



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 02:02 PM BST

PDB ID : 3JA7
EMDB ID: : EMD-6324
Title : Cryo-EM structure of the bacteriophage T4 portal protein assembly at near-atomic resolution
Authors : Sun, L.; Zhang, X.; Gao, S.; Rao, P.A.; Padilla-Sanchez, V.; Chen, Z.; Sun, S.; Xiang, Y.; Subramaniam, S.; Rao, V.B.; Rossmann, M.G.
Deposited on : 2015-04-21
Resolution : 3.63 Å(reported)

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 : unknown
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

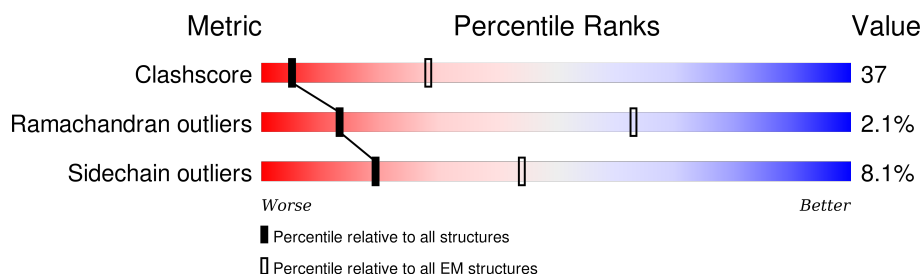
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.63 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	455	46% 42% 6% 6%
1	B	455	46% 42% 6% 6%
1	C	455	47% 41% 6% 6%
1	D	455	46% 42% 6% 6%
1	E	455	46% 42% 6% 6%
1	F	455	46% 42% 6% 6%
1	G	455	46% 42% 6% 6%
1	H	455	45% 43% 6% 6%
1	I	455	46% 42% 6% 6%

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Mol	Chain	Length	Quality of chain
1	J	455	<div><div></div><div>46%42%6%6%</div></div>
1	K	455	<div><div></div><div>47%41%6%6%</div></div>
1	L	455	<div><div></div><div>46%42%6%6%</div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 41280 atoms, of which 0 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 Portal protein gp20.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	B	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	C	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	D	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	E	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	F	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	G	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	H	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	I	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	J	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	K	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		
1	L	428	Total	C	N	O	S	0	0
			3440	2164	608	649	19		

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	ASN	-	EXPRESSION TAG	UNP P13334
A	71	TRP	-	EXPRESSION TAG	UNP P13334
A	72	ILE	-	EXPRESSION TAG	UNP P13334
A	73	SER	-	EXPRESSION TAG	UNP P13334
B	70	ASN	-	EXPRESSION TAG	UNP P13334
B	71	TRP	-	EXPRESSION TAG	UNP P13334

