



wwPDB EM Validation Summary Report ⓘ

Mar 29, 2026 – 05:11 PM UTC

PDB ID : 6XR4 / pdb_00006xr4
EMDB ID : EMD-20825
Title : Integrative in situ structure of Parkinsons disease-linked human LRRK2
Authors : Villa, E.; Lasker, K.; Audagnotto, M.
Deposited on : 2020-07-10
Resolution : 14.00 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

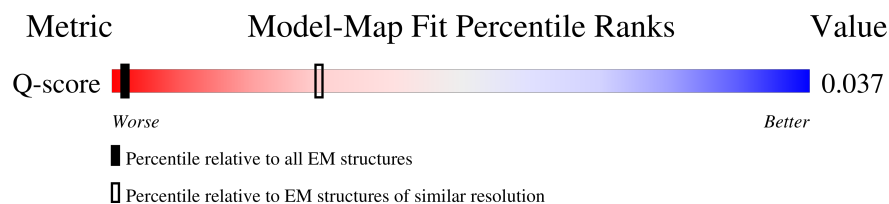
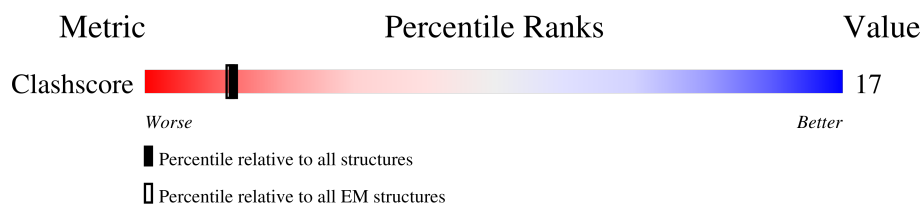
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

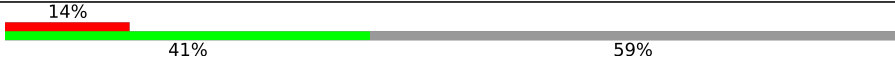
The reported resolution of this entry is 14.00 Å.

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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Q-score	-	25397	61 (13.50 - 14.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	1-A	2527	 10% 41% 59%
1	1-B	2527	 14% 41% 59%
1	10-A	2527	 40% . 59%
1	10-B	2527	 40% . 59%
1	11-A	2527	 41% 59%
1	11-B	2527	 41% 59%

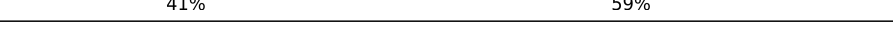
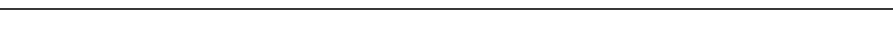
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Mol	Chain	Length	Quality of chain	
1	12-A	2527		
1	12-B	2527		
1	13-A	2527		
1	13-B	2527		
1	14-A	2527		
1	14-B	2527		
1	15-A	2527		
1	15-B	2527		
1	16-A	2527		
1	16-B	2527		
1	17-A	2527		
1	17-B	2527		
1	18-A	2527		
1	18-B	2527		
1	19-A	2527		
1	19-B	2527		
1	2-A	2527		
1	2-B	2527		
1	20-A	2527		
1	20-B	2527		
1	21-A	2527		
1	21-B	2527		
1	22-A	2527		
1	22-B	2527		
1	23-A	2527		

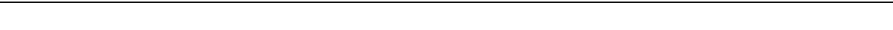
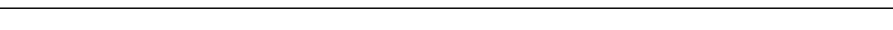
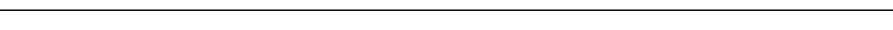
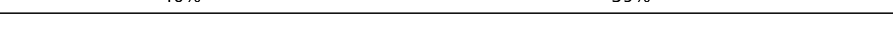
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Mol	Chain	Length	Quality of chain	
1	23-B	2527		
1	24-A	2527		
1	24-B	2527		
1	25-A	2527		
1	25-B	2527		
1	26-A	2527		
1	26-B	2527		
1	27-A	2527		
1	27-B	2527		
1	28-A	2527		
1	28-B	2527		
1	29-A	2527		
1	29-B	2527		
1	3-A	2527		
1	3-B	2527		
1	30-A	2527		
1	30-B	2527		
1	31-A	2527		
1	31-B	2527		
1	32-A	2527		
1	32-B	2527		
1	33-A	2527		
1	33-B	2527		
1	34-A	2527		
1	34-B	2527		

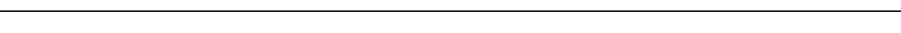
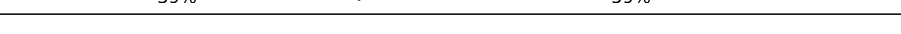
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Mol	Chain	Length	Quality of chain	
1	35-A	2527		
1	35-B	2527		
1	36-A	2527		
1	36-B	2527		
1	37-A	2527		
1	37-B	2527		
1	38-A	2527		
1	38-B	2527		
1	39-A	2527		
1	39-B	2527		
1	4-A	2527		
1	4-B	2527		
1	40-A	2527		
1	40-B	2527		
1	41-A	2527		
1	41-B	2527		
1	42-A	2527		
1	42-B	2527		
1	43-A	2527		
1	43-B	2527		
1	44-A	2527		
1	44-B	2527		
1	45-A	2527		
1	45-B	2527		
1	46-A	2527		

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Mol	Chain	Length	Quality of chain
1	46-B	2527	 40% 59%
1	47-A	2527	 40% 59%
1	47-B	2527	 40% 59%
1	48-A	2527	 40% 59%
1	48-B	2527	 40% 59%
1	49-A	2527	 40% 59%
1	49-B	2527	 40% 59%
1	5-A	2527	 40% 59%
1	5-B	2527	 40% 59%
1	50-A	2527	 39% 59%
1	50-B	2527	 39% 59%
1	51-A	2527	 39% 59%
1	51-B	2527	 39% 59%
1	52-A	2527	 39% 59%
1	52-B	2527	 39% 59%
1	53-A	2527	 39% 59%
1	53-B	2527	 39% 59%
1	6-A	2527	 41% 59%
1	6-B	2527	 41% 59%
1	7-A	2527	 39% 59%
1	7-B	2527	 39% 59%
1	8-A	2527	 40% 59%
1	8-B	2527	 40% 59%
1	9-A	2527	 41% 59%
1	9-B	2527	 40% 59%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 110240 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 Leucine-rich repeat serine/threonine-protein kinase 2.

Mol	Chain	Residues	Atoms	AltConf	Trace
1	1-A	1040	Total C 1040 1040	0	1040
1	2-A	1040	Total C 1040 1040	0	1040
1	3-A	1040	Total C 1040 1040	0	1040
1	4-A	1040	Total C 1040 1040	0	1040
1	5-A	1040	Total C 1040 1040	0	1040
1	6-A	1040	Total C 1040 1040	0	1040
1	7-A	1040	Total C 1040 1040	0	1040
1	8-A	1040	Total C 1040 1040	0	1040
1	9-A	1040	Total C 1040 1040	0	1040
1	10-A	1040	Total C 1040 1040	0	1040
1	11-A	1040	Total C 1040 1040	0	1040
1	12-A	1040	Total C 1040 1040	0	1040
1	13-A	1040	Total C 1040 1040	0	1040
1	14-A	1040	Total C 1040 1040	0	1040
1	15-A	1040	Total C 1040 1040	0	1040
1	16-A	1040	Total C 1040 1040	0	1040
1	17-A	1040	Total C 1040 1040	0	1040

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Mol	Chain	Residues	Atoms	AltConf	Trace
1	18-A	1040	Total C 1040 1040	0	1040
1	19-A	1040	Total C 1040 1040	0	1040
1	20-A	1040	Total C 1040 1040	0	1040
1	21-A	1040	Total C 1040 1040	0	1040
1	22-A	1040	Total C 1040 1040	0	1040
1	23-A	1040	Total C 1040 1040	0	1040
1	24-A	1040	Total C 1040 1040	0	1040
1	25-A	1040	Total C 1040 1040	0	1040
1	26-A	1040	Total C 1040 1040	0	1040
1	27-A	1040	Total C 1040 1040	0	1040
1	28-A	1040	Total C 1040 1040	0	1040
1	29-A	1040	Total C 1040 1040	0	1040
1	30-A	1040	Total C 1040 1040	0	1040
1	31-A	1040	Total C 1040 1040	0	1040
1	32-A	1040	Total C 1040 1040	0	1040
1	33-A	1040	Total C 1040 1040	0	1040
1	34-A	1040	Total C 1040 1040	0	1040
1	35-A	1040	Total C 1040 1040	0	1040
1	36-A	1040	Total C 1040 1040	0	1040
1	37-A	1040	Total C 1040 1040	0	1040
1	38-A	1040	Total C 1040 1040	0	1040

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Mol	Chain	Residues	Atoms	AltConf	Trace
1	39-A	1040	Total C 1040 1040	0	1040
1	40-A	1040	Total C 1040 1040	0	1040
1	41-A	1040	Total C 1040 1040	0	1040
1	42-A	1040	Total C 1040 1040	0	1040
1	43-A	1040	Total C 1040 1040	0	1040
1	44-A	1040	Total C 1040 1040	0	1040
1	45-A	1040	Total C 1040 1040	0	1040
1	46-A	1040	Total C 1040 1040	0	1040
1	47-A	1040	Total C 1040 1040	0	1040
1	48-A	1040	Total C 1040 1040	0	1040
1	49-A	1040	Total C 1040 1040	0	1040
1	50-A	1040	Total C 1040 1040	0	1040
1	51-A	1040	Total C 1040 1040	0	1040
1	52-A	1040	Total C 1040 1040	0	1040
1	53-A	1040	Total C 1040 1040	0	1040
1	1-B	1040	Total C 1040 1040	0	1040
1	2-B	1040	Total C 1040 1040	0	1040
1	3-B	1040	Total C 1040 1040	0	1040
1	4-B	1040	Total C 1040 1040	0	1040
1	5-B	1040	Total C 1040 1040	0	1040
1	6-B	1040	Total C 1040 1040	0	1040

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Mol	Chain	Residues	Atoms	AltConf	Trace
1	7-B	1040	Total C 1040 1040	0	1040
1	8-B	1040	Total C 1040 1040	0	1040
1	9-B	1040	Total C 1040 1040	0	1040
1	10-B	1040	Total C 1040 1040	0	1040
1	11-B	1040	Total C 1040 1040	0	1040
1	12-B	1040	Total C 1040 1040	0	1040
1	13-B	1040	Total C 1040 1040	0	1040
1	14-B	1040	Total C 1040 1040	0	1040
1	15-B	1040	Total C 1040 1040	0	1040
1	16-B	1040	Total C 1040 1040	0	1040
1	17-B	1040	Total C 1040 1040	0	1040
1	18-B	1040	Total C 1040 1040	0	1040
1	19-B	1040	Total C 1040 1040	0	1040
1	20-B	1040	Total C 1040 1040	0	1040
1	21-B	1040	Total C 1040 1040	0	1040
1	22-B	1040	Total C 1040 1040	0	1040
1	23-B	1040	Total C 1040 1040	0	1040
1	24-B	1040	Total C 1040 1040	0	1040
1	25-B	1040	Total C 1040 1040	0	1040
1	26-B	1040	Total C 1040 1040	0	1040
1	27-B	1040	Total C 1040 1040	0	1040

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Mol	Chain	Residues	Atoms	AltConf	Trace
1	28-B	1040	Total C 1040 1040	0	1040
1	29-B	1040	Total C 1040 1040	0	1040
1	30-B	1040	Total C 1040 1040	0	1040
1	31-B	1040	Total C 1040 1040	0	1040
1	32-B	1040	Total C 1040 1040	0	1040
1	33-B	1040	Total C 1040 1040	0	1040
1	34-B	1040	Total C 1040 1040	0	1040
1	35-B	1040	Total C 1040 1040	0	1040
1	36-B	1040	Total C 1040 1040	0	1040
1	37-B	1040	Total C 1040 1040	0	1040
1	38-B	1040	Total C 1040 1040	0	1040
1	39-B	1040	Total C 1040 1040	0	1040
1	40-B	1040	Total C 1040 1040	0	1040
1	41-B	1040	Total C 1040 1040	0	1040
1	42-B	1040	Total C 1040 1040	0	1040
1	43-B	1040	Total C 1040 1040	0	1040
1	44-B	1040	Total C 1040 1040	0	1040
1	45-B	1040	Total C 1040 1040	0	1040
1	46-B	1040	Total C 1040 1040	0	1040
1	47-B	1040	Total C 1040 1040	0	1040
1	48-B	1040	Total C 1040 1040	0	1040

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Mol	Chain	Residues	Atoms		AltConf	Trace
1	49-B	1040	Total 1040	C 1040	0	1040
1	50-B	1040	Total 1040	C 1040	0	1040
1	51-B	1040	Total 1040	C 1040	0	1040
1	52-B	1040	Total 1040	C 1040	0	1040
1	53-B	1040	Total 1040	C 1040	0	1040

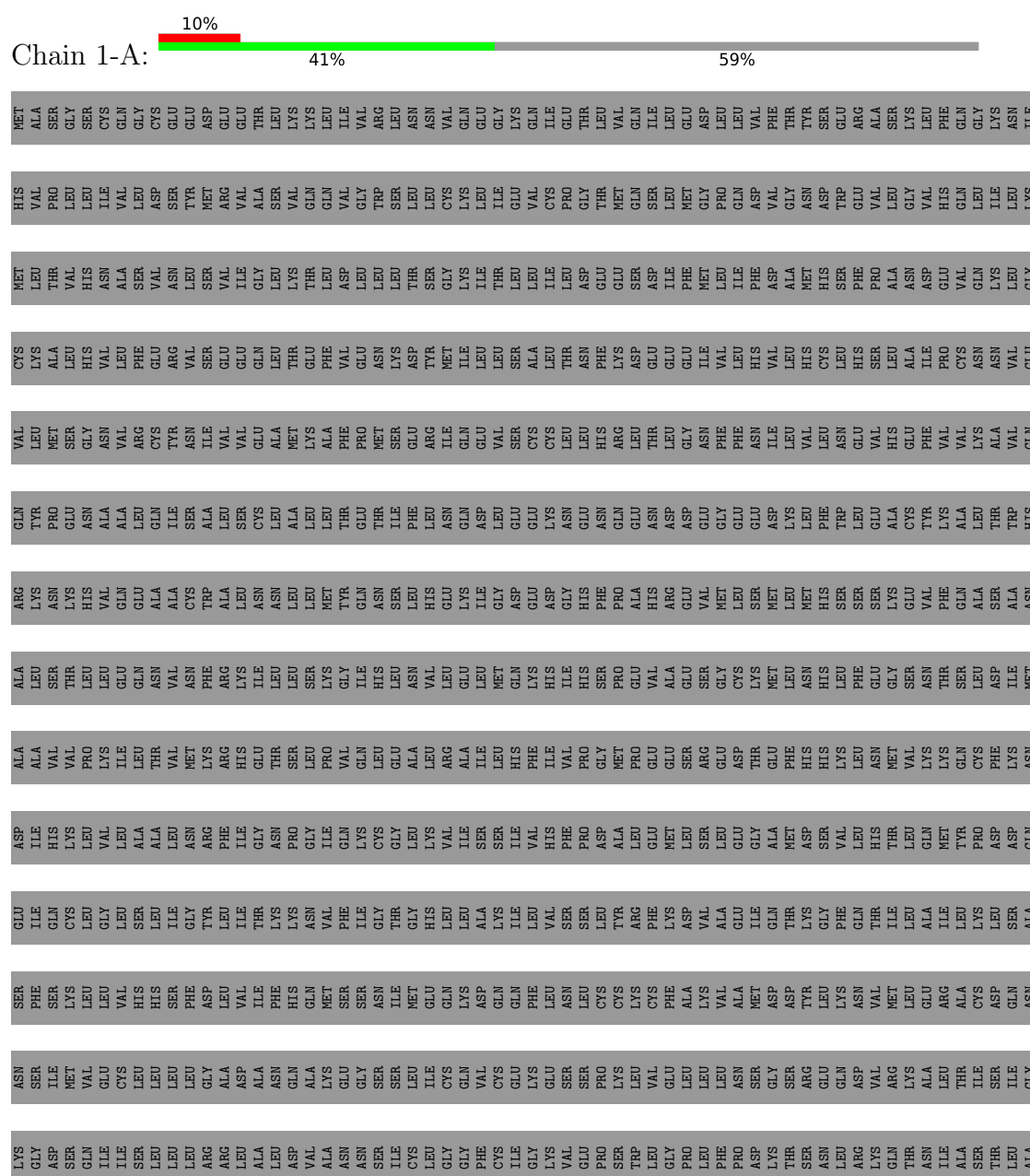
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2020	THR	ILE	variant	UNP Q5S007
B	2020	THR	ILE	variant	UNP Q5S007

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: Leucine-rich repeat serine/threonine-protein kinase 2





