



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 04:56 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 Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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

A user guide is available at

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

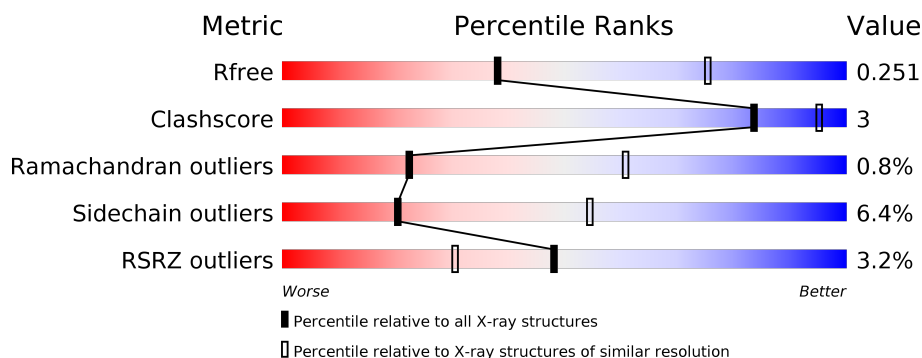
1 Overall quality at a glance

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}	100719	2395 (3.00-2.92)
Clashscore	112137	2773 (3.00-2.92)
Ramachandran outliers	110173	2680 (3.00-2.92)
Sidechain outliers	110143	2683 (3.00-2.92)
RSRZ outliers	101464	2421 (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>0%</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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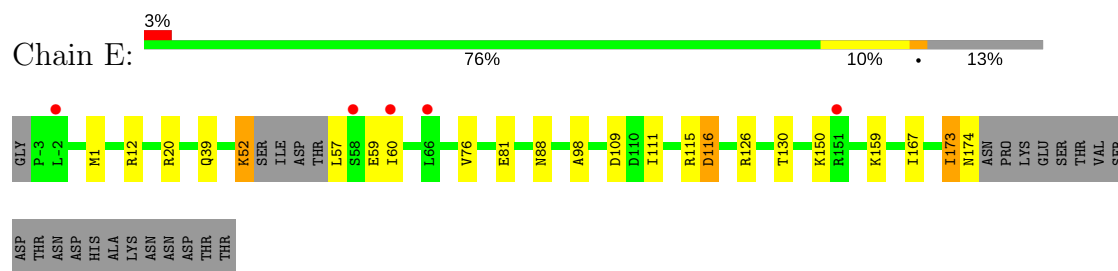
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	11	Total 11	O 11	0	0
3	F	18	Total 18	O 18	0	0
3	G	17	Total 17	O 17	0	0
3	H	16	Total 16	O 16	0	0
3	I	4	Total 4	O 4	0	0
3	J	6	Total 6	O 6	0	0
3	K	3	Total 3	O 3	0	0
3	L	8	Total 8	O 8	0	0
3	M	8	Total 8	O 8	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

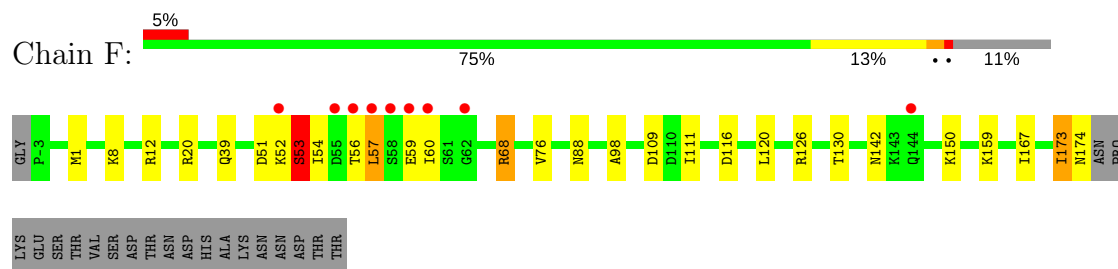
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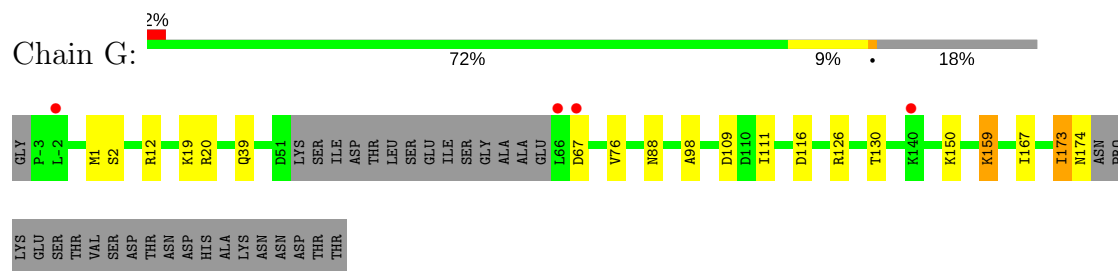
- Molecule 1: MATRIX PROTEIN 2-1