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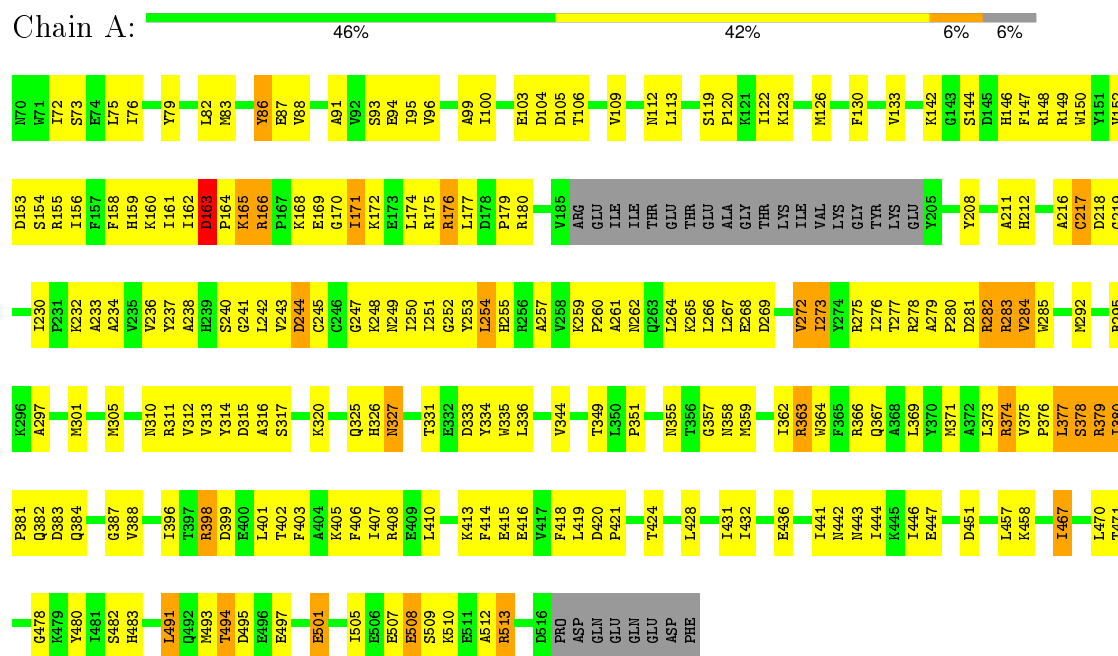
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Chain	Residue	Modelled	Actual	Comment	Reference
B	72	ILE	-	EXPRESSION TAG	UNP P13334
B	73	SER	-	EXPRESSION TAG	UNP P13334
C	70	ASN	-	EXPRESSION TAG	UNP P13334
C	71	TRP	-	EXPRESSION TAG	UNP P13334
C	72	ILE	-	EXPRESSION TAG	UNP P13334
C	73	SER	-	EXPRESSION TAG	UNP P13334
D	70	ASN	-	EXPRESSION TAG	UNP P13334
D	71	TRP	-	EXPRESSION TAG	UNP P13334
D	72	ILE	-	EXPRESSION TAG	UNP P13334
D	73	SER	-	EXPRESSION TAG	UNP P13334
E	70	ASN	-	EXPRESSION TAG	UNP P13334
E	71	TRP	-	EXPRESSION TAG	UNP P13334
E	72	ILE	-	EXPRESSION TAG	UNP P13334
E	73	SER	-	EXPRESSION TAG	UNP P13334
F	70	ASN	-	EXPRESSION TAG	UNP P13334
F	71	TRP	-	EXPRESSION TAG	UNP P13334
F	72	ILE	-	EXPRESSION TAG	UNP P13334
F	73	SER	-	EXPRESSION TAG	UNP P13334
G	70	ASN	-	EXPRESSION TAG	UNP P13334
G	71	TRP	-	EXPRESSION TAG	UNP P13334
G	72	ILE	-	EXPRESSION TAG	UNP P13334
G	73	SER	-	EXPRESSION TAG	UNP P13334
H	70	ASN	-	EXPRESSION TAG	UNP P13334
H	71	TRP	-	EXPRESSION TAG	UNP P13334
H	72	ILE	-	EXPRESSION TAG	UNP P13334
H	73	SER	-	EXPRESSION TAG	UNP P13334
I	70	ASN	-	EXPRESSION TAG	UNP P13334
I	71	TRP	-	EXPRESSION TAG	UNP P13334
I	72	ILE	-	EXPRESSION TAG	UNP P13334
I	73	SER	-	EXPRESSION TAG	UNP P13334
J	70	ASN	-	EXPRESSION TAG	UNP P13334
J	71	TRP	-	EXPRESSION TAG	UNP P13334
J	72	ILE	-	EXPRESSION TAG	UNP P13334
J	73	SER	-	EXPRESSION TAG	UNP P13334
K	70	ASN	-	EXPRESSION TAG	UNP P13334
K	71	TRP	-	EXPRESSION TAG	UNP P13334
K	72	ILE	-	EXPRESSION TAG	UNP P13334
K	73	SER	-	EXPRESSION TAG	UNP P13334
L	70	ASN	-	EXPRESSION TAG	UNP P13334
L	71	TRP	-	EXPRESSION TAG	UNP P13334
L	72	ILE	-	EXPRESSION TAG	UNP P13334
L	73	SER	-	EXPRESSION TAG	UNP P13334