[illegible]

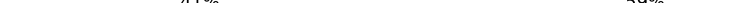
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

59%



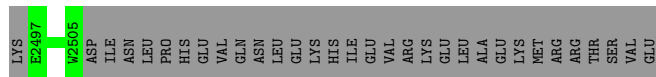
[illegible]

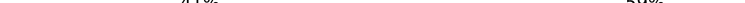
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

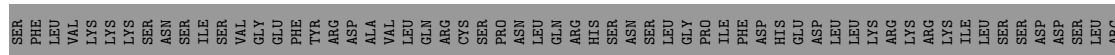
Chain 2-B:  41% 59%

[illegible]





- Chain 3-B:  41% 59%

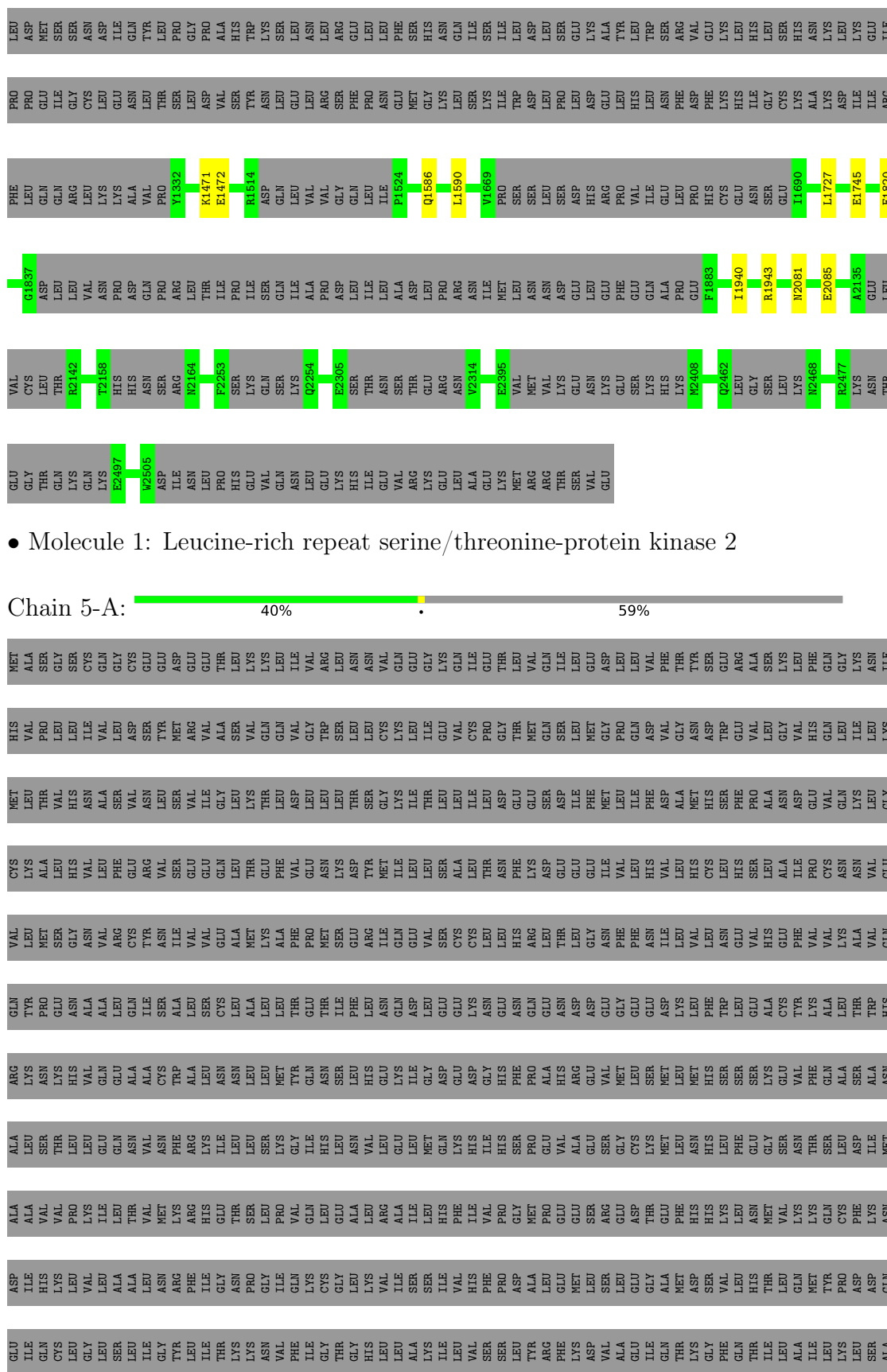




- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

59%

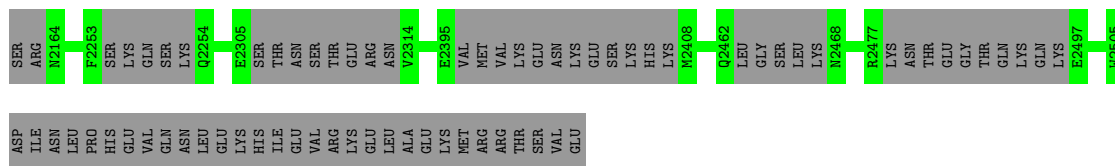




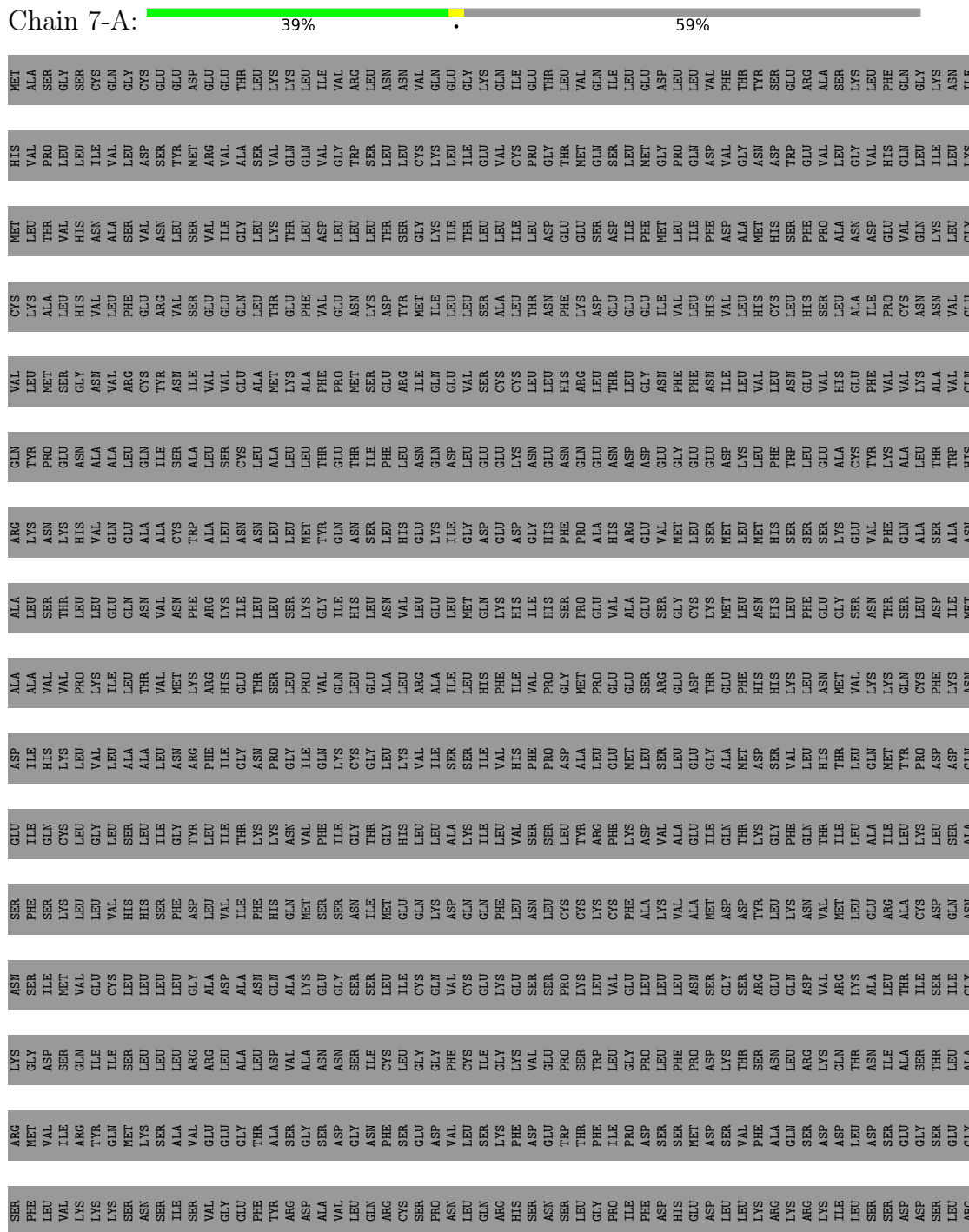






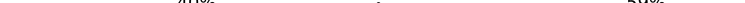


- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2





GLY	THR	GLN	LYS	GLN	LYS	LEU	PRO	HIS	VAL	GLN	ASN	LEU	GLU	HIS	ILE	GLU	VAL	ARG	LYS	GLU	LEU	ALA	GLU	LYS	MET	ARG	ARG	THR	SER	VAL	GLU
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

- Chain 8-A:  40% 59%

SER PHE LEU VAL LYS LYS SER SER SER ILE VAL GLY PHE TYR ARG ASP ALA VAL LEU GLN ARG CYS SER PRO ASN LEU GLN ARG HIS SER ASN SER SER LEU GLY PRO ILE PHE ASP HIS GLU ASP LEU LEU LYS ARG LYS LYS LEU LEU SER SER ASP ASP SER SER LEU





Chain 9-A: 41% 59%





Chain 11-A: 41% 59%



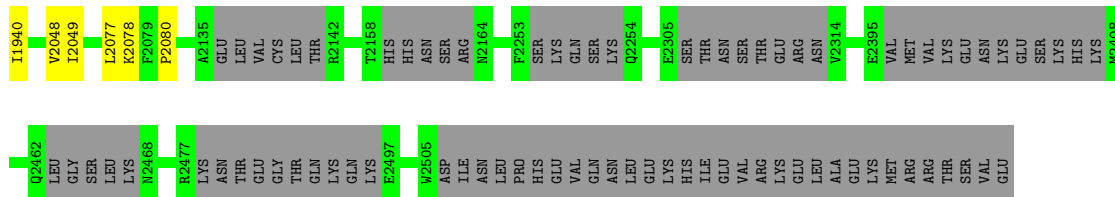
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

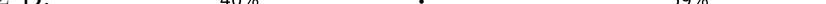
Chain 11-B:



GLU	ILE	ASP	ALA	ALA	ARG	GLN	VAL	CYS	MET	HIS	MET
ILE	GLN	ILE	LEU	LEU	ASN	TYR	TYR	LYS	LEU	VAL	ALA
GLN	HIS	GLN	SER	SER	ASN	PRO	PRO	ALA	THR	PRO	ALA
LEU	LEU	LYS	THR	THR	HIS	GLU	GLY	LEU	VAL	LEU	GLY
GLY	LEU	VAL	LEU	VAL	VAL	ALA	ASN	VAL	ASN	ILE	CYS
LEU	ILE	LEU	GLU	GLN	GLN	ALA	VAL	LEU	VAL	VAL	GLN
SER	LEU	ALA	GLN	GLU	ASN	LEU	ARG	PHE	SER	GLY	GLY
LEU	ALA	THR	ASN	ASN	ALA	GLN	VAL	ARG	VAL	ASP	CYS
GLY	LEU	MET	ASN	ASN	ALA	ILE	TYR	GLU	ASN	ASP	GLU
TYR	ARG	ASN	PHE	PHE	TRP	SER	ALA	VAL	SER	MET	ASP
LEU	PHE	ARG	ARG	ARG	ALA	LEU	VAL	GLU	VAL	ARG	GLU
ILE	ILE	ILE	HIS	ILE	ASN	SER	VAL	GLN	ILE	VAL	THR
THR	GLY	ASN	GLU	ILE	LEU	CYS	GLU	GLN	GLY	ALA	ALA
LYS	ASN	THR	THR	THR	ASN	LEU	ALA	LEU	LEU	SER	LYS
ASN	GLY	PRO	SER	SER	LEU	ALA	MET	THR	LYS	VAL	LYS
VAL	VAL	ILE	PRO	LYS	LEU	LEU	LYS	GLU	THR	GLN	LEU
PHE	PHE	GLN	VAL	VAL	TYR	THR	PHE	VAL	PHE	VAL	ILE
ILE	GLY	LYS	GLN	ILE	ILE	GLU	GLU	GLU	GLU	VAL	GLY
GLY	GLY	CYS	ALA	LEU	ASN	THR	MET	ASN	LEU	TRP	ARG
THR	GLY	GLY	LEU	LEU	LEU	ILE	GLU	LYS	LEU	GLY	GLY
GLY	GLY	GLY	ALA	ASN	LEU	PHE	GLU	ASP	THR	LEU	ASN
HIS	LYS	LYS	LEU	VAL	HIS	GLU	SER	TYR	SER	LEU	ASN
LEU	LEU	VAL	LYS	GLN	GLU	GLU	GLY	ASP	GLY	VAL	VAL
LEU	VAL	ILE	PHE	GLU	GLU	ASN	ILE	MET	GLY	CYS	GLN
SER	SER	PHE	VAL	ILE	ILE	GLN	GLN	ILE	GLY	LYS	VAL
SER	SER	PRO	PRO	GLY	HIS	ASN	GLU	THR	LEU	GLY	THR
LEU	LEU	ASP	GLY	SER	PHE	ASN	HIS	PHE	GLU	MET	ASN
TYR	TYR	ALA	MET	PRO	PRO	GLN	ARG	LYS	GLU	THR	VAL
ARG	ARG	LEU	PRO	LEU	ALA	GLU	LEU	ASP	SER	GLN	GLN
PHE	PHE	GLU	GLU	VAL	ARG	ASN	THR	GLU	ASP	SER	ILE
LYS	LYS	MET	GLU	ALA	HIS	ASN	CYS	LEU	LEU	ILE	ILE
ASP	ASP	LEU	SER	GLU	GLY	ASP	GLY	GLU	PHE	MET	GLY
VAL	VAL	SER	ARG	SER	VAL	GLU	ASN	ILE	MET	GLY	ASP
ALA	ALA	LEU	GLU	GLY	MET	GLY	PHE	VAL	LEU	PRO	LEU
GLU	GLU	GLY	ASP	CYS	LEU	GLU	GLU	VAL	ILE	GLN	LEU
ILE	ILE	GLU	LYS	LYS	SER	GLU	ILE	LEU	ILE	GLN	SER
GLN	GLN	ALA	GLU	MET	MET	LYS	ILE	LEU	ASP	VAL	PHE
THR	THR	MET	PHE	LEU	LEU	LYS	LEU	LEU	ALA	GLY	THR
LYS	LYS	ASP	HIS	ASN	MET	LEU	VAL	VAL	ASN	ASN	TYR
GLY	GLY	SER	HIS	HIS	SER	PHE	LEU	CYS	HIS	ASP	SER
PHE	PHE	VAL	LYS	LYS	SER	TRP	LEU	LEU	TRP	TRP	GLU
GLN	GLN	LEU	PHE	PHE	SER	THR	GLU	HIS	SER	GLU	ARG
THR	THR	LEU	GLU	GLU	SER	GLU	VAL	SER	PRO	VAL	ALA
ILE	ILE	THR	GLY	GLY	LEU	ALA	HIS	LEU	ALA	GLY	SER
ALA	ALA	GLN	ASN	VAL	GLU	CYS	ASN	ALA	ASN	LEU	LYS
LEU	LEU	LEU	SER	SER	GLU	TYR	ASP	ILE	GLY	VAL	GLY
LEU	LEU	ILE	THR	THR	LEU	LEU	THR	THR	THR	VAL	LEU
LEU	LEU	VAL	VAL	PHE	LEU	LEU	LEU	PHE	LEU	VAL	LEU
LEU	LEU	GLY	GLY	GLY	GLY	GL					





Chain 12-B:  40% . 59%



ARG	ARG	THR	SER	VAL	GLU	MET	VAL	LYS	GLU	ASN	LYS	GLU	SER	LYS	HIS	LYS	P2408	P2462	LEU	GLY	SER	SER	LEU	LYS	P2468	P2477	LYS	ASN	THR	GLU	GLY	THR	GLN	LYS	GLN	LYS	P2497	P2505	ASP	ILE	ASN	LEU	PRO	HIS	GLU	VAL	GLN	ASN	LEU	GLY	LYS	HIS	ILE	GLU	VAL	ARG	LYS	GLU	LEU	ALA	GLU	LYS	MET
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- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 13-B: 40% . 59%

