



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:30 PM GMT

PDB ID : 4WHT
Title : Structure of the Hepatitis C virus envelope glycoprotein E2 antigenic region
412-423 bound to the broadly neutralizing antibody 3/11, P1 crystal form
Authors : Krey, T.; Rey, F.A.
Deposited on : 2014-09-23
Resolution : 2.22 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

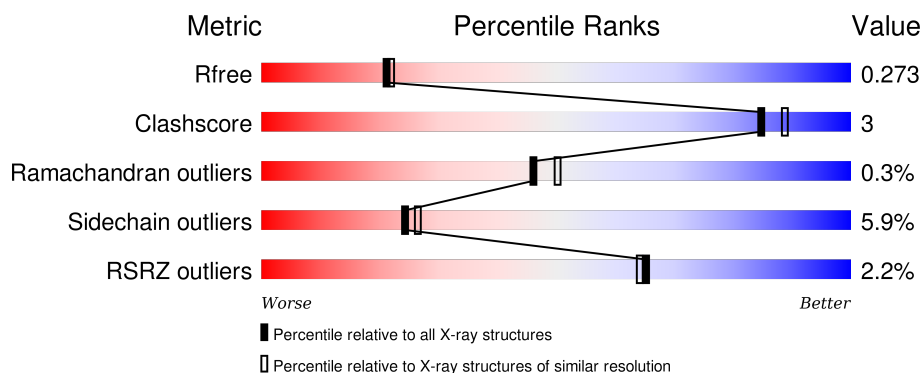
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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}	91344	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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	252	<div> <div></div> <div>73% 9% 18%</div> </div>
1	C	252	<div> <div></div> <div>69% 12% 19%</div> </div>
1	E	252	<div> <div></div> <div>73% 9% 17%</div> </div>
1	G	252	<div> <div></div> <div>73% 9% 18%</div> </div>
1	I	252	<div> <div></div> <div>71% 11% 19%</div> </div>




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Mol	Chain	Length	Quality of chain
1	K	252	
1	M	252	
1	O	252	
1	Q	252	
1	S	252	
1	U	252	
1	X	252	
2	B	220	
2	D	220	
2	F	220	
2	H	220	
2	J	220	
2	L	220	
2	N	220	
2	P	220	
2	R	220	
2	T	220	
2	V	220	
2	Y	220	
3	a	12	
3	c	12	
3	e	12	
3	g	12	
3	i	12	
3	k	12	

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Mol	Chain	Length	Quality of chain
3	m	12	 83%17%
3	o	12	 92%8%
3	q	12	 83%17%
3	s	12	 92%8%
3	u	12	 83%17%
3	x	12	 92%8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 40099 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 Heavy chain of the Fab fragment derived from neutralizing antibody 3/11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1565	988	264	306	7			
1	C	205	Total	C	N	O	S	0	0	0
			1554	982	262	303	7			
1	E	208	Total	C	N	O	S	0	0	0
			1570	991	265	307	7			
1	G	207	Total	C	N	O	S	0	0	0
			1567	989	264	307	7			
1	I	205	Total	C	N	O	S	0	0	0
			1550	980	261	302	7			
1	K	204	Total	C	N	O	S	0	0	0
			1546	978	260	301	7			
1	M	204	Total	C	N	O	S	0	0	0
			1546	978	260	301	7			
1	O	205	Total	C	N	O	S	0	0	0
			1550	980	261	302	7			
1	Q	206	Total	C	N	O	S	0	0	0
			1557	984	262	304	7			
1	S	204	Total	C	N	O	S	0	0	0
			1546	978	260	301	7			
1	U	205	Total	C	N	O	S	0	0	0
			1554	982	262	303	7			
1	X	199	Total	C	N	O	S	0	0	0
			1495	945	250	293	7			

- Molecule 2 is a protein called Light chain of the Fab fragment derived from neutralizing antibody 3/11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	D	216	Total	C	N	O	S	0	0	0
			1642	1018	275	342	7			
2	F	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	H	216	Total	C	N	O	S	0	0	0
			1640	1017	274	342	7			
2	L	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	N	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	P	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	R	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	T	215	Total	C	N	O	S	0	0	0
			1634	1014	273	340	7			
2	V	212	Total	C	N	O	S	0	0	0
			1615	1004	270	334	7			
2	Y	210	Total	C	N	O	S	0	0	0
			1591	987	266	331	7			

- Molecule 3 is a protein called Epitope peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	a	12	Total	C	N	O	0	0	0
			97	60	19	18			
3	c	11	Total	C	N	O	0	0	0
			89	56	17	16			
3	e	11	Total	C	N	O	0	0	0
			89	56	17	16			
3	g	12	Total	C	N	O	0	0	0
			97	60	19	18			
3	i	12	Total	C	N	O	0	0	0
			97	60	19	18			
3	k	11	Total	C	N	O	0	0	0
			89	56	17	16			
3	m	12	Total	C	N	O	0	0	0
			97	60	19	18			
3	o	11	Total	C	N	O	0	0	0
			88	55	17	16			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	q	10	Total	C	N	O	0	0	0
			80	51	15	14			
3	s	11	Total	C	N	O	0	0	0
			89	56	17	16			
3	u	10	Total	C	N	O	0	0	0
			80	51	15	14			
3	x	12	Total	C	N	O	0	0	0
			97	60	19	18			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	38	Total	O	0	0
			38	38		
4	C	44	Total	O	0	0
			44	44		
4	E	34	Total	O	0	0
			34	34		
4	G	29	Total	O	0	0
			29	29		
4	I	31	Total	O	0	0
			31	31		
4	J	55	Total	O	0	0
			55	55		
4	K	13	Total	O	0	0
			13	13		
4	M	18	Total	O	0	0
			18	18		
4	O	26	Total	O	0	0
			26	26		
4	Q	20	Total	O	0	0
			20	20		
4	S	27	Total	O	0	0
			27	27		
4	U	14	Total	O	0	0
			14	14		
4	X	42	Total	O	0	0
			42	42		
4	B	47	Total	O	0	0
			47	47		
4	D	62	Total	O	0	0
			62	62		