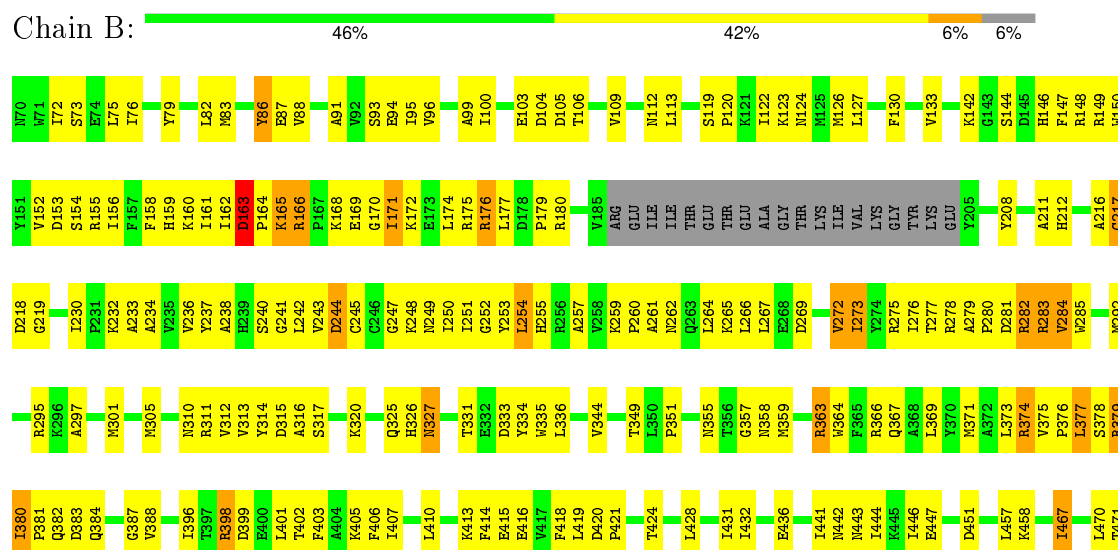
3 Residue-property plots

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. 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. 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: Portal protein gp20