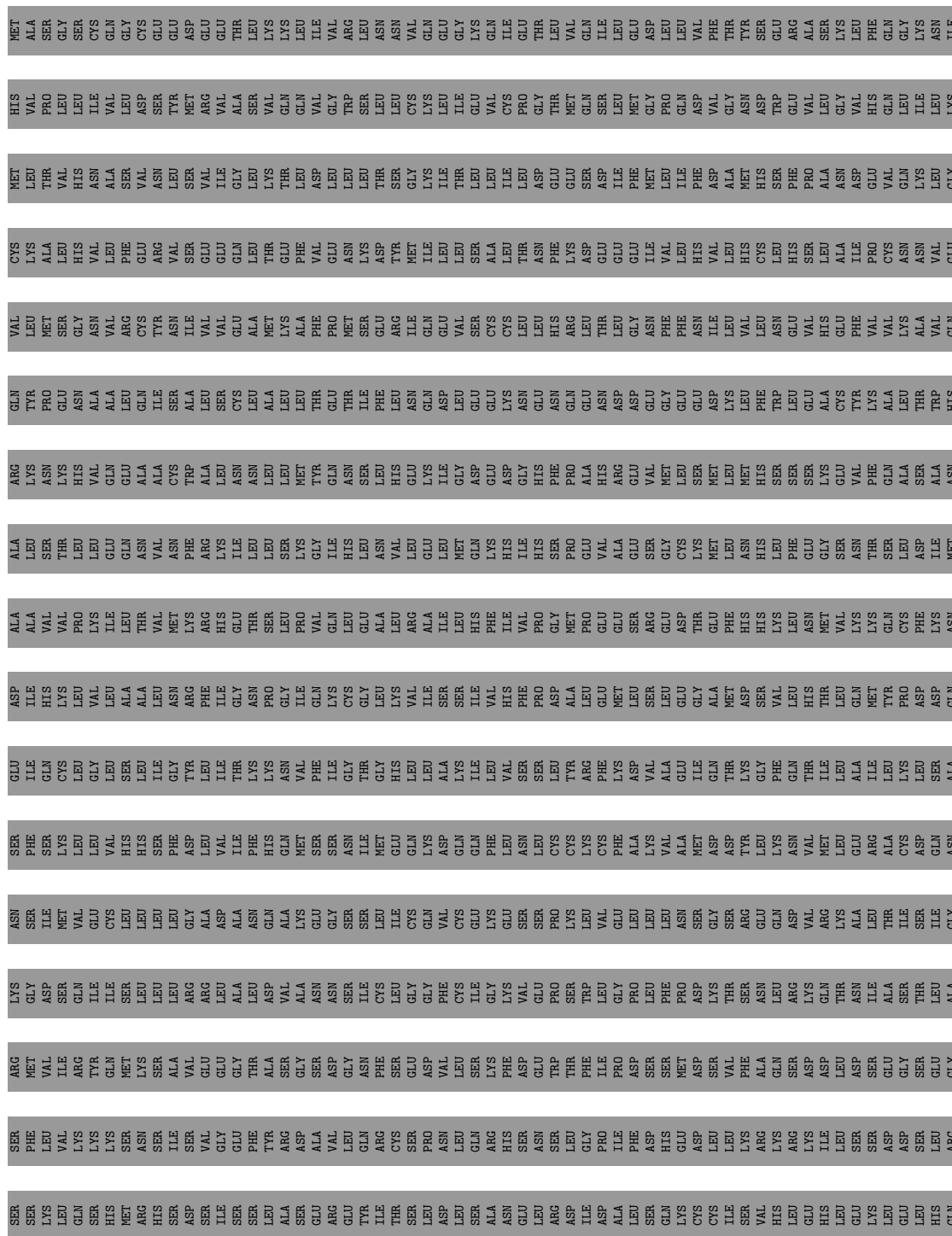
[illegible]





- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2



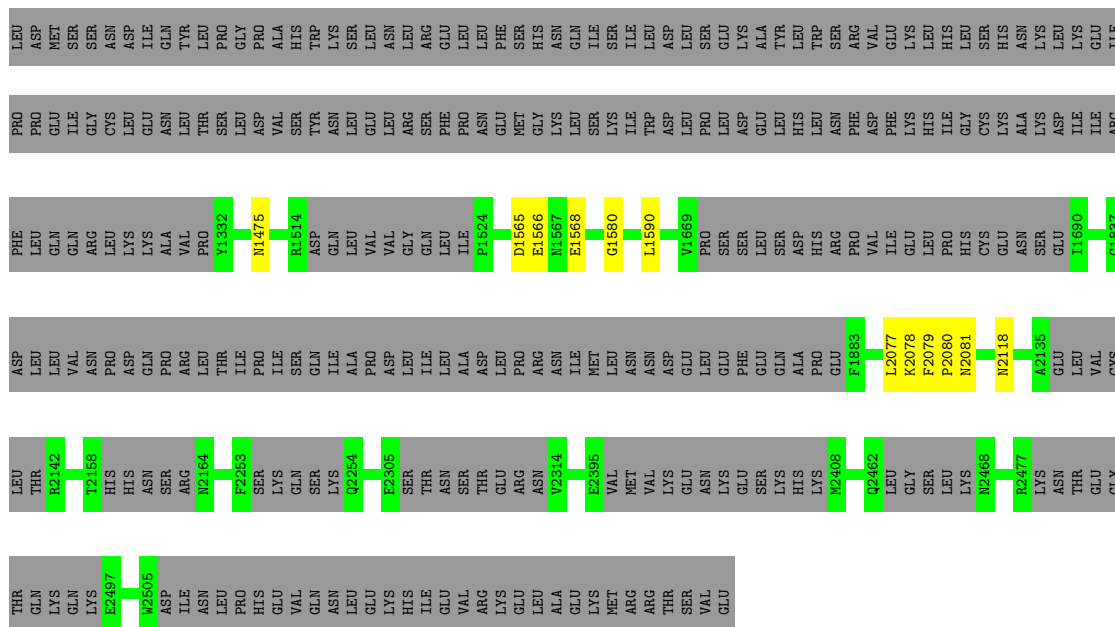




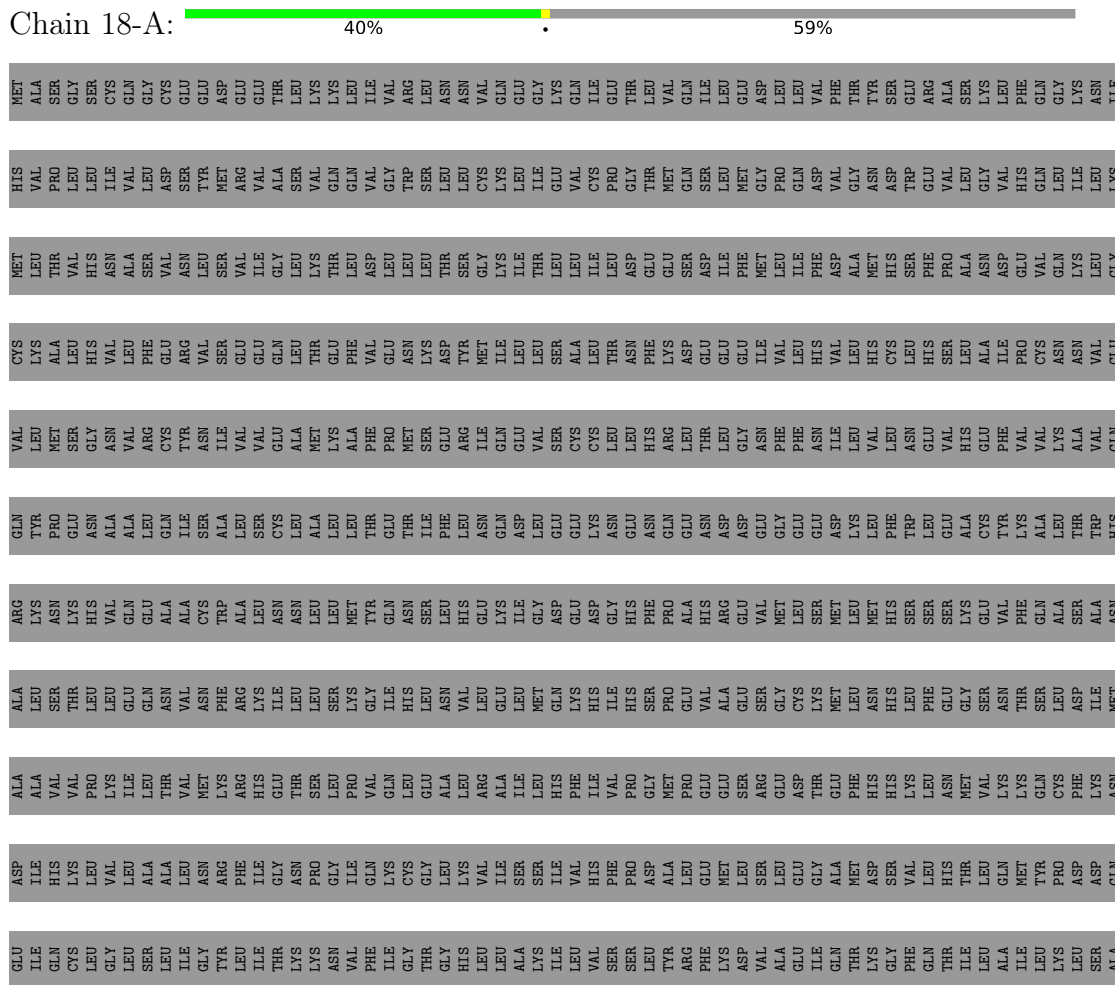
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2



WORLDWIDE
PDB
PROTEIN DATA BANK

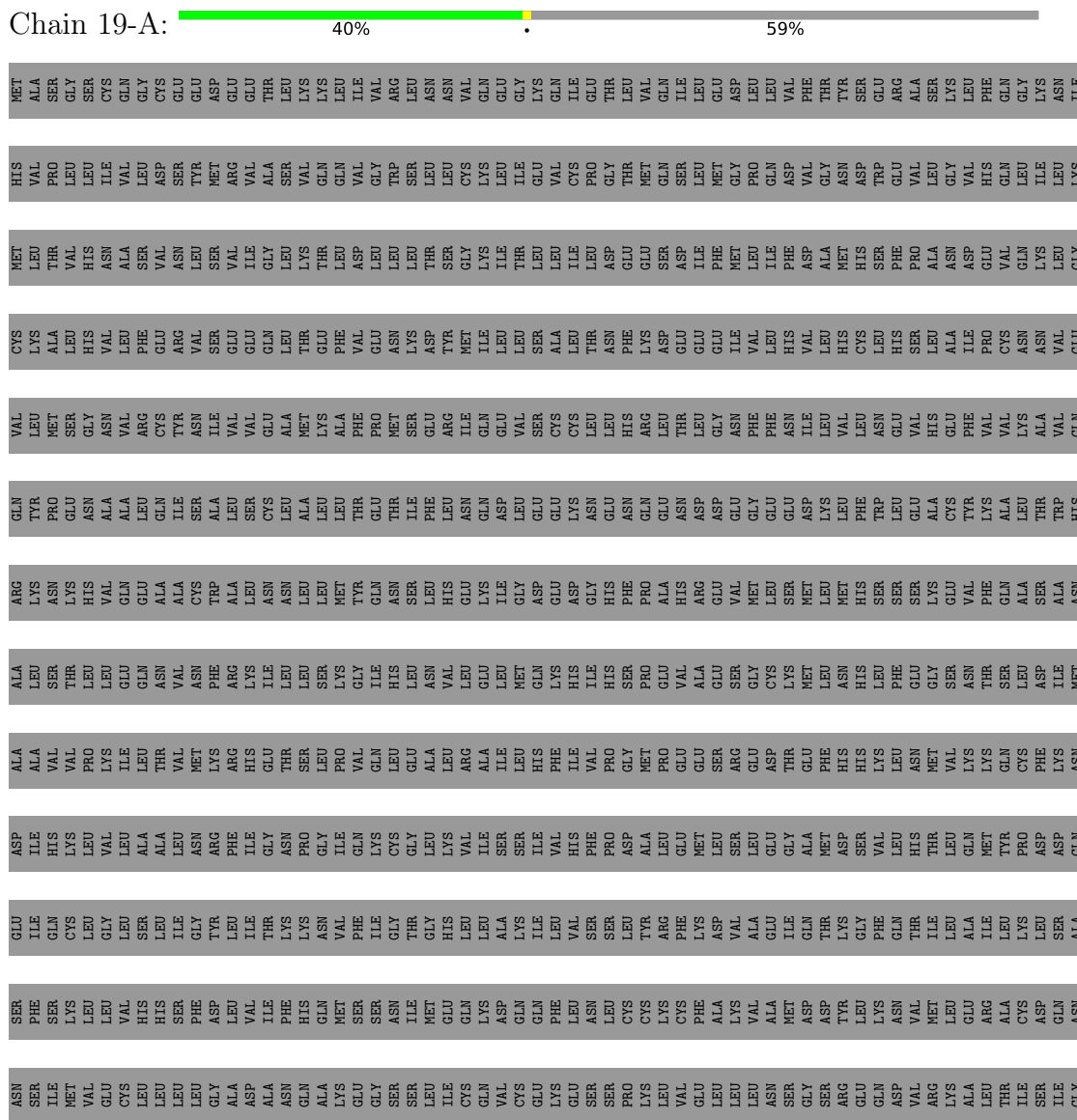


- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2



WORLDWIDE
PDB
PROTEIN DATA BANK

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2



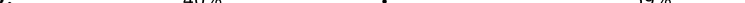


[illegible]

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

[illegible]

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 20-B:  40% 59%

[illegible]



- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

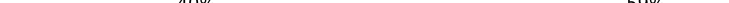
Chain 21-A:

[illegible]



MET	VAL
ARG	MET
ARG	VAL
THR	LYS
SER	GLU
VAL	ASN
GLU	LYS
	GLU
	SER
	LYS
	HIS
	LYS
	E2408
	G2462
	LEU
	GLY
	SER
	LEU
	LYS
	E2468
	E2477
	LYS
	ASN
	THR
	GLU
	GLY
	THR
	GLN
	LYS
	GLN
	LYS
	E2497
	E2505
	ASP
	ILE
	ASN
	LEU
	PRO
	HIS
	GLU
	VAL
	GLN
	ASN
	LEU
	GLU
	LYS
	HIS
	ILE
	GLU
	VAL
	ARG
	LYS
	GLU
	LEU
	ALA
	GLU
	LYS

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 22-A:  40% 59%

[illegible]

