



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2018 – 08:34 am GMT

PDB ID : 4C3B
Title : HRSV M2-1, P21 crystal form
Authors : Tanner, S.J.; Ariza, A.; Richard, C.A.; Wu, W.; Trincao, J.; Hiscox, J.A.;
Carroll, M.W.; Silman, N.J.; Eleouet, J.F.; Edwards, T.A.; Barr, J.N.
Deposited on : 2013-08-22
Resolution : 2.95 Å(reported)

This is a wwPDB X-ray Structure 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/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : trunk30967
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk30967

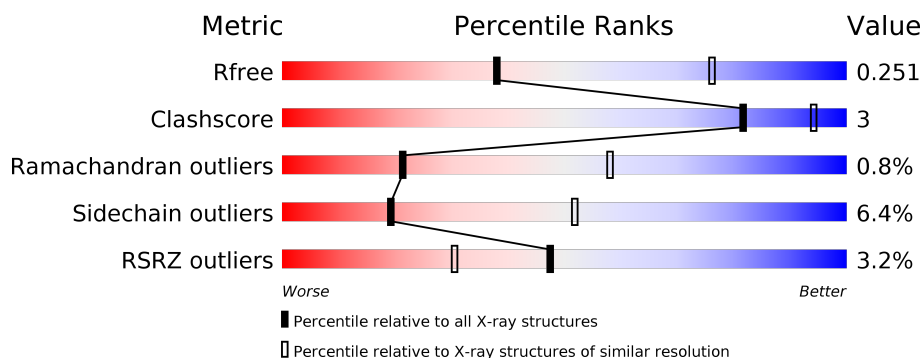
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	111664	2641 (3.00-2.92)
Clashscore	122126	2988 (3.00-2.92)
Ramachandran outliers	120053	2892 (3.00-2.92)
Sidechain outliers	120020	2895 (3.00-2.92)
RSRZ outliers	108989	2527 (3.00-2.92)

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

Mol	Chain	Length	Quality of chain
1	A	199	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>10%</div> <div>•</div> <div>17%</div> </div> </div>
1	B	199	<div> <div>2%</div> <div> <div></div> <div>71%</div> <div>10%</div> <div>•</div> <div>18%</div> </div> </div>
1	C	199	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>14%</div> <div>••</div> <div>11%</div> </div> </div>
1	D	199	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>11%</div> <div>•</div> <div>13%</div> </div> </div>
1	E	199	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>10%</div> <div>•</div> <div>13%</div> </div> </div>
1	F	199	<div> <div>5%</div> <div> <div></div> <div>75%</div> <div>13%</div> <div>••</div> <div>11%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	199	<p>2% 72% 9% 18%</p>
1	H	199	<p>4% 73% 12% 13%</p>
1	I	199	<p>% 69% 8% 22%</p>
1	J	199	<p>4% 71% 9% 19%</p>
1	K	199	<p>4% 73% 10% 16%</p>
1	L	199	<p>4% 72% 7% 20%</p>
1	M	199	<p>2% 71% 10% 19%</p>
1	N	199	<p>% 67% 10% 22%</p>
1	O	199	<p>5% 74% 10% 15%</p>
1	P	199	<p>4% 72% 8% 20%</p>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21688 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called MATRIX PROTEIN 2-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	1	0
			1344	844	248	244	8			
1	B	164	Total	C	N	O	S	0	0	0
			1326	833	245	240	8			
1	C	178	Total	C	N	O	S	0	0	0
			1425	893	259	265	8			
1	D	174	Total	C	N	O	S	0	0	0
			1396	876	255	257	8			
1	E	174	Total	C	N	O	S	0	0	0
			1396	876	255	257	8			
1	F	178	Total	C	N	O	S	0	0	0
			1425	893	259	265	8			
1	G	164	Total	C	N	O	S	0	0	0
			1327	834	244	241	8			
1	H	174	Total	C	N	O	S	0	0	0
			1405	882	255	260	8			
1	I	156	Total	C	N	O	S	0	0	0
			1272	800	236	229	7			
1	J	161	Total	C	N	O	S	0	0	0
			1311	824	242	238	7			
1	K	167	Total	C	N	O	S	0	0	0
			1344	844	244	249	7			
1	L	160	Total	C	N	O	S	0	0	0
			1300	818	238	237	7			
1	M	161	Total	C	N	O	S	0	0	0
			1305	821	239	238	7			
1	N	155	Total	C	N	O	S	0	0	0
			1259	794	230	228	7			
1	O	169	Total	C	N	O	S	0	0	0
			1363	855	250	251	7			
1	P	160	Total	C	N	O	S	0	0	0
			1296	815	237	237	7			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP P04545
A	-3	PRO	-	expression tag	UNP P04545
A	-2	LEU	-	expression tag	UNP P04545
A	-1	GLY	-	expression tag	UNP P04545
A	0	SER	-	expression tag	UNP P04545
B	-4	GLY	-	expression tag	UNP P04545
B	-3	PRO	-	expression tag	UNP P04545
B	-2	LEU	-	expression tag	UNP P04545
B	-1	GLY	-	expression tag	UNP P04545
B	0	SER	-	expression tag	UNP P04545
C	-4	GLY	-	expression tag	UNP P04545
C	-3	PRO	-	expression tag	UNP P04545
C	-2	LEU	-	expression tag	UNP P04545
C	-1	GLY	-	expression tag	UNP P04545
C	0	SER	-	expression tag	UNP P04545
D	-4	GLY	-	expression tag	UNP P04545
D	-3	PRO	-	expression tag	UNP P04545
D	-2	LEU	-	expression tag	UNP P04545
D	-1	GLY	-	expression tag	UNP P04545
D	0	SER	-	expression tag	UNP P04545
E	-4	GLY	-	expression tag	UNP P04545
E	-3	PRO	-	expression tag	UNP P04545
E	-2	LEU	-	expression tag	UNP P04545
E	-1	GLY	-	expression tag	UNP P04545
E	0	SER	-	expression tag	UNP P04545
F	-4	GLY	-	expression tag	UNP P04545
F	-3	PRO	-	expression tag	UNP P04545
F	-2	LEU	-	expression tag	UNP P04545
F	-1	GLY	-	expression tag	UNP P04545
F	0	SER	-	expression tag	UNP P04545
G	-4	GLY	-	expression tag	UNP P04545
G	-3	PRO	-	expression tag	UNP P04545
G	-2	LEU	-	expression tag	UNP P04545
G	-1	GLY	-	expression tag	UNP P04545
G	0	SER	-	expression tag	UNP P04545
H	-4	GLY	-	expression tag	UNP P04545
H	-3	PRO	-	expression tag	UNP P04545
H	-2	LEU	-	expression tag	UNP P04545
H	-1	GLY	-	expression tag	UNP P04545
H	0	SER	-	expression tag	UNP P04545
I	-4	GLY	-	expression tag	UNP P04545
I	-3	PRO	-	expression tag	UNP P04545