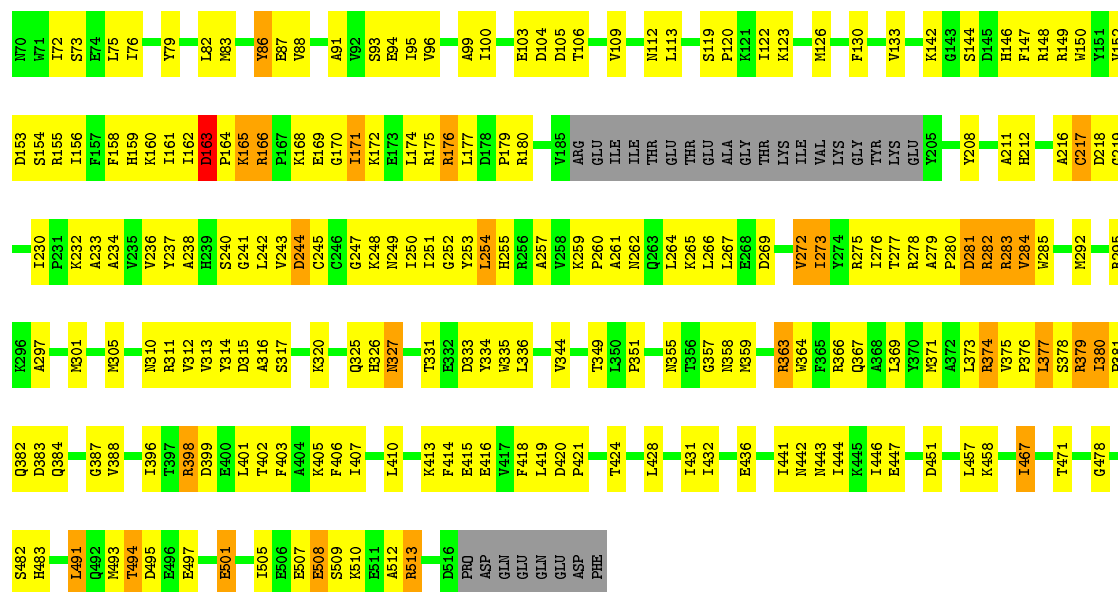
• Molecule 1: Portal protein gp20





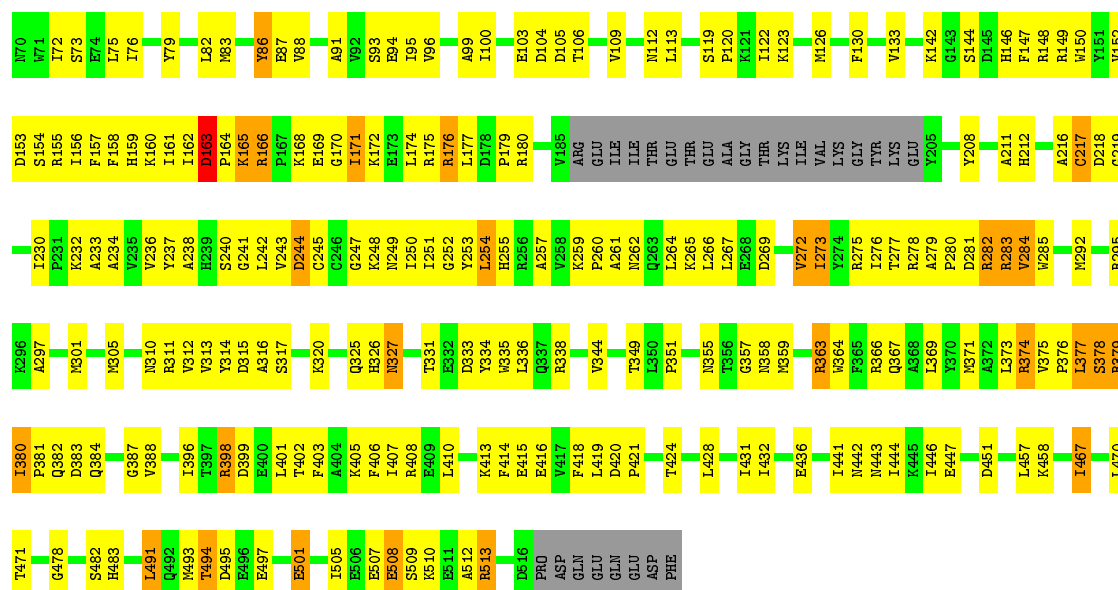
• Molecule 1: Portal protein gp20

Chain C: 47% 41% 6% 6%



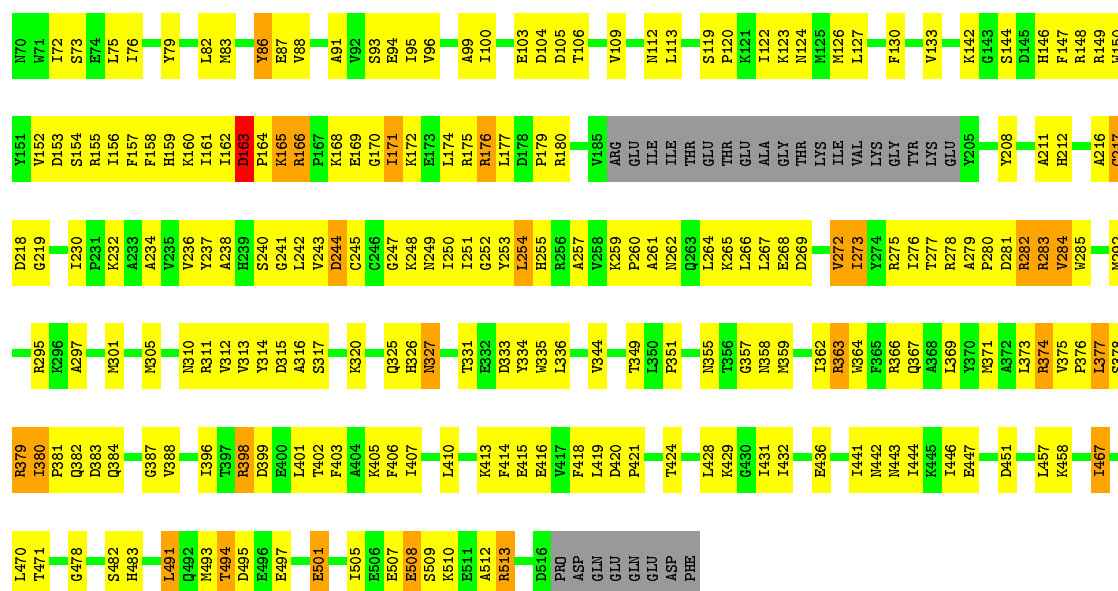
• Molecule 1: Portal protein gp20

Chain D: 46% 42% 6% 6%



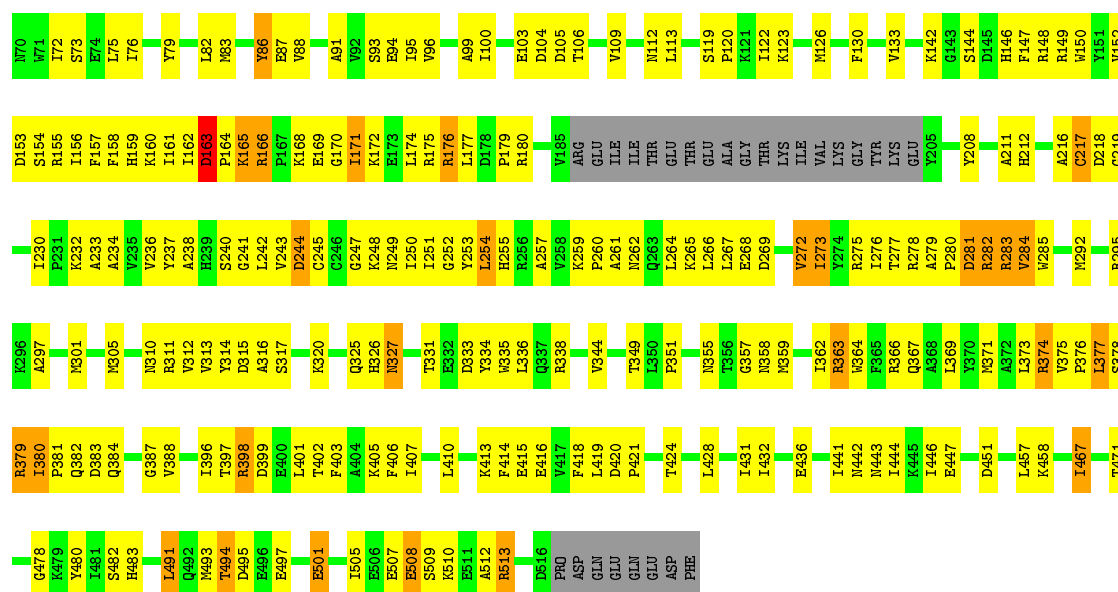
• Molecule 1: Portal protein gp20

Chain E: 46% 42% 6% 6%



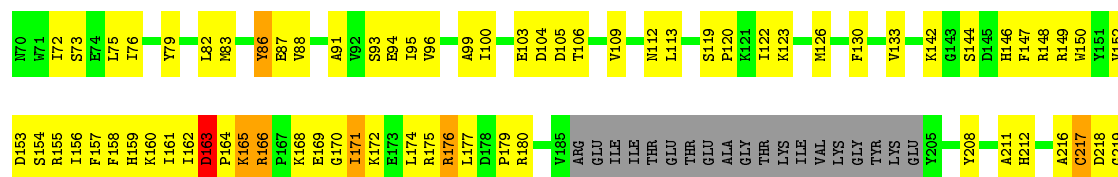
• Molecule 1: Portal protein gp20

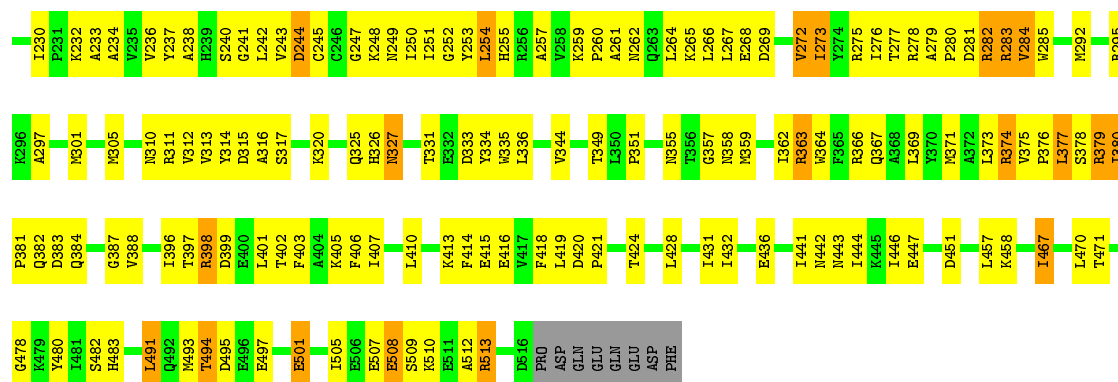
Chain F: 46% 42% 6% 6%



• Molecule 1: Portal protein gp20

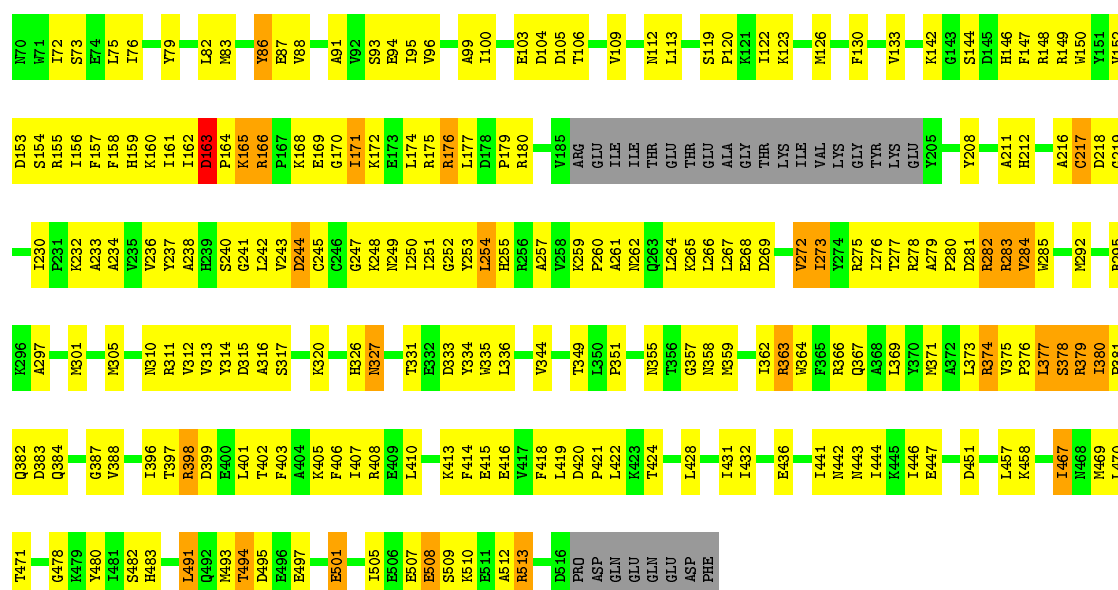
Chain G: 46% 42% 6% 6%





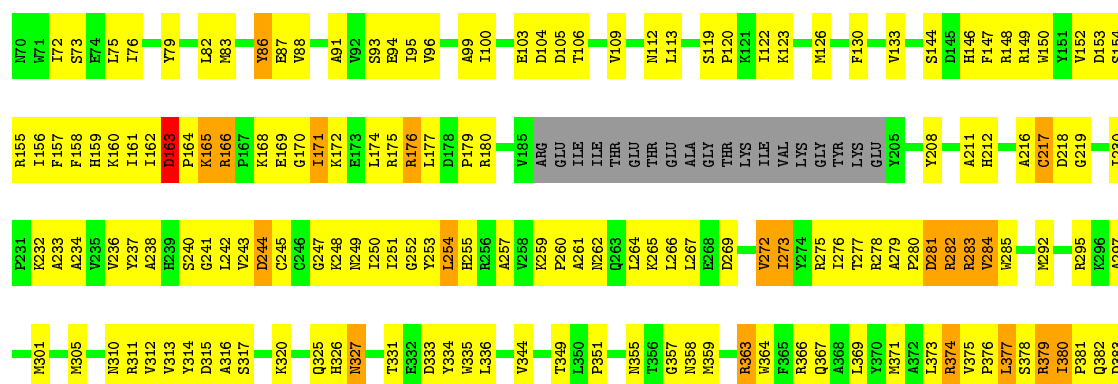
• Molecule 1: Portal protein gp20

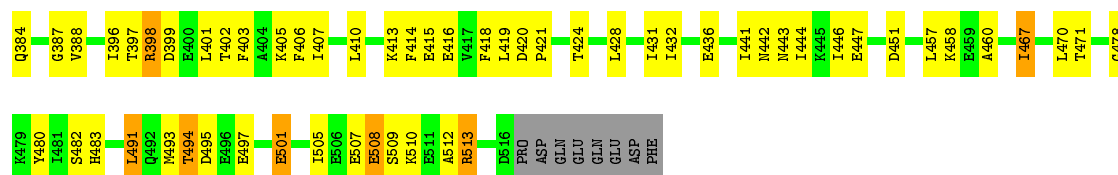
Chain H: 45% 43% 6% 6%



• Molecule 1: Portal protein gp20

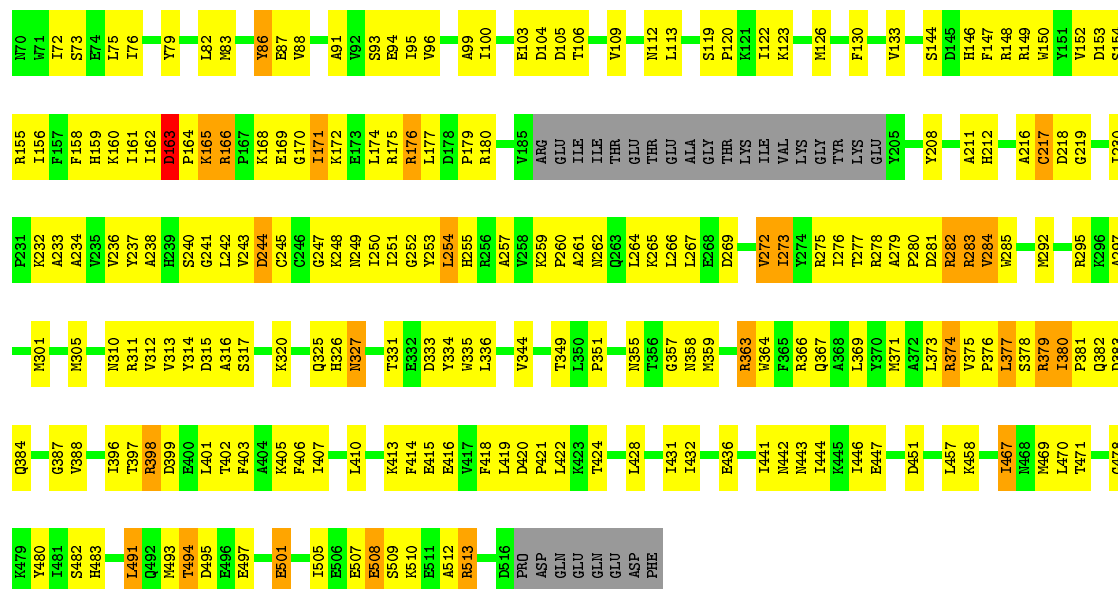
Chain I: 46% 42% 6% 6%





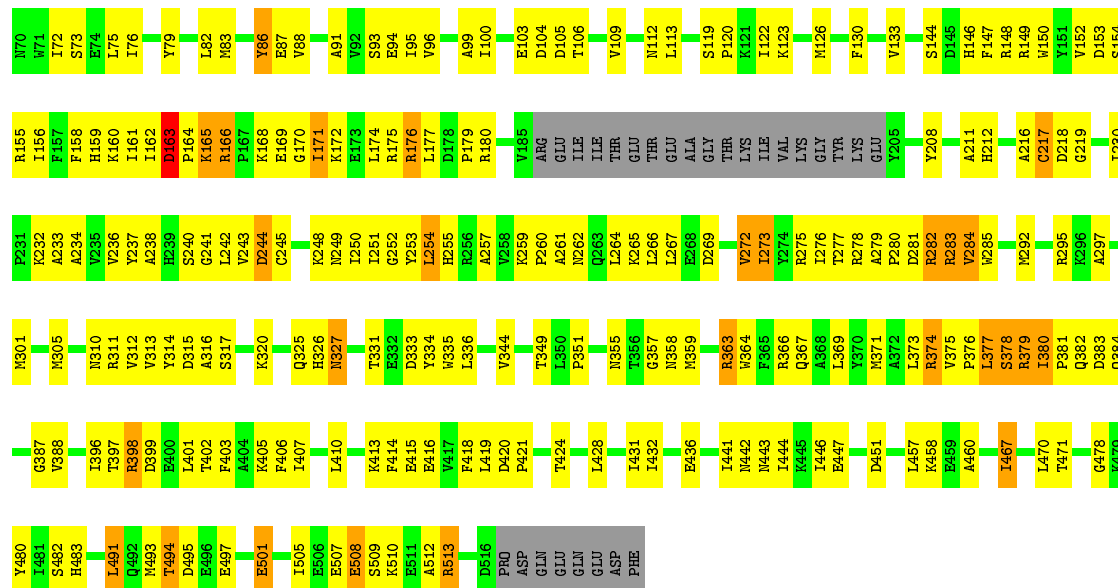
• Molecule 1: Portal protein gp20

Chain J: 46% 42% 6% 6%