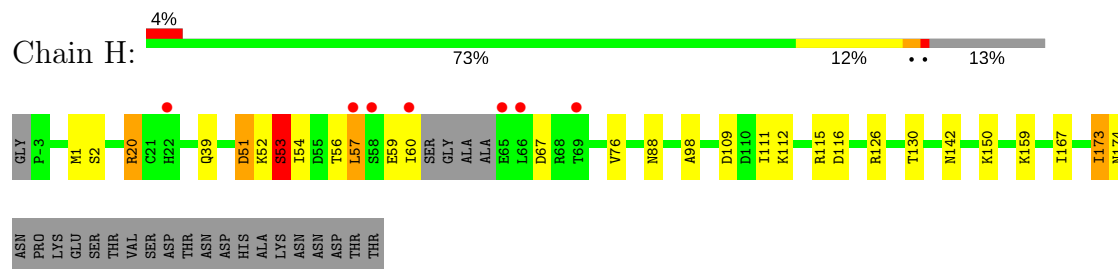
- Molecule 1: MATRIX PROTEIN 2-1



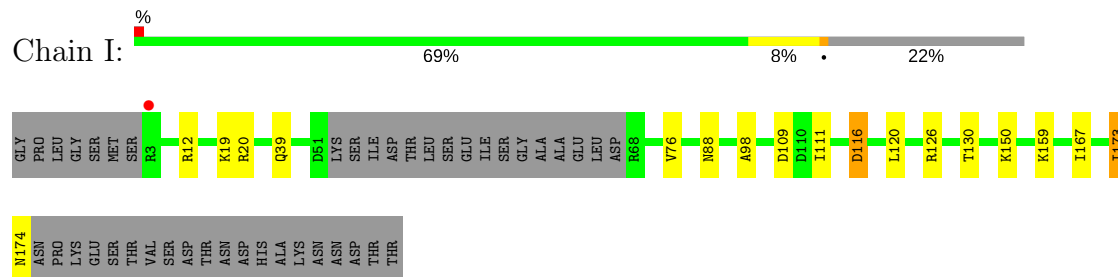
- Molecule 1: MATRIX PROTEIN 2-1



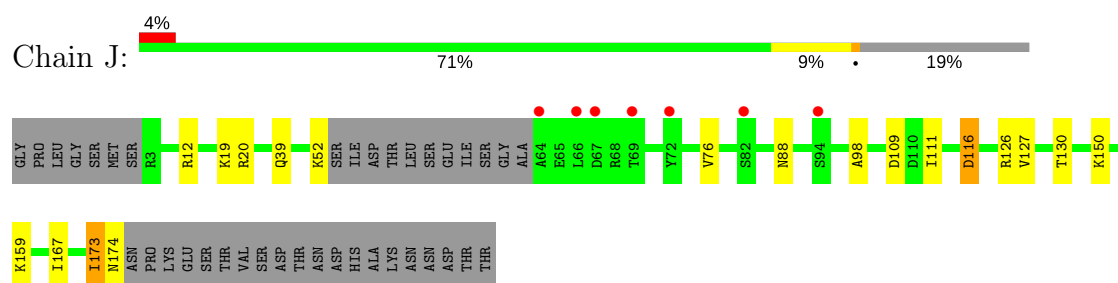
- Molecule 1: MATRIX PROTEIN 2-1



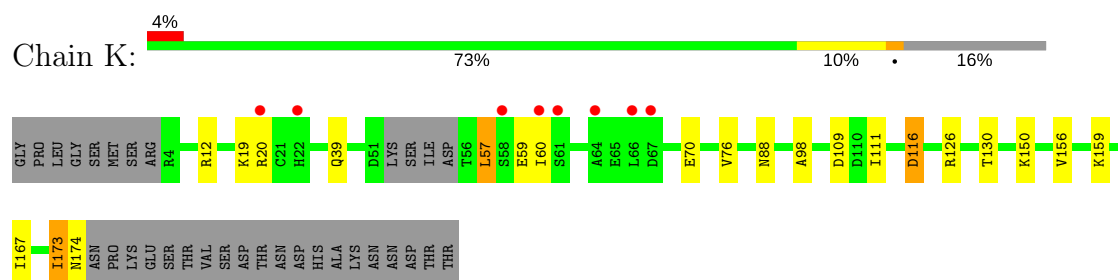
- Molecule 1: MATRIX PROTEIN 2-1



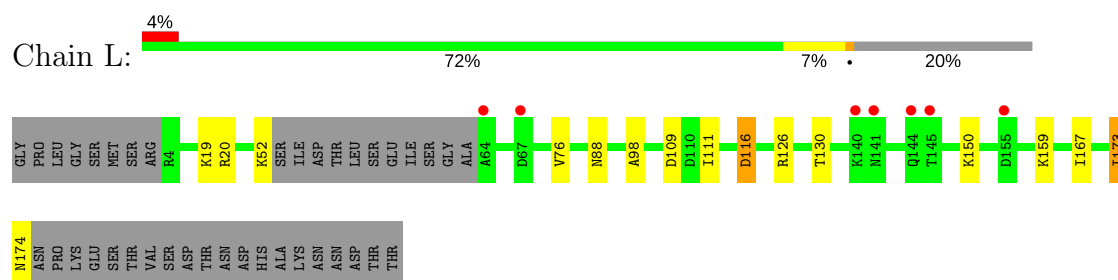
- Molecule 1: MATRIX PROTEIN 2-1



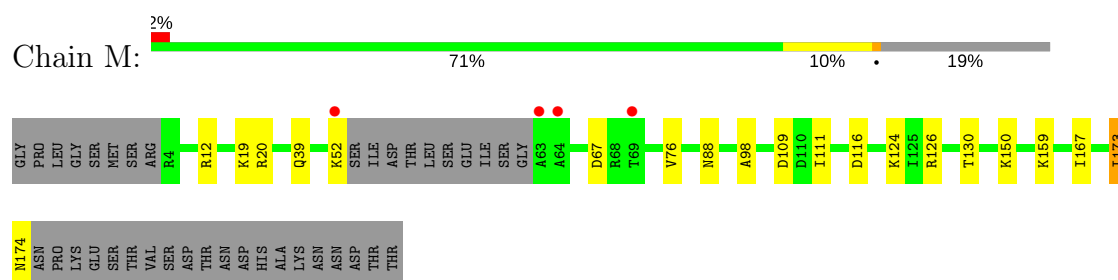
- Molecule 1: MATRIX PROTEIN 2-1



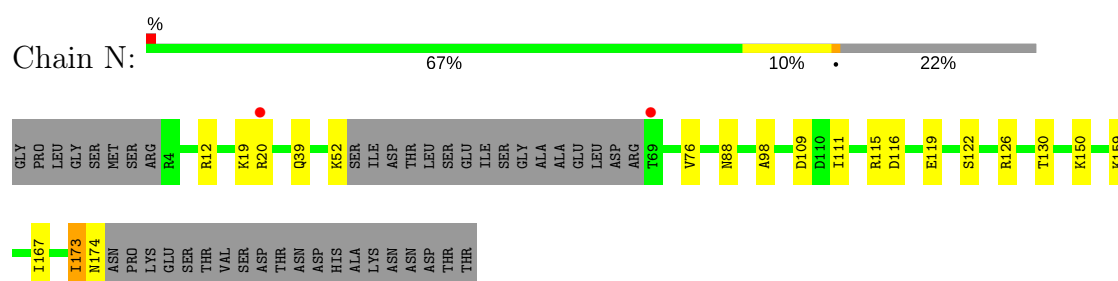
- Molecule 1: MATRIX PROTEIN 2-1



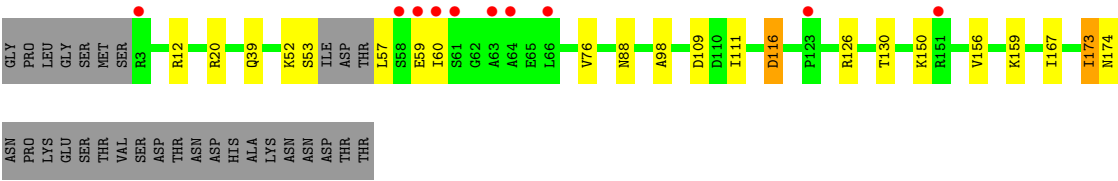
- Molecule 1: MATRIX PROTEIN 2-1



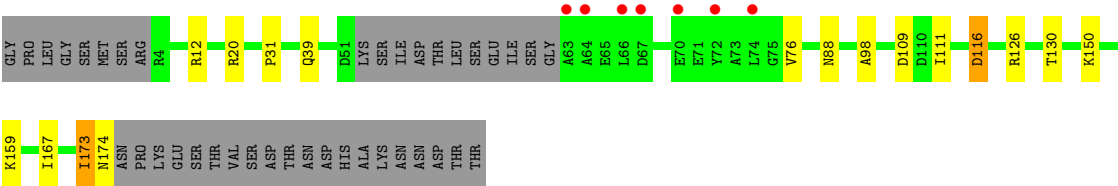
- Molecule 1: MATRIX PROTEIN 2-1



- 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 (5.29%)	DCC
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.