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-2	LEU	-	expression tag	UNP P04545
I	-1	GLY	-	expression tag	UNP P04545
I	0	SER	-	expression tag	UNP P04545
J	-4	GLY	-	expression tag	UNP P04545
J	-3	PRO	-	expression tag	UNP P04545
J	-2	LEU	-	expression tag	UNP P04545
J	-1	GLY	-	expression tag	UNP P04545
J	0	SER	-	expression tag	UNP P04545
K	-4	GLY	-	expression tag	UNP P04545
K	-3	PRO	-	expression tag	UNP P04545
K	-2	LEU	-	expression tag	UNP P04545
K	-1	GLY	-	expression tag	UNP P04545
K	0	SER	-	expression tag	UNP P04545
L	-4	GLY	-	expression tag	UNP P04545
L	-3	PRO	-	expression tag	UNP P04545
L	-2	LEU	-	expression tag	UNP P04545
L	-1	GLY	-	expression tag	UNP P04545
L	0	SER	-	expression tag	UNP P04545
M	-4	GLY	-	expression tag	UNP P04545
M	-3	PRO	-	expression tag	UNP P04545
M	-2	LEU	-	expression tag	UNP P04545
M	-1	GLY	-	expression tag	UNP P04545
M	0	SER	-	expression tag	UNP P04545
N	-4	GLY	-	expression tag	UNP P04545
N	-3	PRO	-	expression tag	UNP P04545
N	-2	LEU	-	expression tag	UNP P04545
N	-1	GLY	-	expression tag	UNP P04545
N	0	SER	-	expression tag	UNP P04545
O	-4	GLY	-	expression tag	UNP P04545
O	-3	PRO	-	expression tag	UNP P04545
O	-2	LEU	-	expression tag	UNP P04545
O	-1	GLY	-	expression tag	UNP P04545
O	0	SER	-	expression tag	UNP P04545
P	-4	GLY	-	expression tag	UNP P04545
P	-3	PRO	-	expression tag	UNP P04545
P	-2	LEU	-	expression tag	UNP P04545
P	-1	GLY	-	expression tag	UNP P04545
P	0	SER	-	expression tag	UNP P04545

