



wwPDB X-ray Structure Validation Summary Report

Mar 1, 2014 – 12:30 AM GMT

PDB ID : 1PCQ
Title : Crystal structure of groEL-groES
Authors : Chaudhry, C.; Farr, G.W.; Todd, M.J.; Rye, H.S.; Brunger, A.T.; Adams, P.D.; Horwich, A.L.; Sigler, P.B.
Deposited on : 2003-05-16
Resolution : 2.81 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

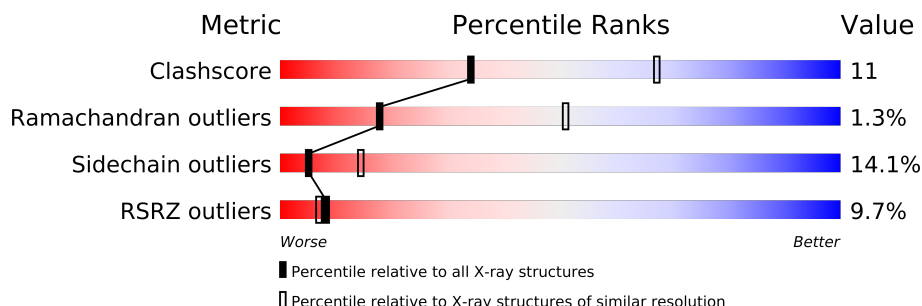
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.81 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))
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	524	
1	B	524	
1	C	524	
1	D	524	
1	E	524	
1	F	524	
1	G	524	
1	H	524	
1	I	524	
1	J	524	
1	K	524	
1	L	524	
1	M	524	
1	N	524	
2	O	97	
2	P	97	

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Mol	Chain	Length	Quality of chain
2	Q	97	
2	R	97	
2	S	97	
2	T	97	
2	U	97	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 59304 atoms, of which 0 are hydrogen and 0 are deuterium.

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 groEL protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	B	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	C	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	D	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	E	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	F	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	G	524	Total	C	N	O	S	0	0	0
			3855	2397	665	773	20			
1	H	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	I	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	J	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	K	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	L	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	M	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			
1	N	524	Total	C	N	O	S	0	0	0
			3856	2397	665	774	20			

- Molecule 2 is a protein called groES protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	O	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	P	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	Q	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	R	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	S	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	T	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			
2	U	97	Total	C	N	O	S	0	0	0
			728	454	127	145	2			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	A	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is POTASSIUM ION (three-letter code: K) (formula: K).

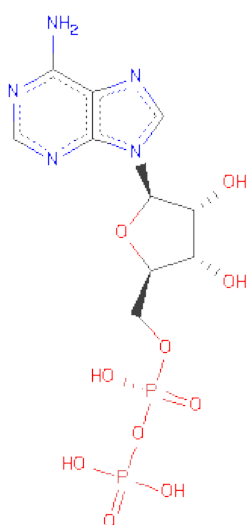
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total	K	0	0
			1	1		
4	D	1	Total	K	0	0
			1	1		
4	E	1	Total	K	0	0
			1	1		
4	B	1	Total	K	0	0
			1	1		

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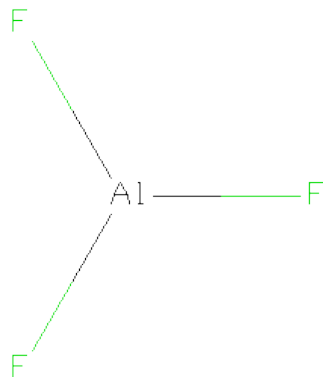
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	K	0	0
			1	1		
4	A	1	Total	K	0	0
			1	1		
4	F	1	Total	K	0	0
			1	1		

- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	C	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	D	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	E	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	F	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	G	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 6 is ALUMINUM FLUORIDE (three-letter code: AF3) (formula: AlF_3).



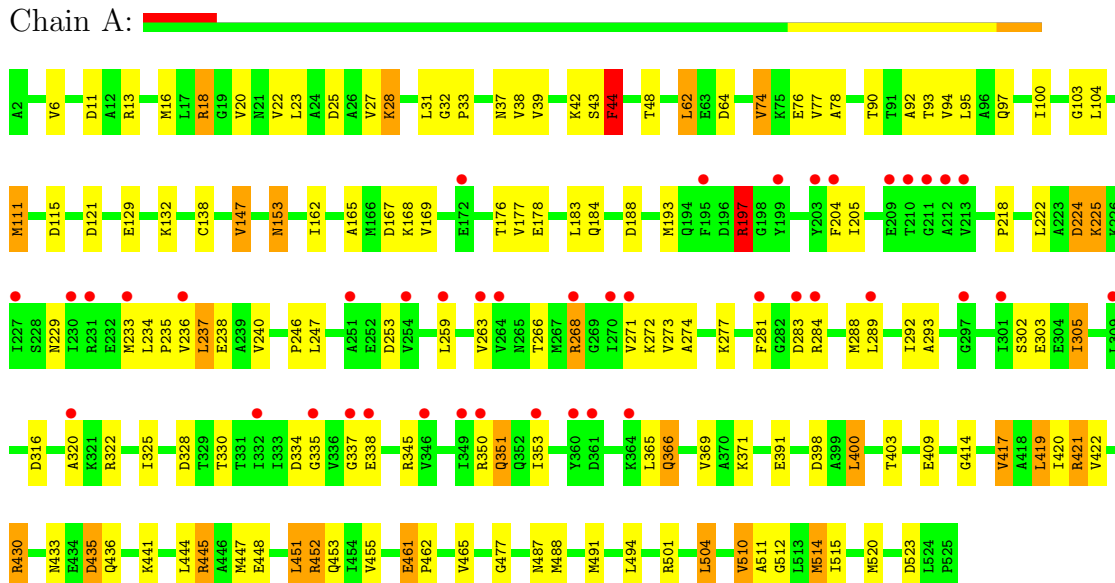
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	Al	F	0	0
			4	1	3		
6	B	1	Total	Al	F	0	0
			4	1	3		
6	C	1	Total	Al	F	0	0
			4	1	3		
6	D	1	Total	Al	F	0	0
			4	1	3		
6	E	1	Total	Al	F	0	0
			4	1	3		
6	F	1	Total	Al	F	0	0
			4	1	3		
6	G	1	Total	Al	F	0	0
			4	1	3		

3 Residue-property plots

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.