All (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
1	A	68[A]	ARG	CG-CD-NE	-7.59	95.86	111.80
1	A	68[B]	ARG	CG-CD-NE	-7.59	95.86	111.80
1	H	57	LEU	CA-CB-CG	7.33	132.16	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	12	ARG	NE-CZ-NH2	-7.02	116.79	120.30
1	J	12	ARG	NE-CZ-NH2	-6.77	116.92	120.30
1	C	12	ARG	NE-CZ-NH2	-6.43	117.09	120.30
1	B	12	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	G	12	ARG	NE-CZ-NH2	-6.33	117.14	120.30
1	O	12	ARG	NE-CZ-NH2	-6.05	117.27	120.30
1	A	12	ARG	NE-CZ-NH2	-6.00	117.30	120.30
1	F	51	ASP	CB-CG-OD1	5.93	123.64	118.30
1	C	51	ASP	CB-CG-OD1	5.89	123.60	118.30
1	F	12	ARG	NE-CZ-NH2	-5.79	117.41	120.30
1	K	12	ARG	NE-CZ-NH2	-5.78	117.41	120.30
1	N	12	ARG	NE-CZ-NH2	-5.54	117.53	120.30
1	D	12	ARG	NE-CZ-NH2	-5.45	117.58	120.30
1	H	51	ASP	CB-CG-OD1	5.37	123.13	118.30
1	P	12	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	H	57	LEU	CB-CA-C	5.28	120.23	110.20
1	M	12	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	I	12	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	F	68	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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
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.