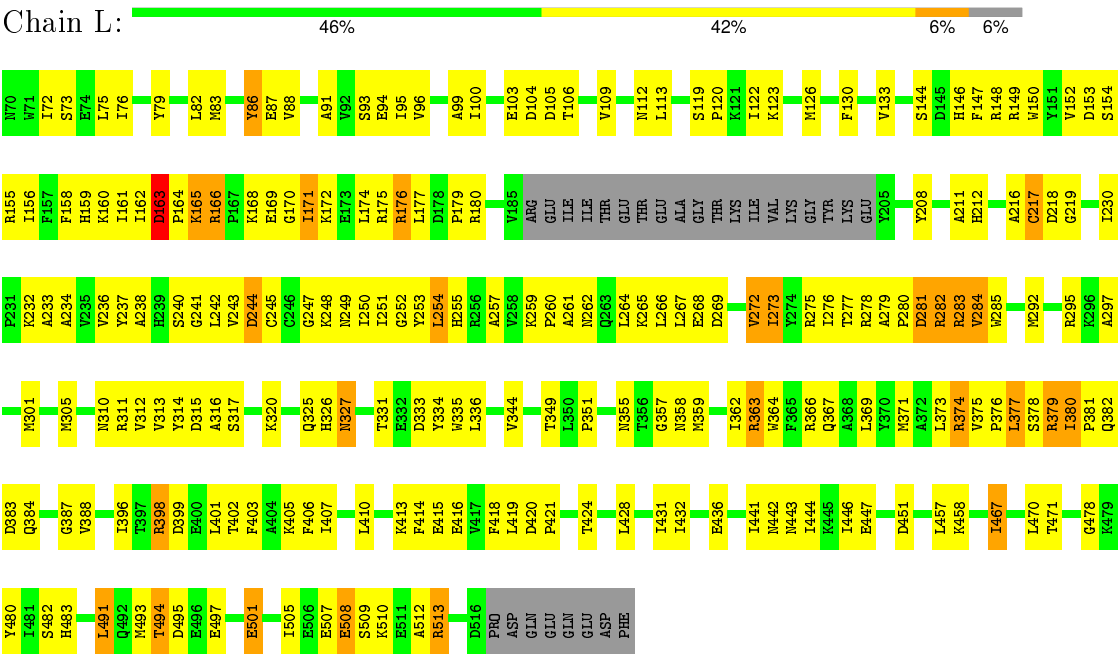


• Molecule 1: Portal protein gp20

Chain K: 47% 41% 6% 6%



• Molecule 1: Portal protein gp20



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	98000	Depositor
Resolution determination method	FSC 0.143	Depositor
CTF correction method	each particle	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	2	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	29000	Depositor
Image detector	GATAN K2 (4k x 4 k)	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.46	0/3504	0.64	2/4730 (0.0%)
1	B	0.46	0/3504	0.64	2/4730 (0.0%)
1	C	0.46	0/3504	0.64	2/4730 (0.0%)
1	D	0.46	0/3504	0.64	2/4730 (0.0%)
1	E	0.46	0/3504	0.64	2/4730 (0.0%)
1	F	0.46	0/3504	0.64	2/4730 (0.0%)
1	G	0.46	0/3504	0.64	2/4730 (0.0%)
1	H	0.46	0/3504	0.64	2/4730 (0.0%)
1	I	0.46	0/3504	0.64	2/4730 (0.0%)
1	J	0.46	0/3504	0.64	2/4730 (0.0%)
1	K	0.46	0/3504	0.64	2/4730 (0.0%)
1	L	0.46	0/3504	0.64	2/4730 (0.0%)
All	All	0.46	0/42048	0.64	24/56760 (0.0%)

