



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 12:41 PM UTC

PDB ID : 4FQV / pdb_00004fqv
Title : Crystal structure of broadly neutralizing antibody CR9114 bound to H7 influenza hemagglutinin
Authors : Ekiert, D.C.; Dreyfus, C.; Wilson, I.A.
Deposited on : 2012-06-25
Resolution : 5.75 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 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:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

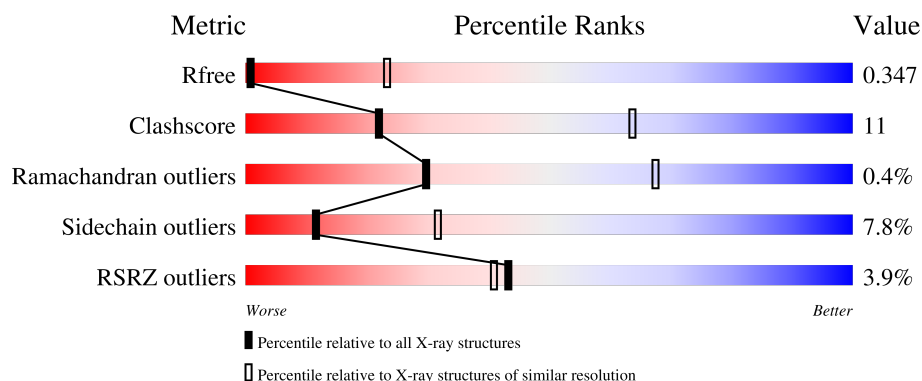
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 5.75 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}	180053	1119 (7.50-4.00)
Clashscore	190562	1182 (7.50-4.00)
Ramachandran outliers	187476	1014 (7.50-4.00)
Sidechain outliers	187428	1006 (7.54-3.98)
RSRZ outliers	180081	1112 (7.50-4.00)

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 of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	327	4% 74% 17% 5% • •
1	C	327	3% 73% 18% 5% •
1	E	327	4% 77% 16% • •
2	B	176	7% 64% 26% 7% •
2	D	176	8% 63% 29% 5% •

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Mol	Chain	Length	Quality of chain
2	F	176	10% 65% 26% 5%
3	H	224	% 75% 18%
3	I	224	2% 84% 11%
3	J	224	6% 87% 8%
4	L	216	76% 19%
4	M	216	3% 85% 10%
4	N	216	80% 16%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 20938 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	0	0
			2423	1503	435	471	14			
1	C	316	Total	C	N	O	S	0	0	0
			2423	1503	435	471	14			
1	E	317	Total	C	N	O	S	0	0	0
			2427	1505	436	472	14			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	ALA	-	expression tag	UNP Q6VMK1
A	8	ASP	-	expression tag	UNP Q6VMK1
A	9	PRO	-	expression tag	UNP Q6VMK1
A	10	GLY	-	expression tag	UNP Q6VMK1
A	254	LEU	PRO	conflict	UNP Q6VMK1
C	7	ALA	-	expression tag	UNP Q6VMK1
C	8	ASP	-	expression tag	UNP Q6VMK1
C	9	PRO	-	expression tag	UNP Q6VMK1
C	10	GLY	-	expression tag	UNP Q6VMK1
C	254	LEU	PRO	conflict	UNP Q6VMK1
E	7	ALA	-	expression tag	UNP Q6VMK1
E	8	ASP	-	expression tag	UNP Q6VMK1
E	9	PRO	-	expression tag	UNP Q6VMK1
E	10	GLY	-	expression tag	UNP Q6VMK1
E	254	LEU	PRO	conflict	UNP Q6VMK1

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	170	Total	C	N	O	S	0	0	0
			1380	851	243	279	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	171	Total	C	N	O	S	0	0	0
			1388	857	244	280	7			
2	F	169	Total	C	N	O	S	0	0	0
			1369	845	239	278	7			

- Molecule 3 is a protein called Antibody CR9114 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			
3	I	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			
3	J	215	Total	C	N	O	S	0	2	0
			1608	1013	268	320	7			