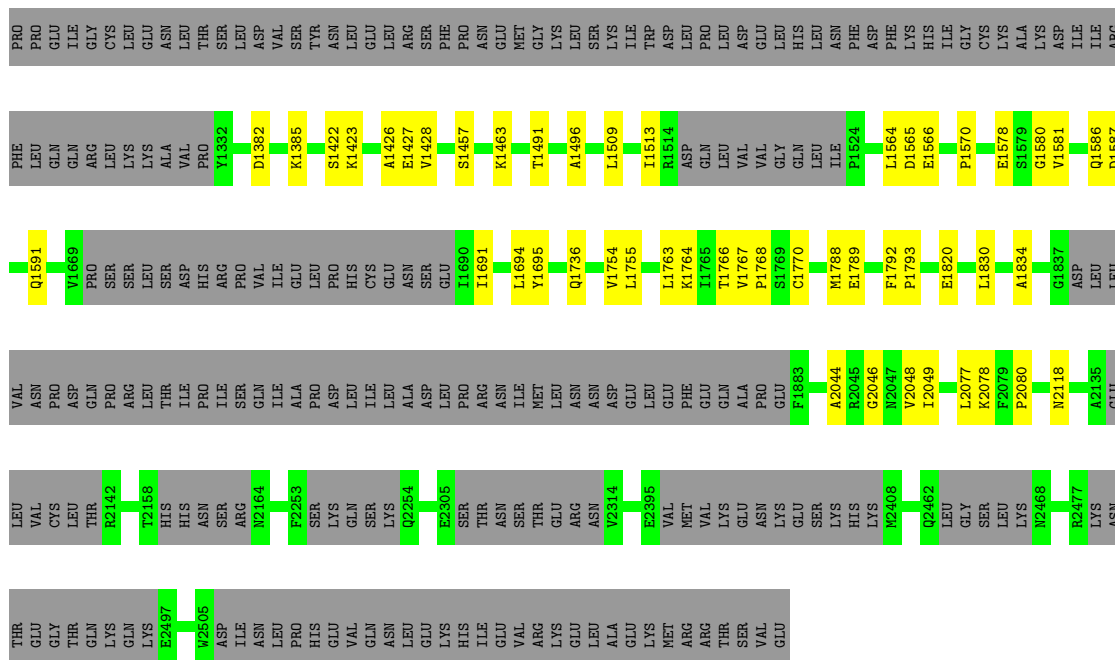
Chain 23-A: 41% 59%



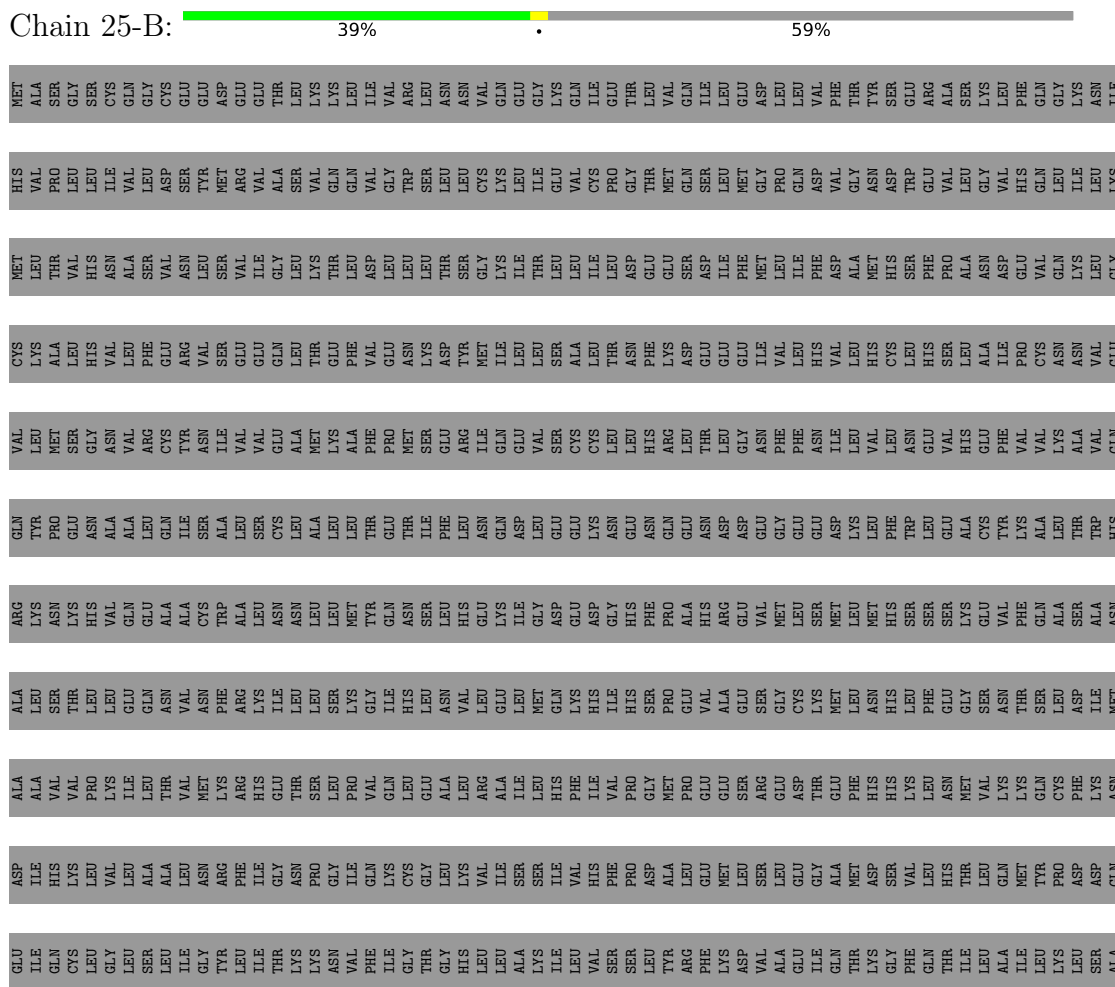








- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2













VAL GLU																												
	ASN LYS GLU SER LYS HIS LYS M2408 Q2462 LEU GLY SER LEU LYS M2468 R2477 LYS ASN THR GLU GLY THR GLN LYS GLN LYS E2497 Q2505 ASP ILE ASN LEU PRO HIS GLU VAL GLN ASN LEU GLY LYS HIS ILE GLU VAL ARG LYS GLU LEU ALA LYS MET ARG THR SER	PHE GLU GLN ALA PRO GLU F1983 R1918 L2077 K2078 F2079 P2080 E2117 N2118 A2135 GLU LEU VAL CYS LEU THR R2142 HIS HIS ASN SER ARG M2164 F2253 SER LYS GLN SER LYS Q2254 E2305 SER THR ASN SER THR ASN GLU ARG ASN V2314 E2395 VAL MET VAL LYS GLU																										

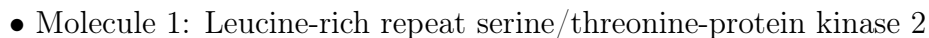
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 28-B:

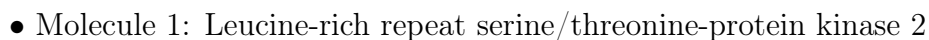
40%59%

ARG	MET	VAL	ILE	GLN	ILE	GLY	LEU	LEU	LEU	ASN	SER	PHE	GLY	ILE	ASP	ILE	ALA	ALA	LEU	LEU	ARG	LYS	THR	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY
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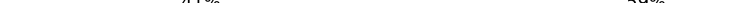


[illegible]



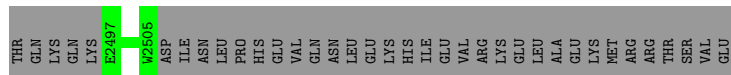
[illegible]

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

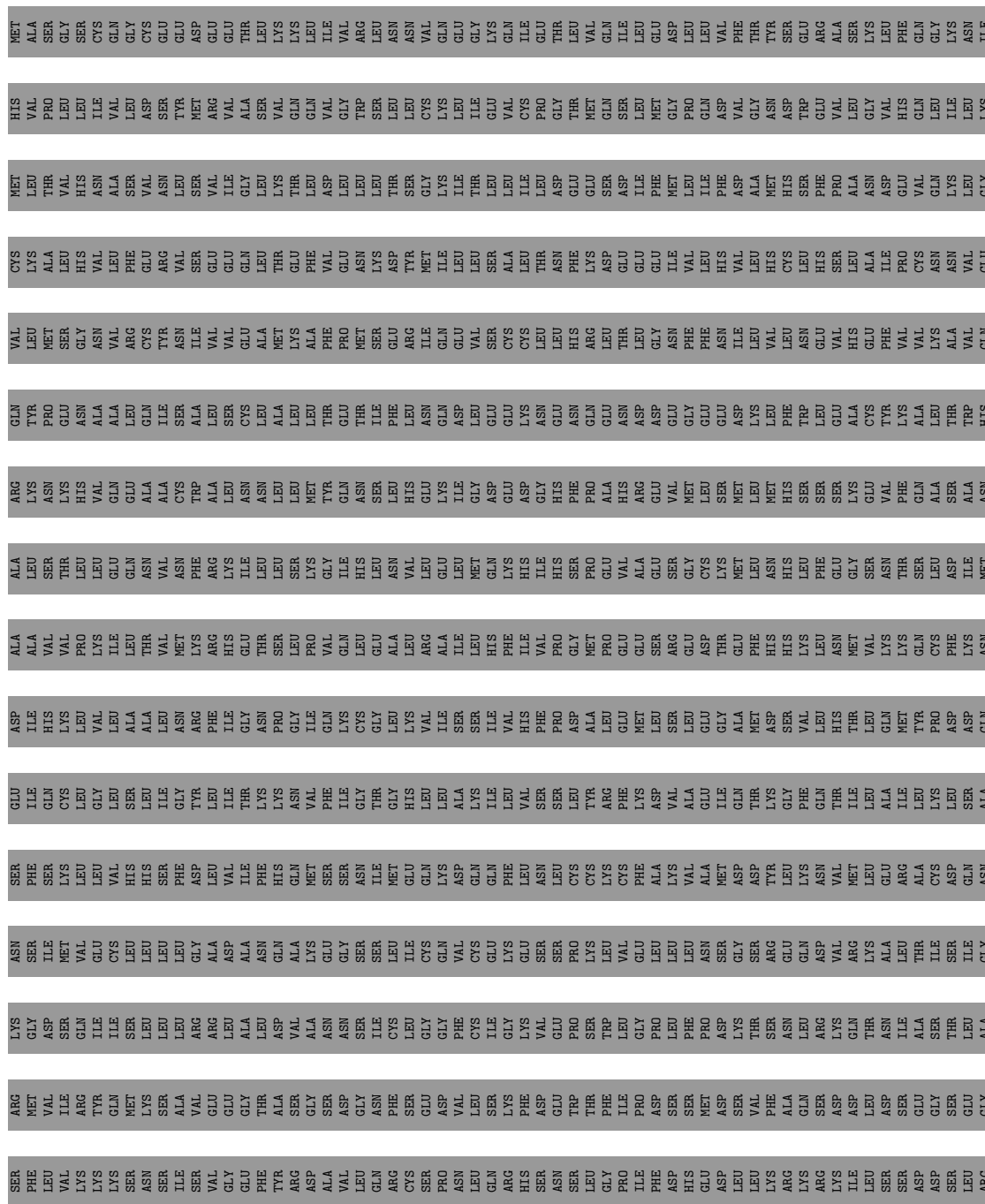
Chain 31-A:  41% 59%

[illegible]





- Chain 31-B: 41% 59%













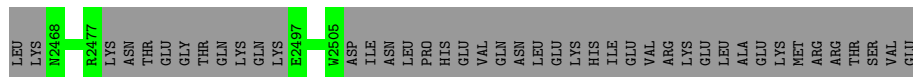
Q2462	LEU	VAL	PHE	PRO
LEU	GLU	ILE	LEU	PRO
GLY	PHE	GLN	GLN	ILE
SER	GLN	LEU	ARG	GLY
LEU	ALA	HIS	LEU	CYS
LYS	PRO	CYS	LYS	LEU
M2468	GLU	GLY	GLU	GLU
F1883	F1883	ASN	ALA	ASN
A2135	A2135	SER	VAL	LEU
GLU	GLU	GLU	PRO	THR
LEU	LEU	I1690	Y1332	THR
VAL	VAL			ASP
CYS	P1793		S1444	LEU
LEU	G1794			ASP
THR	L1795		R1477	VAL
R2142	A1812		G1478	SER
T2158	L1813		F1479	TVR
HIS	Q1823		P1480	ASN
ASN	G1837		R1483	GLU
SER	ASP			LEU
ARG	LEU		R1514	ARG
N2164	LEU		GLN	SER
F2253	VAL		ASP	PHE
SER	ASN		LEU	PRO
LYS	PRO		VAL	ASN
GLN	GLN		VAL	GLU
VAL	ASP		GLY	GLY
GLN	GLN		LEU	LYS
ASN	PRO		ILE	LEU
Q2254	ARG		P1524	SER
GLU	LEU			LYS
LYS	THR		D1565	ILE
HIS	ILE			TRP
SER	PRO			ASP
THR	ILE			LEU
F2305	PRO		Q1586	PRO
ASN	ILE		D1587	LEU
VAL	SER		P1588	LEU
ARG	GLN		A1589	ASP
LYS	ILE		L1590	GLU
GLU	ALA		Q1591	LEU
ASN	PRO			HIS
V2314	ASP		G1617	LEU
F2395	LEU			ASN
VAL	ALA		P1622	PHE
MET	ASP		K1623	ASP
ARG	ASP		Q1648	PHE
THR	LEU			LYS
GLU	PRO		V1669	ILE
VAL	ARG			HIS
SER	ASN		PRO	GLY
VAL	ASN		SER	CYS
GLU	ILE		SER	LYS
LYS	MET		LEU	ALA
THR	LEU			LYS
THR	PRO			LYS
SER	ARG			ASP
VAL	ASN			ASP
GLU	ASN			ILE
LYS	HIS			ASP
M2408	LYS			TRP
	GLN			ARG

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

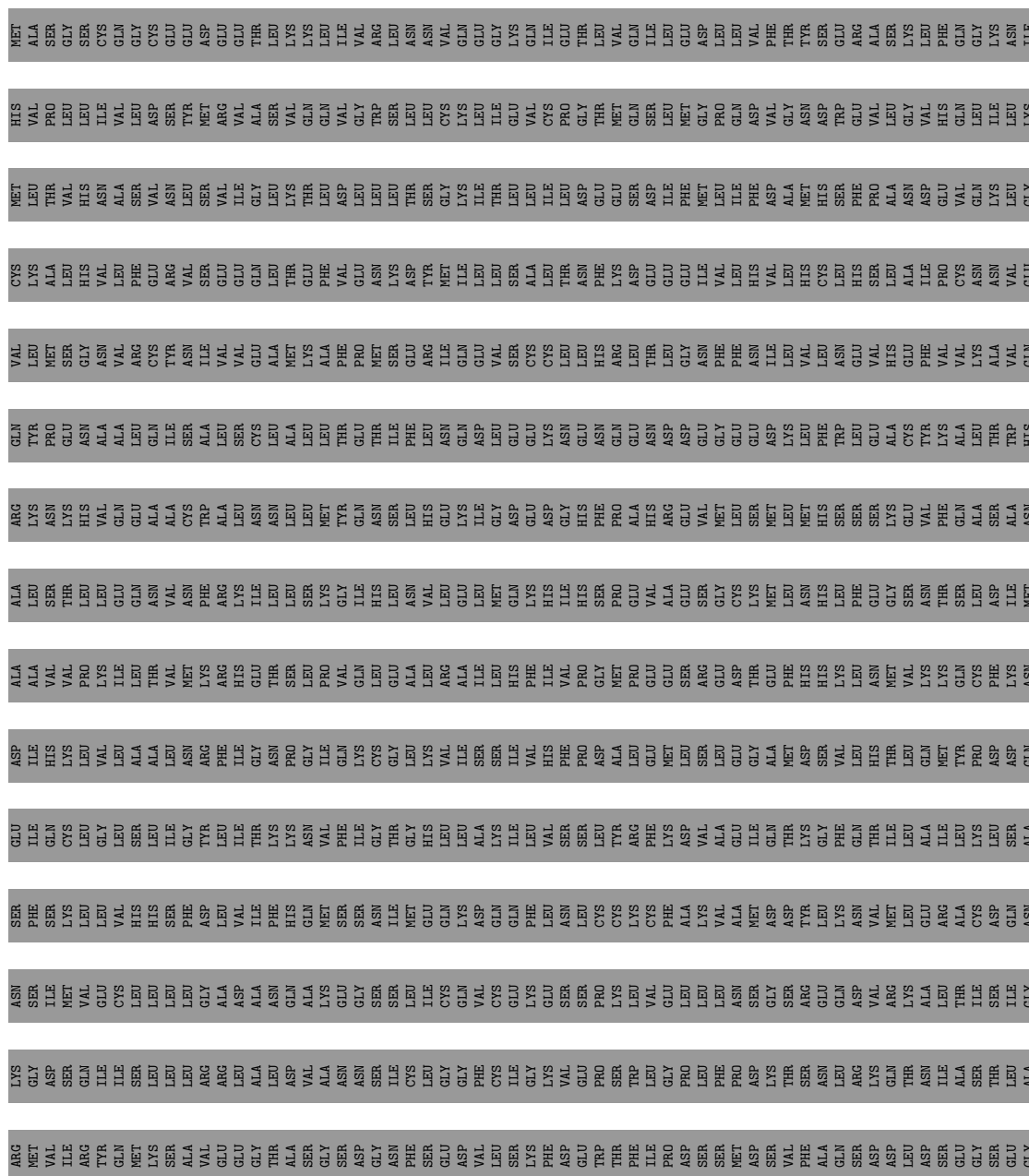
Chain 34-A: 40% 59%

[illegible]