All (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
1:A:67:ASP:O	1:B:115:ARG:NH2	2.26	0.67
1:E:52:LYS:CG	3:E:2007:HOH:O	2.31	0.66
1:D:173:ILE:O	1:D:174:ASN:CG	2.33	0.65
1:C:173:ILE:O	1:C:174:ASN:CG	2.36	0.64
1:P:173:ILE:O	1:P:174:ASN:CG	2.36	0.64
1:G:159:LYS:CE	3:G:2015:HOH:O	2.41	0.64
1:N:173:ILE:O	1:N:174:ASN:CG	2.37	0.63
1:O:173:ILE:O	1:O:174:ASN:CG	2.37	0.63
1:M:173:ILE:O	1:M:174:ASN:CG	2.38	0.62
1:E:173:ILE:O	1:E:174:ASN:CG	2.38	0.62
1:I:19:LYS:NZ	1:K:116:ASP:OD2	2.29	0.62
1:G:173:ILE:O	1:G:174:ASN:CG	2.38	0.62
1:J:173:ILE:O	1:J:174:ASN:CG	2.38	0.62
1:A:173:ILE:O	1:A:174:ASN:CG	2.38	0.61
1:C:119:GLU:HB2	1:I:120:LEU:HD11	1.80	0.61
1:F:54:ILE:HA	1:F:57:LEU:HB2	1.81	0.61
1:I:173:ILE:O	1:I:174:ASN:CG	2.39	0.61
1:F:173:ILE:O	1:F:174:ASN:CG	2.39	0.61
1:H:173:ILE:O	1:H:174:ASN:CG	2.40	0.60
1:C:54:ILE:HA	1:C:57:LEU:HB2	1.83	0.60
1:K:173:ILE:O	1:K:174:ASN:CG	2.40	0.60
1:B:173:ILE:O	1:B:174:ASN:CG	2.39	0.60
1:L:173:ILE:O	1:L:174:ASN:CG	2.40	0.60
1:B:20:ARG:NH2	1:C:81:GLU:OE1	2.34	0.59
1:B:76:VAL:HG13	1:B:98:ALA:HB1	1.88	0.55
1:G:76:VAL:HG13	1:G:98:ALA:HB1	1.89	0.55
1:H:57:LEU:O	1:H:60:ILE:HG22	2.07	0.54
1:H:76:VAL:HG13	1:H:98:ALA:HB1	1.89	0.54
1:C:57:LEU:O	1:C:60:ILE:HG22	2.08	0.54
1:P:76:VAL:HG13	1:P:98:ALA:HB1	1.89	0.54
1:K:57:LEU:O	1:K:60:ILE:HG22	2.08	0.54
1:J:76:VAL:HG13	1:J:98:ALA:HB1	1.90	0.54
1:M:19:LYS:NZ	1:O:116:ASP:OD2	2.34	0.54
1:F:57:LEU:O	1:F:60:ILE:HG22	2.08	0.54
1:M:76:VAL:HG13	1:M:98:ALA:HB1	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:76:VAL:HG13	1:L:98:ALA:HB1	1.90	0.54
1:N:76:VAL:HG13	1:N:98:ALA:HB1	1.91	0.53
1:A:76:VAL:HG13	1:A:98:ALA:HB1	1.90	0.53
1:D:57:LEU:O	1:D:60:ILE:HG22	2.08	0.53
1:P:31:PRO:HB3	3:P:2004:HOH:O	2.08	0.53
1:E:57:LEU:O	1:E:60:ILE:HG22	2.08	0.53
1:O:57:LEU:O	1:O:60:ILE:HG22	2.08	0.53
1:O:76:VAL:HG13	1:O:98:ALA:HB1	1.91	0.52
1:E:76:VAL:HG13	1:E:98:ALA:HB1	1.92	0.52
1:K:76:VAL:HG13	1:K:98:ALA:HB1	1.92	0.52
1:B:70:GLU:CG	3:B:2013:HOH:O	2.55	0.51
1:I:76:VAL:HG13	1:I:98:ALA:HB1	1.91	0.51
1:C:76:VAL:HG13	1:C:98:ALA:HB1	1.91	0.51
1:I:39:GLN:OE1	1:J:39:GLN:OE1	2.28	0.51
1:F:76:VAL:HG13	1:F:98:ALA:HB1	1.92	0.51
1:F:120:LEU:HD12	1:N:122:SER:HB2	1.93	0.50
1:A:68[B]:ARG:HD2	3:A:2008:HOH:O	2.11	0.50
1:G:67:ASP:O	1:H:115:ARG:NH2	2.36	0.50
1:D:76:VAL:HG13	1:D:98:ALA:HB1	1.93	0.50
1:G:67:ASP:OD1	1:H:112:LYS:NZ	2.44	0.49
1:K:70:GLU:HG2	3:K:2002:HOH:O	2.12	0.49
1:C:111:ILE:CD1	1:C:167:ILE:HG21	2.43	0.49