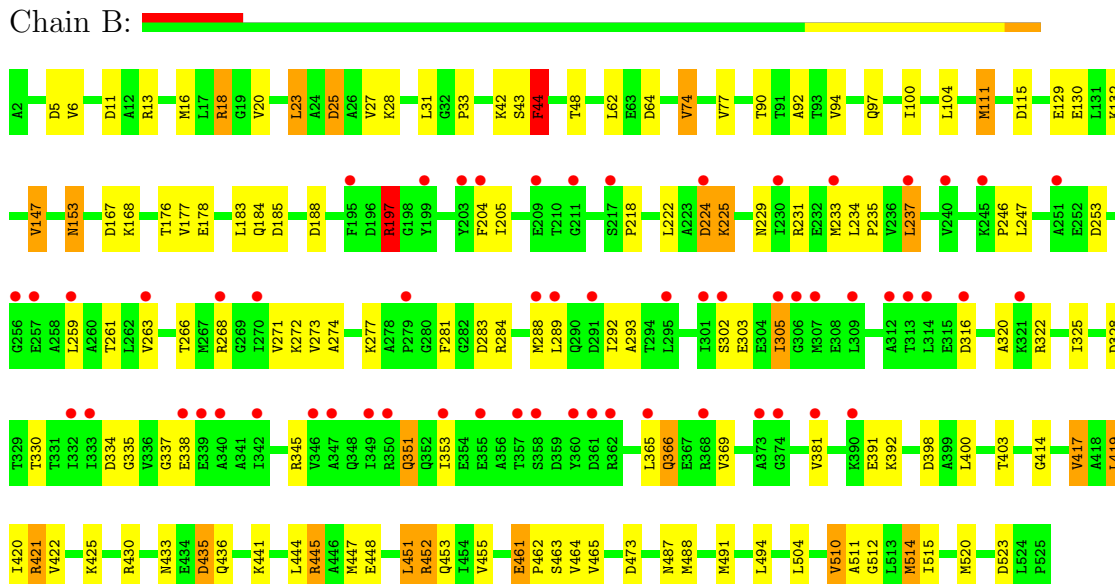
- Molecule 1: groEL protein

Chain A:

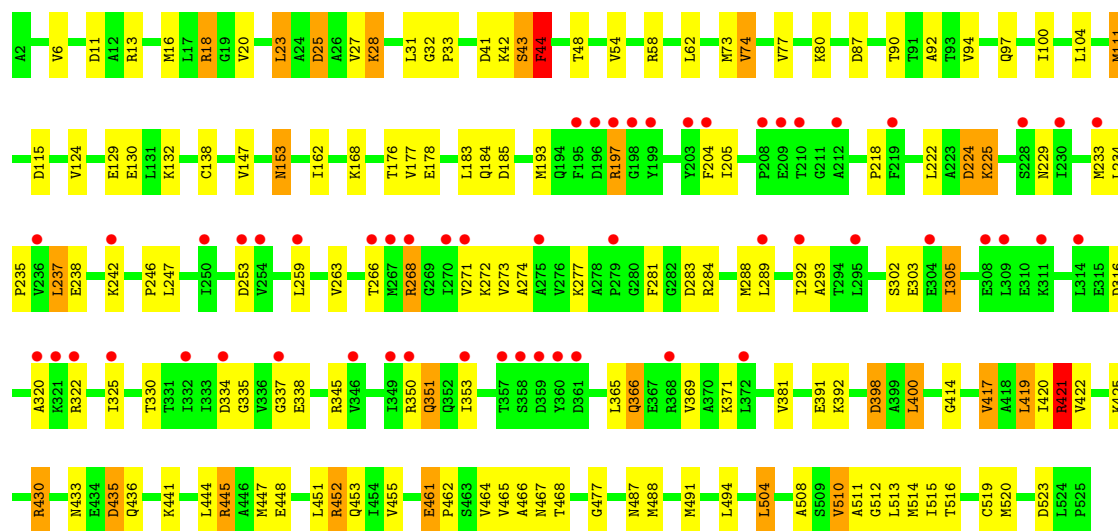


- Molecule 1: groEL protein

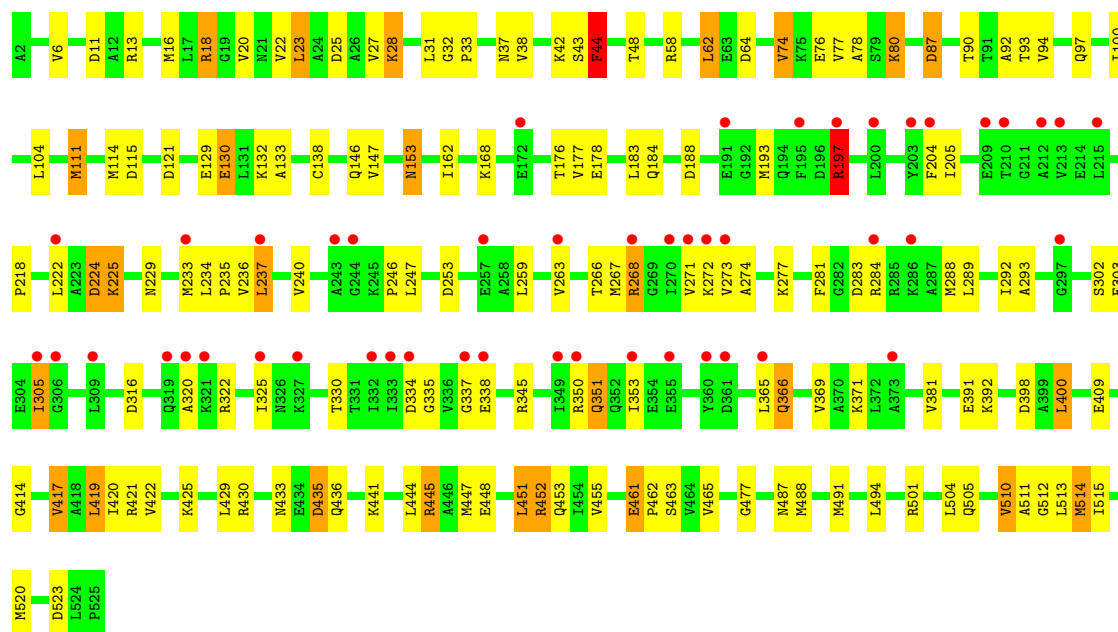
Chain B:



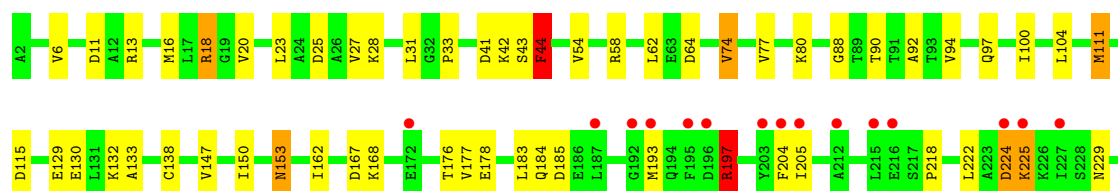
- Molecule 1: groEL protein

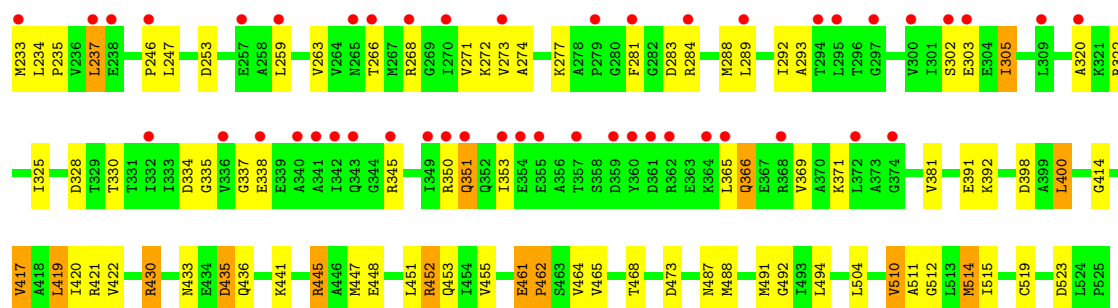
Chain C: 