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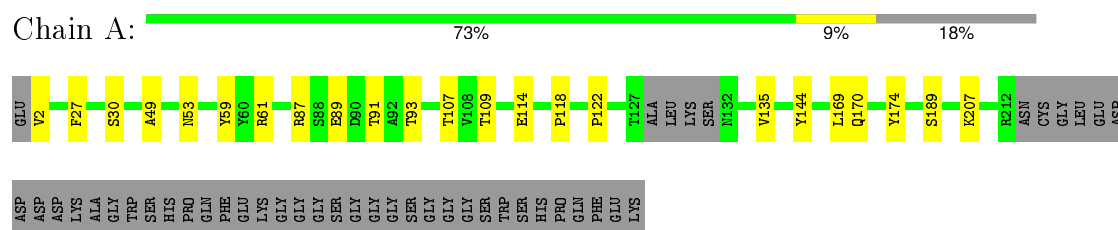
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	F	48	Total 48	O 48	0	0
4	H	51	Total 51	O 51	0	0
4	L	46	Total 46	O 46	0	0
4	N	24	Total 24	O 24	0	0
4	P	57	Total 57	O 57	0	0
4	R	29	Total 29	O 29	0	0
4	T	32	Total 32	O 32	0	0
4	V	21	Total 21	O 21	0	0
4	Y	40	Total 40	O 40	0	0
4	a	1	Total 1	O 1	0	0
4	s	1	Total 1	O 1	0	0

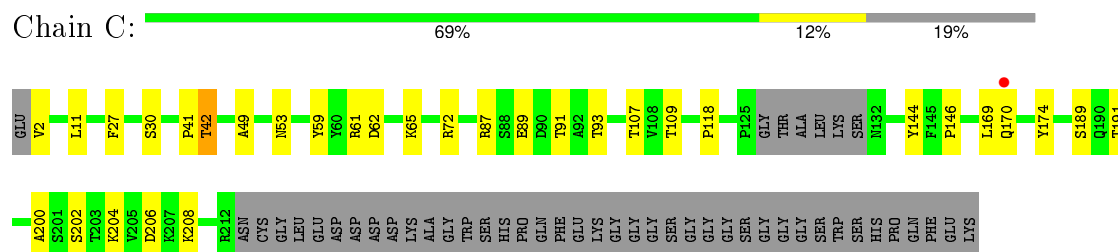
3 Residue-property plots [i](#)

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

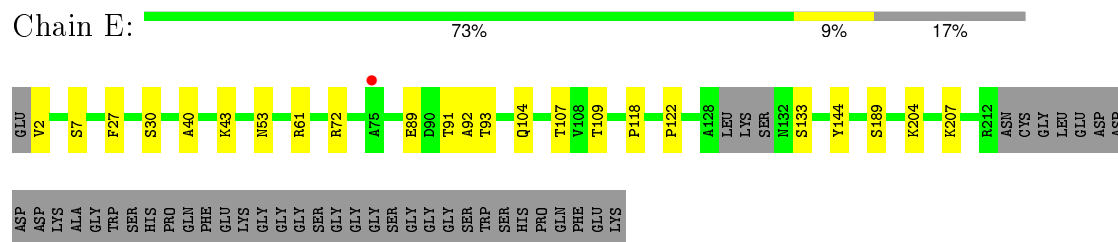
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



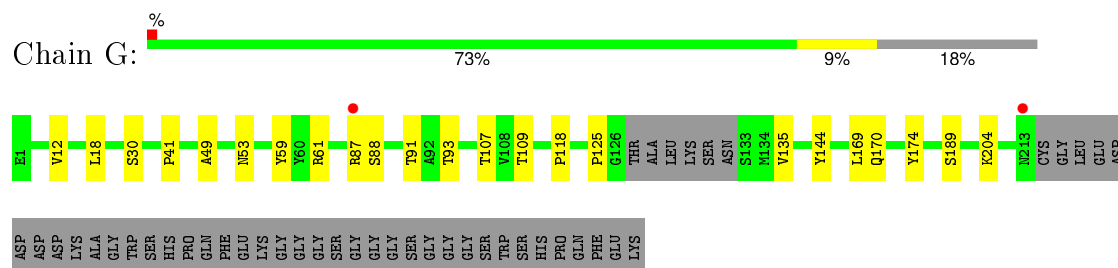
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



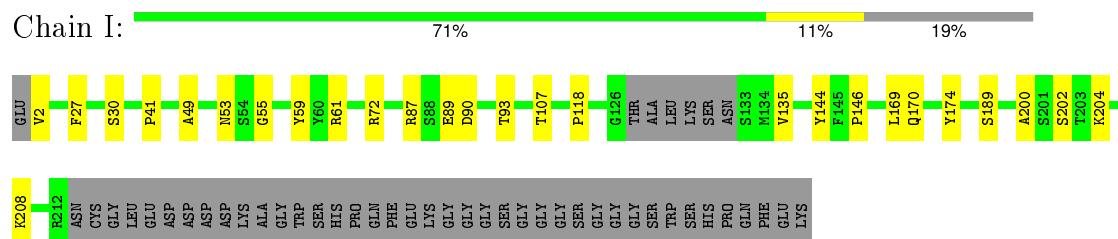
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



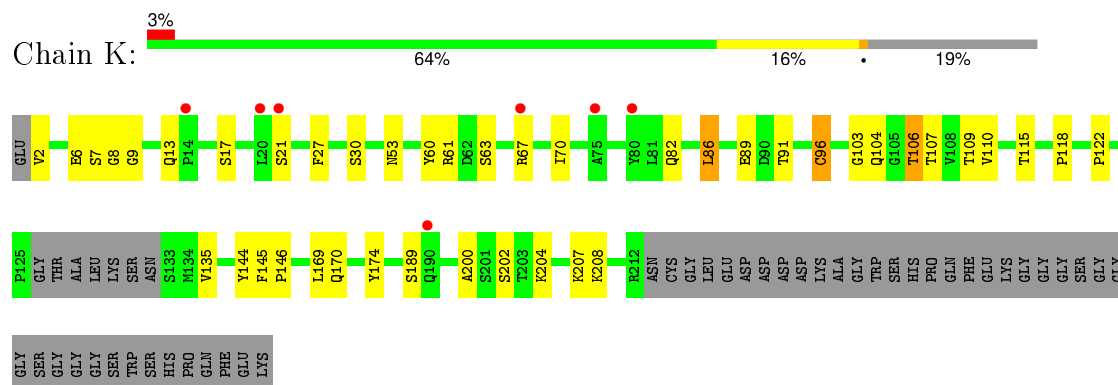
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