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	P	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0
2	J	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	K	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	I	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	A	1	Total Zn 1 1	0	0
2	N	1	Total Zn 1 1	0	0
2	O	1	Total Zn 1 1	0	0
2	L	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	M	1	Total Zn 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	15	Total O 15 15	0	0
3	B	20	Total O 20 20	0	0
3	C	15	Total O 15 15	0	0
3	D	11	Total O 11 11	0	0

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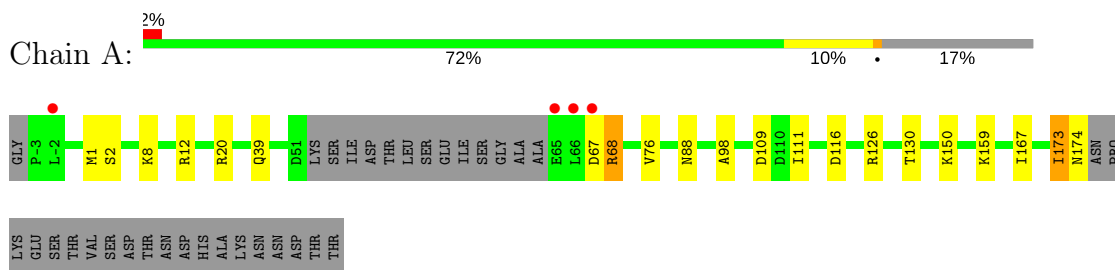
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	11	Total 11	O 11	0	0
3	F	18	Total 18	O 18	0	0
3	G	17	Total 17	O 17	0	0
3	H	16	Total 16	O 16	0	0
3	I	4	Total 4	O 4	0	0
3	J	6	Total 6	O 6	0	0
3	K	3	Total 3	O 3	0	0
3	L	8	Total 8	O 8	0	0
3	M	8	Total 8	O 8	0	0
3	N	8	Total 8	O 8	0	0
3	O	4	Total 4	O 4	0	0
3	P	14	Total 14	O 14	0	0

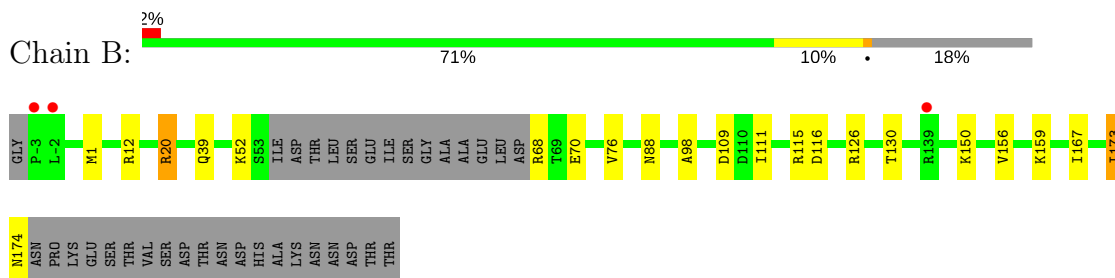
3 Residue-property plots [i](#)