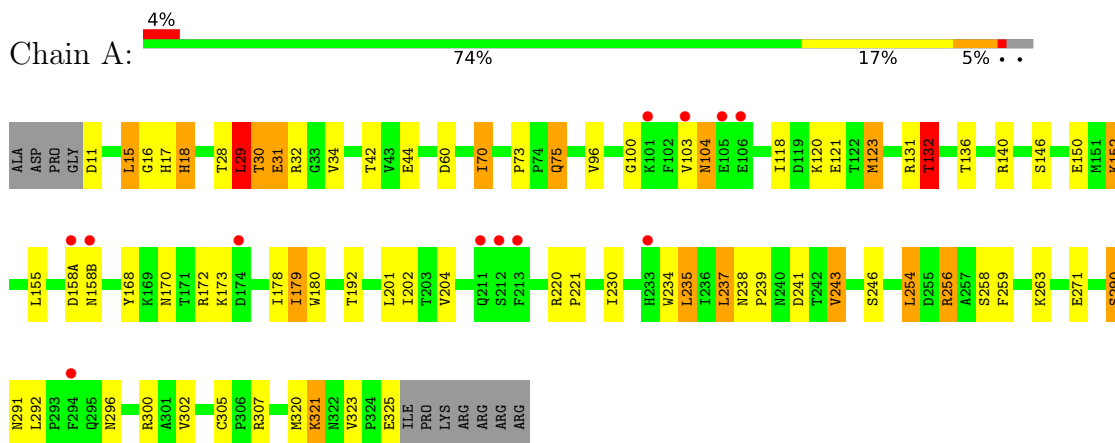
- Molecule 4 is a protein called Antibody CR9114 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			
4	M	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			
4	N	211	Total	C	N	O	S	0	0	0
			1568	978	266	320	4			

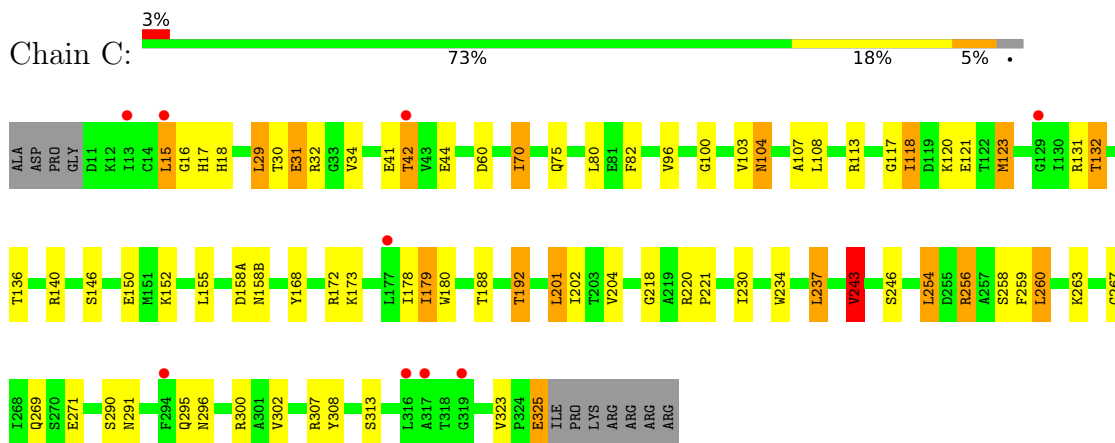
3 Residue-property plots [i](#)

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 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.