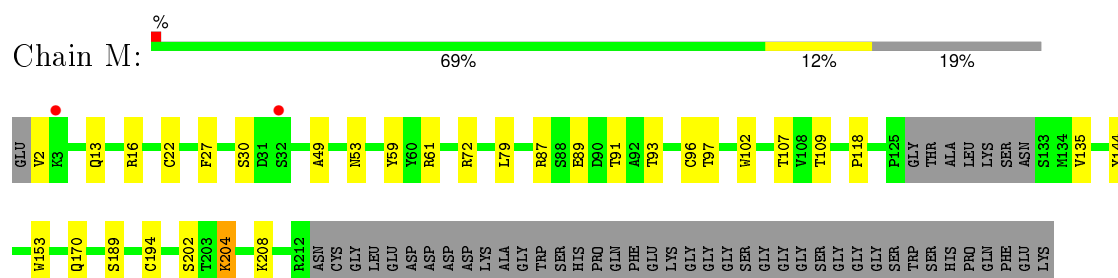
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



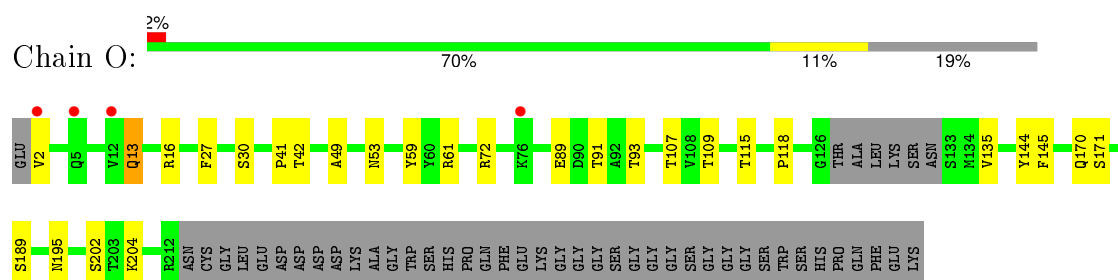
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



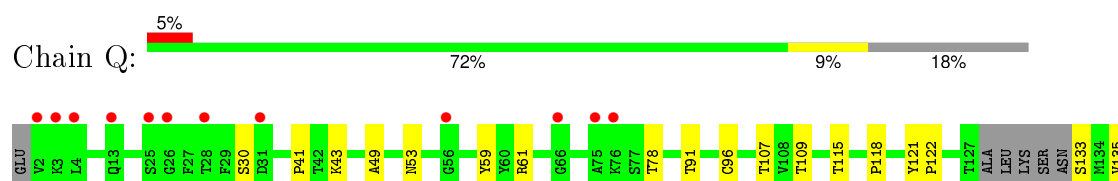
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11

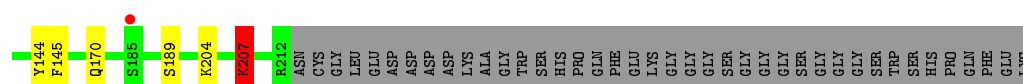


- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11

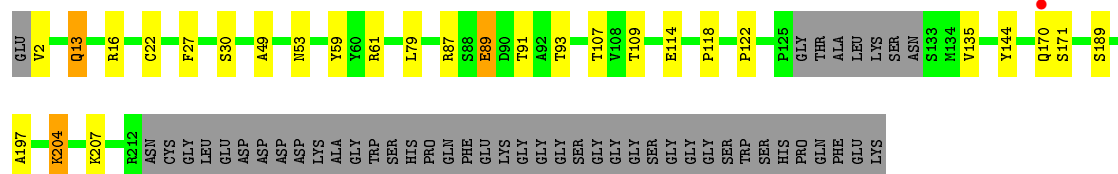


- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11

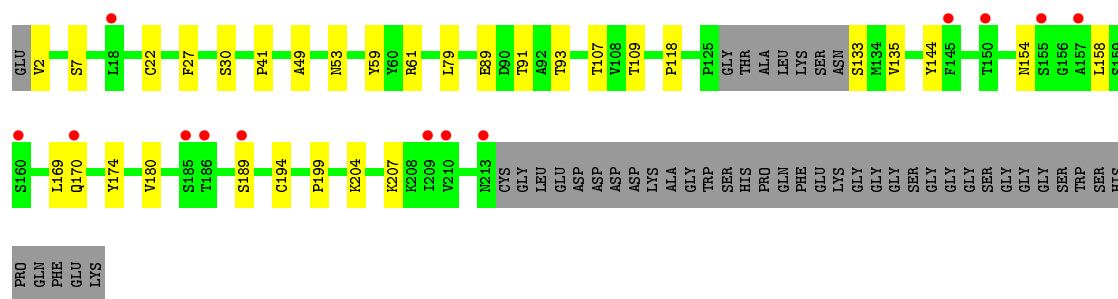




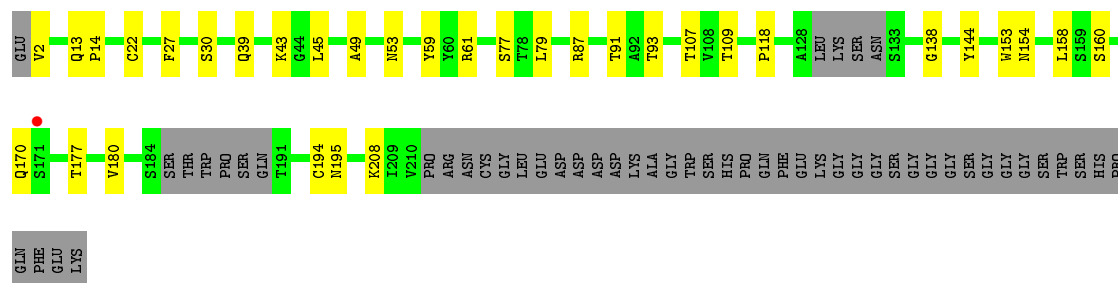
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



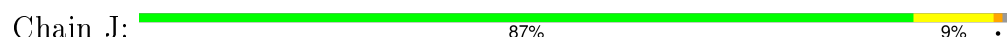
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