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

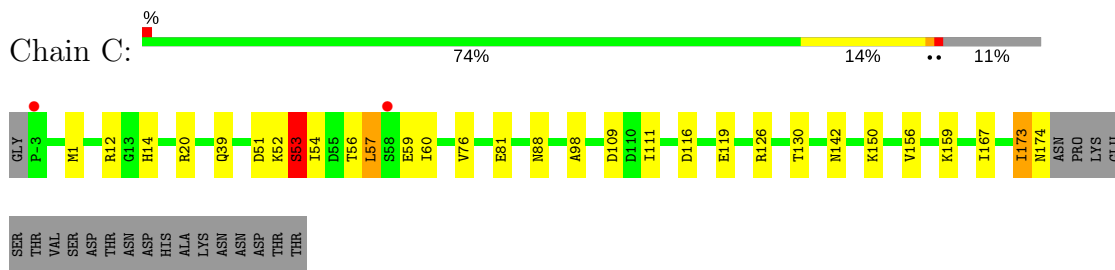
• Molecule 1: MATRIX PROTEIN 2-1



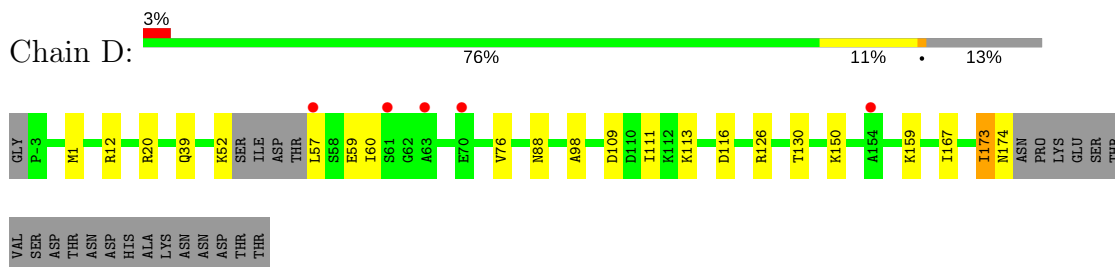
• Molecule 1: MATRIX PROTEIN 2-1



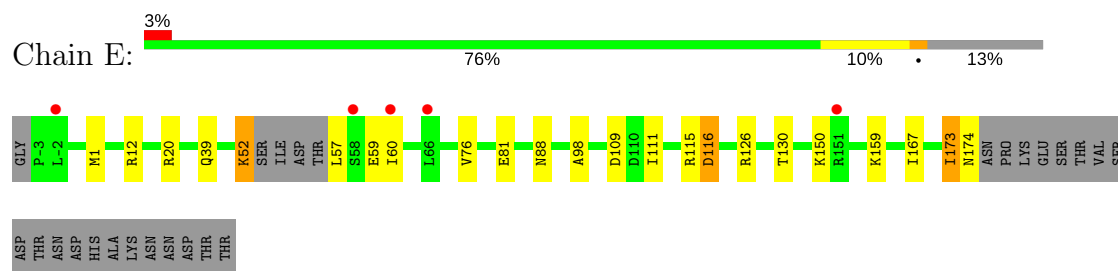
• Molecule 1: MATRIX PROTEIN 2-1



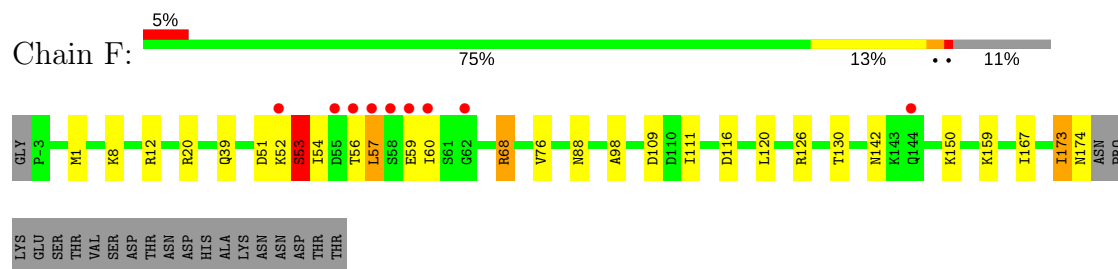
• Molecule 1: MATRIX PROTEIN 2-1



