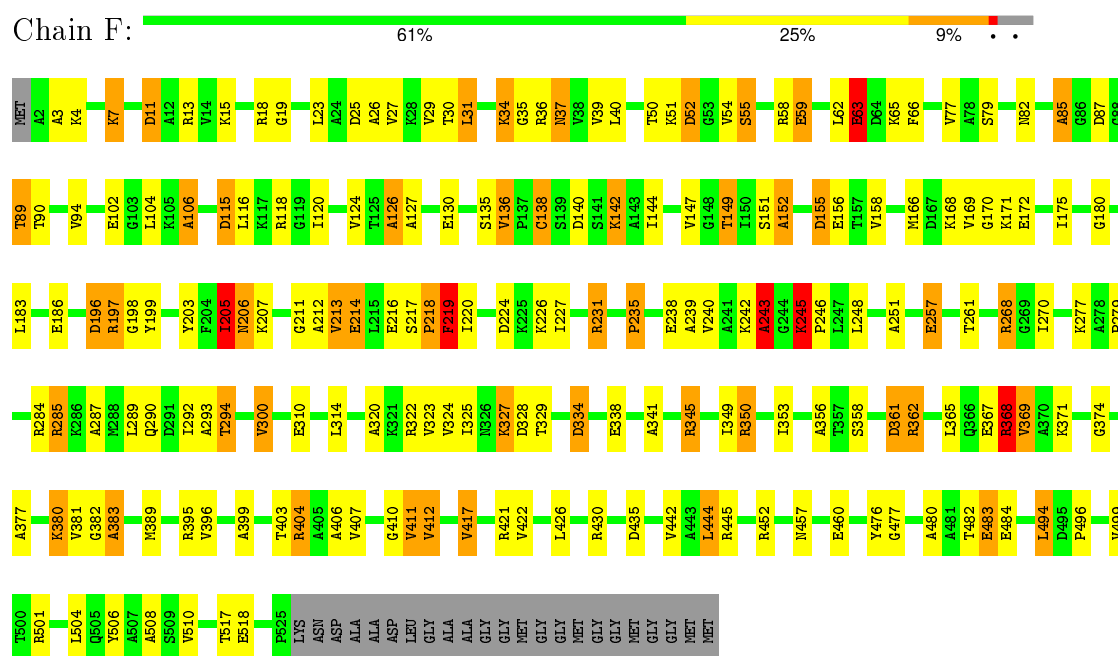


• Molecule 1: 60 KDA CHAPERONIN

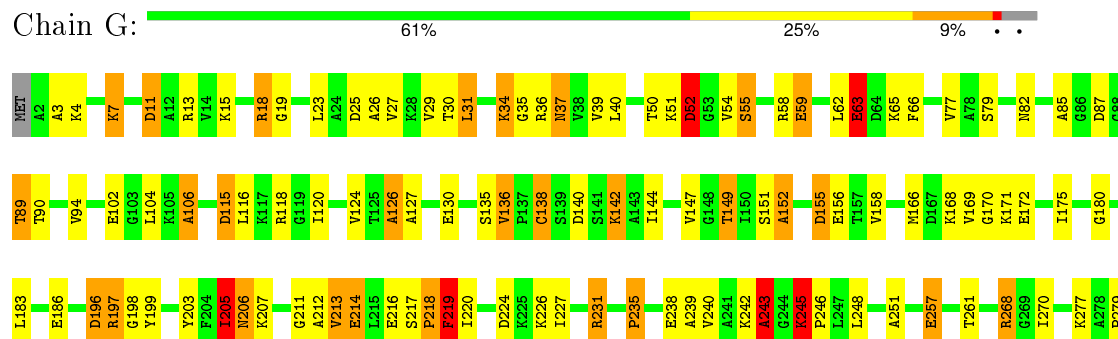
Chain E: 61% 25% 9%

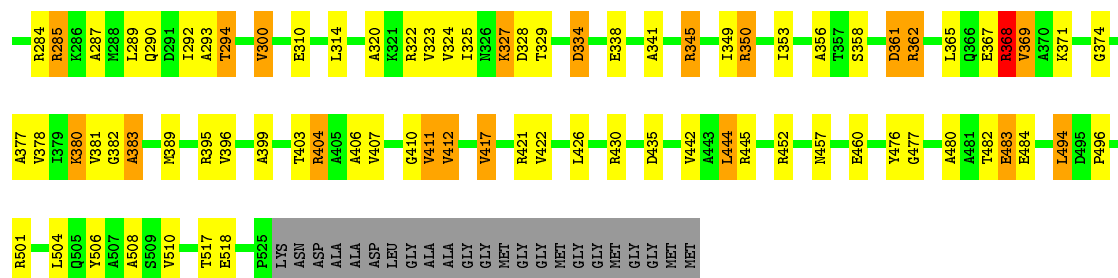


• Molecule 1: 60 KDA CHAPERONIN



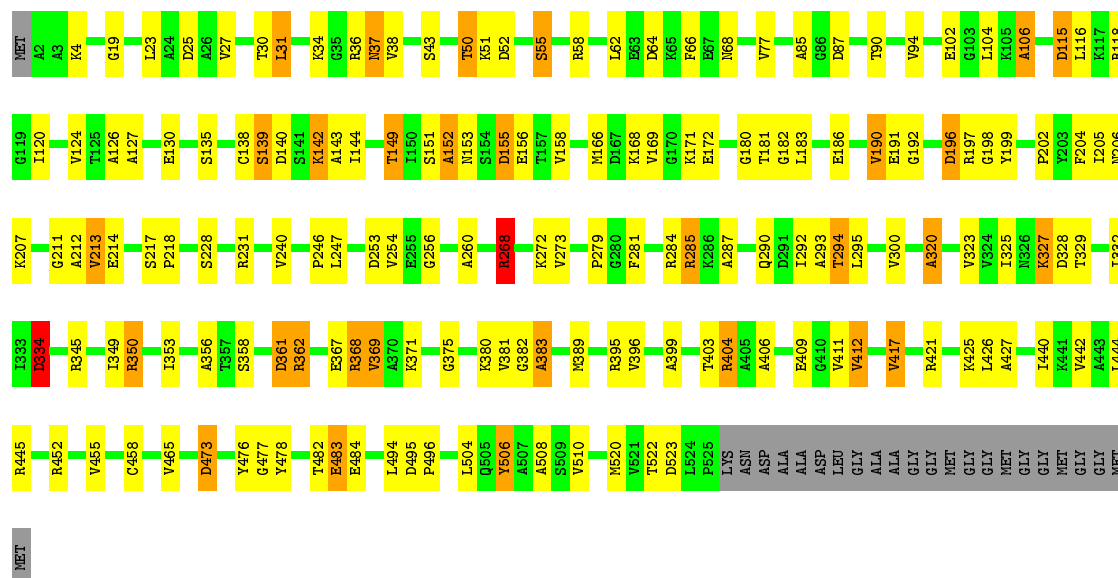
• Molecule 1: 60 KDA CHAPERONIN