1:I:111:ILE:CD1	1:I:167:ILE:HG21	2.43	0.49
1:D:111:ILE:CD1	1:D:167:ILE:HG21	2.43	0.49
1:K:111:ILE:CD1	1:K:167:ILE:HG21	2.43	0.49
1:H:54:ILE:HA	1:H:57:LEU:HB3	1.94	0.48
1:L:111:ILE:CD1	1:L:167:ILE:HG21	2.43	0.48
1:A:111:ILE:CD1	1:A:167:ILE:HG21	2.44	0.48
1:O:111:ILE:CD1	1:O:167:ILE:HG21	2.43	0.48
1:F:111:ILE:CD1	1:F:167:ILE:HG21	2.43	0.48
1:M:67:ASP:O	1:N:115:ARG:NH2	2.43	0.48
1:G:111:ILE:CD1	1:G:167:ILE:HG21	2.44	0.48
1:C:14:HIS:HB3	3:C:2006:HOH:O	2.14	0.48
1:P:111:ILE:CD1	1:P:167:ILE:HG21	2.44	0.47
1:E:111:ILE:CD1	1:E:167:ILE:HG21	2.44	0.47
1:J:127:VAL:HG11	3:J:2005:HOH:O	2.14	0.47
1:O:57:LEU:O	1:O:59:GLU:N	2.48	0.47
1:E:116:ASP:OD2	1:G:19:LYS:NZ	2.40	0.47
1:J:111:ILE:CD1	1:J:167:ILE:HG21	2.44	0.47
1:H:53:SER:O	1:H:56:THR:HG22	2.15	0.46
1:B:111:ILE:CD1	1:B:167:ILE:HG21	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:111:ILE:CD1	1:M:167:ILE:HG21	2.45	0.46
1:N:111:ILE:CD1	1:N:167:ILE:HG21	2.45	0.46
1:F:53:SER:O	1:F:56:THR:HG22	2.16	0.46
1:E:115:ARG:NH2	1:H:67:ASP:O	2.48	0.46
1:M:39:GLN:OE1	1:N:39:GLN:OE1	2.34	0.46
1:E:57:LEU:O	1:E:59:GLU:N	2.50	0.45
1:F:57:LEU:O	1:F:59:GLU:N	2.50	0.45
1:H:51:ASP:OD1	1:H:53:SER:OG	2.30	0.45
1:J:19:LYS:NZ	1:L:116:ASP:OD2	2.45	0.45
1:M:124:LYS:NZ	3:M:2006:HOH:O	2.50	0.45
1:C:57:LEU:O	1:C:59:GLU:N	2.50	0.44
1:K:57:LEU:O	1:K:59:GLU:N	2.50	0.44
1:H:57:LEU:O	1:H:59:GLU:N	2.49	0.44
1:D:113:LYS:HE2	3:D:2007:HOH:O	2.17	0.44
1:D:57:LEU:O	1:D:59:GLU:N	2.51	0.44
1:A:68[B]:ARG:HE	1:A:68[B]:ARG:HB3	1.79	0.44
1:H:111:ILE:CD1	1:H:167:ILE:HG21	2.47	0.44
1:I:116:ASP:OD2	1:K:19:LYS:NZ	2.45	0.43
1:B:68:ARG:HB3	3:B:2012:HOH:O	2.18	0.43
1:C:53:SER:O	1:C:56:THR:HG22	2.17	0.43
1:E:81:GLU:OE1	1:H:20:ARG:NH2	2.49	0.43
1:N:39:GLN:OE1	1:O:39:GLN:OE1	2.37	0.43
1:D:57:LEU:O	1:D:60:ILE:N	2.51	0.42
1:J:116:ASP:HB3	1:L:19:LYS:NZ	2.33	0.42
1:F:53:SER:HB3	1:F:142:ASN:ND2	2.34	0.42
1:E:39:GLN:OE1	1:F:39:GLN:OE1	2.38	0.42
1:G:39:GLN:OE1	1:H:39:GLN:OE1	2.38	0.41
1:N:19:LYS:NZ	1:P:116:ASP:OD2	2.46	0.41
1:B:39:GLN:OE1	1:C:39:GLN:OE1	2.38	0.41
1:H:53:SER:HB3	1:H:142:ASN:ND2	2.36	0.41
1:K:57:LEU:O	1:K:60:ILE:N	2.53	0.41
1:F:8:LYS:HB3	3:F:2003:HOH:O	2.20	0.41
1:O:39:GLN:OE1	1:P:39:GLN:OE1	2.39	0.41
1:A:8:LYS:NZ	3:A:2003:HOH:O	2.54	0.41
1:C:53:SER:HB3	1:C:142:ASN:ND2	2.36	0.41
1:E:57:LEU:O	1:E:60:ILE:N	2.53	0.41
1:J:39:GLN:OE1	1:K:39:GLN:OE1	2.39	0.41
1:A:39:GLN:OE1	1:D:39:GLN:OE1	2.39	0.40
1:H:51:ASP:OD1	1:H:53:SER:CB	2.69	0.40