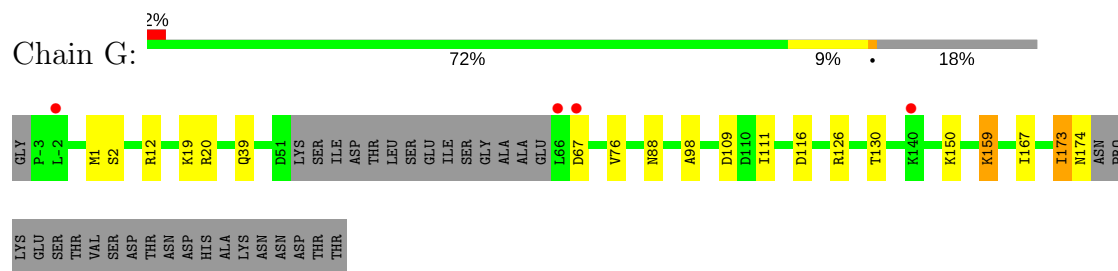
- Molecule 1: MATRIX PROTEIN 2-1



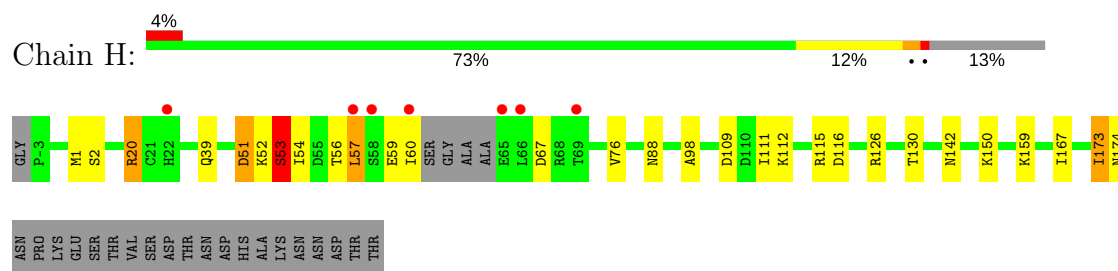
- Molecule 1: MATRIX PROTEIN 2-1



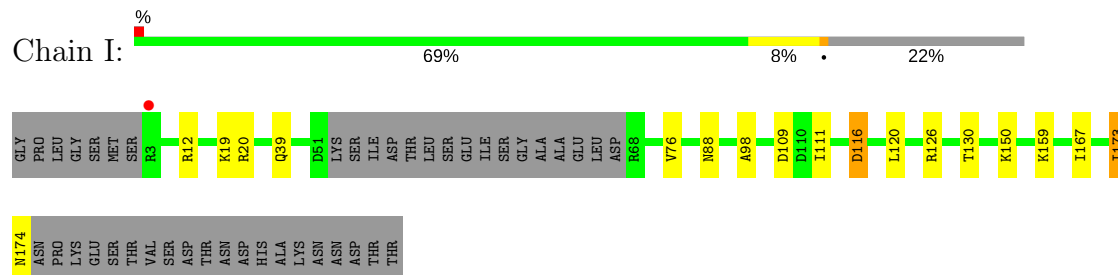
- Molecule 1: MATRIX PROTEIN 2-1



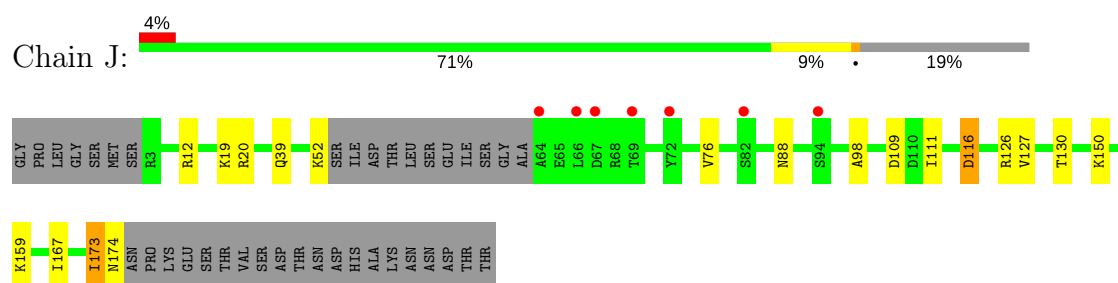
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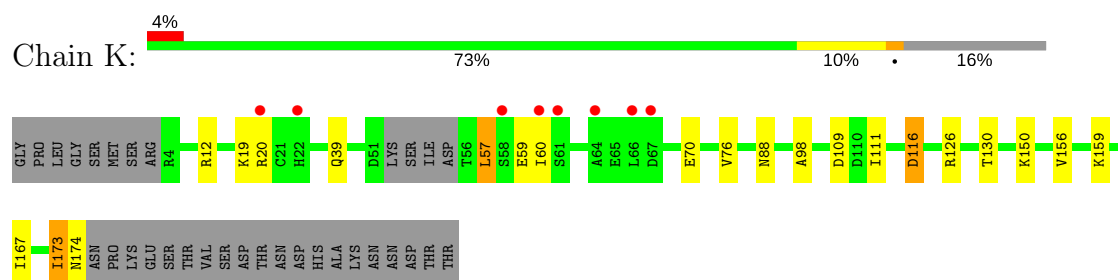
- Molecule 1: MATRIX PROTEIN 2-1