• Molecule 1: groEL protein

Chain D: 

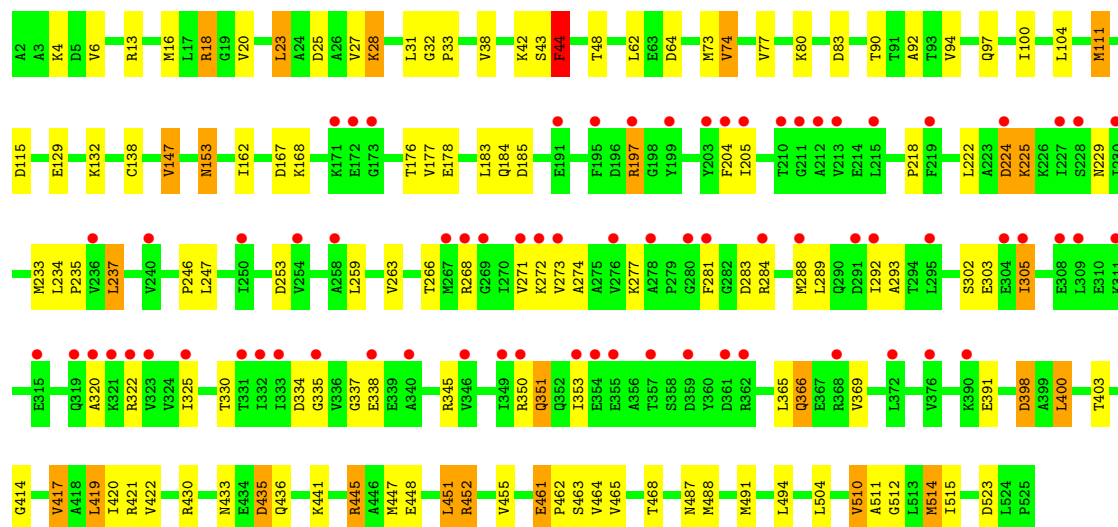
• Molecule 1: groEL protein

Chain E: 



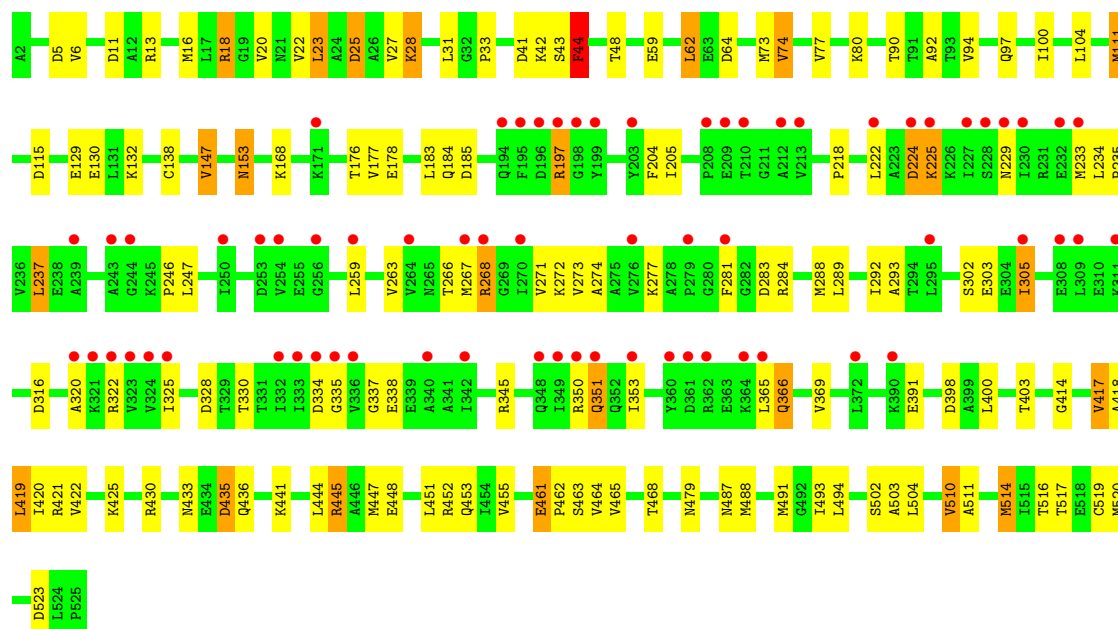
• Molecule 1: groEL protein

Chain F:



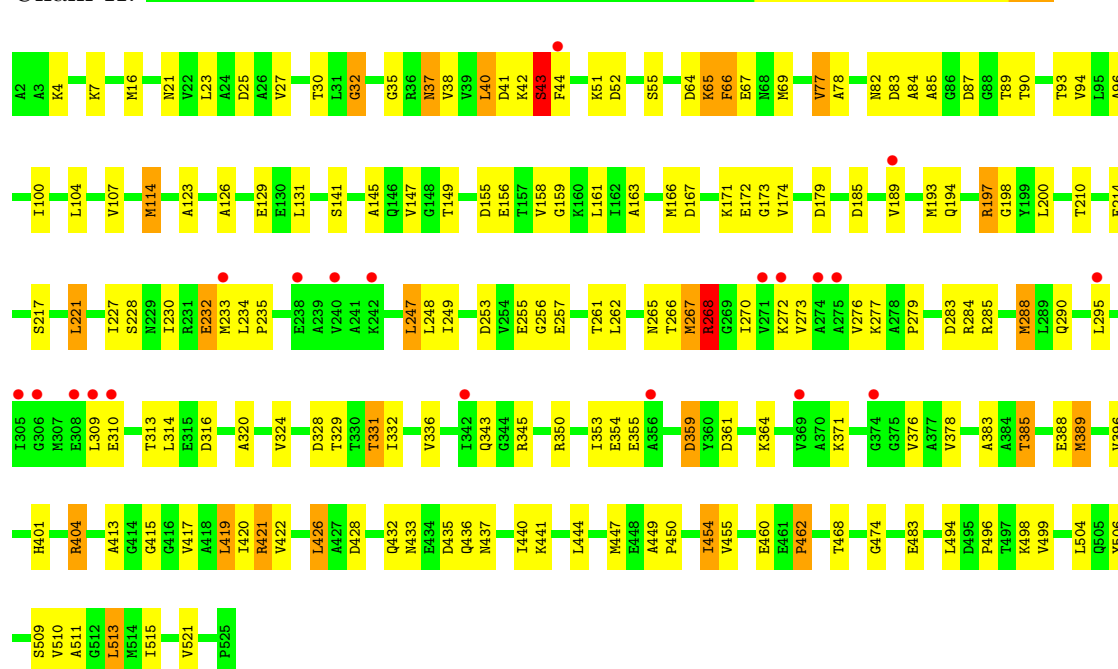
• Molecule 1: groEL protein

Chain G:



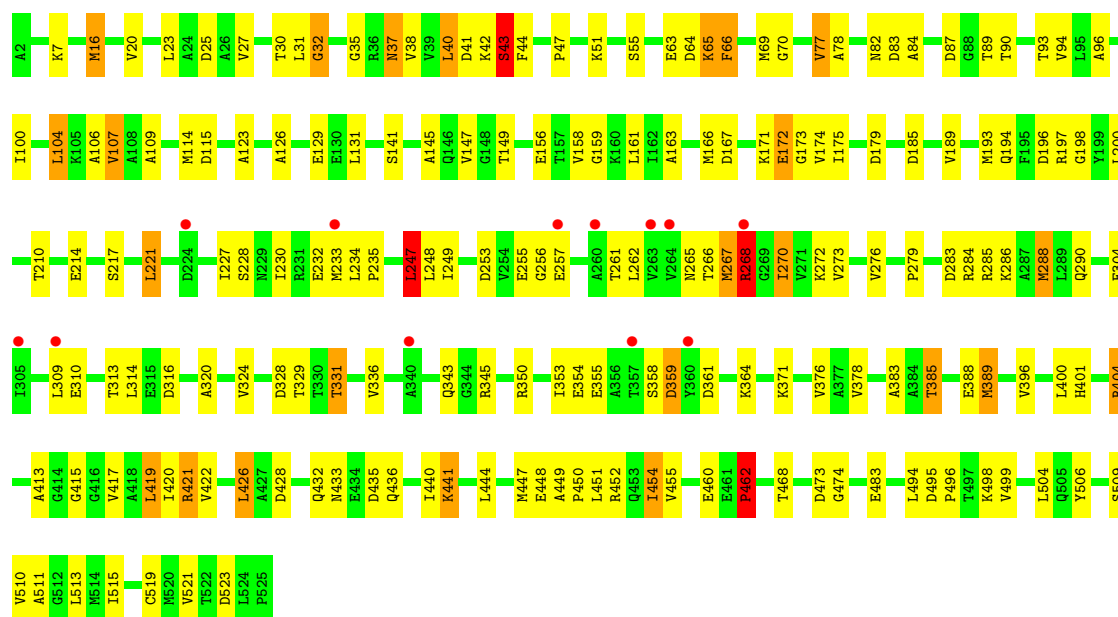
- Molecule 1: groEL protein

Chain H:



- Molecule 1: groEL protein

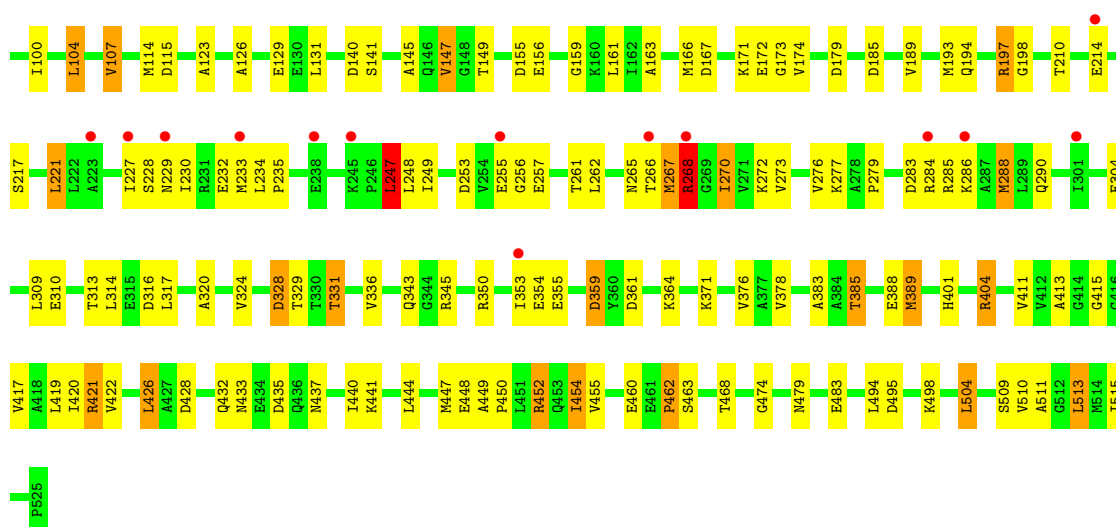
Chain I:



- Molecule 1: groEL protein

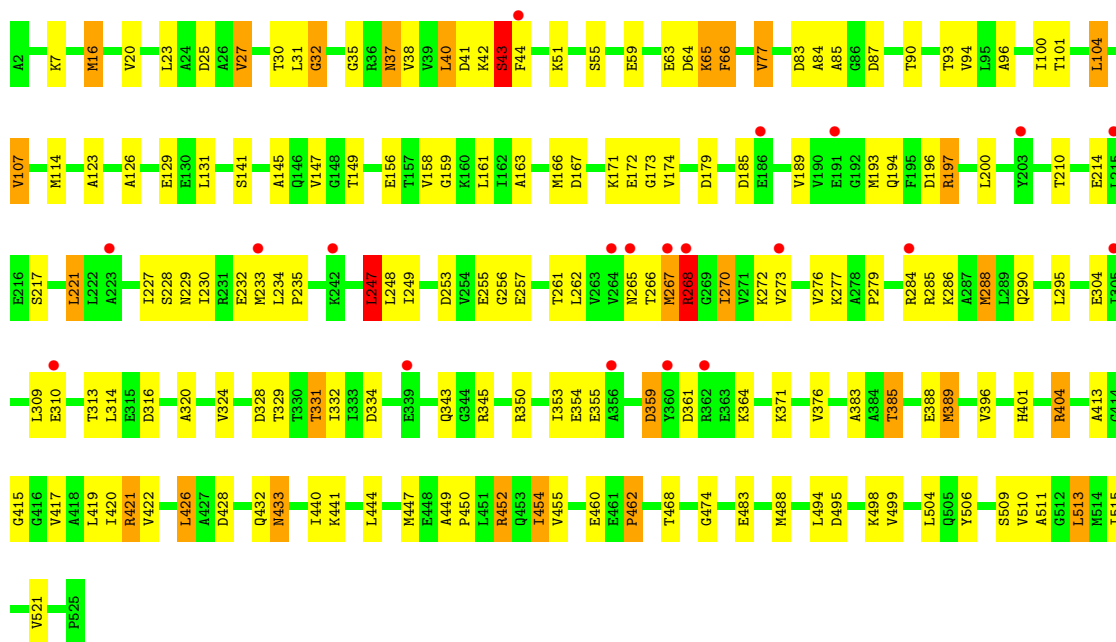
Chain J:





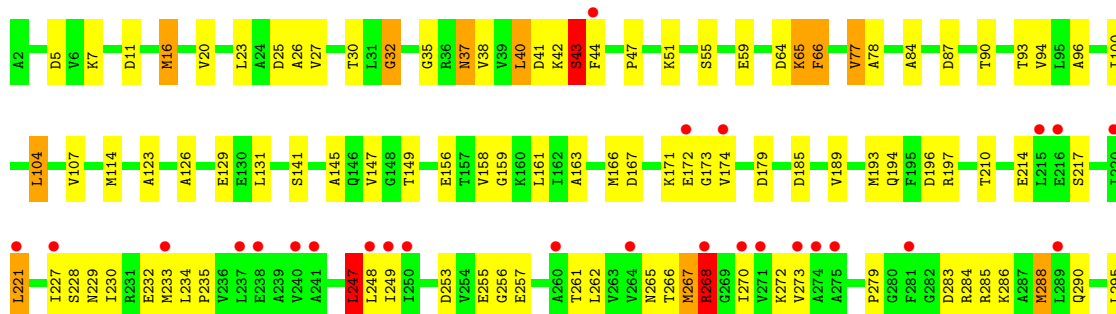
• Molecule 1: groEL protein

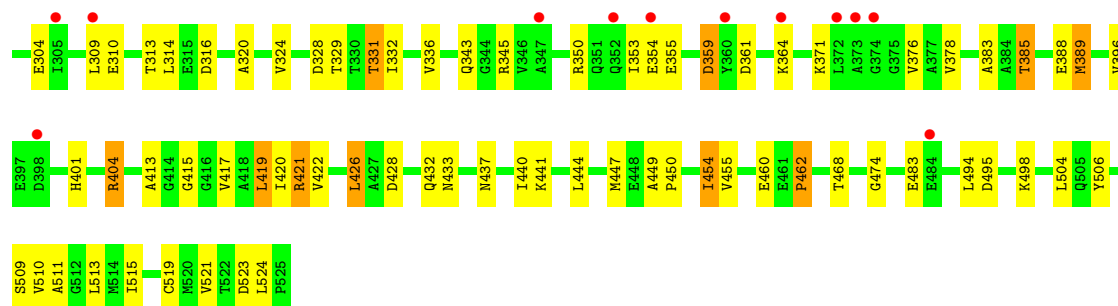
Chain K:



• Molecule 1: groEL protein

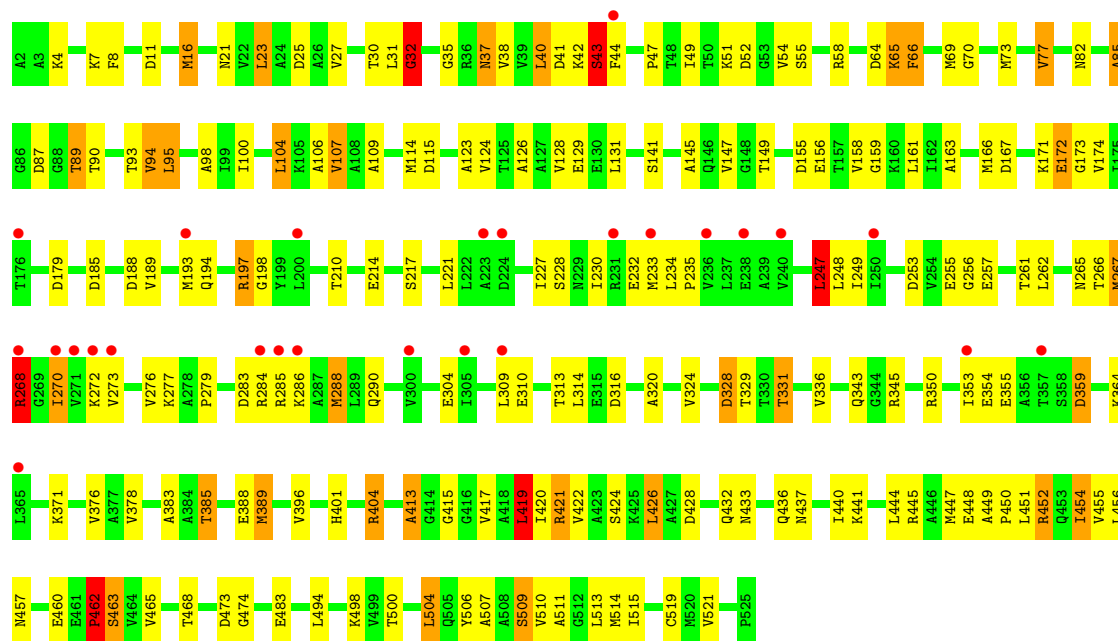
Chain L:





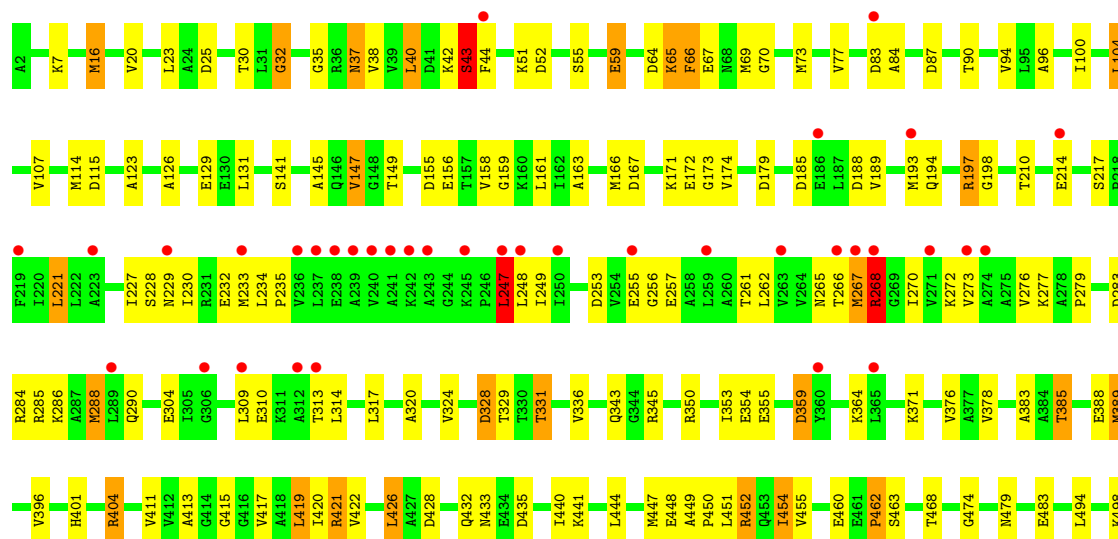
• Molecule 1: groEL protein

Chain M:



• Molecule 1: groEL protein

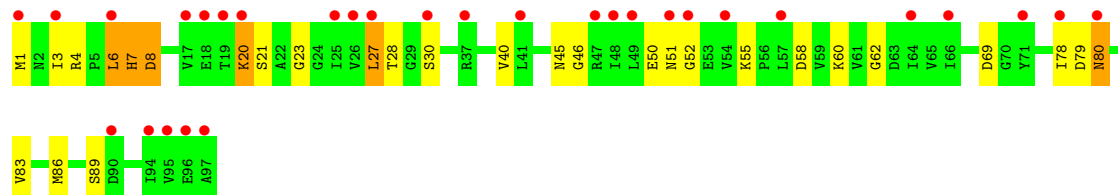
Chain N:





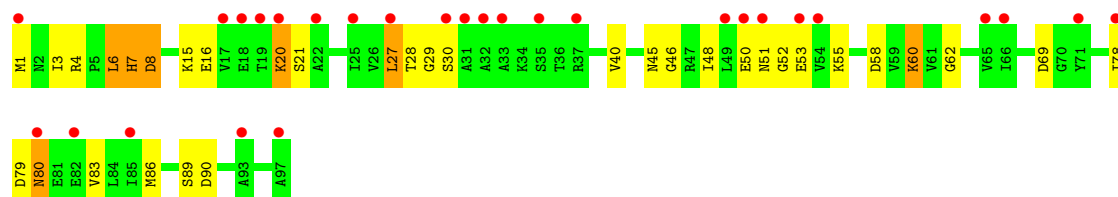
• Molecule 2: groES protein

Chain O:



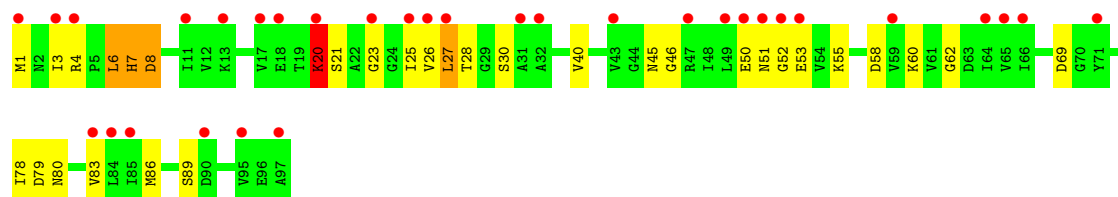
• Molecule 2: groES protein

Chain P:



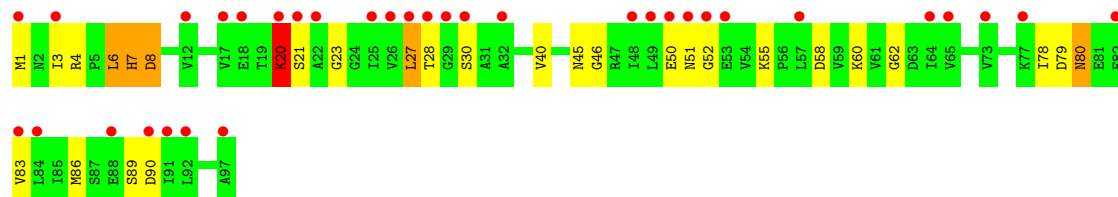
• Molecule 2: groES protein

Chain Q:



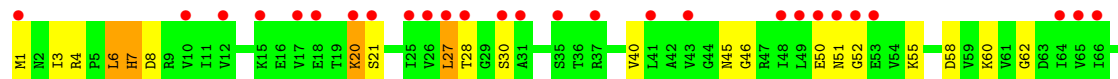
• Molecule 2: groES protein

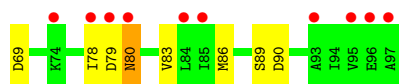
Chain R:



• Molecule 2: groES protein

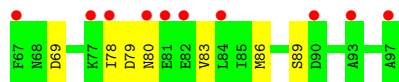
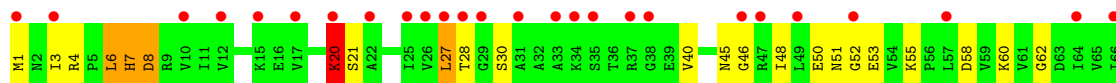
Chain S:





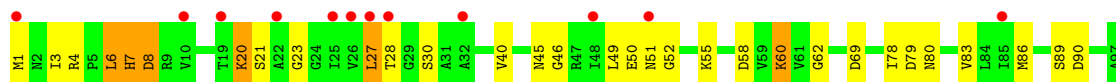
- Molecule 2: groES protein

Chain T:



- Molecule 2: groES protein

Chain U:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	255.55Å 266.86Å 187.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.39 – 2.81 49.44 – 2.31	Depositor EDS
% Data completeness (in resolution range)	(Not available) (49.39-2.81) 57.0 (49.44-2.31)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.32Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.262 , 0.278 (Not available) , (Not available)	Depositor DCC
R_{free} test set	NotAvailable	DCC
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.958	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 43.1	EDS
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	6 of 345474 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	59304	wwPDB-VP
Average B, all atoms (Å ²)	2.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, MG, ADP, AF3

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	0.70	1/3883 (0.0%)	0.93	19/5243 (0.4%)
1	B	0.61	0/3883	0.96	23/5243 (0.4%)
1	C	0.66	0/3883	0.91	19/5243 (0.4%)
1	D	0.68	1/3883 (0.0%)	0.92	19/5243 (0.4%)
1	E	0.54	0/3883	0.86	17/5243 (0.3%)
1	F	0.54	0/3883	0.87	17/5243 (0.3%)
1	G	0.62	0/3883	0.92	20/5243 (0.4%)
1	H	0.60	0/3884	0.83	15/5243 (0.3%)
1	I	0.64	0/3884	0.87	19/5243 (0.4%)
1	J	0.60	1/3884 (0.0%)	0.85	20/5243 (0.4%)
1	K	0.52	0/3884	0.86	16/5243 (0.3%)
1	L	0.49	0/3884	0.81	17/5243 (0.3%)
1	M	0.83	4/3884 (0.1%)	0.99	24/5243 (0.5%)
1	N	0.60	0/3884	0.85	16/5243 (0.3%)
2	O	0.35	0/732	0.69	4/983 (0.4%)
2	P	0.35	0/732	0.69	5/983 (0.5%)
2	Q	0.34	0/732	0.69	4/983 (0.4%)
2	R	0.34	0/732	0.69	4/983 (0.4%)
2	S	0.35	0/732	0.69	4/983 (0.4%)
2	T	0.35	0/732	0.69	4/983 (0.4%)
2	U	0.35	0/732	0.69	5/983 (0.5%)
All	All	0.60	7/59493 (0.0%)	0.87	291/80283 (0.4%)