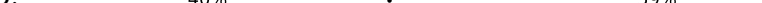


Chain 35-A: 40% 59%



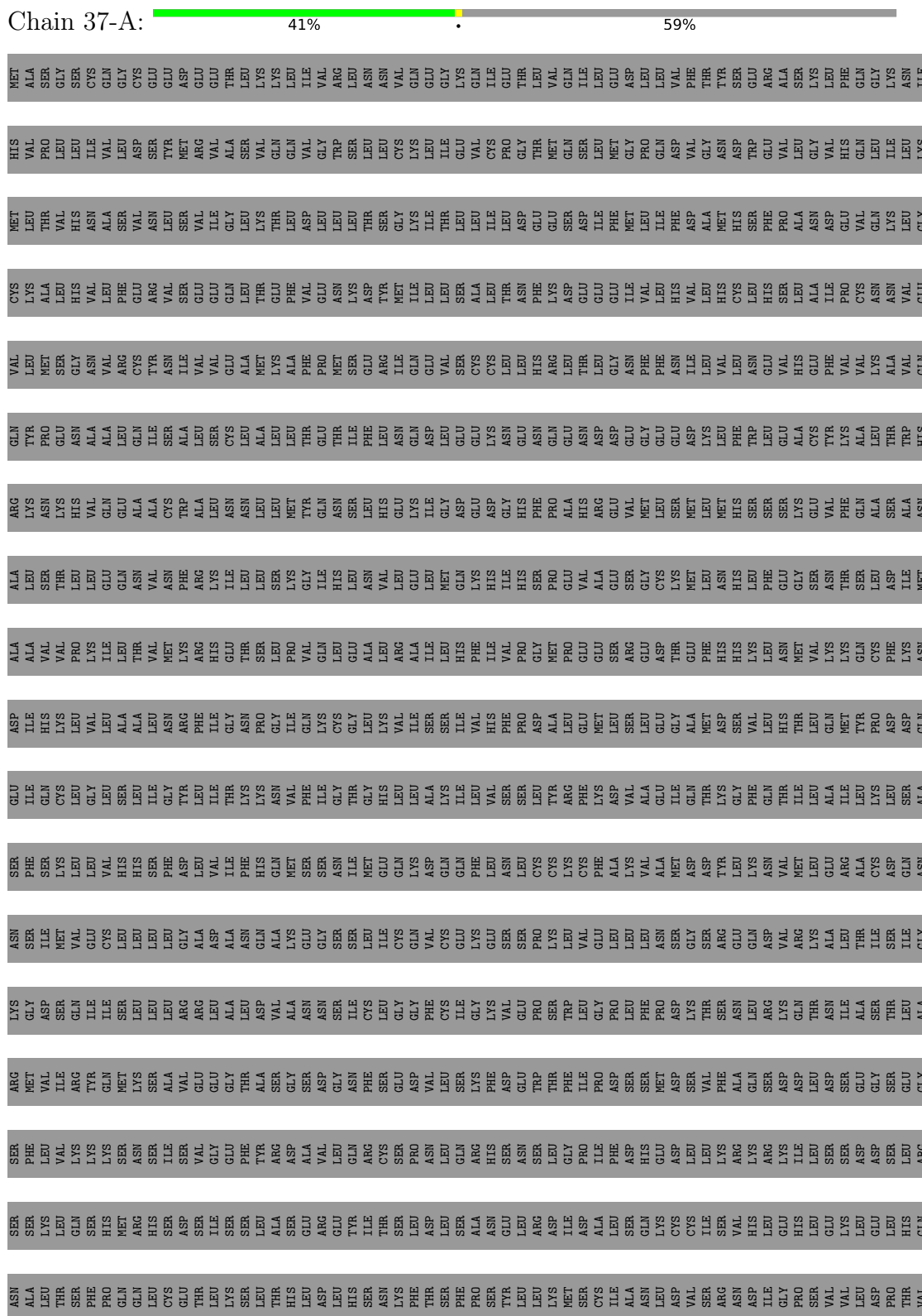


- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 36-B:  40% 59%

ARG	GLN	VAL	CYS	MET	HIS	MET	ALA
LYS	TYR	LEU	LYS	VAL	VAL	VAL	ALA
ASN	PRO	MET	LEU	THR	VAL	THR	PRO
LYS	GLN	SER	GLY	HIS	LEU	GLY	GLY
VAL	ALA	ASN	VAL	ASN	ILE	ASN	CYS
GLN	LEU	VAL	VAL	ALA	VAL	GLN	GLN
GLU	ALA	ARG	PHE	SER	LEU	SER	CYS
ALA	CYS	GLN	GLU	VAL	VAL	GLY	GLY
ILE	ILE	TYR	ARG	ASN	SER	ASN	ASN
CYS	SER	ASN	VAL	LEU	TYR	LEU	GLU
TRP	ALA	ILE	SER	SER	MET	MET	ASP
ALA	LEU	VAL	GLU	VAL	ARG	GLU	GLU
ASN	SER	VAL	GLN	ILE	VAL	VAL	GLU
ASN	CYS	GLU	GLN	GLY	ALA	THR	THR
ASN	LEU	ALA	LEU	LEU	SER	SER	LYS
LEU	ALA	MET	THR	THR	VAL	VAL	LYS
LEU	LYS	GLU	GLU	THR	GLN	GLN	LYS
MET	ALA	PHE	PHE	LEU	GLN	ILE	ILE
THR	THR	PHE	VAL	GLY	VAL	VAL	ILE
GLN	GLU	PRO	GLU	LEU	GLY	VAL	VAL
ASN	THR	MET	ASN	LEU	TRP	VAL	ARG
ILE	ILE	SER	LYS	LEU	SER	VAL	LYS
LEU	PHE	GLU	ASP	THR	LEU	GLY	LYS
LEU	LEU	ARG	ASP	THR	LEU	ASN	ASN
HIS	LEU	ARG	TYR	GLY	LEU	CYS	ASN
GLU	ASN	ILE	MET	GLY	SER	VAL	VAL
LYS	GLN	GLN	ILE	ILE	LYS	GLN	GLN
ILE	ILE	GLU	LEU	ILE	ILE	GLU	GLU
GLY	LEU	VAL	LEU	THR	ILE	GLY	GLY
ASP	GLU	SER	SER	LEU	GLU	VAL	LYS
GLU	GLU	CYS	ALA	LEU	VAL	GLN	GLN
ASP	LYS	CYS	LEU	ILE	CYS	ILE	ILE
GLY	ASN	LEU	GLU	LEU	PRO	GLU	GLU
HIS	GLU	LEU	ASN	LEU	GLY	THR	THR
PHE	ASN	HIS	PHE	GLU	THR	LEU	LEU
PRO	GLN	ARG	LYS	GLU	MET	VAL	VAL
ALA	GLU	ALA	ASP	SER	GLN	GLN	GLN
HIS	ASN	THR	GLU	ASN	SER	ILE	ILE
ARG	ASP	LEU	GLU	ASP	LEU	LEU	LEU
GLU	ASP	GLY	GLU	PHE	MET	GLY	GLU
VAL	GLU	ASN	ILE	MET	GLY	ASP	LEU
MET	GLY	PHE	VAL	LEU	PRO	PRO	LEU
LEU	GLU	PHE	ASN	ILE	GLN	VAL	LEU
SER	GLU	ASN	HIS	MET	ASN	ASP	VAL
GLU	THR	ASN	CYS	HIS	ASP	TRP	GLY
SER	LEU	GLU	HIS	SER	TRP	GLU	ARG
SER	GLU	VAL	SER	PHE	GLU	VAL	ALA
LYS	ALA	HIS	LEU	ALA	LEU	SER	SER
GLU	CYS	GLU	ALA	ASN	GLY	LYS	LYS
VAL	TYR	PHE	ILE	ASP	VAL	VAL	LEU
PHE	LYS	VAL	PRO	ASP	HIS	PHI	PHI
GLN	ALA	VAL	CYS	VAL	GLN	GLN	GLY
ALA	LEU	LYS	ASN	GLN	LEU	ILE	LYS
SER	THR	ALA	ASN	LYS	ILE	LYS	LYS
THR	THR	ALA	VAL	THR	LEU	LEU	ASN
ASP	TRP	VAL	CTR	GLY	LYS	CTR	CTR





- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2





- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

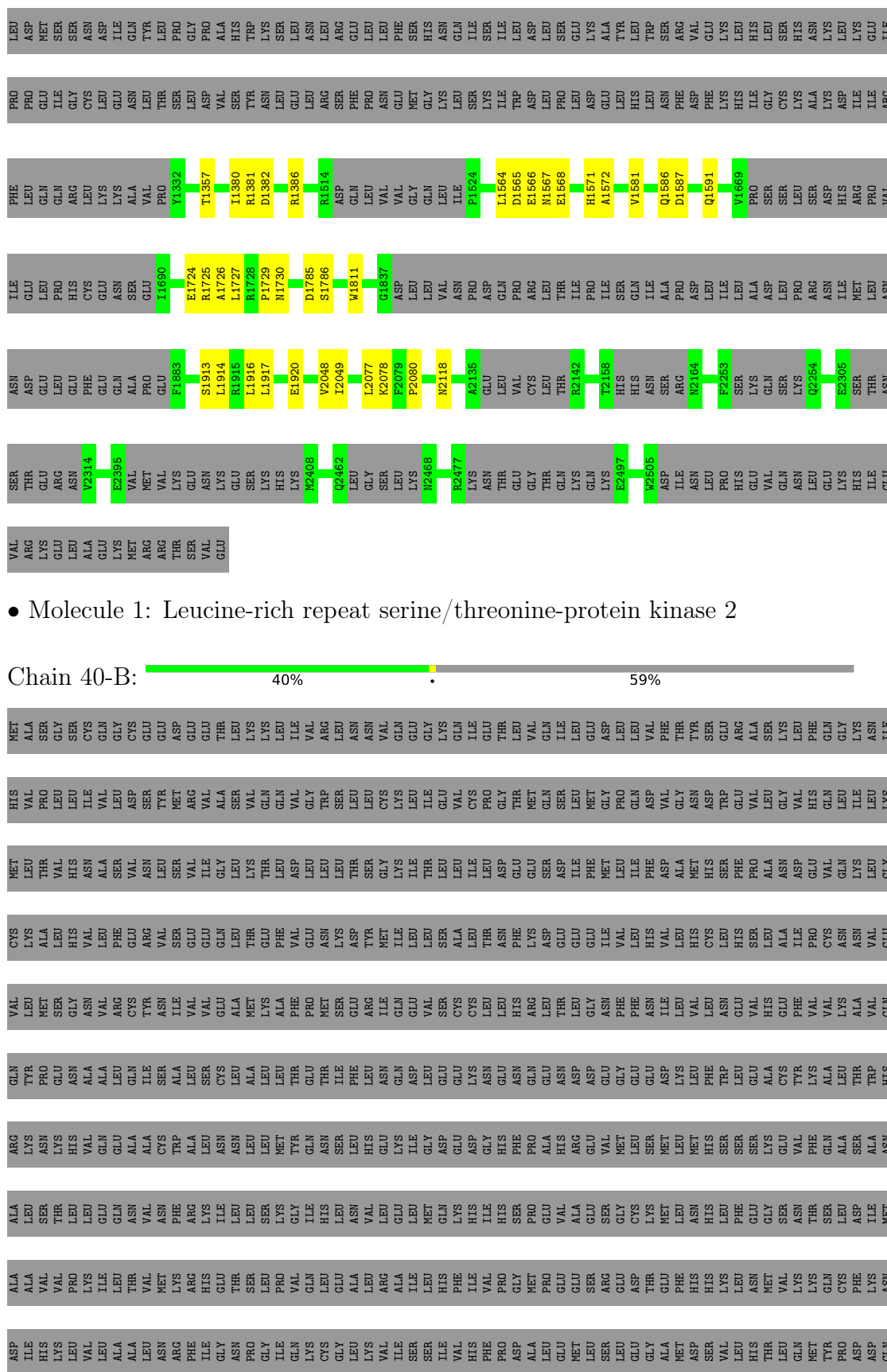


WORLDWIDE
PDB
PROTEIN DATA BANK

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2



WORLDWIDE
PDB
PROTEIN DATA BANK

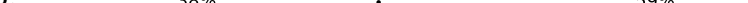




WORLDWIDE
PDB
PROTEIN DATA BANK

THR	LYS	GLN		L1590	PHE	PRO	LEU
	GLU	ALA	T1806		LEU	PRO	ASP
SER	ASN		L1807	P1622	GLN	GLU	MET
	LYS		L1808		GLN	ILE	SER
GLU		F1883	L1809	V1669	ARG	GLY	SER
	SER		K1810		LEU	CYS	ASN
LYS		G2028	W1811		LEU	LEU	ASP
	HIS				LYS	LEU	ILE
M2408	LYS	R2044	G1819		LYS	GLU	GLN
		G2045			VAL	LEU	ASN
G2482		N2047	L1827		ASP	THR	TYR
	LEU				PRO	THR	LEU
GLY	SER	L2077	L1830		H1332	SER	PRO
	LEU	K2078				LEU	GLY
LEU	F2079	A1834	K1833	VAL	S1422	ASP	PRO
	LYS	P2080	E1835	ILE	K1423	VAL	ALA
W2468			E1836	GLU	Q1424	SER	HIS
			E1837	LEU	Q1425	TYR	TRP
R2477	LYS	N2118		PRO	A1426	ASN	LYS
	ASN			HIS	E1427	LEU	SER
THR	GLN	A2135	LEU	CYS	V1428	GLU	LEU
	THR	LEU	LEU	GLY	D1429	LEU	ASN
GLY	VAL	VAL	VAL	ASN	F1436	ARG	LEU
	CYS	VAL	ASN	SER		SER	ARG
THR	LEU	LEU	PRO	GLU	H1453	PHE	GLU
	GLN	R2142	ASP			PRO	LEU
LYS	GLN	T2158	ARG	L1693	V1456	ASN	LEU
			LEU		S1457	MET	PHE
E2487	LYS	HIS	THR	L1708	C1465	LYS	ASN
		HIS	ILE			LEU	GLN
W2505	ASP	ASN	PRO	G1722	K1471	SER	ILE
	ILE	SER	ILE	Y1733	E1472	LYS	SER
ASN	ASN	ARG	SER	W1734	R1477	TRP	LEU
	LEU	W2164	GLN			ASP	ASP
PRO	HIS	F2253	ALA	G1737	R1514	LEU	LEU
	GLU	LYS	SER	L1738	ASP	PRO	SER
VAL	GLN	LYS	LEU	L1740	GLN	LEU	GLU
	GLN	SER	ILE	M1741	VAL	GLY	LYS
ASN	LYS	Q2254	LEU	P1744	VAL	LEU	ALA
	GLU	E2305	ASP	E1745	GLN	HIS	TYR
HIS	THR	E2305	LEU	Y1747	GLY	LEU	TRP
	ILE	THR	ARG		ILE	ASN	ARG
GLU	VAL	ASN	ASN	S1752	P1524	PHE	VAL
	ARG	THR	ILE	L1755	D1565	GLY	LYS
LEU	LYS	GLU	MET		E1566	HIS	LEU
	ASN	ARG	LEU	F1761	W1567	ILE	HIS
ALA	ASN	ASN	ASN		E1568	CYS	SER
	GLU	ASN	ASN		L1569	GLY	SER
GLU	ASN	ASN	ASN		P1570	LYS	HIS
	VAL	ASN	ASP	M1788		ALA	ASN
MET	VAL	E2395	GLU	E1789	H1574	LYS	LYS
	ARG		LEU	E1790	F1575	ASP	LEU
THR	VAL		GLU	W1791		ILE	LYS
	VAL		PHE		G1580	ILE	GLU
ARG	VAL		THR	F1792		ASP	THR
	THR		GLN				GLY

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 41-B:  38% 1% 59%

[illegible]



- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Response	Percentage
Yes, the U.S. is a democracy	40%
No, the U.S. is not a democracy	59%



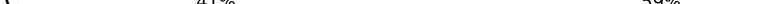
PRO	HIS	F2253	GLN	D1630	PHE	PRO	LEU
GLU	GLY	LEU	ILE	Y1631	LEU	PRO	ASP
VAL	GLN	LYS	ALA	E1632	GLN	GLU	MET
GLN	SER	ASP	PRO	K1643	ARG	ILE	SER
ASN	LYS	LEU	LEU		LEU	CYS	ASN
	Q2254	ILE	ILE	Q1648	LYS	ASP	ASN
GLU		LEU	LEU	Y1649	LYS	GLU	ILE
LEU	E2305	ALA	ASP	F1650	ALA	ASN	GLN
HIS	SER	ASP	ASP	K1651	VAL	LEU	TYR
ILE	THR	LEU	LEU		PRO	THR	LEU
GLU	ASN	PRO	PRO	G1663	Y11332	SER	PRO
VAL	SER	ARG	ARG			LEU	GLY
ARG	THR	ASN	ASN	Y1669	P1377	ASP	PRO
LYS	GLU	ILE	ILE	PRO	I1378	VAL	ALA
GLU	ARG	MET	MET	SER	Q1379	SER	HIS
LEU	ASN	LEU	LEU	SER		TYR	TRP
ALA	V2314	ASN	ASN	LEU	D1382	ASN	LYS
GLU		ASP	ASP	SER	K1383	LEU	SER
LYS	E2395	ASP	ASP	ASP	R1384	GLU	LEU
MET	VAL	GLU	GLU	HIS	K1385	LEU	ASN
ARG	MET	LEU	LEU	ARG	R1396	ARG	LEU
ARG	VAL	GLU	GLU	PRO		SER	ARG
THR	VAL	PHE	PHE	VAL	D1458	PHE	GLU
SER	GLU	GLU	GLU	ILE		PRO	LEU
VAL	ASN	GLN	GLN	LEU	E1493	ASN	LEU
GLU	LYS	ALA	ALA	LEU	S1494	GLU	PHE
	GLU	PRO	PRO	PRO	D1495	MET	SER
	SER	GLU	GLU	HIS		GLY	HIS
LYS	LYS	F1883		CYS	K1499	LYS	ASN
HIS	LYS			GLU	L1500	LEU	GLN
	E2408	S1913		ASN	R1501	SER	ILE
GLU				SER	K1502	LYS	GLN
		R1918		GLU		ILE	ILE
						TRP	LEU
	Q2462	V1922		I1690	R1514	ASP	ASP
LEU						LEU	LEU
GLY		M1730			GLN	PRO	SER
SER		R1731			VAL	LEU	GLU
LEU		M1732			VAL	GLU	GLU
LYS		N2047			GLY	ASP	LYS
	N2468	Y1733			VAL	GLY	ALA
		Y2048			GLN	LEU	TYR
		I2049				HIS	LEU
	E2477	A2135			ILE	LEU	TRP
LYS	ASN	GLU	GLU	ASP	P1524	LEU	ASP
THR	THR	LEU	LEU	LEU		PHE	ARG
GLU	GLY	VAL	VAL	LEU	R1550	ASP	VAL
GLY		CYS	CYS	VAL	K1551	PHE	GLU
THR	THR	LEU	LEU	ASN		LYS	LYS
GLN	GLN	R2142		ASN	L1554	HIS	LEU
				PRO		ILE	HIS
GLN	GLN	T2158		GLN	E1566	GLY	LEU
LYS	LYS			PRO		CYS	SER
	E2497	HIS	HIS	ARG	L1569	LYS	HIS
		HIS	ASN	THR		ALA	ASN
	N2505	LEU	ASN	THR	P1588	LYS	LYS
ASP	ASP	ILE	SER	PRO	A1589	ASP	LEU
ILE	ILE	ARG	ARG	PRO	L1590	ILE	LYS
ASN	ASN	N2164		ILE	Q1591	ILE	GLU
LEU	LEU	SER				ASP	GLY