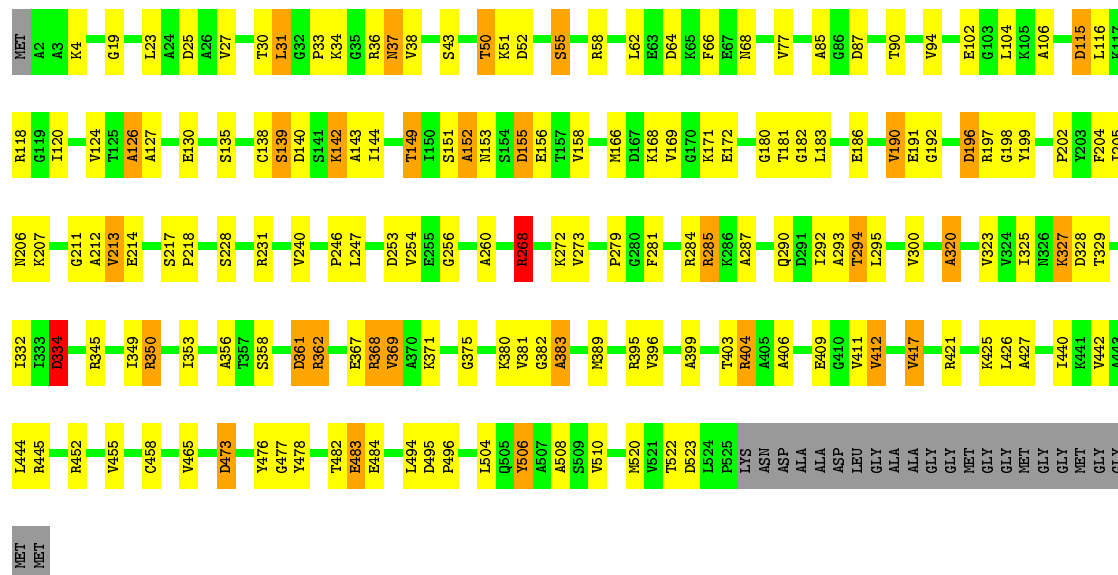
• Molecule 1: 60 KDA CHAPERONIN

Chain H: 65% 25% 5%

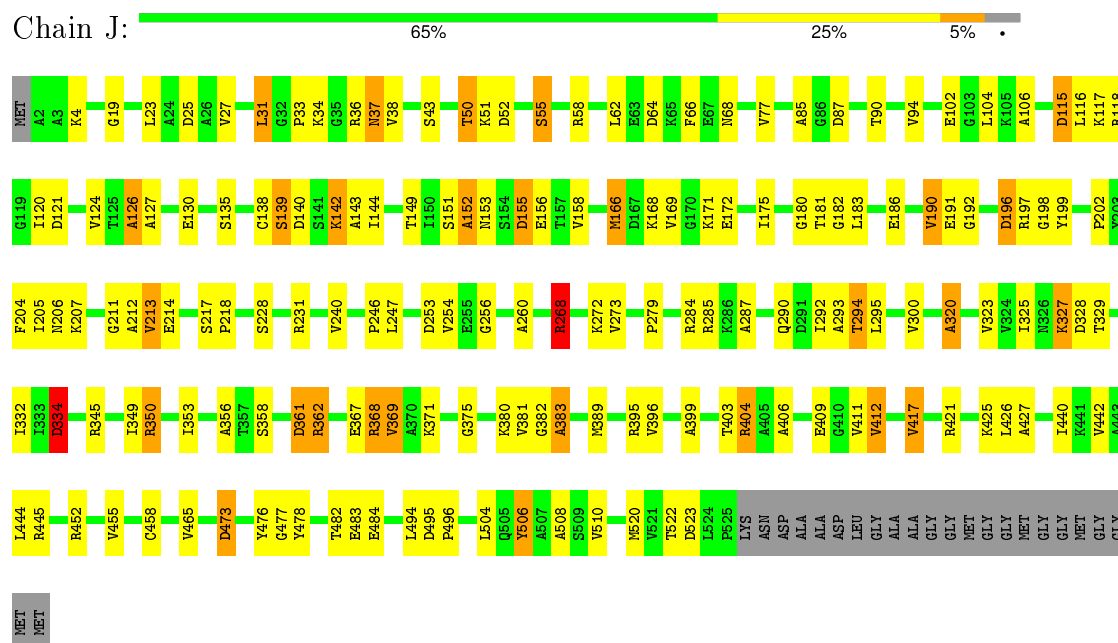


• Molecule 1: 60 KDA CHAPERONIN

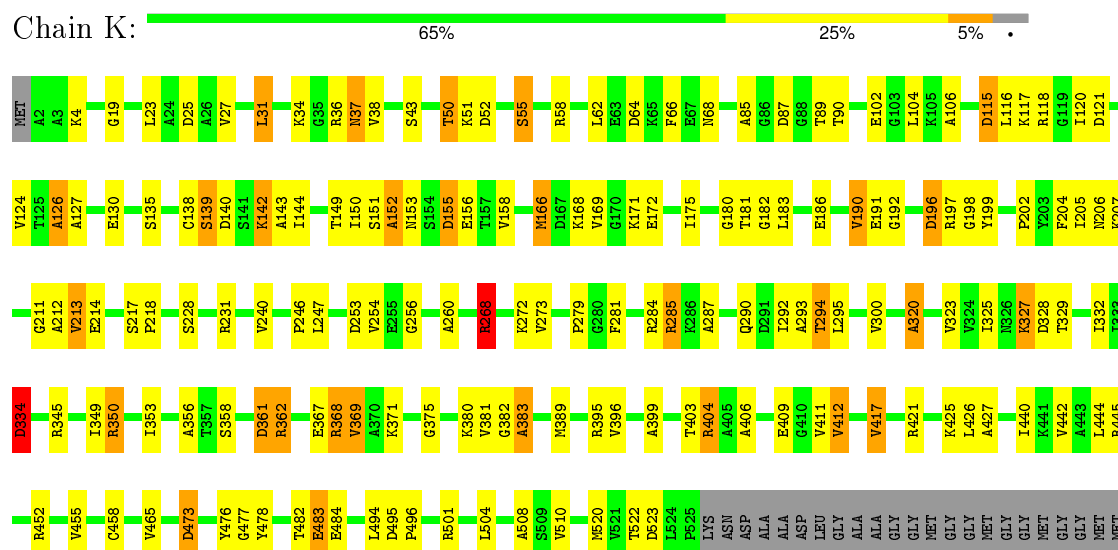
Chain I: 65% 25% 5%



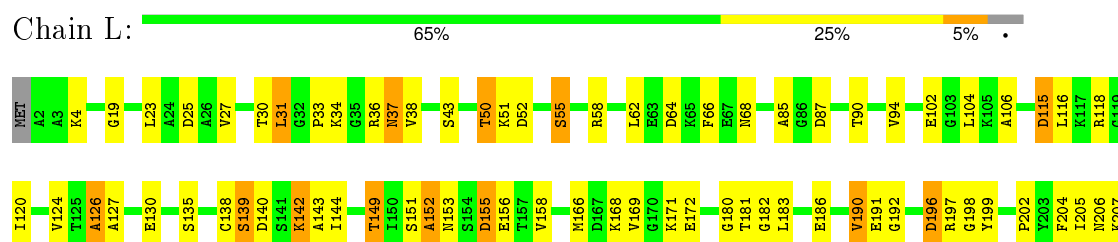