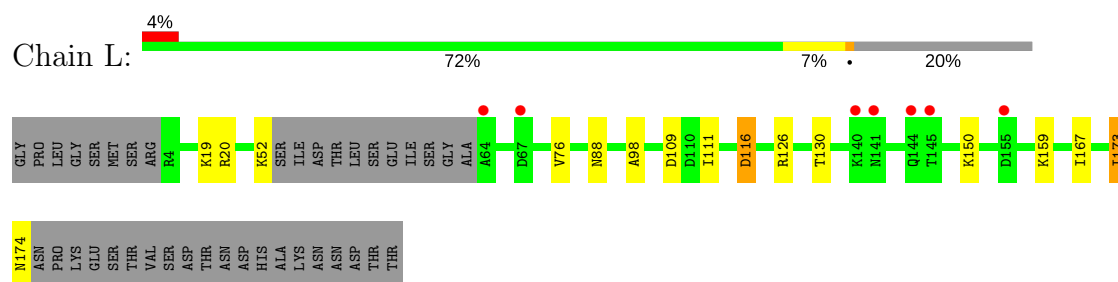
- Molecule 1: MATRIX PROTEIN 2-1



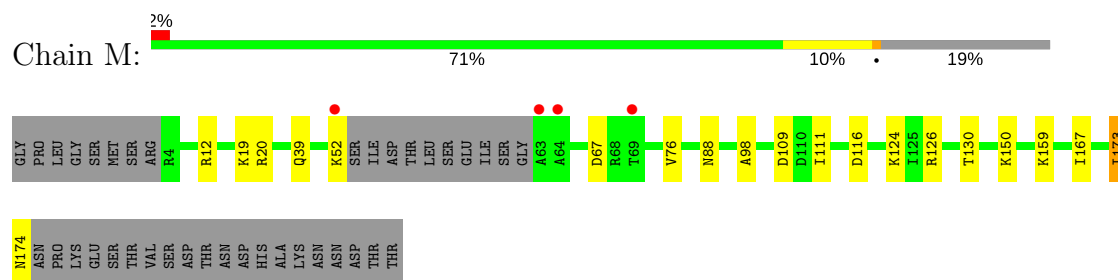
- Molecule 1: MATRIX PROTEIN 2-1



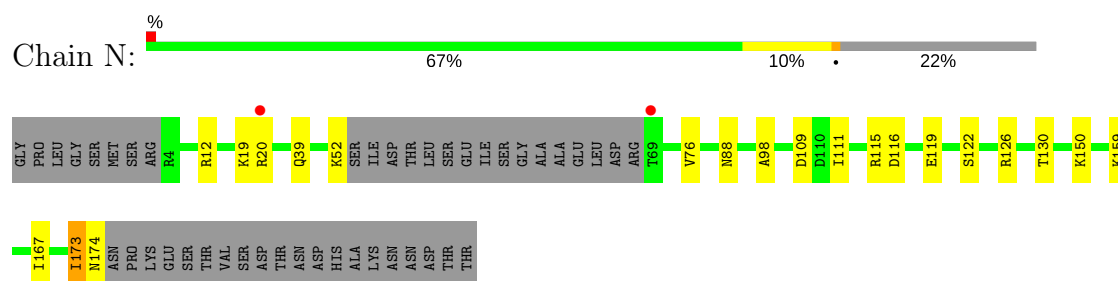
- Molecule 1: MATRIX PROTEIN 2-1



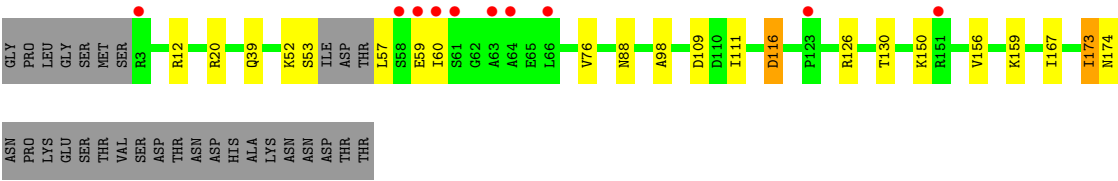
- Molecule 1: MATRIX PROTEIN 2-1



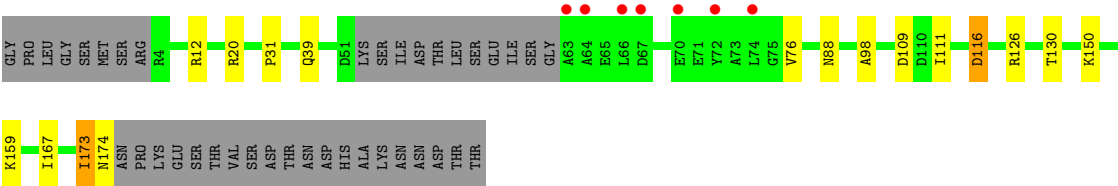
- Molecule 1: MATRIX PROTEIN 2-1



- Molecule 1: MATRIX PROTEIN 2-1



• Molecule 1: MATRIX PROTEIN 2-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.85Å 141.62Å 141.69Å 90.00° 96.14° 90.00°	Depositor
Resolution (Å)	140.88 – 2.95 140.88 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.5 (140.88-2.95) 99.5 (140.88-2.95)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.96Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.205 , 0.255 0.207 , 0.251	Depositor DCC
R_{free} test set	3797 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	53.0	Xtriage
Anisotropy	0.896	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 59.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	21688	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.74	0/1368	1.08	7/1838 (0.4%)
1	B	0.70	0/1347	0.81	1/1809 (0.1%)
1	C	0.74	0/1447	0.83	2/1947 (0.1%)
1	D	0.70	0/1417	0.79	1/1904 (0.1%)
1	E	0.70	0/1417	0.82	1/1904 (0.1%)
1	F	0.73	0/1447	0.87	4/1947 (0.2%)
1	G	0.72	0/1348	0.80	1/1812 (0.1%)
1	H	0.71	0/1426	0.85	3/1917 (0.2%)
1	I	0.62	0/1292	0.76	1/1737 (0.1%)
1	J	0.64	0/1331	0.78	1/1789 (0.1%)
1	K	0.62	0/1364	0.77	1/1836 (0.1%)
1	L	0.62	0/1320	0.76	0/1775
1	M	0.63	0/1325	0.75	1/1782 (0.1%)
1	N	0.64	0/1279	0.77	1/1720 (0.1%)
1	O	0.65	0/1383	0.79	1/1859 (0.1%)
1	P	0.60	0/1316	0.75	1/1771 (0.1%)
All	All	0.68	0/21827	0.82	27/29347 (0.1%)

There are no bond length outliers.