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 43-B: 40% . 59%

[illegible]

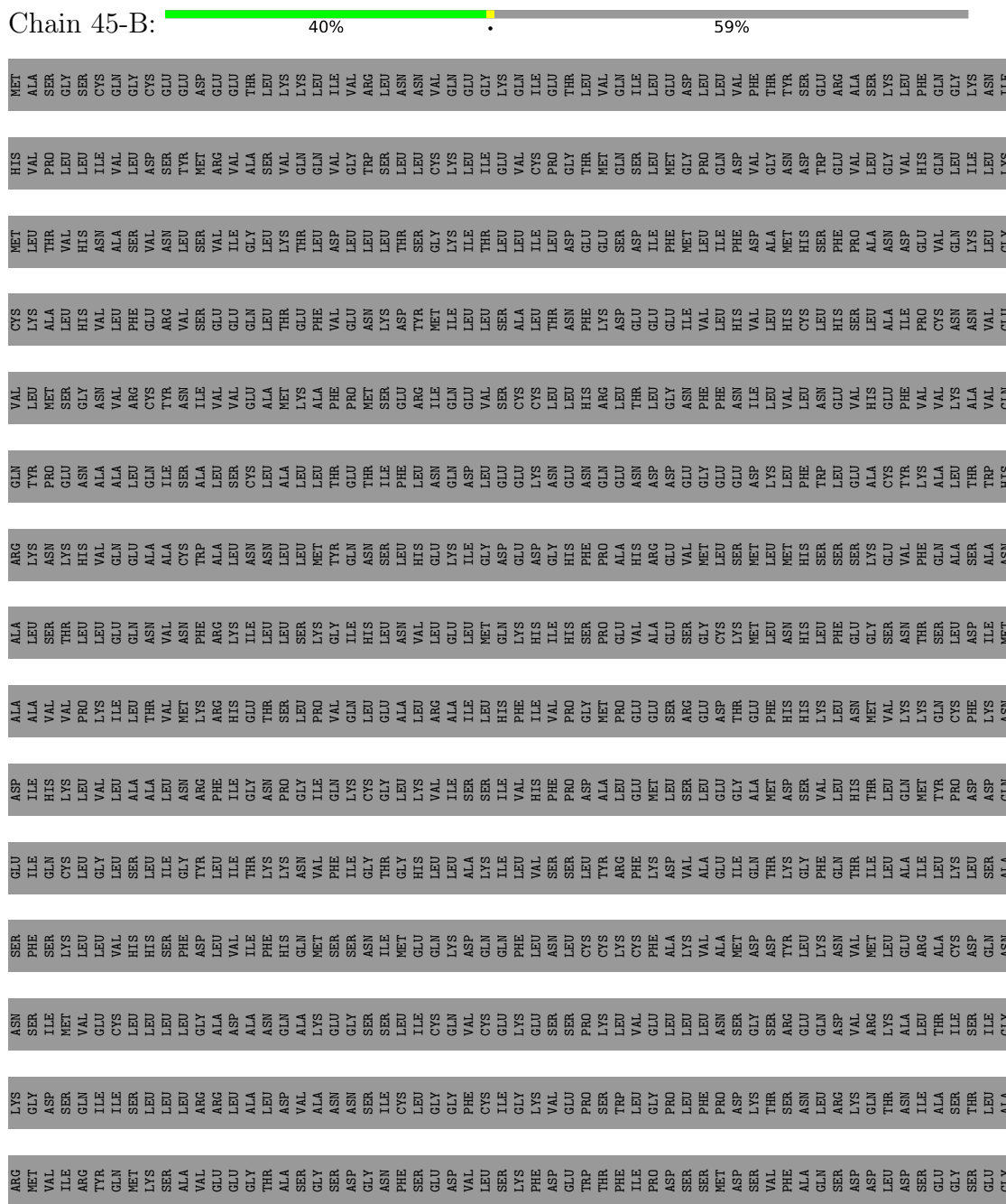
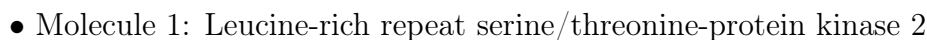
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 44-A:  41% 59%

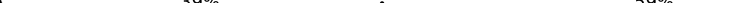








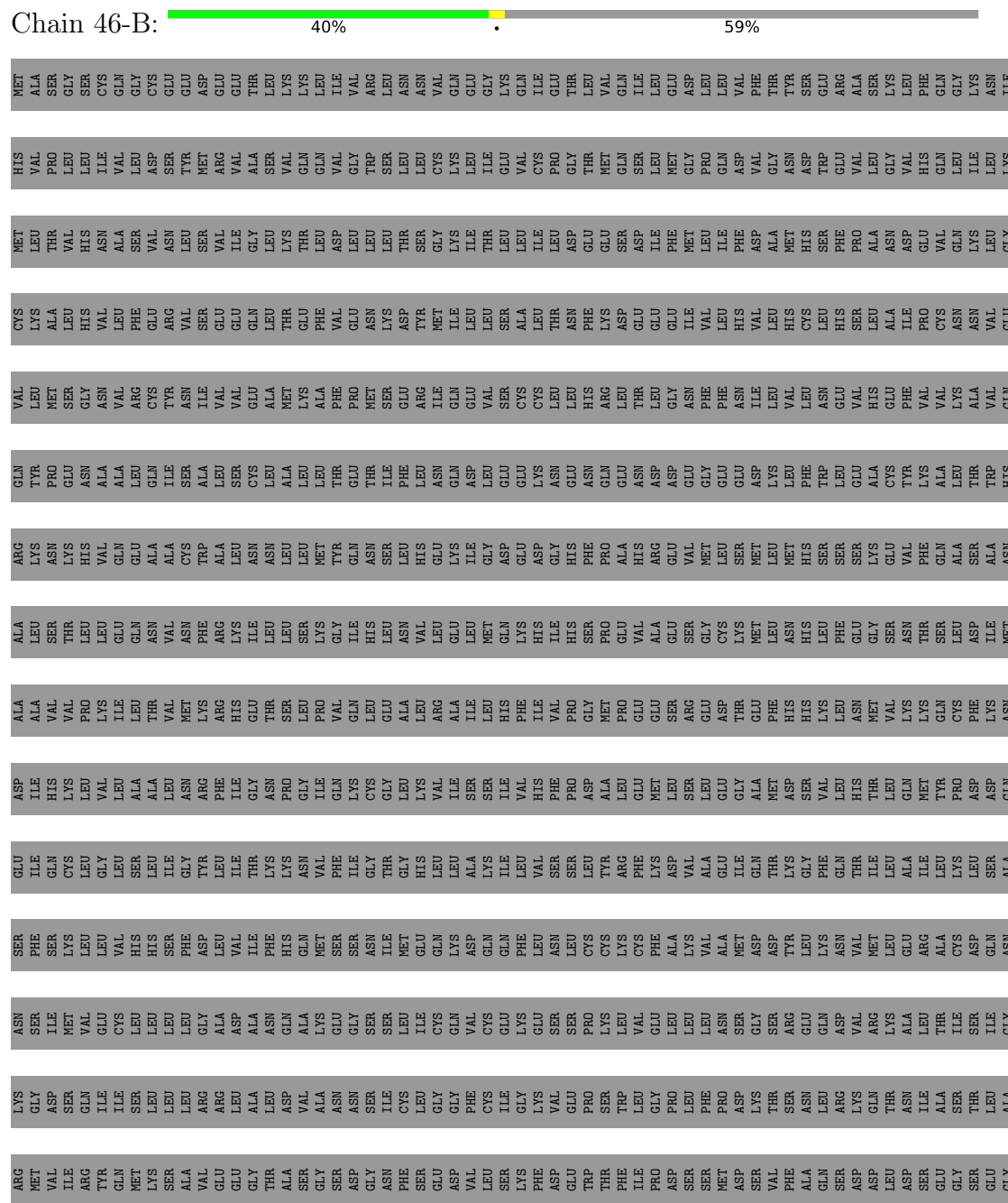
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 46-A:  39% 1% 59%

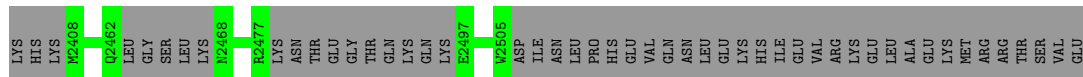
[illegible]



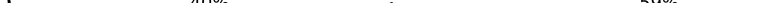
- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

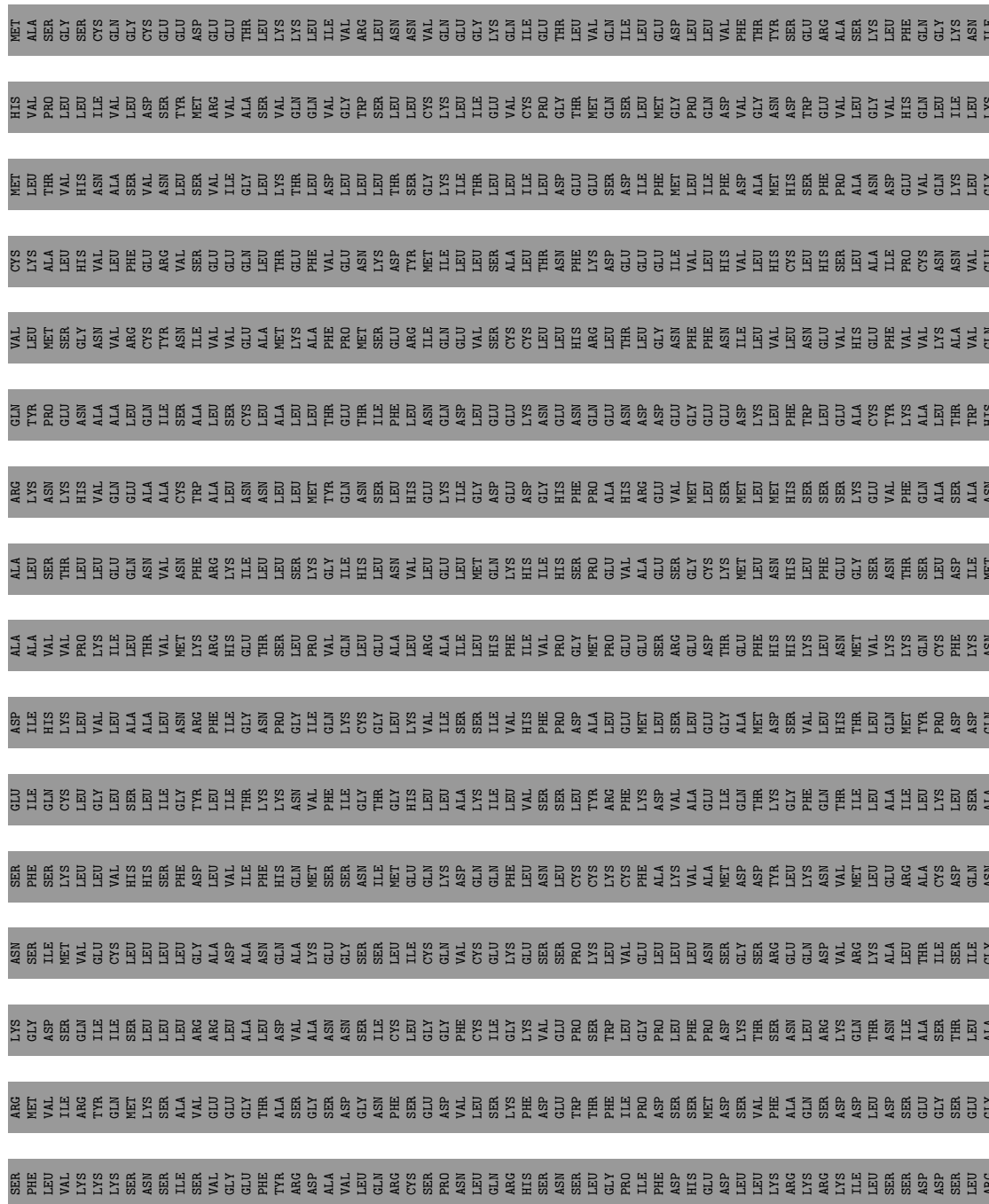






- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Chain 47-B:  40% 59%



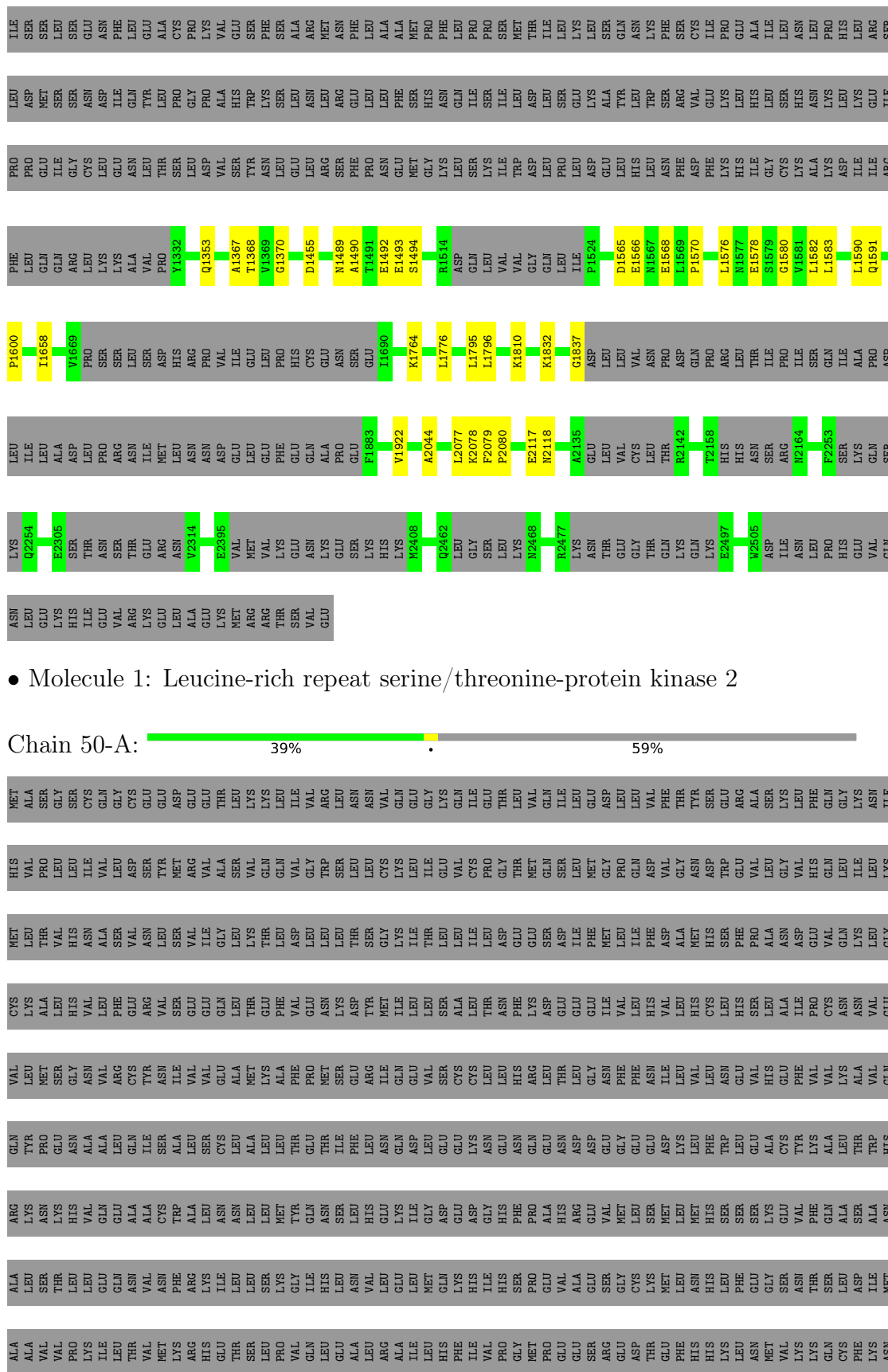






Chain 49-B: 40% 1% 59%





- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Government	Percentage
Current government	39%
Previous government	59%



- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

[illegible]

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Opinion	Percentage
Doing a good job	39%
Doing a bad job	59%

WORLDWIDE
PDB
PROTEIN DATA BANK

[illegible]



WORLDWIDE
PDB
PROTEIN DATA BANK

- Molecule 1: Leucine-rich repeat serine/threonine-protein kinase 2

Response	Percentage
Democracy	39%
Dictatorship	59%

WORLDWIDE
PDB
PROTEIN DATA BANK



4 Experimental information

Property	Value	Source
EM reconstruction method	SUBTOMOGRAM AVERAGING	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of subtomograms used	4307	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI POLARA 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	2	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.209	Depositor
Minimum map value	-0.128	Depositor
Average map value	0.003	Depositor
Map value standard deviation	0.031	Depositor
Recommended contour level	0.05	Depositor
Map size ($\text{\AA}$)	281.6, 281.6, 281.6	wwPDB
Map dimensions	128, 128, 128	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	2.2, 2.2, 2.2	Depositor

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).