- Molecule 1: 60 KDA CHAPERONIN

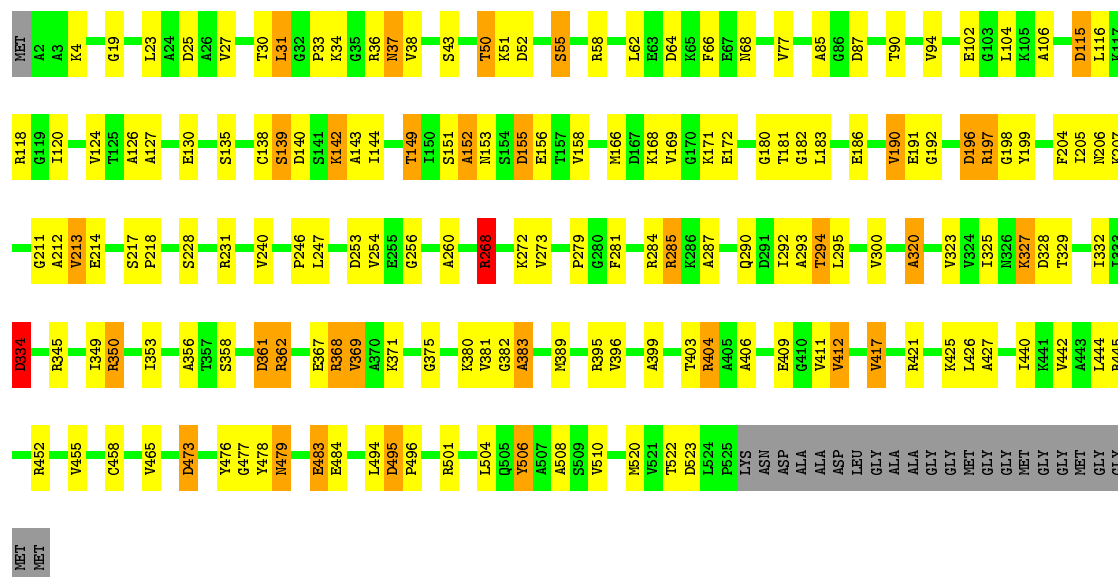


- Molecule 1: 60 KDA CHAPERONIN



- Molecule 1: 60 KDA CHAPERONIN





4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	EACH PARTICLE WAS PHASE FLIPPED, Not provided	Depositor
Microscope	FEI TECNAI F20	Depositor
Voltage (kV)	120	Depositor
Electron dose ($e^-/\text{\AA}^2$)	15	Depositor
Minimum defocus (nm)	700	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	Depositor
Image detector	GATAN ULTRASCAN 4000 4K CCD CAM- ERA	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, MG, ATP