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%)	28	67
1	B	160/199 (80%)	151 (94%)	8 (5%)	1 (1%)	28	67
1	C	176/199 (88%)	163 (93%)	10 (6%)	3 (2%)	11	41
1	D	170/199 (85%)	162 (95%)	7 (4%)	1 (1%)	28	67
1	E	170/199 (85%)	160 (94%)	9 (5%)	1 (1%)	28	67
1	F	176/199 (88%)	163 (93%)	10 (6%)	3 (2%)	11	41
1	G	160/199 (80%)	152 (95%)	7 (4%)	1 (1%)	28	67
1	H	170/199 (85%)	157 (92%)	11 (6%)	2 (1%)	15	50
1	I	152/199 (76%)	143 (94%)	8 (5%)	1 (1%)	25	64
1	J	157/199 (79%)	148 (94%)	8 (5%)	1 (1%)	28	67
1	K	163/199 (82%)	152 (93%)	9 (6%)	2 (1%)	15	50
1	L	156/199 (78%)	147 (94%)	8 (5%)	1 (1%)	28	67
1	M	157/199 (79%)	147 (94%)	9 (6%)	1 (1%)	28	67
1	N	151/199 (76%)	142 (94%)	8 (5%)	1 (1%)	25	64
1	O	165/199 (83%)	154 (93%)	10 (6%)	1 (1%)	28	67
1	P	156/199 (78%)	147 (94%)	8 (5%)	1 (1%)	28	67
All	All	2601/3184 (82%)	2441 (94%)	138 (5%)	22 (1%)	22	61

All (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
1	C	88	ASN

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Mol	Chain	Res	Type
1	D	88	ASN
1	E	88	ASN
1	F	88	ASN
1	J	88	ASN
1	L	88	ASN
1	O	88	ASN
1	C	57	LEU
1	F	57	LEU
1	H	88	ASN
1	K	57	LEU
1	K	88	ASN
1	M	88	ASN
1	N	88	ASN
1	P	88	ASN
1	G	88	ASN
1	I	88	ASN

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%)	20	53
1	B	153/183 (84%)	142 (93%)	11 (7%)	17	48
1	C	164/183 (90%)	152 (93%)	12 (7%)	16	48
1	D	160/183 (87%)	150 (94%)	10 (6%)	21	55
1	E	160/183 (87%)	150 (94%)	10 (6%)	21	55
1	F	164/183 (90%)	152 (93%)	12 (7%)	16	48
1	G	153/183 (84%)	143 (94%)	10 (6%)	20	53
1	H	163/183 (89%)	151 (93%)	12 (7%)	16	47
1	I	146/183 (80%)	138 (94%)	8 (6%)	25	61
1	J	150/183 (82%)	141 (94%)	9 (6%)	22	57
1	K	154/183 (84%)	145 (94%)	9 (6%)	23	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	L	149/183 (81%)	140 (94%)	9 (6%)	22	57
1	M	149/183 (81%)	140 (94%)	9 (6%)	22	57
1	N	145/183 (79%)	136 (94%)	9 (6%)	21	56
1	O	156/183 (85%)	146 (94%)	10 (6%)	20	54
1	P	148/183 (81%)	140 (95%)	8 (5%)	26	61
All	All	2469/2928 (84%)	2311 (94%)	158 (6%)	20	54

All (158) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	SER
1	A	20	ARG
1	A	109	ASP
1	A	116	ASP
1	A	126	ARG
1	A	130	THR
1	A	150	LYS
1	A	159	LYS
1	A	173	ILE
1	B	1	MET
1	B	20	ARG
1	B	52	LYS
1	B	109	ASP
1	B	116	ASP
1	B	126	ARG
1	B	130	THR
1	B	150	LYS
1	B	156	VAL
1	B	159	LYS
1	B	173	ILE
1	C	1	MET
1	C	20	ARG
1	C	52	LYS
1	C	53	SER
1	C	109	ASP
1	C	116	ASP
1	C	126	ARG
1	C	130	THR
1	C	150	LYS