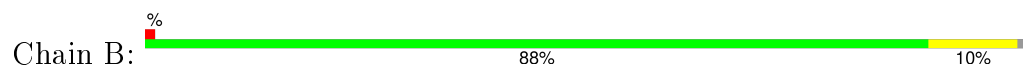
- Molecule 1: Heavy chain of the Fab fragment derived from neutralizing antibody 3/11



- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11

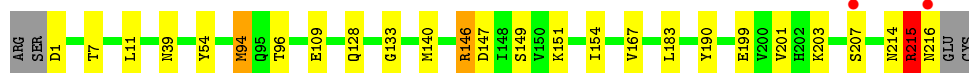
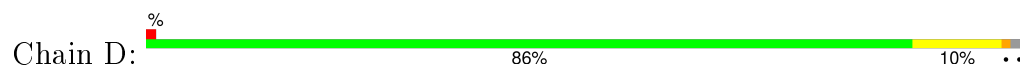


- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11

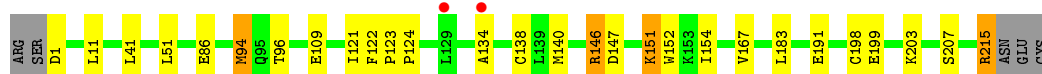
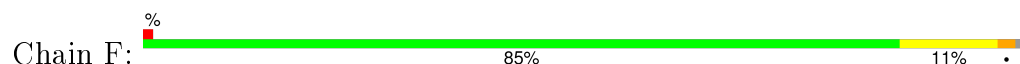




- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11



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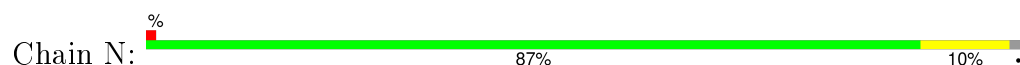
- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11



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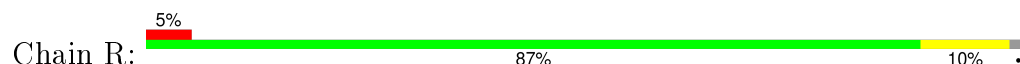
- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11

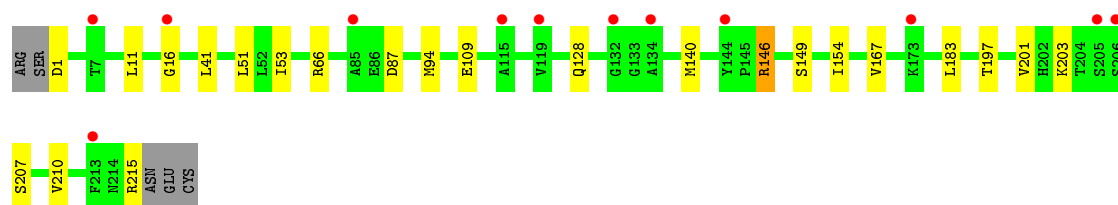


- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11

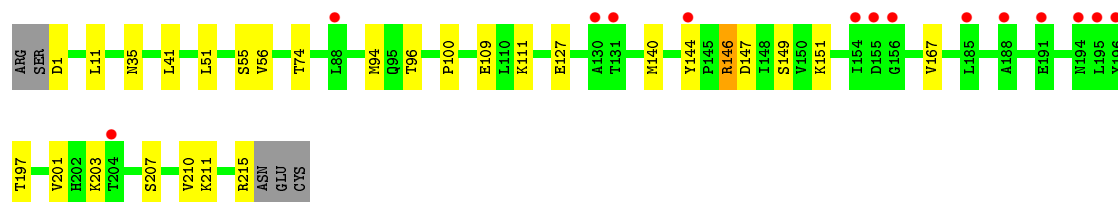
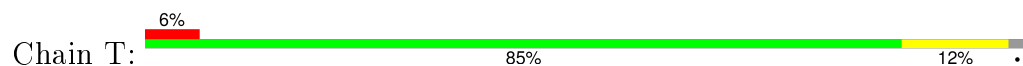


- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11

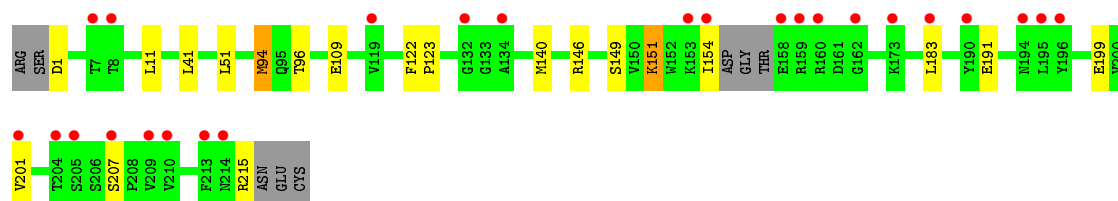
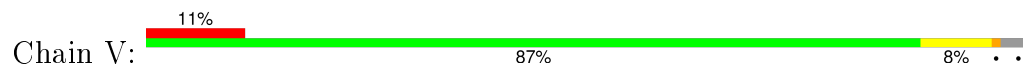




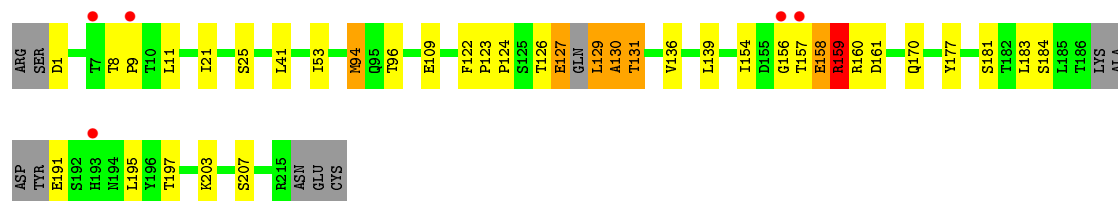
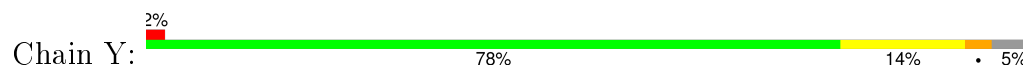
- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11



- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11



- Molecule 2: Light chain of the Fab fragment derived from neutralizing antibody 3/11



- Molecule 3: Epitope peptide



- Molecule 3: Epitope peptide





- Molecule 3: Epitope peptide

Chain e: 83% 8% 8%



- Molecule 3: Epitope peptide

Chain g: 100%

There are no outlier residues recorded for this chain.