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	M	0	1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	95	LEU	CG-CD1	7.61	1.80	1.51
1	D	130	GLU	CD-OE1	5.75	1.31	1.25
1	M	94	VAL	C-O	5.54	1.33	1.23
1	M	85	ALA	CA-CB	5.50	1.64	1.52
1	A	76	GLU	CD-OE1	5.26	1.31	1.25

The worst 5 of 291 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	231	ARG	NE-CZ-NH1	-19.61	110.50	120.30
1	B	231	ARG	NE-CZ-NH2	18.76	129.68	120.30
1	K	268	ARG	NE-CZ-NH2	-14.05	113.28	120.30
1	M	268	ARG	NE-CZ-NH2	-13.54	113.53	120.30
1	K	268	ARG	NE-CZ-NH1	13.45	127.03	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	M	32	GLY	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3855	0	3976	91	0
1	B	3855	0	3976	75	0
1	C	3855	0	3976	95	0
1	D	3855	0	3976	90	0
1	E	3855	0	3976	75	0
1	F	3855	0	3976	72	0
1	G	3855	0	3976	83	0
1	H	3856	0	3976	92	0
1	I	3856	0	3976	97	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	J	3856	0	3976	98	0
1	K	3856	0	3976	92	0
1	L	3856	0	3976	81	0
1	M	3856	0	3976	129	0
1	N	3856	0	3976	102	0
2	O	728	0	762	13	0
2	P	728	0	762	19	0
2	Q	728	0	762	16	0
2	R	728	0	762	12	0
2	S	728	0	762	11	0
2	T	728	0	762	13	0
2	U	728	0	762	14	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
5	A	27	0	12	4	0
5	B	27	0	12	1	0
5	C	27	0	12	5	0
5	D	27	0	12	2	0
5	E	27	0	12	1	0
5	F	27	0	12	3	0
5	G	27	0	12	1	0
6	A	4	0	0	1	0
6	B	4	0	0	0	0
6	C	4	0	0	1	0
6	D	4	0	0	0	0
6	E	4	0	0	0	0
6	F	4	0	0	0	0
6	G	4	0	0	1	0
All	All	59304	0	61082	1315	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 11.

The worst 5 of 1315 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:M:95:LEU:CG	1:M:95:LEU:CD1	1.80	1.60
1:F:73:MET:CE	1:F:73:MET:SD	2.01	1.48
1:H:114:MET:CE	1:H:114:MET:SD	2.02	1.48
1:G:73:MET:SD	1:G:73:MET:CE	2.01	1.48
1:N:16:MET:SD	1:N:16:MET:CE	2.01	1.47

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	522/524 (100%)	486 (93%)	32 (6%)	4 (1%)	27	65
1	B	522/524 (100%)	486 (93%)	32 (6%)	4 (1%)	27	65
1	C	522/524 (100%)	484 (93%)	34 (6%)	4 (1%)	27	65
1	D	522/524 (100%)	486 (93%)	31 (6%)	5 (1%)	22	60
1	E	522/524 (100%)	488 (94%)	30 (6%)	4 (1%)	27	65
1	F	522/524 (100%)	487 (93%)	31 (6%)	4 (1%)	27	65
1	G	522/524 (100%)	489 (94%)	29 (6%)	4 (1%)	27	65
1	H	522/524 (100%)	483 (92%)	33 (6%)	6 (1%)	21	57
1	I	522/524 (100%)	480 (92%)	36 (7%)	6 (1%)	21	57
1	J	522/524 (100%)	477 (91%)	39 (8%)	6 (1%)	21	57
1	K	522/524 (100%)	483 (92%)	33 (6%)	6 (1%)	21	57
1	L	522/524 (100%)	485 (93%)	31 (6%)	6 (1%)	21	57
1	M	522/524 (100%)	468 (90%)	42 (8%)	12 (2%)	10	31
1	N	522/524 (100%)	479 (92%)	37 (7%)	6 (1%)	21	57
2	O	95/97 (98%)	74 (78%)	18 (19%)	3 (3%)	6	20
2	P	95/97 (98%)	74 (78%)	17 (18%)	4 (4%)	4	13

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	Q	95/97 (98%)	74 (78%)	16 (17%)	5 (5%)	3	9
2	R	95/97 (98%)	74 (78%)	17 (18%)	4 (4%)	4	13
2	S	95/97 (98%)	74 (78%)	18 (19%)	3 (3%)	6	20
2	T	95/97 (98%)	74 (78%)	16 (17%)	5 (5%)	3	9
2	U	95/97 (98%)	73 (77%)	18 (19%)	4 (4%)	4	13
All	All	7973/8015 (100%)	7278 (91%)	590 (7%)	105 (1%)	18	51

5 of 105 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	PHE
1	A	337	GLY
1	B	44	PHE
1	B	337	GLY
1	C	44	PHE

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	404/404 (100%)	352 (87%)	52 (13%)	6	18
1	B	404/404 (100%)	354 (88%)	50 (12%)	7	20
1	C	404/404 (100%)	353 (87%)	51 (13%)	7	19
1	D	404/404 (100%)	351 (87%)	53 (13%)	6	17
1	E	404/404 (100%)	352 (87%)	52 (13%)	6	18
1	F	404/404 (100%)	352 (87%)	52 (13%)	6	18
1	G	404/404 (100%)	352 (87%)	52 (13%)	6	18
1	H	404/404 (100%)	344 (85%)	60 (15%)	4	12
1	I	404/404 (100%)	343 (85%)	61 (15%)	4	12
1	J	404/404 (100%)	344 (85%)	60 (15%)	4	12
1	K	404/404 (100%)	342 (85%)	62 (15%)	4	12
1	L	404/404 (100%)	345 (85%)	59 (15%)	5	13

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	M	404/404 (100%)	342 (85%)	62 (15%)	4	12
1	N	404/404 (100%)	345 (85%)	59 (15%)	5	13
2	O	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	P	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	Q	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	R	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	S	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	T	80/80 (100%)	67 (84%)	13 (16%)	3	10
2	U	80/80 (100%)	67 (84%)	13 (16%)	3	10
All	All	6216/6216 (100%)	5340 (86%)	876 (14%)	5	14

5 of 876 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	329	THR
1	J	55	SER
2	P	86	MET
1	H	404	ARG
1	I	230	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 126 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	153	ASN
1	I	37	ASN
2	R	45	ASN
1	G	348	GLN
1	H	37	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 28 ligands modelled in this entry, 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
5	ADP	A	600	3,4,6	29,29,29	1.01	2 (6%)	45,45,45	2.68	14 (31%)
6	AF3	A	602	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	B	700	3,4,6	29,29,29	1.01	2 (6%)	45,45,45	2.45	12 (26%)
6	AF3	B	702	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	C	800	3,4,6	29,29,29	1.16	5 (17%)	45,45,45	2.50	12 (26%)
6	AF3	C	802	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	D	900	3,4,6	29,29,29	1.07	4 (13%)	45,45,45	2.76	14 (31%)
6	AF3	D	902	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	E	1000	3,4,6	29,29,29	1.06	3 (10%)	45,45,45	2.63	14 (31%)
6	AF3	E	1002	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	F	1100	3,4,6	29,29,29	1.11	3 (10%)	45,45,45	2.38	10 (22%)
6	AF3	F	1102	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-
5	ADP	G	1200	3,4,6	29,29,29	1.18	3 (10%)	45,45,45	2.69	15 (33%)
6	AF3	G	1202	1,3,5,4	0,3,3	0.00	-	0,3,3	0.00	-