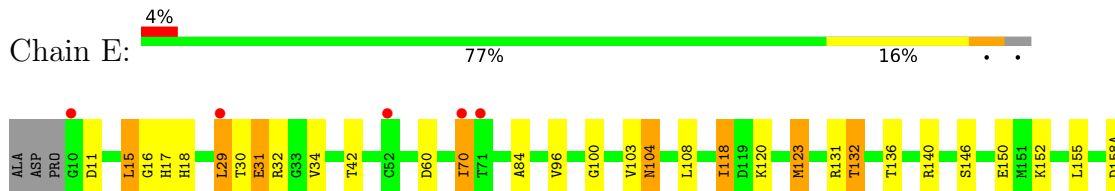
- Molecule 1: Hemagglutinin HA1

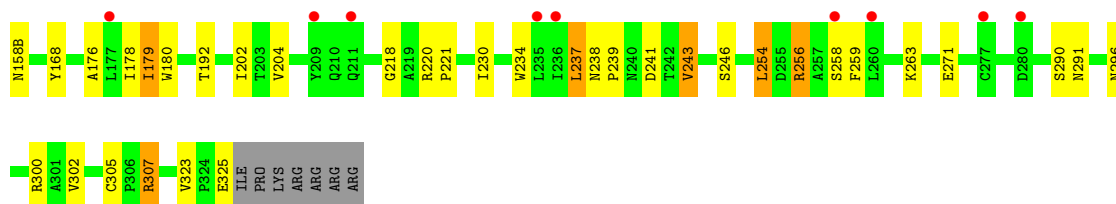


- Molecule 1: Hemagglutinin HA1

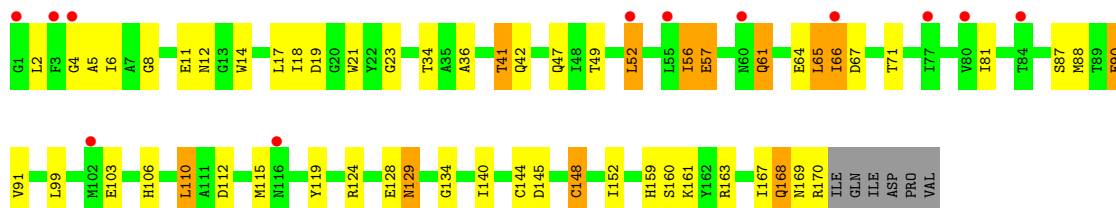


- Molecule 1: Hemagglutinin HA1

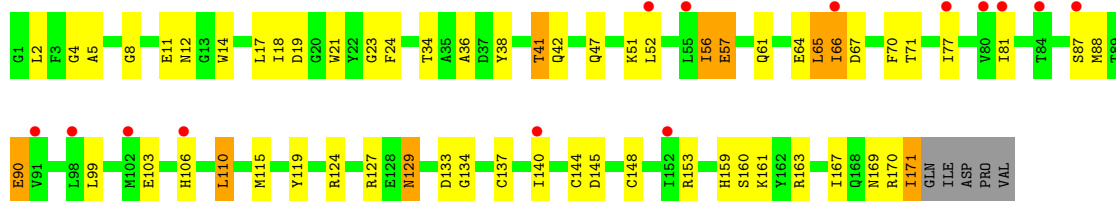




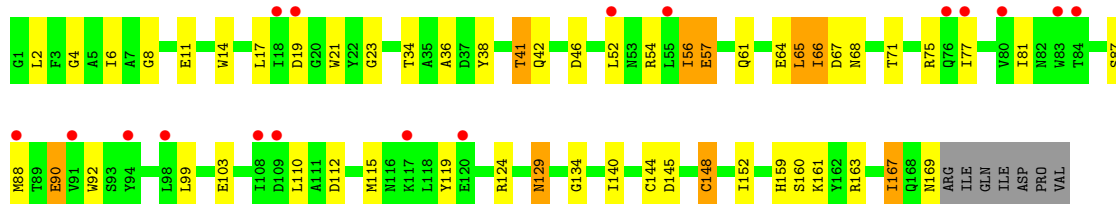
• Molecule 2: Hemagglutinin HA2



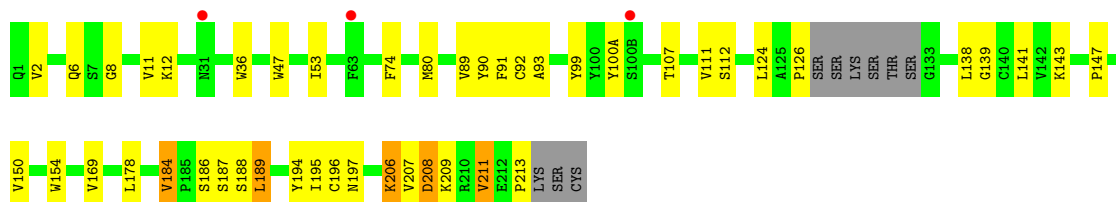
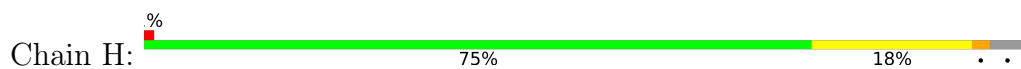
• Molecule 2: Hemagglutinin HA2