There are no bond length outliers.

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	379	ARG	N-CA-C	-7.33	91.21	111.00
1	D	379	ARG	N-CA-C	-7.33	91.21	111.00
1	G	379	ARG	N-CA-C	-7.33	91.21	111.00
1	J	379	ARG	N-CA-C	-7.33	91.21	111.00
1	C	379	ARG	N-CA-C	-7.33	91.21	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3440	0	3344	285	0
1	B	3440	0	3344	282	0
1	C	3440	0	3344	272	0
1	D	3440	0	3344	280	0
1	E	3440	0	3344	278	0
1	F	3440	0	3344	281	0
1	G	3440	0	3344	276	0
1	H	3440	0	3344	285	0
1	I	3440	0	3344	285	0
1	J	3440	0	3344	278	0
1	K	3440	0	3344	285	0
1	L	3440	0	3344	278	0
All	All	41280	0	40128	3003	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 37.

The worst 5 of 3003 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:162:ILE:CG2	1:G:171:ILE:HG13	1.39	1.52
1:F:162:ILE:CG2	1:F:171:ILE:HG13	1.39	1.52
1:A:162:ILE:CG2	1:A:171:ILE:HG13	1.39	1.51
1:H:162:ILE:CG2	1:H:171:ILE:HG13	1.39	1.51
1:E:162:ILE:CG2	1:E:171:ILE:HG13	1.39	1.51