- Molecule 3: Epitope peptide

Chain i: 92% 8%



- Molecule 3: Epitope peptide

Chain k: 92% 8%



- Molecule 3: Epitope peptide

Chain m: 83% 17%



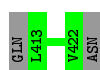
- Molecule 3: Epitope peptide

Chain o: 92% 8%



- Molecule 3: Epitope peptide

Chain q: 83% 17%




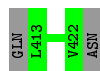
- Molecule 3: Epitope peptide

Chain s:  92% 8%



- Molecule 3: Epitope peptide

Chain u:  83% 17%



- Molecule 3: Epitope peptide

Chain x:  92% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	64.79Å 128.22Å 163.63Å 88.79° 94.36° 96.15°	Depositor
Resolution (Å)	32.50 – 2.22 48.74 – 2.22	Depositor EDS
% Data completeness (in resolution range)	95.4 (32.50-2.22) 95.2 (48.74-2.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.22Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.209 , 0.243 0.232 , 0.273	Depositor DCC
R_{free} test set	12374 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	33.1	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 246060 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	40099	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.51	0/1603	0.76	0/2188
1	C	0.52	0/1592	0.75	0/2173
1	E	0.51	0/1608	0.74	0/2195
1	G	0.50	0/1605	0.75	0/2190
1	I	0.50	0/1588	0.75	1/2167 (0.0%)
1	K	0.51	0/1584	0.74	0/2162
1	M	0.47	0/1584	0.73	0/2162
1	O	0.49	0/1588	0.74	0/2167
1	Q	0.48	0/1595	0.73	1/2177 (0.0%)
1	S	0.53	0/1584	0.75	0/2162
1	U	0.49	0/1592	0.75	0/2173
1	X	0.56	0/1528	0.82	0/2082
2	B	0.53	0/1664	0.74	0/2264
2	D	0.52	0/1672	0.74	0/2275
2	F	0.51	0/1664	0.72	0/2264
2	H	0.51	0/1670	0.73	0/2272
2	J	0.51	0/1664	0.73	0/2264
2	L	0.51	0/1664	0.72	0/2264
2	N	0.44	0/1664	0.71	0/2264
2	P	0.48	0/1664	0.73	1/2264 (0.0%)
2	R	0.45	0/1664	0.71	0/2264
2	T	0.50	0/1664	0.73	0/2264
2	V	0.47	0/1644	0.69	0/2235
2	Y	0.61	0/1618	0.88	2/2199 (0.1%)
3	a	0.51	0/99	0.67	0/135
3	c	0.48	0/91	0.71	0/124
3	e	0.57	0/91	0.71	0/124
3	g	0.56	0/99	0.70	0/135
3	i	0.50	0/99	0.75	0/135
3	k	0.55	0/91	0.78	0/124
3	m	0.45	0/99	0.75	0/135
3	o	0.45	0/90	0.64	0/123
3	q	0.46	0/82	0.73	0/112
3	s	0.52	0/91	0.74	0/124

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
3	u	0.49	0/82	0.73	0/112
3	x	0.67	0/99	0.82	0/135
All	All	0.51	0/40080	0.74	5/54609 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Y	130	ALA	C-N-CA	5.87	136.37	121.70
2	P	94	MET	N-CA-CB	-5.52	100.66	110.60
1	I	55	GLY	N-CA-C	5.43	126.67	113.10
1	Q	207	LYS	CG-CD-CE	-5.22	96.25	111.90
2	Y	158	GLU	C-N-CA	5.04	134.30	121.70