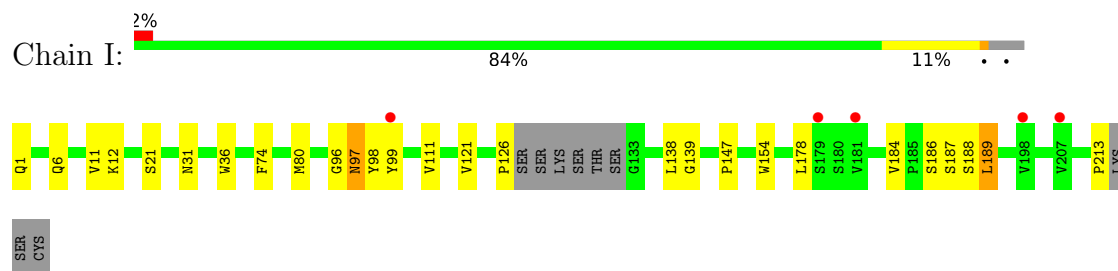
• Molecule 2: Hemagglutinin HA2



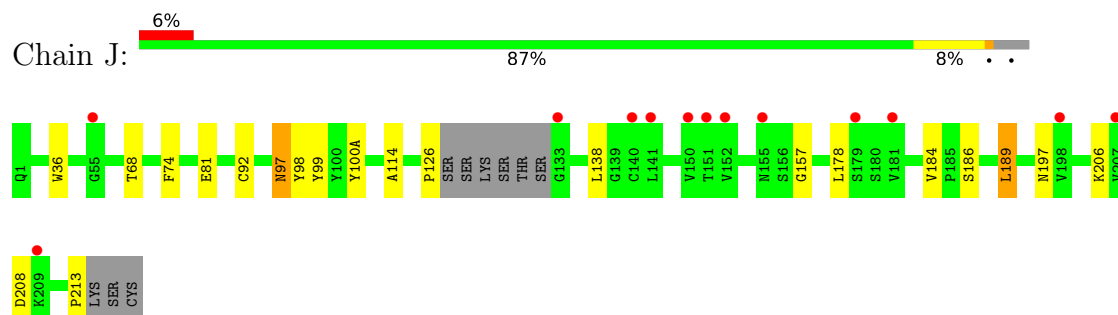
• Molecule 3: Antibody CR9114 heavy chain



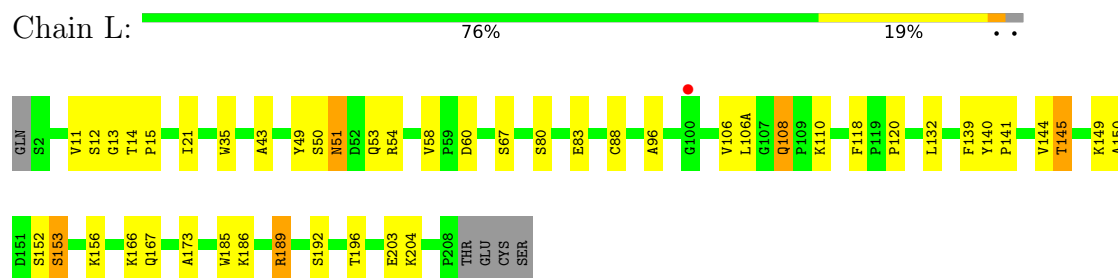
- Molecule 3: Antibody CR9114 heavy chain



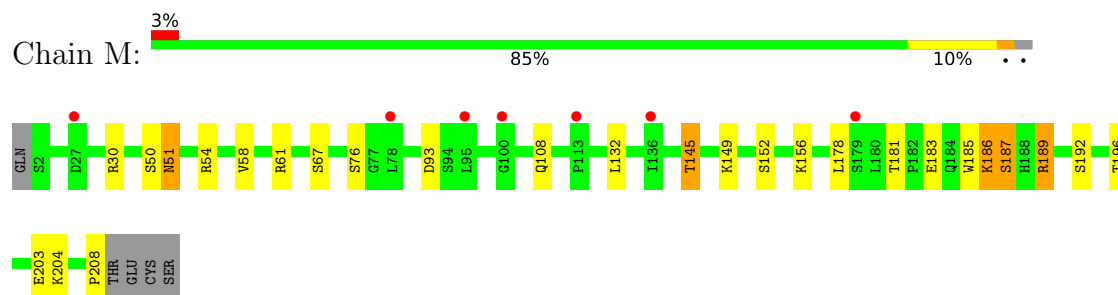
- Molecule 3: Antibody CR9114 heavy chain