There are no protein, RNA or DNA chains available to summarize Z scores of covalent bonds and angles.

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	1-A	1040	0	0	2	0
1	1-B	1040	0	0	2	0
1	2-A	1040	0	0	0	0
1	2-B	1040	0	0	2	0
1	3-A	1040	0	0	7	0
1	3-B	1040	0	0	2	0
1	4-A	1040	0	0	12	0
1	4-B	1040	0	0	7	0
1	5-A	1040	0	0	17	0
1	5-B	1040	0	0	16	0
1	6-A	1040	0	0	4	0
1	6-B	1040	0	0	4	0
1	7-A	1040	0	0	42	0
1	7-B	1040	0	0	42	0
1	8-A	1040	0	0	19	0
1	8-B	1040	0	0	19	0
1	9-A	1040	0	0	6	0
1	9-B	1040	0	0	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	10-A	1040	0	0	20	0
1	10-B	1040	0	0	20	0
1	11-A	1040	0	0	1	0
1	11-B	1040	0	0	0	0
1	12-A	1040	0	0	11	0
1	12-B	1040	0	0	11	0
1	13-A	1040	0	0	18	0
1	13-B	1040	0	0	18	0
1	14-A	1040	0	0	32	0
1	14-B	1040	0	0	32	0
1	15-A	1040	0	0	14	0
1	15-B	1040	0	0	14	0
1	16-A	1040	0	0	31	0
1	16-B	1040	0	0	31	0
1	17-A	1040	0	0	7	0
1	17-B	1040	0	0	7	0
1	18-A	1040	0	0	11	0
1	18-B	1040	0	0	11	0
1	19-A	1040	0	0	34	0
1	19-B	1040	0	0	34	0
1	20-A	1040	0	0	10	0
1	20-B	1040	0	0	10	0
1	21-A	1040	0	0	18	0
1	21-B	1040	0	0	18	0
1	22-A	1040	0	0	13	0
1	22-B	1040	0	0	13	0
1	23-A	1040	0	0	7	0
1	23-B	1040	0	0	7	0
1	24-A	1040	0	0	39	0
1	24-B	1040	0	0	39	0
1	25-A	1040	0	0	42	0
1	25-B	1040	0	0	42	0
1	26-A	1040	0	0	7	0
1	26-B	1040	0	0	7	0
1	27-A	1040	0	0	26	0
1	27-B	1040	0	0	26	0
1	28-A	1040	0	0	16	0
1	28-B	1040	0	0	16	0
1	29-A	1040	0	0	22	0
1	29-B	1040	0	0	22	0
1	30-A	1040	0	0	17	0
1	30-B	1040	0	0	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	31-A	1040	0	0	7	0
1	31-B	1040	0	0	7	0
1	32-A	1040	0	0	12	0
1	32-B	1040	0	0	12	0
1	33-A	1040	0	0	17	0
1	33-B	1040	0	0	17	0
1	34-A	1040	0	0	15	0
1	34-B	1040	0	0	15	0
1	35-A	1040	0	0	13	0
1	35-B	1040	0	0	14	0
1	36-A	1040	0	0	23	0
1	36-B	1040	0	0	23	0
1	37-A	1040	0	0	10	0
1	37-B	1040	0	0	11	0
1	38-A	1040	0	0	49	0
1	38-B	1040	0	0	49	0
1	39-A	1040	0	0	31	0
1	39-B	1040	0	0	31	0
1	40-A	1040	0	0	26	0
1	40-B	1040	0	0	26	0
1	41-A	1040	0	0	62	0
1	41-B	1040	0	0	62	0
1	42-A	1040	0	0	20	0
1	42-B	1040	0	0	21	0
1	43-A	1040	0	0	24	0
1	43-B	1040	0	0	24	0
1	44-A	1040	0	0	4	0
1	44-B	1040	0	0	0	0
1	45-A	1040	0	0	17	0
1	45-B	1040	0	0	17	0
1	46-A	1040	0	0	38	0
1	46-B	1040	0	0	38	0
1	47-A	1040	0	0	13	0
1	47-B	1040	0	0	13	0
1	48-A	1040	0	0	34	0
1	48-B	1040	0	0	34	0
1	49-A	1040	0	0	23	0
1	49-B	1040	0	0	24	0
1	50-A	1040	0	0	50	0
1	50-B	1040	0	0	50	0
1	51-A	1040	0	0	36	0
1	51-B	1040	0	0	36	0

Continued on next page...

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	52-A	1040	0	0	46	0
1	52-B	1040	0	0	46	0
1	53-A	1040	0	0	34	0
1	53-B	1040	0	0	34	0
All	All	110240	0	0	1840	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 17.

The worst 5 of 1840 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:1591:GLN:CA	1:A:2118:ASN:CA	1.75	1.65
1:A:1428:VAL:CA	1:A:1811:TRP:CA	1.75	1.64
1:A:1458:ASP:CA	1:A:1601:LYS:CA	1.77	1.63
1:B:1426:ALA:CA	1:B:1807:LEU:CA	1.76	1.63
1:A:1740:LEU:CA	1:B:1754:VAL:CA	1.76	1.62

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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.

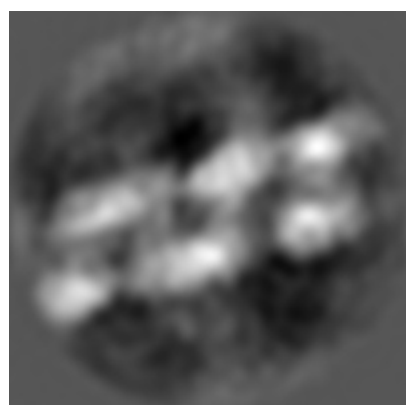
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-20825. These allow visual inspection of the internal detail of the map and identification of artifacts.

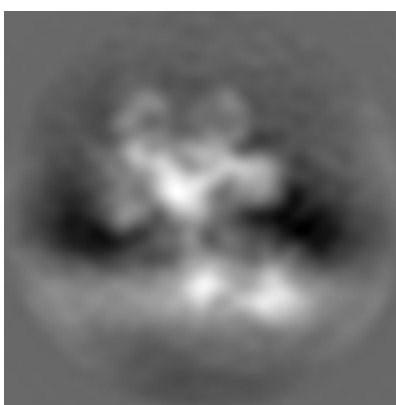
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

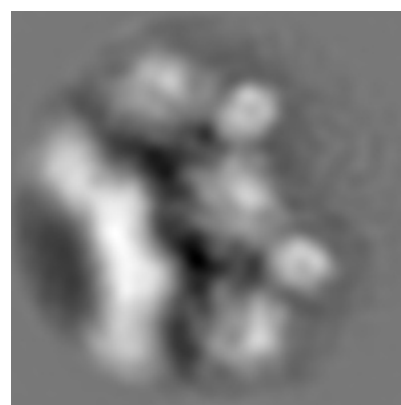
6.1.1 Primary map



X



Y

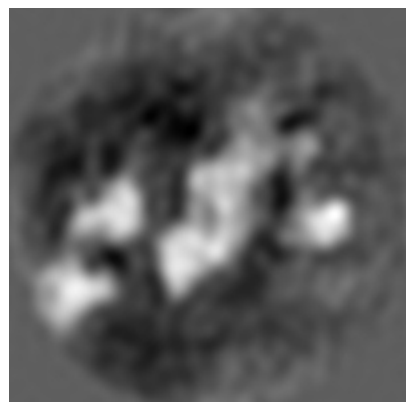


Z

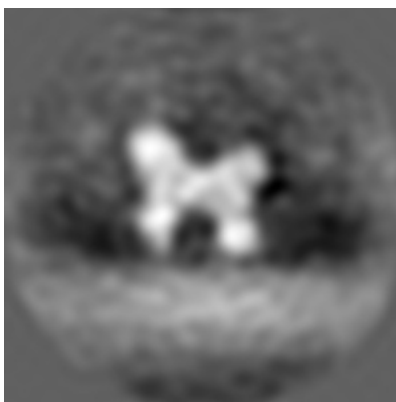
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

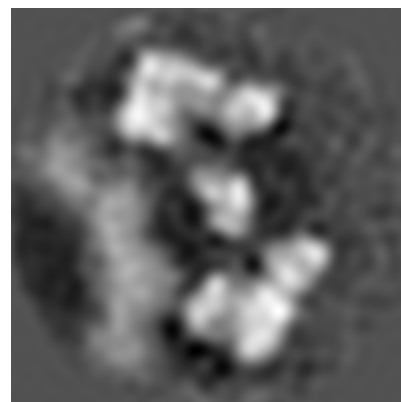
6.2.1 Primary map



X Index: 64



Y Index: 64

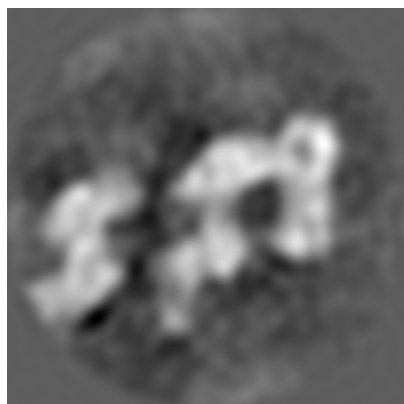


Z Index: 64

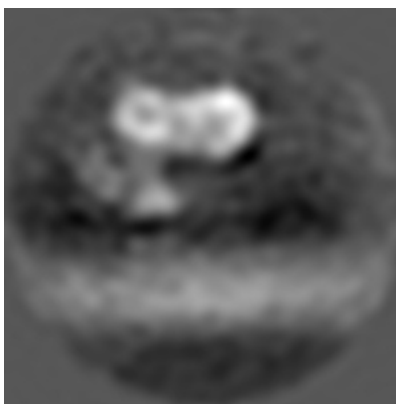
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

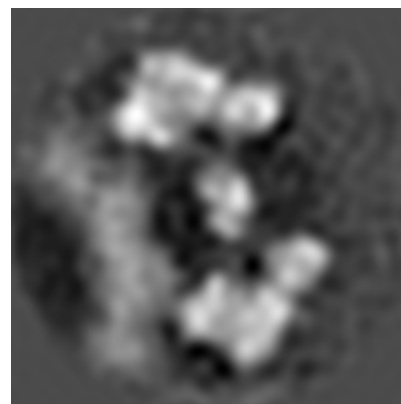
6.3.1 Primary map



X Index: 76



Y Index: 49

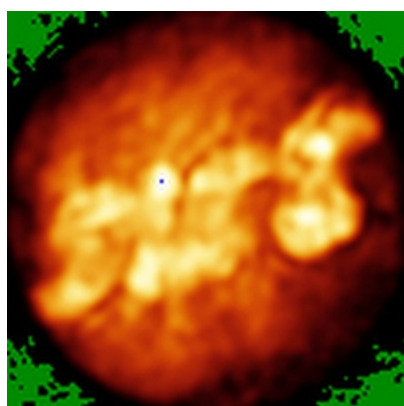


Z Index: 62

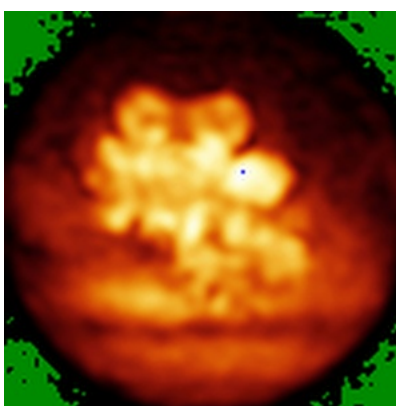
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

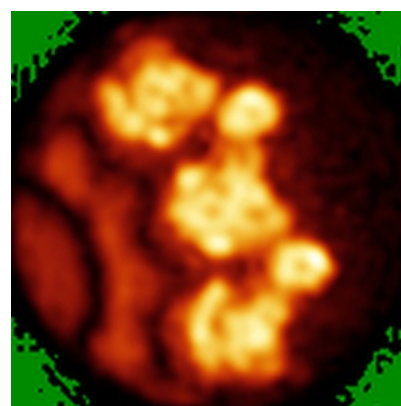
6.4.1 Primary map



X



Y

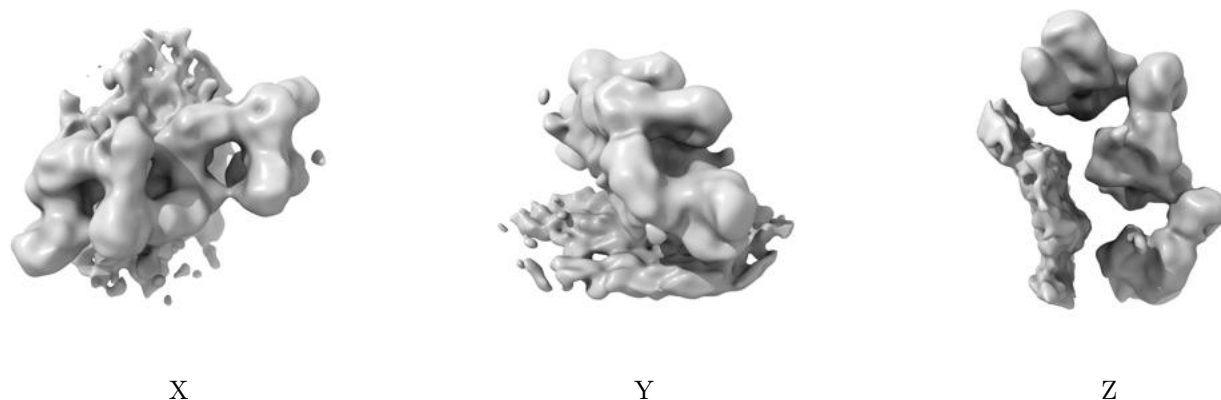


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.05. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

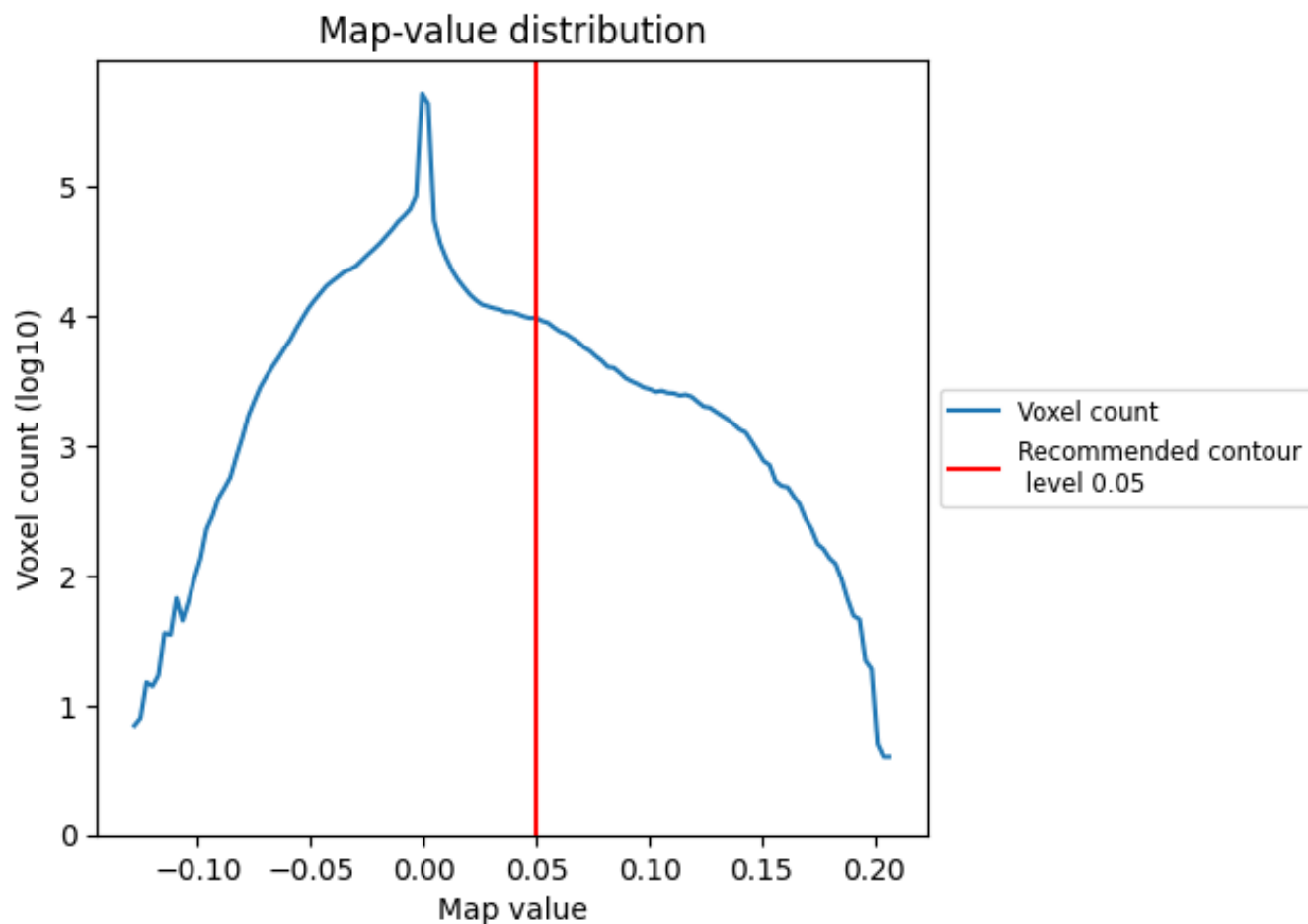
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