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Mol	Chain	Res	Type
1	C	156	VAL
1	C	159	LYS
1	C	173	ILE
1	D	1	MET
1	D	20	ARG
1	D	52	LYS
1	D	109	ASP
1	D	116	ASP
1	D	126	ARG
1	D	130	THR
1	D	150	LYS
1	D	159	LYS
1	D	173	ILE
1	E	1	MET
1	E	20	ARG
1	E	52	LYS
1	E	109	ASP
1	E	116	ASP
1	E	126	ARG
1	E	130	THR
1	E	150	LYS
1	E	159	LYS
1	E	173	ILE
1	F	1	MET
1	F	20	ARG
1	F	52	LYS
1	F	53	SER
1	F	68	ARG
1	F	109	ASP
1	F	116	ASP
1	F	126	ARG
1	F	130	THR
1	F	150	LYS
1	F	159	LYS
1	F	173	ILE
1	G	1	MET
1	G	2	SER
1	G	20	ARG
1	G	109	ASP
1	G	116	ASP
1	G	126	ARG
1	G	130	THR

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Mol	Chain	Res	Type
1	G	150	LYS
1	G	159	LYS
1	G	173	ILE
1	H	1	MET
1	H	2	SER
1	H	20	ARG
1	H	52	LYS
1	H	53	SER
1	H	109	ASP
1	H	116	ASP
1	H	126	ARG
1	H	130	THR
1	H	150	LYS
1	H	159	LYS
1	H	173	ILE
1	I	20	ARG
1	I	109	ASP
1	I	116	ASP
1	I	126	ARG
1	I	130	THR
1	I	150	LYS
1	I	159	LYS
1	I	173	ILE
1	J	20	ARG
1	J	52	LYS
1	J	109	ASP
1	J	116	ASP
1	J	126	ARG
1	J	130	THR
1	J	150	LYS
1	J	159	LYS
1	J	173	ILE
1	K	20	ARG
1	K	109	ASP
1	K	116	ASP
1	K	126	ARG
1	K	130	THR
1	K	150	LYS
1	K	156	VAL
1	K	159	LYS
1	K	173	ILE
1	L	20	ARG

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Mol	Chain	Res	Type
1	L	52	LYS
1	L	109	ASP
1	L	116	ASP
1	L	126	ARG
1	L	130	THR
1	L	150	LYS
1	L	159	LYS
1	L	173	ILE
1	M	20	ARG
1	M	52	LYS
1	M	109	ASP
1	M	116	ASP
1	M	126	ARG
1	M	130	THR
1	M	150	LYS
1	M	159	LYS
1	M	173	ILE
1	N	20	ARG
1	N	52	LYS
1	N	109	ASP
1	N	116	ASP
1	N	126	ARG
1	N	130	THR
1	N	150	LYS
1	N	159	LYS
1	N	173	ILE
1	O	20	ARG
1	O	52	LYS
1	O	109	ASP
1	O	116	ASP
1	O	126	ARG
1	O	130	THR
1	O	150	LYS
1	O	156	VAL
1	O	159	LYS
1	O	173	ILE
1	P	20	ARG
1	P	109	ASP
1	P	116	ASP
1	P	126	ARG
1	P	130	THR
1	P	150	LYS

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Mol	Chain	Res	Type
1	P	159	LYS
1	P	173	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	GLN
1	B	138	ASN
1	B	144	GLN
1	C	138	ASN
1	C	144	GLN
1	D	138	ASN
1	D	144	GLN
1	E	144	GLN
1	F	138	ASN
1	F	144	GLN
1	H	144	GLN
1	I	138	ASN
1	I	144	GLN
1	J	138	ASN
1	J	144	GLN
1	K	138	ASN
1	K	144	GLN
1	L	138	ASN
1	L	144	GLN
1	M	144	GLN
1	N	138	ASN
1	N	144	GLN
1	O	138	ASN
1	O	144	GLN
1	P	138	ASN
1	P	144	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [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%) 69 50	25, 56, 107, 149	0
1	C	178/199 (89%)	0.15	2 (1%) 80 63	22, 55, 124, 144	0
1	D	174/199 (87%)	0.21	6 (3%) 46 29	22, 58, 119, 170	0
1	E	174/199 (87%)	0.21	5 (2%) 52 33	23, 59, 109, 169	0
1	F	178/199 (89%)	0.24	9 (5%) 29 18	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%) 39 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%) 36 22	38, 68, 126, 153	0
1	K	167/199 (83%)	0.35	8 (4%) 31 19	37, 70, 128, 175	0
1	L	160/199 (80%)	0.33	7 (4%) 35 21	38, 75, 123, 157	0
1	M	161/199 (80%)	0.08	4 (2%) 58 39	37, 71, 118, 141	0
1	N	155/199 (77%)	0.12	2 (1%) 77 59	36, 66, 115, 139	0
1	O	169/199 (84%)	0.28	10 (5%) 23 14	36, 65, 131, 152	0
1	P	160/199 (80%)	0.31	7 (4%) 35 21	36, 77, 135, 152	0
All	All	2660/3184 (83%)	0.18	86 (3%) 48 30	22, 64, 122, 175	0