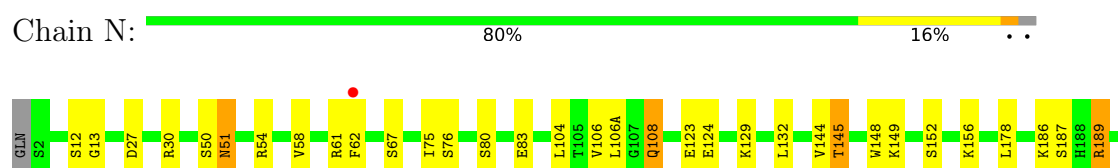
- Molecule 4: Antibody CR9114 light chain



- Molecule 4: Antibody CR9114 light chain



- Molecule 4: Antibody CR9114 light chain



S192					
T196					
H197					
E203					
K204					
P208					
THR					
GLU					
CYS					
SER					

4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	197.57Å 197.57Å 223.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.35 – 5.75 49.35 – 5.75	Depositor EDS
% Data completeness (in resolution range)	99.4 (49.35-5.75) 98.7 (49.35-5.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 5.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.314 , 0.339 0.323 , 0.347	Depositor DCC
R_{free} test set	642 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	319.3	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 390.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	20938	wwPDB-VP
Average B, all atoms (Å ²)	244.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.52% 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

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	0.44	0/2468	0.82	3/3336 (0.1%)
1	C	0.45	0/2468	0.85	2/3336 (0.1%)
1	E	0.44	0/2472	0.83	3/3341 (0.1%)
2	B	0.41	0/1404	0.76	1/1892 (0.1%)
2	D	0.47	0/1412	0.80	0/1903
2	F	0.43	0/1393	0.79	2/1878 (0.1%)
3	H	0.49	0/1647	1.12	12/2245 (0.5%)
3	I	0.43	0/1647	0.82	4/2245 (0.2%)
3	J	0.49	0/1647	0.84	2/2245 (0.1%)
4	L	0.45	0/1606	0.90	5/2193 (0.2%)
4	M	0.48	0/1606	0.83	1/2193 (0.0%)
4	N	0.43	0/1606	0.87	3/2193 (0.1%)
All	All	0.45	0/21376	0.86	38/29000 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
1	E	0	1
2	B	0	1
2	D	0	1
2	F	0	1
4	M	0	1
All	All	0	7

There are no bond length outliers.

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	153	SER	CA-C-N	-11.41	105.58	119.84
4	L	153	SER	C-N-CA	-11.41	105.58	119.84
3	H	92	CYS	N-CA-C	-10.22	93.14	109.59
3	I	97	ASN	N-CA-C	9.60	120.54	108.45
4	N	108	GLN	CA-C-N	9.09	128.84	119.76

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	LEU	Peptide
2	B	56	ILE	Peptide
1	C	29	LEU	Peptide
2	D	56	ILE	Peptide
1	E	29	LEU	Peptide

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	2423	0	2373	65	0
1	C	2423	0	2373	57	0
1	E	2427	0	2376	53	0
2	B	1380	0	1284	53	0
2	D	1388	0	1295	58	0
2	F	1369	0	1271	49	0
3	H	1608	0	1556	63	0
3	I	1608	0	1556	35	0
3	J	1608	0	1556	36	2
4	L	1568	0	1517	37	2
4	M	1568	0	1517	15	1
4	N	1568	0	1517	30	0
All	All	20938	0	20191	439	3

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 11.