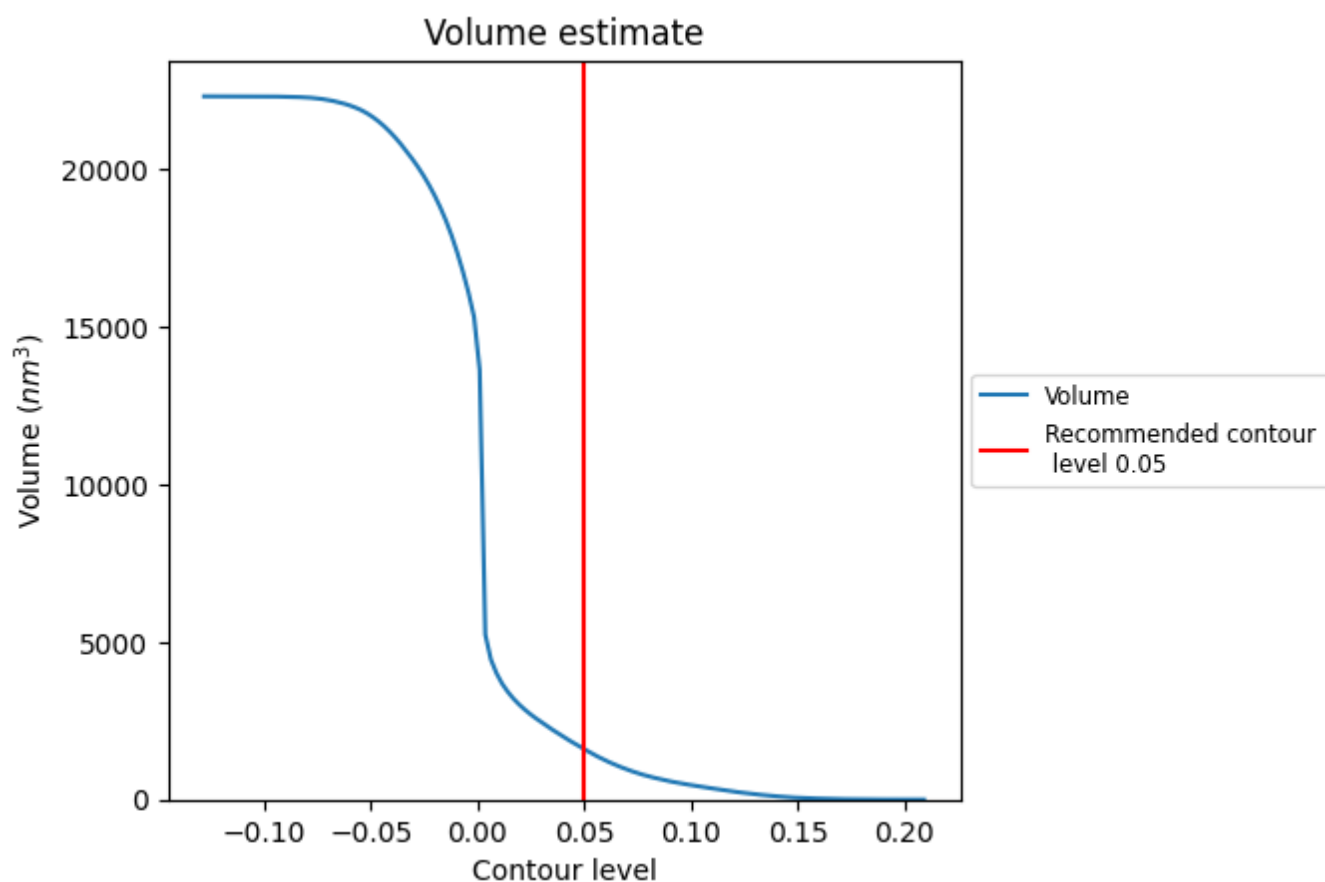
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

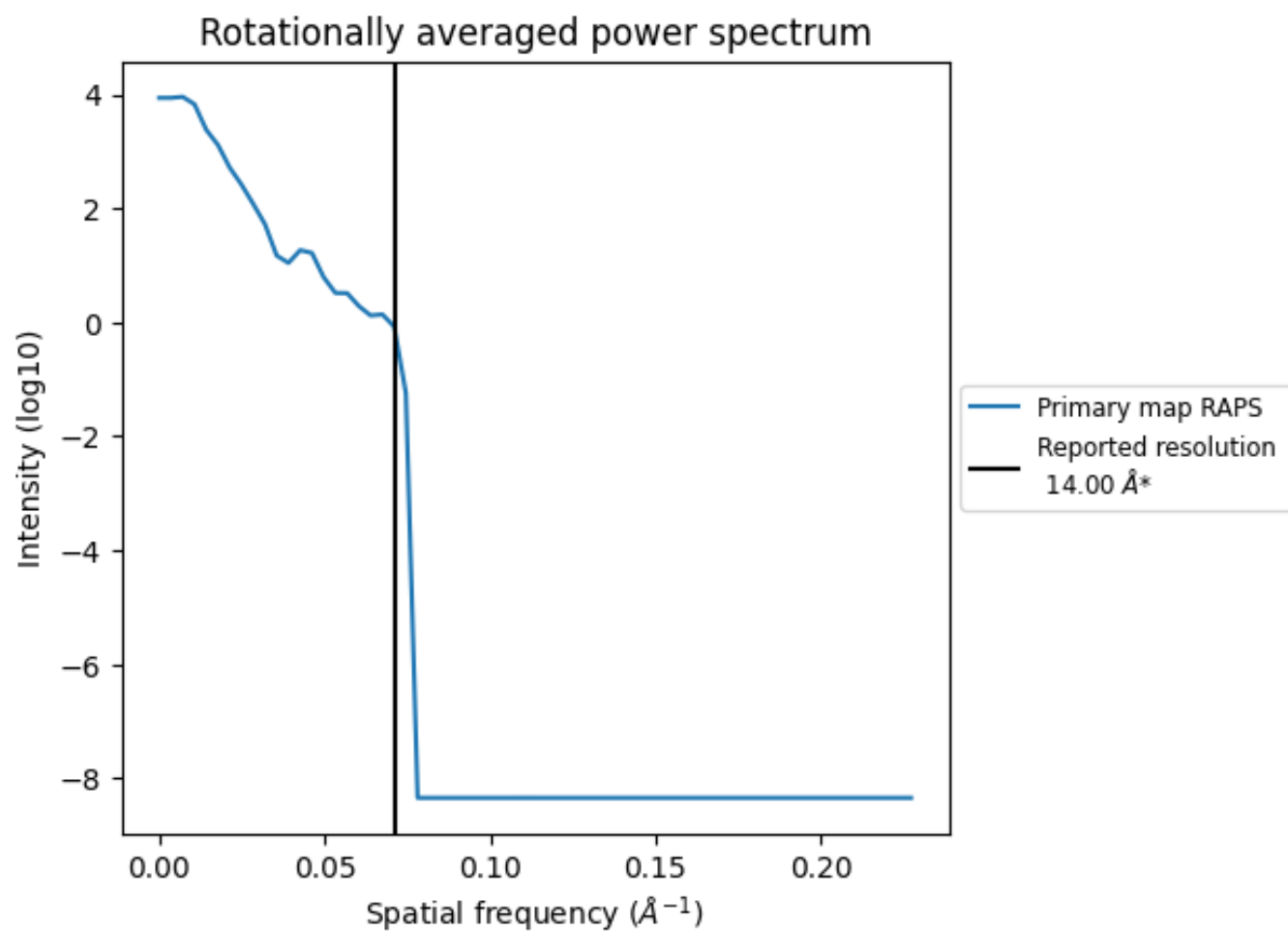
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1604 nm³; this corresponds to an approximate mass of 1449 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ



*Reported resolution corresponds to spatial frequency of 0.071 Å⁻¹

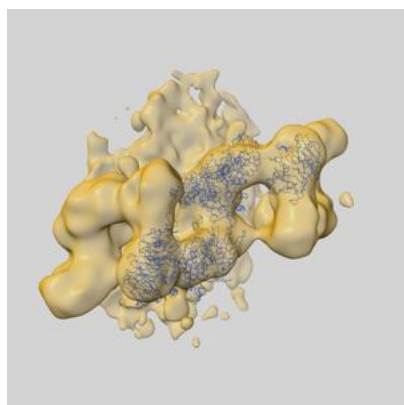
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

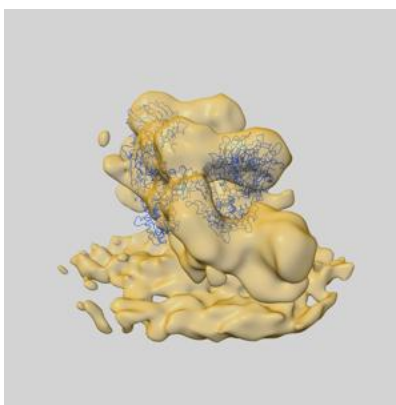
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-20825 and PDB model 6XR4. Per-residue inclusion information can be found in section [3](#) on page [13](#).

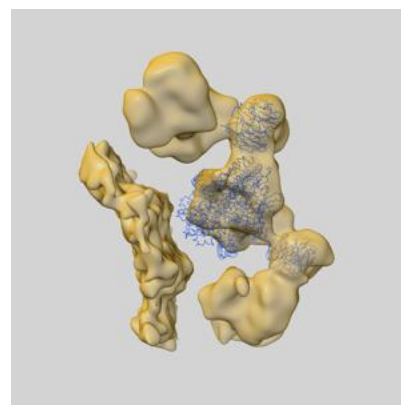
9.1 Map-model overlay [i](#)



X



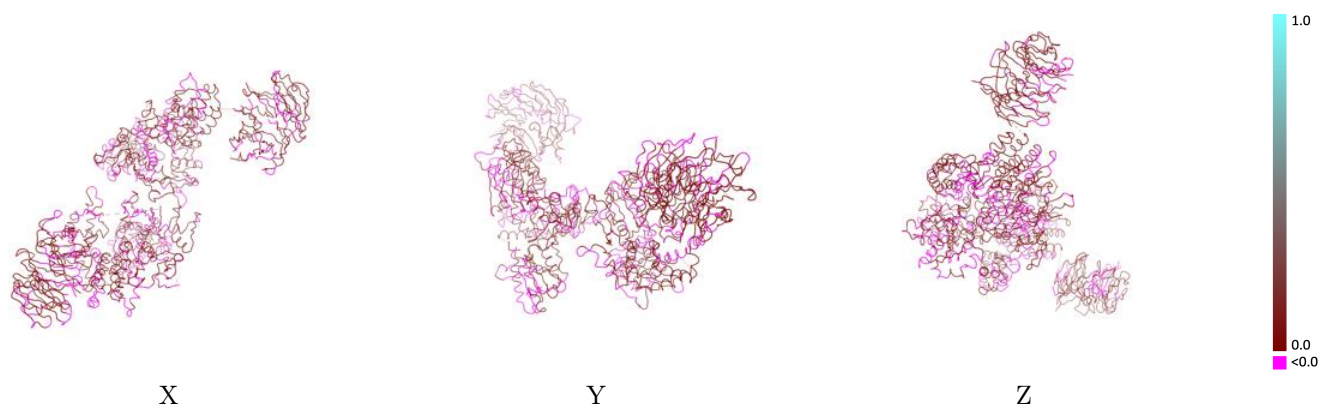
Y



Z

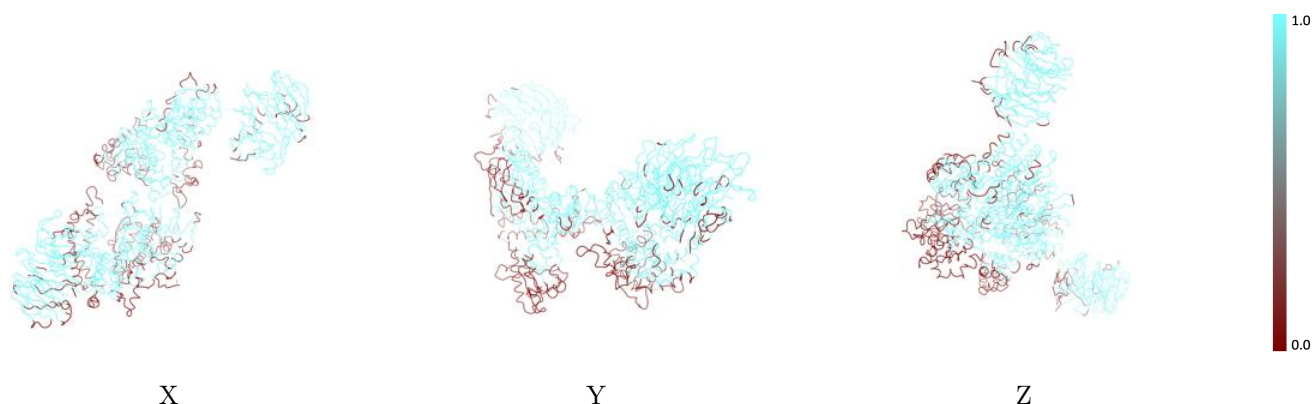
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



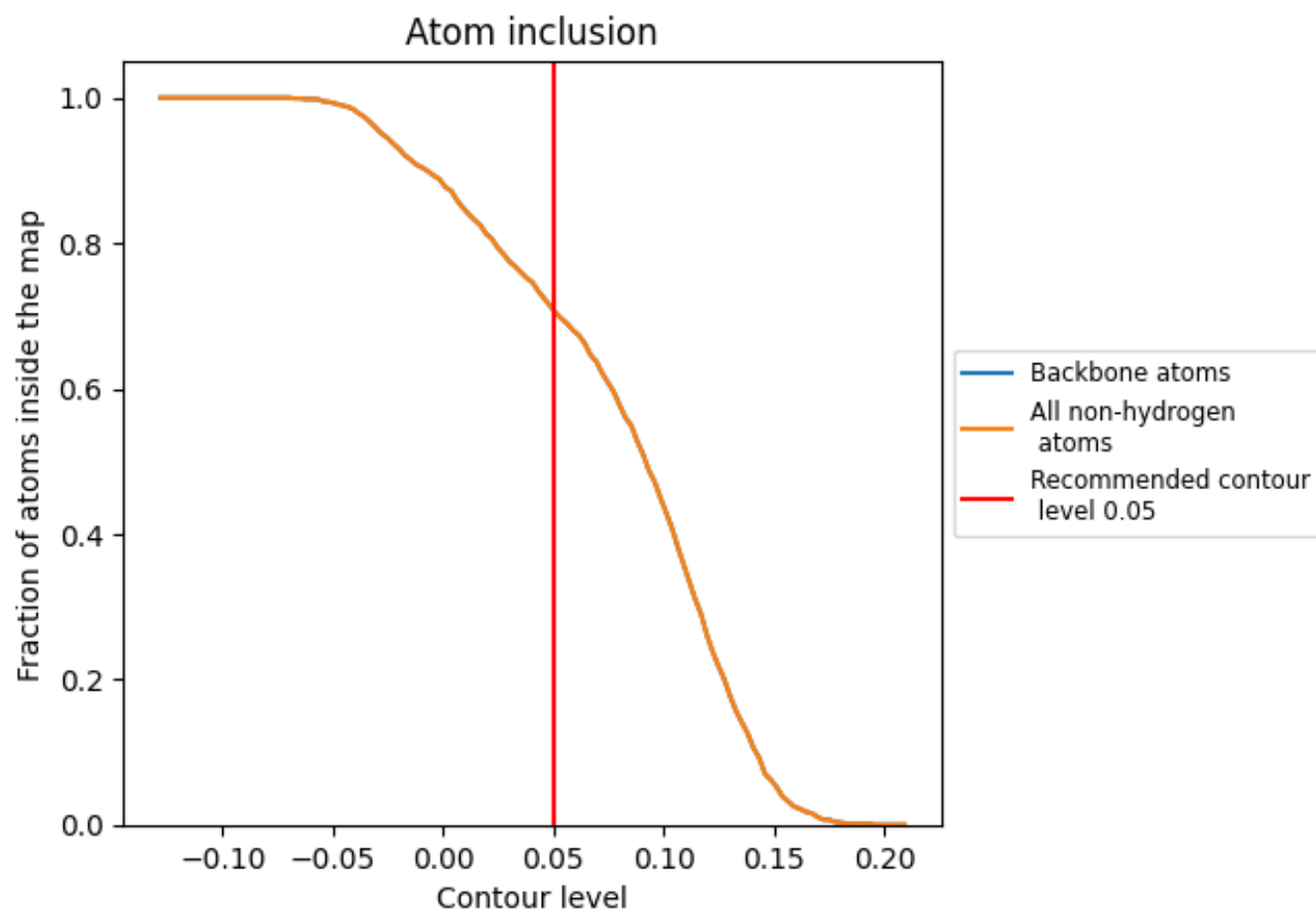
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.05).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 71% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7080	0.0370
A	0.7670	0.0390
B	0.6480	0.0340