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	1.01	0/3873	1.46	68/5229 (1.3%)
1	B	1.01	0/3873	1.46	67/5229 (1.3%)
1	C	1.01	0/3873	1.46	66/5229 (1.3%)
1	D	1.01	0/3873	1.46	67/5229 (1.3%)
1	E	1.01	0/3873	1.45	67/5229 (1.3%)
1	F	1.01	0/3873	1.46	67/5229 (1.3%)
1	G	1.01	0/3873	1.46	68/5229 (1.3%)
1	H	1.00	0/3871	1.40	57/5223 (1.1%)
1	I	1.00	0/3871	1.39	55/5223 (1.1%)
1	J	1.00	0/3871	1.39	53/5223 (1.0%)
1	K	1.00	0/3871	1.40	55/5223 (1.1%)
1	L	1.00	0/3871	1.39	56/5223 (1.1%)
1	M	1.08	1/3871 (0.0%)	1.40	58/5223 (1.1%)
1	N	1.00	0/3871	1.40	57/5223 (1.1%)
All	All	1.01	1/54208 (0.0%)	1.43	861/73164 (1.2%)

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	10
1	B	0	10
1	C	0	10
1	D	0	10
1	E	0	10
1	F	0	10
1	G	0	10
1	H	0	10
1	I	0	10

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Mol	Chain	#Chirality outliers	#Planarity outliers
1	J	0	10
1	K	0	10
1	L	0	10
1	M	0	10
1	N	1	10
All	All	1	140

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	493	ILE	CG1-CD1	26.52	3.33	1.50

The worst 5 of 861 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	245	LYS	N-CA-CB	-14.08	85.26	110.60
1	D	245	LYS	N-CA-CB	-14.05	85.30	110.60
1	B	245	LYS	N-CA-CB	-14.05	85.31	110.60
1	A	245	LYS	N-CA-CB	-14.05	85.31	110.60
1	G	245	LYS	N-CA-CB	-14.04	85.33	110.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	N	479	ASN	CA

5 of 140 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	118	ARG	Sidechain
1	A	18	ARG	Sidechain
1	A	197	ARG	Sidechain
1	A	345	ARG	Sidechain
1	A	350	ARG	Sidechain

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	3846	0	3970	127	0
1	B	3846	0	3970	126	0
1	C	3846	0	3970	125	0
1	D	3846	0	3970	122	0
1	E	3846	0	3970	122	0
1	F	3846	0	3970	123	0
1	G	3846	0	3970	123	0
1	H	3846	0	3968	86	0
1	I	3846	0	3968	89	0
1	J	3846	0	3968	88	0
1	K	3846	0	3968	89	0
1	L	3846	0	3968	87	0
1	M	3846	0	3968	108	0
1	N	3846	0	3968	87	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0
2	N	1	0	0	0	0
3	A	1	0	0	5	0
3	B	1	0	0	5	0
3	C	1	0	0	5	0
3	D	1	0	0	5	0
3	E	1	0	0	4	0
3	F	1	0	0	5	0
3	G	1	0	0	4	0
3	H	1	0	0	5	0
3	I	1	0	0	5	0
3	J	1	0	0	5	0
3	K	1	0	0	5	0
3	L	1	0	0	5	0
3	M	1	0	0	5	0
3	N	1	0	0	4	0
4	A	31	12	12	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	31	12	12	4	0
4	C	31	12	12	4	0
4	D	31	12	12	4	0
4	E	31	12	12	3	0
4	F	31	12	12	4	0
4	G	31	12	12	3	0
4	H	31	12	12	4	0
4	I	31	12	12	5	0
4	J	31	12	12	5	0
4	K	31	12	12	5	0
4	L	31	12	12	5	0
4	M	31	12	12	34	0
4	N	31	12	12	5	0
All	All	54306	168	55734	1536	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 14.

The worst 5 of 1536 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)
4:M:1527:ATP:C4	4:M:1527:ATP:C5	1.93	1.55
4:M:1527:ATP:C8	4:M:1527:ATP:N7	1.72	1.54
4:M:1527:ATP:N7	4:M:1527:ATP:C5	1.93	1.37
4:M:1527:ATP:C8	4:M:1527:ATP:N9	1.94	1.36
4:M:1527:ATP:C4	4:M:1527:ATP:N9	2.02	1.26

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 PDB entries followed by that with respect to all EM entries.

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	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	B	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	C	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	D	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	E	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	F	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	G	522/548 (95%)	518 (99%)	2 (0%)	2 (0%)	39	80
1	H	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	I	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	J	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	K	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	L	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	M	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
1	N	518/548 (94%)	513 (99%)	5 (1%)	0	100	100
All	All	7280/7672 (95%)	7217 (99%)	49 (1%)	14 (0%)	56	86

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	ALA
1	B	243	ALA
1	C	243	ALA
1	D	243	ALA
1	E	243	ALA

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 PDB entries followed by that with respect to all EM entries.