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	A	1565	0	1536	7	0
1	C	1554	0	1526	14	0
1	E	1570	0	1541	7	0
1	G	1567	0	1538	6	0
1	I	1550	0	1523	10	0
1	K	1546	0	1520	18	0
1	M	1546	0	1520	12	0
1	O	1550	0	1523	9	0
1	Q	1557	0	1530	7	0
1	S	1546	0	1520	12	0
1	U	1554	0	1526	10	0
1	X	1495	0	1472	13	0
2	B	1634	0	1593	6	0
2	D	1642	0	1599	15	0
2	F	1634	0	1593	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	1640	0	1598	6	0
2	J	1634	0	1593	10	0
2	L	1634	0	1593	7	0
2	N	1634	0	1593	8	0
2	P	1634	0	1593	10	0
2	R	1634	0	1593	8	0
2	T	1634	0	1593	9	0
2	V	1615	0	1578	8	0
2	Y	1591	0	1552	11	0
3	a	97	0	88	0	0
3	c	89	0	82	0	0
3	e	89	0	82	0	0
3	g	97	0	88	0	0
3	i	97	0	88	0	0
3	k	89	0	82	0	0
3	m	97	0	88	0	0
3	o	88	0	80	0	0
3	q	80	0	74	0	0
3	s	89	0	82	0	0
3	u	80	0	74	0	0
3	x	97	0	88	0	0
4	A	38	0	0	0	0
4	B	47	0	0	0	0
4	C	44	0	0	2	0
4	D	62	0	0	1	0
4	E	34	0	0	0	0
4	F	48	0	0	0	0
4	G	29	0	0	0	0
4	H	51	0	0	0	0
4	I	31	0	0	0	0
4	J	55	0	0	0	0
4	K	13	0	0	0	0
4	L	46	0	0	0	0
4	M	18	0	0	1	0
4	N	24	0	0	0	0
4	O	26	0	0	0	0
4	P	57	0	0	0	0
4	Q	20	0	0	0	0
4	R	29	0	0	1	0
4	S	27	0	0	1	0
4	T	32	0	0	1	0
4	U	14	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	V	21	0	0	0	0
4	X	42	0	0	0	0
4	Y	40	0	0	0	0
4	a	1	0	0	0	0
4	s	1	0	0	0	0
All	All	40099	0	38342	223	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 (223) 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:M:97:THR:HG21	1:M:102:TRP:CE3	2.13	0.83
1:C:202:SER:HB3	1:I:208:LYS:HB3	1.65	0.76
2:J:44:ARG:NH1	2:J:86:GLU:O	2.23	0.71
1:K:208:LYS:HB3	1:M:202:SER:HB3	1.75	0.66
2:D:214:ASN:O	2:D:215:ARG:HB2	1.95	0.66
2:L:154:ILE:HD11	2:L:183:LEU:HD21	1.78	0.66
2:J:160:ARG:HH21	2:J:163:VAL:HG11	1.62	0.65
2:F:191:GLU:HA	2:F:215:ARG:HH22	1.62	0.65
2:V:191:GLU:HA	2:V:215:ARG:HH22	1.62	0.64
2:Y:127:GLU:O	2:Y:129:LEU:N	2.31	0.63
1:G:30:SER:O	1:G:53:ASN:HB2	2.00	0.62
1:U:118:PRO:HB3	1:U:144:TYR:HB3	1.83	0.61
1:M:97:THR:HG21	1:M:102:TRP:CD2	2.37	0.60
1:G:118:PRO:HB3	1:G:144:TYR:HB3	1.84	0.59
1:I:118:PRO:HB3	1:I:144:TYR:HB3	1.84	0.59
1:S:118:PRO:HB3	1:S:144:TYR:HB3	1.83	0.59
1:S:171:SER:HB3	2:Y:159:ARG:HG3	1.84	0.58
1:Q:30:SER:O	1:Q:53:ASN:HB2	2.04	0.57
1:M:30:SER:O	1:M:53:ASN:HB2	2.04	0.57
1:O:30:SER:O	1:O:53:ASN:HB2	2.03	0.57
1:Q:118:PRO:HB3	1:Q:144:TYR:HB3	1.87	0.57
1:X:118:PRO:HB3	1:X:144:TYR:HB3	1.86	0.56
2:T:197:THR:HG23	2:T:210:VAL:HG13	1.88	0.56
2:N:154:ILE:HD11	2:N:183:LEU:HD21	1.88	0.56
1:U:30:SER:O	1:U:53:ASN:HB2	2.06	0.56
1:A:118:PRO:HB3	1:A:144:TYR:HB3	1.88	0.55
1:O:118:PRO:HB3	1:O:144:TYR:HB3	1.88	0.55
1:S:30:SER:O	1:S:53:ASN:HB2	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:118:PRO:HB3	1:K:144:TYR:HB3	1.89	0.55
1:I:30:SER:O	1:I:53:ASN:HB2	2.06	0.55
1:C:208:LYS:HB3	1:I:202:SER:HB3	1.88	0.55
1:E:118:PRO:HB3	1:E:144:TYR:HB3	1.87	0.55
1:X:30:SER:O	1:X:53:ASN:HB2	2.06	0.55
1:X:22:CYS:HB3	1:X:79:LEU:HB3	1.89	0.54
2:N:146:ARG:HD3	2:N:167:VAL:HG11	1.89	0.54
2:B:146:ARG:HD3	2:B:167:VAL:HG11	1.90	0.54
1:O:202:SER:HB3	1:X:208:LYS:HB3	1.89	0.54
1:C:206:ASP:HB2	1:I:204:LYS:HB2	1.88	0.54
2:D:94:MET:CE	2:D:96:THR:HG23	2.38	0.54
2:D:154:ILE:HD11	2:D:183:LEU:HD21	1.89	0.54
1:C:118:PRO:HB3	1:C:144:TYR:HB3	1.89	0.54
2:T:55:SER:O	2:T:56:VAL:HG12	2.07	0.54
1:A:30:SER:O	1:A:53:ASN:HB2	2.07	0.53
2:D:190:TYR:CE2	2:D:216:ASN:CB	2.91	0.53
2:R:197:THR:HG23	2:R:210:VAL:HG13	1.91	0.53
2:D:190:TYR:CE2	2:D:216:ASN:HB3	2.44	0.53
1:K:146:PRO:HD2	1:K:200:ALA:CB	2.39	0.53
1:M:118:PRO:HB3	1:M:144:TYR:HB3	1.89	0.53
2:R:146:ARG:HD3	2:R:167:VAL:HG11	1.90	0.52
1:U:22:CYS:HB3	1:U:79:LEU:HB3	1.92	0.52
2:H:146:ARG:HG2	2:H:177:TYR:CE1	2.45	0.52
2:J:146:ARG:HD3	2:J:167:VAL:HG11	1.92	0.52
2:F:121:ILE:HD12	2:F:198:CYS:HB2	1.90	0.52
2:T:111:LYS:HA	2:T:144:TYR:OH	2.10	0.52