The worst 5 of 27 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	68[A]	ARG	NE-CZ-NH2	-15.96	112.32	120.30
1	A	68[B]	ARG	NE-CZ-NH2	-15.96	112.32	120.30
1	A	68[A]	ARG	NE-CZ-NH1	10.58	125.59	120.30
1	A	68[B]	ARG	NE-CZ-NH1	10.58	125.59	120.30
1	F	68	ARG	CG-CD-NE	9.53	131.81	111.80

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	1344	0	1387	8	0
1	B	1326	0	1371	9	0
1	C	1425	0	1466	12	0
1	D	1396	0	1438	8	0
1	E	1396	0	1438	12	0
1	F	1425	0	1466	12	0
1	G	1327	0	1368	9	0
1	H	1405	0	1447	15	0
1	I	1272	0	1310	7	0
1	J	1311	0	1349	8	0
1	K	1344	0	1376	10	0
1	L	1300	0	1336	5	0
1	M	1305	0	1341	7	0
1	N	1259	0	1297	9	0
1	O	1363	0	1400	9	0
1	P	1296	0	1328	6	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0
2	N	1	0	0	0	0
2	O	1	0	0	0	0
2	P	1	0	0	0	0
3	A	15	0	0	2	0
3	B	20	0	0	3	0
3	C	15	0	0	1	0
3	D	11	0	0	1	0
3	E	11	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	18	0	0	1	0
3	G	17	0	0	2	0
3	H	16	0	0	0	0
3	I	4	0	0	0	0
3	J	6	0	0	1	0
3	K	3	0	0	1	0
3	L	8	0	0	0	0
3	M	8	0	0	1	0
3	N	8	0	0	0	0
3	O	4	0	0	0	0
3	P	14	0	0	1	0
All	All	21688	0	22118	120	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

The worst 5 of 120 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:E:52:LYS:HG2	3:E:2007:HOH:O	1.59	1.00
1:G:159:LYS:HE2	3:G:2015:HOH:O	1.78	0.83
1:O:53:SER:HA	1:O:57:LEU:HB2	1.62	0.81
1:F:120:LEU:HD11	1:N:119:GLU:HB2	1.67	0.76
1:B:70:GLU:HG2	3:B:2013:HOH:O	1.91	0.70

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	162/199 (81%)	153 (94%)	8 (5%)	1 (1%)	27 64

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	160/199 (80%)	151 (94%)	8 (5%)	1 (1%)	27 64
1	C	176/199 (88%)	163 (93%)	10 (6%)	3 (2%)	10 38
1	D	170/199 (85%)	162 (95%)	7 (4%)	1 (1%)	27 64
1	E	170/199 (85%)	160 (94%)	9 (5%)	1 (1%)	27 64
1	F	176/199 (88%)	163 (93%)	10 (6%)	3 (2%)	10 38
1	G	160/199 (80%)	152 (95%)	7 (4%)	1 (1%)	27 64
1	H	170/199 (85%)	157 (92%)	11 (6%)	2 (1%)	14 48
1	I	152/199 (76%)	143 (94%)	8 (5%)	1 (1%)	24 61
1	J	157/199 (79%)	148 (94%)	8 (5%)	1 (1%)	27 64
1	K	163/199 (82%)	152 (93%)	9 (6%)	2 (1%)	14 48
1	L	156/199 (78%)	147 (94%)	8 (5%)	1 (1%)	27 64
1	M	157/199 (79%)	147 (94%)	9 (6%)	1 (1%)	27 64
1	N	151/199 (76%)	142 (94%)	8 (5%)	1 (1%)	24 61
1	O	165/199 (83%)	154 (93%)	10 (6%)	1 (1%)	27 64
1	P	156/199 (78%)	147 (94%)	8 (5%)	1 (1%)	27 64
All	All	2601/3184 (82%)	2441 (94%)	138 (5%)	22 (1%)	21 58

5 of 22 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	53	SER
1	H	53	SER
1	A	88	ASN
1	B	88	ASN
1	C	53	SER

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	155/183 (85%)	145 (94%)	10 (6%)	19 51