All (86) 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

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Mol	Chain	Res	Type	RSRZ
1	G	67	ASP	4.5
1	D	63	ALA	4.3
1	H	58	SER	4.2
1	P	70	GLU	4.1
1	L	64	ALA	4.1
1	D	57	LEU	4.0
1	P	66	LEU	4.0
1	F	57	LEU	3.9
1	H	60	ILE	3.8
1	A	67	ASP	3.7
1	F	62	GLY	3.6
1	O	59	GLU	3.6
1	L	67	ASP	3.5
1	L	144	GLN	3.5
1	O	58	SER	3.5
1	O	151	ARG	3.4
1	E	66	LEU	3.4
1	D	61	SER	3.3
1	J	66	LEU	3.2
1	C	58	SER	3.2
1	N	69	THR	3.1
1	P	64	ALA	3.1
1	B	-2	LEU	3.0
1	A	-2	LEU	2.9
1	F	58	SER	2.9
1	O	66	LEU	2.9
1	C	-3	PRO	2.9
1	E	-2	LEU	2.9
1	I	3	ARG	2.9
1	G	-2	LEU	2.8
1	H	69	THR	2.8
1	M	63	ALA	2.7
1	L	140	LYS	2.7
1	M	52	LYS	2.7
1	J	64	ALA	2.7
1	O	60	ILE	2.7
1	D	154	ALA	2.7
1	F	52	LYS	2.6
1	F	55	ASP	2.6
1	B	-3	PRO	2.6
1	F	60	ILE	2.6
1	F	59	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	O	3	ARG	2.5
1	K	66	LEU	2.5
1	L	155	ASP	2.5
1	K	64	ALA	2.5
1	H	66	LEU	2.5
1	L	141	ASN	2.5
1	F	56	THR	2.5
1	G	140	LYS	2.4
1	A	65	GLU	2.4
1	E	60	ILE	2.4
1	N	20	ARG	2.4
1	A	66	LEU	2.4
1	J	67	ASP	2.4
1	H	65	GLU	2.3
1	O	61	SER	2.3
1	P	74	LEU	2.3
1	K	61	SER	2.3
1	L	145	THR	2.3
1	O	123	PRO	2.3
1	B	139	ARG	2.3
1	F	144	GLN	2.3
1	J	82	SER	2.2
1	E	151	ARG	2.2
1	P	72	TYR	2.2
1	J	94	SER	2.2
1	M	69	THR	2.2
1	K	22	HIS	2.1
1	P	67	ASP	2.1
1	K	60	ILE	2.1
1	K	67	ASP	2.1
1	M	64	ALA	2.1
1	O	63	ALA	2.1
1	G	66	LEU	2.1
1	D	70	GLU	2.1
1	H	22	HIS	2.1
1	H	57	LEU	2.1
1	J	72	TYR	2.0
1	O	64	ALA	2.0
1	D	69	THR	2.0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	ZN	L	201	1/1	0.96	0.04	-1.85	58,58,58,58	0
2	ZN	J	201	1/1	0.97	0.04	-1.98	53,53,53,53	0
2	ZN	I	201	1/1	0.99	0.05	-2.11	56,56,56,56	0
2	ZN	A	201	1/1	0.99	0.04	-2.23	42,42,42,42	0
2	ZN	G	201	1/1	0.99	0.06	-2.60	39,39,39,39	0
2	ZN	K	201	1/1	0.99	0.04	-2.65	54,54,54,54	0
2	ZN	M	201	1/1	0.99	0.05	-2.74	59,59,59,59	0
2	ZN	B	201	1/1	0.99	0.05	-2.75	32,32,32,32	0
2	ZN	N	201	1/1	0.99	0.02	-2.76	45,45,45,45	0
2	ZN	O	201	1/1	0.99	0.03	-2.87	50,50,50,50	0
2	ZN	D	201	1/1	0.97	0.05	-3.10	33,33,33,33	0
2	ZN	E	201	1/1	0.98	0.04	-3.30	31,31,31,31	0
2	ZN	C	201	1/1	0.98	0.05	-3.70	30,30,30,30	0
2	ZN	H	201	1/1	0.99	0.03	-3.92	37,37,37,37	0
2	ZN	F	201	1/1	0.99	0.03	-6.67	35,35,35,35	0
2	ZN	P	201	1/1	0.98	0.04	-	47,47,47,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.