2:Y:124:PRO:HD3	2:Y:136:VAL:HG22	1.92	0.52
1:Q:122:PRO:HD3	1:Q:207:LYS:HE3	1.92	0.51
2:D:146:ARG:HD3	2:D:167:VAL:HG11	1.91	0.51
2:P:154:ILE:HD11	2:P:183:LEU:HD21	1.92	0.51
1:S:13:GLN:HG2	1:S:16:ARG:HD3	1.92	0.51
1:X:177:THR:HG21	2:Y:139:LEU:HD21	1.93	0.51
2:D:39:ASN:OD1	2:D:54:TYR:HA	2.09	0.51
1:K:6:GLU:OE1	1:K:103:GLY:HA3	2.10	0.51
2:H:154:ILE:HD11	2:H:183:LEU:HD21	1.92	0.51
1:U:158:LEU:HD23	1:U:180:VAL:HG21	1.93	0.51
2:V:151:LYS:HB3	2:V:199:GLU:HB3	1.93	0.51
1:S:87:ARG:HG3	1:S:89:GLU:HB3	1.93	0.51
2:B:154:ILE:HD11	2:B:183:LEU:HD21	1.93	0.51
1:A:91:THR:HG23	1:A:109:THR:HA	1.91	0.51
1:X:154:ASN:HB2	1:X:158:LEU:HD13	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:151:LYS:HB3	2:F:199:GLU:HB3	1.91	0.50
2:R:154:ILE:HD11	2:R:183:LEU:HD21	1.93	0.50
2:N:42:LEU:HB2	2:N:52:LEU:HD11	1.94	0.50
2:L:124:PRO:HD3	2:L:136:VAL:HG22	1.94	0.50
1:C:30:SER:O	1:C:53:ASN:HB2	2.11	0.50
1:K:2:VAL:HG13	1:K:27:PHE:CD1	2.48	0.49
2:R:16:GLY:HA2	4:R:320:HOH:O	2.12	0.49
1:X:39:GLN:HB2	1:X:45:LEU:HD23	1.95	0.49
1:O:13:GLN:HG3	1:O:16:ARG:HD3	1.95	0.49
2:F:154:ILE:HD11	2:F:183:LEU:HD21	1.95	0.49
1:O:41:PRO:O	1:O:42:THR:HB	2.13	0.48
2:D:190:TYR:CE2	2:D:216:ASN:HB2	2.49	0.48
2:V:154:ILE:HD11	2:V:183:LEU:HD21	1.95	0.48
2:J:154:ILE:HD11	2:J:183:LEU:HD21	1.95	0.48
1:K:63:SER:O	1:K:67:ARG:NH2	2.46	0.48
1:S:22:CYS:HB3	1:S:79:LEU:HB3	1.95	0.48
1:U:2:VAL:HG13	1:U:27:PHE:CD1	2.49	0.48
2:T:35:ASN:HB3	4:T:327:HOH:O	2.14	0.48
2:D:94:MET:HE3	2:D:96:THR:HG23	1.96	0.47
1:E:91:THR:HG23	1:E:109:THR:HA	1.97	0.47
1:K:30:SER:O	1:K:53:ASN:HB2	2.14	0.47
1:E:2:VAL:HG13	1:E:27:PHE:CD1	2.49	0.47
1:S:2:VAL:HG13	1:S:27:PHE:CD1	2.49	0.47
2:D:151:LYS:HB3	2:D:199:GLU:HB3	1.96	0.47
1:K:115:THR:HA	1:K:145:PHE:O	2.14	0.47
2:F:146:ARG:HD3	2:F:167:VAL:HG11	1.95	0.47
1:I:2:VAL:HG13	1:I:27:PHE:CD1	2.49	0.47
1:E:30:SER:O	1:E:53:ASN:HB2	2.13	0.47
2:J:151:LYS:HB3	2:J:199:GLU:HB3	1.96	0.47
1:X:91:THR:HG23	1:X:109:THR:HA	1.97	0.47
1:A:2:VAL:HG13	1:A:27:PHE:CD1	2.50	0.47
1:X:2:VAL:HG13	1:X:27:PHE:CD1	2.50	0.46
1:O:2:VAL:HG13	1:O:27:PHE:CD1	2.51	0.46
1:A:49:ALA:HA	1:A:59:TYR:O	2.15	0.46
1:E:122:PRO:HD3	1:E:207:LYS:HE3	1.98	0.46
2:L:151:LYS:HB3	2:L:199:GLU:HB3	1.97	0.46
2:L:149:SER:HB3	2:L:201:VAL:HB	1.97	0.46
1:C:49:ALA:HA	1:C:59:TYR:O	2.16	0.46
1:G:49:ALA:HA	1:G:59:TYR:O	2.15	0.46
1:U:199:PRO:HD2	4:U:306:HOH:O	2.16	0.46
1:M:2:VAL:HG13	1:M:27:PHE:CD1	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:147:ASP:O	2:P:202:HIS:CD2	2.68	0.46
1:U:91:THR:HG23	1:U:109:THR:HA	1.97	0.45
2:R:41:LEU:HD23	2:R:51:LEU:HA	1.98	0.45
1:S:91:THR:HG23	1:S:109:THR:HA	1.98	0.45
1:C:91:THR:HG23	1:C:109:THR:HA	1.98	0.45
2:Y:170:GLN:HB2	2:Y:177:TYR:CZ	2.52	0.45
1:X:49:ALA:HA	1:X:59:TYR:O	2.17	0.45
2:H:146:ARG:HD3	2:H:167:VAL:HG11	1.97	0.45
1:K:91:THR:HG23	1:K:109:THR:HA	1.98	0.45
1:Q:49:ALA:HA	1:Q:59:TYR:O	2.17	0.45
1:M:204:LYS:HG2	4:M:303:HOH:O	2.16	0.45
1:O:91:THR:HG23	1:O:109:THR:HA	1.99	0.45
1:Q:91:THR:HG23	1:Q:109:THR:HA	1.98	0.45
1:G:91:THR:HG23	1:G:109:THR:HA	1.98	0.45
2:P:214:ASN:O	2:P:215:ARG:HB2	2.17	0.45
2:P:147:ASP:O	2:P:202:HIS:HD2	2.00	0.45
1:K:9:GLY:HA3	1:K:106:THR:HG23	1.98	0.45
1:K:122:PRO:HD3	1:K:207:LYS:HE3	1.99	0.45
2:Y:94:MET:CE	2:Y:96:THR:HG23	2.47	0.45
1:C:11:LEU:HD11	4:C:325:HOH:O	2.17	0.45
1:I:49:ALA:HA	1:I:59:TYR:O	2.17	0.45
2:P:149:SER:HB3	2:P:201:VAL:HB	1.99	0.45
2:T:96:THR:HG23	2:T:100:PRO:HG3	1.99	0.45
2:N:151:LYS:HB3	2:N:199:GLU:HB3	1.99	0.44
1:X:138:GLY:HA2	1:X:153:TRP:CH2	2.53	0.44
2:L:94:MET:HE3	2:L:96:THR:HG23	1.99	0.44
1:M:91:THR:HG23	1:M:109:THR:HA	2.00	0.44
2:B:194:ASN:O	2:B:214:ASN:HA	2.17	0.44
1:C:2:VAL:HG13	1:C:27:PHE:CD1	2.53	0.44
2:P:151:LYS:HB3	2:P:199:GLU:HB3	2.00	0.44
2:B:149:SER:HB3	2:B:201:VAL:HB	1.99	0.44