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	153/183 (84%)	142 (93%)	11 (7%)	16	46
1	C	164/183 (90%)	152 (93%)	12 (7%)	15	45
1	D	160/183 (87%)	150 (94%)	10 (6%)	20	52
1	E	160/183 (87%)	150 (94%)	10 (6%)	20	52
1	F	164/183 (90%)	152 (93%)	12 (7%)	15	45
1	G	153/183 (84%)	143 (94%)	10 (6%)	19	51
1	H	163/183 (89%)	151 (93%)	12 (7%)	15	45
1	I	146/183 (80%)	138 (94%)	8 (6%)	24	58
1	J	150/183 (82%)	141 (94%)	9 (6%)	21	54
1	K	154/183 (84%)	145 (94%)	9 (6%)	22	56
1	L	149/183 (81%)	140 (94%)	9 (6%)	21	54
1	M	149/183 (81%)	140 (94%)	9 (6%)	21	54
1	N	145/183 (79%)	136 (94%)	9 (6%)	20	53
1	O	156/183 (85%)	146 (94%)	10 (6%)	19	52
1	P	148/183 (81%)	140 (95%)	8 (5%)	24	59
All	All	2469/2928 (84%)	2311 (94%)	158 (6%)	19	52

5 of 158 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	130	THR
1	I	109	ASP
1	O	150	LYS
1	G	159	LYS
1	H	109	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 26 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	138	ASN
1	J	144	GLN
1	P	138	ASN
1	I	144	GLN
1	J	138	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 16 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	165/199 (82%)	0.10	4 (2%)	59	40	25, 52, 105, 133	0
1	B	164/199 (82%)	0.01	3 (1%)	68	50	25, 56, 107, 149	0
1	C	178/199 (89%)	0.15	2 (1%)	80	64	22, 55, 124, 144	0
1	D	174/199 (87%)	0.21	5 (2%)	51	34	22, 58, 119, 170	0
1	E	174/199 (87%)	0.21	5 (2%)	51	34	23, 59, 109, 169	0
1	F	178/199 (89%)	0.24	9 (5%)	28	17	23, 57, 128, 167	0
1	G	164/199 (82%)	0.03	4 (2%)	59	40	24, 52, 110, 129	0
1	H	174/199 (87%)	0.07	7 (4%)	38	24	26, 56, 125, 170	0
1	I	156/199 (78%)	0.11	1 (0%)	89	77	37, 69, 120, 145	0
1	J	161/199 (80%)	0.21	7 (4%)	35	22	38, 68, 126, 153	0
1	K	167/199 (83%)	0.35	8 (4%)	30	19	37, 70, 128, 175	0
1	L	160/199 (80%)	0.33	7 (4%)	34	21	38, 75, 123, 157	0
1	M	161/199 (80%)	0.08	4 (2%)	57	39	37, 71, 118, 141	0
1	N	155/199 (77%)	0.12	2 (1%)	77	60	36, 66, 115, 139	0
1	O	169/199 (84%)	0.28	10 (5%)	22	13	36, 65, 131, 152	0
1	P	160/199 (80%)	0.31	7 (4%)	34	21	36, 77, 135, 152	0
All	All	2660/3184 (83%)	0.18	85 (3%)	47	30	22, 64, 122, 175	0

The worst 5 of 85 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	58	SER	7.4
1	J	69	THR	6.9
1	K	58	SER	5.8
1	K	20	ARG	4.8
1	P	63	ALA	4.8

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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. 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	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	L	201	1/1	0.96	0.03	58,58,58,58	0
2	ZN	D	201	1/1	0.97	0.05	33,33,33,33	0
2	ZN	J	201	1/1	0.97	0.04	53,53,53,53	0
2	ZN	E	201	1/1	0.98	0.04	31,31,31,31	0
2	ZN	P	201	1/1	0.98	0.04	47,47,47,47	0
2	ZN	C	201	1/1	0.98	0.05	30,30,30,30	0
2	ZN	H	201	1/1	0.99	0.03	37,37,37,37	0
2	ZN	N	201	1/1	0.99	0.02	45,45,45,45	0
2	ZN	A	201	1/1	0.99	0.04	42,42,42,42	0
2	ZN	G	201	1/1	0.99	0.06	39,39,39,39	0
2	ZN	K	201	1/1	0.99	0.04	54,54,54,54	0
2	ZN	M	201	1/1	0.99	0.05	59,59,59,59	0
2	ZN	B	201	1/1	0.99	0.05	32,32,32,32	0
2	ZN	I	201	1/1	0.99	0.05	56,56,56,56	0
2	ZN	O	201	1/1	0.99	0.03	50,50,50,50	0
2	ZN	F	201	1/1	0.99	0.03	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.