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	402/414 (97%)	358 (89%)	44 (11%)	8	35
1	B	402/414 (97%)	358 (89%)	44 (11%)	8	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	402/414 (97%)	359 (89%)	43 (11%)	8	36
1	D	402/414 (97%)	357 (89%)	45 (11%)	7	33
1	E	402/414 (97%)	359 (89%)	43 (11%)	8	36
1	F	402/414 (97%)	359 (89%)	43 (11%)	8	36
1	G	402/414 (97%)	359 (89%)	43 (11%)	8	36
1	H	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	I	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	J	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	K	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	L	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	M	402/414 (97%)	370 (92%)	32 (8%)	15	50
1	N	402/414 (97%)	369 (92%)	33 (8%)	14	49
All	All	5628/5796 (97%)	5098 (91%)	530 (9%)	15	42

5 of 530 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	219	PHE
1	G	358	SER
1	M	350	ARG
1	F	268	ARG
1	G	79	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	453	GLN
1	H	453	GLN
1	L	453	GLN
1	F	453	GLN
1	K	453	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 42 ligands modelled in this entry, 14 are modelled with single atom and 14 are monoatomic - leaving 14 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
4	ATP	A	1528	2	26,33,33	0.89	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	B	1528	2	26,33,33	0.90	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	C	1528	2	26,33,33	0.90	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	D	1528	2	26,33,33	0.88	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	E	1528	2	26,33,33	0.90	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	F	1528	2	26,33,33	0.90	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	G	1528	2	26,33,33	0.89	1 (3%)	26,52,52	1.27	4 (15%)
4	ATP	H	1527	2	26,33,33	0.98	1 (3%)	26,52,52	1.34	5 (19%)
4	ATP	I	1527	2	26,33,33	0.82	0	26,52,52	1.08	2 (7%)
4	ATP	J	1527	2	26,33,33	0.83	1 (3%)	26,52,52	1.11	2 (7%)
4	ATP	K	1527	2	26,33,33	0.89	0	26,52,52	1.42	6 (23%)
4	ATP	L	1527	2	26,33,33	0.82	1 (3%)	26,52,52	1.08	2 (7%)
4	ATP	M	1527	2	26,33,33	6.77	5 (19%)	26,52,52	2.99	6 (23%)
4	ATP	N	1527	2	26,33,33	1.00	1 (3%)	26,52,52	1.39	6 (23%)

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
4	ATP	A	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	B	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	C	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	D	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	E	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	F	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	G	1528	2	-	0/18/38/38	0/3/3/3
4	ATP	H	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	I	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	J	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	K	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	L	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	M	1527	2	-	0/18/38/38	0/3/3/3
4	ATP	N	1527	2	-	0/18/38/38	0/3/3/3

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	N	1527	ATP	C2'-C1'	-3.47	1.48	1.53
4	H	1527	ATP	C2'-C1'	-3.27	1.48	1.53
4	C	1528	ATP	C2'-C1'	-2.78	1.49	1.53
4	F	1528	ATP	C2'-C1'	-2.77	1.49	1.53
4	E	1528	ATP	C2'-C1'	-2.77	1.49	1.53

The worst 5 of 57 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	M	1527	ATP	O3G-PG-O2G	-3.79	93.53	107.44
4	K	1527	ATP	C4'-O4'-C1'	-3.54	105.89	109.64
4	N	1527	ATP	C1'-N9-C4	-2.94	123.53	126.81
4	H	1527	ATP	C4'-O4'-C1'	-2.53	106.96	109.64
4	N	1527	ATP	C4'-O4'-C1'	-2.41	107.08	109.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

14 monomers are involved in 89 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1528	ATP	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1528	ATP	4	0
4	C	1528	ATP	4	0
4	D	1528	ATP	4	0
4	E	1528	ATP	3	0
4	F	1528	ATP	4	0
4	G	1528	ATP	3	0
4	H	1527	ATP	4	0
4	I	1527	ATP	5	0
4	J	1527	ATP	5	0
4	K	1527	ATP	5	0
4	L	1527	ATP	5	0
4	M	1527	ATP	34	0
4	N	1527	ATP	5	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.