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	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	B	424/455 (93%)	369 (87%)	46 (11%)	9 (2%)	9	53
1	C	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	D	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	E	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	F	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	G	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	H	424/455 (93%)	369 (87%)	46 (11%)	9 (2%)	9	53
1	I	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	J	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	K	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
1	L	424/455 (93%)	368 (87%)	47 (11%)	9 (2%)	9	53
All	All	5088/5460 (93%)	4418 (87%)	562 (11%)	108 (2%)	13	53

5 of 108 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	ILE
1	A	380	ILE
1	B	171	ILE
1	B	380	ILE
1	C	171	ILE

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	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	B	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	C	364/407 (89%)	334 (92%)	30 (8%)	14	53
1	D	364/407 (89%)	335 (92%)	29 (8%)	15	55

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	F	364/407 (89%)	334 (92%)	30 (8%)	14	53
1	G	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	H	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	I	364/407 (89%)	334 (92%)	30 (8%)	14	53
1	J	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	K	364/407 (89%)	335 (92%)	29 (8%)	15	55
1	L	364/407 (89%)	334 (92%)	30 (8%)	14	53
All	All	4368/4884 (89%)	4016 (92%)	352 (8%)	19	54

5 of 352 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	312	VAL
1	G	457	LEU
1	L	176	ARG
1	F	378	SER
1	G	166	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 132 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	262	ASN
1	G	358	ASN
1	L	90	ASN
1	F	355	ASN
1	G	146	HIS

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 [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.