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
5	ADP	A	600	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	A	602	1,3,5,4	-	0/0/0/0	0/0/0/0
5	ADP	B	700	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	B	702	1,3,5,4	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ADP	C	800	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	C	802	1,3,5,4	-	0/0/0/0	0/0/0/0
5	ADP	D	900	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	D	902	1,3,5,4	-	0/0/0/0	0/0/0/0
5	ADP	E	1000	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	E	1002	1,3,5,4	-	0/0/0/0	0/0/0/0
5	ADP	F	1100	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	F	1102	1,3,5,4	-	0/0/0/0	0/0/0/0
5	ADP	G	1200	3,4,6	-	0/16/32/32	0/1/3/3
6	AF3	G	1202	1,3,5,4	-	0/0/0/0	0/0/0/0

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	1200	ADP	C2-N1	3.08	1.40	1.33
5	A	600	ADP	C2-N3	3.06	1.38	1.32
5	F	1100	ADP	C2-N3	2.94	1.38	1.32
5	E	1000	ADP	C2-N3	2.81	1.37	1.32
5	D	900	ADP	C2-N3	2.71	1.37	1.32

The worst 5 of 91 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	1000	ADP	N3-C2-N1	-11.39	119.18	128.71
5	D	900	ADP	N3-C2-N1	-11.11	119.42	128.71
5	F	1100	ADP	N3-C2-N1	-10.43	119.99	128.71
5	G	1200	ADP	O4'-C1'-N9	10.40	118.12	108.44
5	B	700	ADP	N3-C2-N1	-10.36	120.05	128.71

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	524/524 (100%)	0.42	42 (8%)	12 11	2, 2, 4, 6	0
1	B	524/524 (100%)	0.55	59 (11%)	6 5	2, 2, 4, 6	0
1	C	524/524 (100%)	0.49	54 (10%)	7 6	2, 2, 4, 6	0
1	D	524/524 (100%)	0.39	48 (9%)	9 7	2, 2, 4, 6	0
1	E	524/524 (100%)	0.52	62 (11%)	5 4	2, 2, 4, 6	0
1	F	524/524 (100%)	0.68	72 (13%)	4 3	2, 2, 4, 6	0
1	G	524/524 (100%)	0.61	67 (12%)	4 3	2, 2, 4, 6	0
1	H	524/524 (100%)	0.18	20 (3%)	38 38	2, 2, 4, 6	0
1	I	524/524 (100%)	0.15	12 (2%)	57 58	2, 2, 4, 6	0
1	J	524/524 (100%)	0.16	14 (2%)	52 52	2, 2, 4, 6	0
1	K	524/524 (100%)	0.28	20 (3%)	38 38	2, 2, 4, 6	0
1	L	524/524 (100%)	0.34	38 (7%)	15 13	2, 2, 4, 6	0
1	M	524/524 (100%)	0.26	26 (4%)	28 28	2, 2, 4, 6	0
1	N	524/524 (100%)	0.31	37 (7%)	16 14	2, 2, 4, 6	0
2	O	97/97 (100%)	1.59	30 (30%)	1 1	2, 2, 2, 3	0
2	P	97/97 (100%)	1.57	28 (28%)	1 1	2, 2, 2, 3	0
2	Q	97/97 (100%)	1.39	32 (32%)	1 1	2, 2, 2, 2	0
2	R	97/97 (100%)	1.79	34 (35%)	1 1	2, 2, 2, 3	0
2	S	97/97 (100%)	1.83	37 (38%)	1 0	2, 2, 2, 3	0
2	T	97/97 (100%)	1.70	36 (37%)	1 0	2, 2, 2, 3	0
2	U	97/97 (100%)	1.05	12 (12%)	5 4	2, 2, 2, 2	0
All	All	8015/8015 (100%)	0.48	780 (9%)	8 7	2, 2, 4, 6	0

The worst 5 of 780 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	353	ILE	11.5
1	E	361	ASP	10.7
1	A	204	PHE	10.2
1	C	212	ALA	9.9
1	G	350	ARG	9.8

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	AF3	F	1102	4/4	0.19	0.25	5,7,8,12	0
4	K	D	903	1/1	0.15	-1.58	24,24,24,24	0
5	ADP	F	1100	27/27	0.11	-1.60	2,2,5,8	0
4	K	F	1103	1/1	0.09	-1.67	24,24,24,24	0
4	K	E	1003	1/1	0.14	-1.67	24,24,24,24	0
5	ADP	G	1200	27/27	0.11	-2.43	2,2,6,8	0
5	ADP	E	1000	27/27	0.11	-2.48	2,2,5,8	0
6	AF3	E	1002	4/4	0.10	-2.65	5,7,8,12	0
5	ADP	B	700	27/27	0.10	-2.73	2,2,5,8	0
5	ADP	D	900	27/27	0.10	-2.81	2,2,5,8	0
6	AF3	D	902	4/4	0.12	-3.15	5,7,8,12	0
4	K	A	603	1/1	0.13	-3.15	24,24,24,24	0
5	ADP	A	600	27/27	0.09	-3.17	2,2,6,8	0
4	K	B	703	1/1	0.10	-3.19	24,24,24,24	0
3	MG	A	601	1/1	0.06	-3.20	2,2,2,2	0
6	AF3	C	802	4/4	0.09	-3.28	5,7,8,12	0
6	AF3	A	602	4/4	0.09	-3.42	5,7,8,12	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	AF3	G	1202	4/4	0.10	-3.50	5,7,8,12	0
3	MG	E	1001	1/1	0.06	-3.52	2,2,2,2	0
5	ADP	C	800	27/27	0.10	-3.66	2,2,5,8	0
4	K	C	803	1/1	0.11	-3.71	24,24,24,24	0
6	AF3	B	702	4/4	0.07	-3.91	5,7,8,12	0
4	K	G	1203	1/1	0.06	-4.58	24,24,24,24	0
3	MG	G	1201	1/1	0.07	-4.93	2,2,2,2	0
3	MG	B	701	1/1	0.06	-5.23	2,2,2,2	0
3	MG	C	801	1/1	0.06	-5.68	2,2,2,2	0
3	MG	D	901	1/1	0.08	-5.92	2,2,2,2	0
3	MG	F	1101	1/1	0.06	-8.55	2,2,2,2	0

6.5 Other polymers ⓘ

There are no such residues in this entry.