The worst 5 of 439 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)
2:B:19:ASP:OD2	3:H:100(A):TYR:OH	1.65	1.11
1:C:123:MET:SD	1:C:254:LEU:HD13	2.00	1.01
3:H:11:VAL:CG2	3:H:147:PRO:HG3	1.94	0.97
1:C:123:MET:HE3	1:C:168:TYR:CG	2.01	0.96
1:E:123:MET:SD	1:E:254:LEU:CD1	2.55	0.95

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:189:ARG:NH2	3:J:157:GLY:O[6_554]	1.99	0.21
4:L:189:ARG:NE	3:J:157:GLY:O[6_554]	2.10	0.10
4:M:187:SER:OG	4:M:187:SER:OG[2_465]	2.14	0.06

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	314/327 (96%)	295 (94%)	19 (6%)	0	100	100
1	C	314/327 (96%)	297 (95%)	17 (5%)	0	100	100
1	E	315/327 (96%)	296 (94%)	19 (6%)	0	100	100
2	B	168/176 (96%)	151 (90%)	14 (8%)	3 (2%)	6	34
2	D	169/176 (96%)	152 (90%)	14 (8%)	3 (2%)	6	34
2	F	167/176 (95%)	151 (90%)	14 (8%)	2 (1%)	10	43
3	H	213/224 (95%)	210 (99%)	3 (1%)	0	100	100
3	I	213/224 (95%)	212 (100%)	1 (0%)	0	100	100
3	J	213/224 (95%)	212 (100%)	1 (0%)	0	100	100
4	L	209/216 (97%)	204 (98%)	4 (2%)	1 (0%)	24	63
4	M	209/216 (97%)	206 (99%)	2 (1%)	1 (0%)	24	63

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	N	209/216 (97%)	205 (98%)	3 (1%)	1 (0%)	24	63
All	All	2713/2829 (96%)	2591 (96%)	111 (4%)	11 (0%)	30	67

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	57	GLU
2	D	57	GLU
2	F	57	GLU
2	B	119	TYR
2	D	119	TYR

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	268/277 (97%)	232 (87%)	36 (13%)	4	14
1	C	268/277 (97%)	233 (87%)	35 (13%)	4	15
1	E	268/277 (97%)	236 (88%)	32 (12%)	5	17
2	B	145/151 (96%)	127 (88%)	18 (12%)	4	16
2	D	146/151 (97%)	128 (88%)	18 (12%)	4	17
2	F	144/151 (95%)	127 (88%)	17 (12%)	5	18
3	H	180/187 (96%)	176 (98%)	4 (2%)	45	64
3	I	180/187 (96%)	178 (99%)	2 (1%)	65	75
3	J	180/187 (96%)	178 (99%)	2 (1%)	65	75
4	L	175/180 (97%)	169 (97%)	6 (3%)	32	54
4	M	175/180 (97%)	170 (97%)	5 (3%)	37	57
4	N	175/180 (97%)	170 (97%)	5 (3%)	37	57
All	All	2304/2385 (97%)	2124 (92%)	180 (8%)	11	32

5 of 180 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	70	ILE
2	F	52	LEU
1	E	123	MET
1	E	243	VAL
2	F	110	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
4	N	37	GLN
3	J	192	GLN
3	H	204	ASN
4	M	197	HIS
3	H	192	GLN

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 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	316/327 (96%)	0.08	12 (3%) 44 41	178, 369, 432, 457	0
1	C	316/327 (96%)	0.07	9 (2%) 55 47	145, 357, 410, 426	0
1	E	317/327 (96%)	0.09	14 (4%) 39 38	166, 309, 394, 402	0
2	B	170/176 (96%)	0.22	12 (7%) 22 27	175, 222, 260, 276	0
2	D	171/176 (97%)	0.45	14 (8%) 17 24	168, 211, 267, 300	0
2	F	169/176 (96%)	0.38	17 (10%) 12 19	166, 217, 262, 274	0
3	H	215/224 (95%)	-0.07	3 (1%) 73 62	110, 233, 322, 341	2 (0%)
3	I	215/224 (95%)	-0.01	5 (2%) 61 52	65, 159, 190, 206	2 (0%)
3	J	215/224 (95%)	0.09	13 (6%) 27 32	52, 143, 198, 232	2 (0%)
4	L	211/216 (97%)	-0.30	1 (0%) 87 77	218, 272, 365, 402	0
4	M	211/216 (97%)	-0.05	7 (3%) 49 44	159, 188, 242, 274	0
4	N	211/216 (97%)	-0.33	1 (0%) 87 77	116, 159, 194, 204	0
All	All	2737/2829 (96%)	0.04	108 (3%) 43 41	52, 223, 397, 457	6 (0%)

The worst 5 of 108 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	212	SER	7.4
2	D	98	LEU	7.4
1	E	177	LEU	7.2
2	D	80	VAL	7.1
1	C	317	ALA	6.7

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.