2:N:151:LYS:HG2	2:N:158:GLU:HG3	2.00	0.44
1:M:49:ALA:HA	1:M:59:TYR:O	2.18	0.44
1:X:13:GLN:HG3	1:X:14:PRO:HD2	2.00	0.44
2:Y:122:PHE:HA	2:Y:123:PRO:HD3	1.92	0.43
1:K:86:LEU:HG	1:K:110:VAL:HG21	2.00	0.43
1:C:41:PRO:O	1:C:42:THR:HG22	2.18	0.43
2:V:149:SER:HB3	2:V:201:VAL:HB	2.01	0.43
1:K:60:TYR:CE1	1:K:70:ILE:HG22	2.54	0.43
2:B:151:LYS:HB3	2:B:199:GLU:HB3	2.00	0.43
1:I:87:ARG:O	1:I:90:ASP:HB2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:149:SER:HB3	2:H:201:VAL:HB	2.00	0.43
1:Q:121:TYR:CD2	2:R:128:GLN:HG3	2.54	0.43
2:F:124:PRO:HG3	2:F:134:ALA:HB1	2.01	0.43
1:O:49:ALA:HA	1:O:59:TYR:O	2.19	0.43
2:V:94:MET:CE	2:V:96:THR:HG23	2.48	0.43
1:S:197:ALA:HB2	1:S:204:LYS:HE2	2.01	0.43
2:N:41:LEU:HD23	2:N:51:LEU:HA	2.00	0.43
1:U:49:ALA:HA	1:U:59:TYR:O	2.19	0.43
1:C:146:PRO:HD2	1:C:200:ALA:CB	2.49	0.43
1:E:40:ALA:HB3	1:E:43:LYS:HB2	2.00	0.43
2:J:94:MET:CE	2:J:96:THR:HG23	2.49	0.43
2:P:197:THR:HG1	2:P:212:SER:HG	1.67	0.43
1:X:153:TRP:CZ3	1:X:194:CYS:HB3	2.53	0.42
2:B:41:LEU:HD23	2:B:51:LEU:HA	2.00	0.42
2:D:149:SER:HB3	2:D:201:VAL:HB	2.00	0.42
2:J:151:LYS:HG2	2:J:158:GLU:HG3	2.01	0.42
2:D:146:ARG:HB2	4:D:340:HOH:O	2.19	0.42
2:T:41:LEU:HD23	2:T:51:LEU:HA	2.02	0.42
2:T:146:ARG:HD3	2:T:167:VAL:HG11	2.02	0.42
2:D:94:MET:HE1	2:D:96:THR:HG23	2.01	0.42
2:L:94:MET:CE	2:L:96:THR:HG23	2.49	0.42
1:K:202:SER:HB3	1:M:208:LYS:HB3	2.00	0.42
1:A:122:PRO:HD3	1:A:207:LYS:HE3	2.02	0.42
1:K:7:SER:OG	1:K:8:GLY:N	2.53	0.42
2:J:94:MET:HE2	2:J:96:THR:HG23	2.02	0.42
1:G:12:VAL:CG2	1:G:18:LEU:HD13	2.48	0.42
1:S:49:ALA:HA	1:S:59:TYR:O	2.20	0.42
2:V:41:LEU:HD23	2:V:51:LEU:HA	2.02	0.42
2:J:41:LEU:HD23	2:J:51:LEU:HA	2.00	0.42
2:F:94:MET:CE	2:F:96:THR:HG23	2.48	0.42
2:P:41:LEU:HD23	2:P:51:LEU:HA	2.01	0.42
2:P:107:LYS:HE3	2:P:109:GLU:HB3	2.02	0.42
1:U:154:ASN:HB2	1:U:158:LEU:HD13	2.02	0.42
2:R:66:ARG:NE	2:R:87:ASP:OD2	2.46	0.42
2:T:149:SER:HB3	2:T:201:VAL:HB	2.02	0.42
1:A:169:LEU:HG	1:A:174:TYR:CE1	2.54	0.41
1:K:169:LEU:HG	1:K:174:TYR:CE1	2.55	0.41
4:S:304:HOH:O	2:T:127:GLU:HG3	2.19	0.41
1:C:62:ASP:HA	1:C:65:LYS:HD2	2.00	0.41
1:C:169:LEU:HG	1:C:174:TYR:CE1	2.55	0.41
2:N:149:SER:HB3	2:N:201:VAL:HB	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:115:THR:HA	1:O:145:PHE:O	2.20	0.41
2:L:41:LEU:HD23	2:L:51:LEU:HA	2.02	0.41
1:G:169:LEU:HG	1:G:174:TYR:CE1	2.55	0.41
2:Y:154:ILE:HD11	2:Y:183:LEU:HD21	2.01	0.41
2:H:94:MET:CE	2:H:96:THR:HG23	2.51	0.41
2:Y:129:LEU:O	2:Y:131:THR:N	2.53	0.41
1:K:6:GLU:HG3	1:K:96:CYS:HB3	2.03	0.41
1:E:91:THR:O	1:E:92:ALA:HB2	2.21	0.41
2:P:151:LYS:HG2	2:P:158:GLU:HG3	2.02	0.41
2:R:149:SER:HB3	2:R:201:VAL:HB	2.03	0.41
2:D:190:TYR:HE2	2:D:216:ASN:HB3	1.86	0.41
2:F:122:PHE:HA	2:F:123:PRO:HD3	1.97	0.41
2:J:149:SER:HB3	2:J:201:VAL:HB	2.02	0.41
2:V:94:MET:HE3	2:V:96:THR:HG23	2.03	0.41
2:H:41:LEU:HD23	2:H:51:LEU:HA	2.02	0.41
1:S:122:PRO:HD3	1:S:207:LYS:HE3	2.03	0.41
1:U:169:LEU:HG	1:U:174:TYR:CE1	2.56	0.40
1:C:191:THR:HA	4:C:327:HOH:O	2.21	0.40
2:D:128:GLN:HG2	2:D:133:GLY:O	2.21	0.40
1:M:153:TRP:CH2	1:M:194:CYS:HB3	2.56	0.40
2:F:138:CYS:HB2	2:F:152:TRP:CH2	2.56	0.40
1:I:146:PRO:HD2	1:I:200:ALA:CB	2.51	0.40
2:V:122:PHE:HA	2:V:123:PRO:HD3	1.98	0.40
1:M:22:CYS:HB3	1:M:79:LEU:HB3	2.03	0.40
2:N:122:PHE:HA	2:N:123:PRO:HD3	2.00	0.40
2:F:41:LEU:HD23	2:F:51:LEU:HA	2.03	0.40
1:S:171:SER:CB	2:Y:159:ARG:HG3	2.51	0.40
2:Y:94:MET:HE2	2:Y:96:THR:HG23	2.04	0.40
1:Q:115:THR:HA	1:Q:145:PHE:O	2.22	0.40
1:I:169:LEU:HG	1:I:174:TYR:CE1	2.56	0.40
1:K:17:SER:HB3	1:K:82:GLN:HE22	1.87	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	203/252 (81%)	195 (96%)	8 (4%)	0	100	100
1	C	201/252 (80%)	195 (97%)	6 (3%)	0	100	100
1	E	204/252 (81%)	197 (97%)	7 (3%)	0	100	100
1	G	203/252 (81%)	197 (97%)	4 (2%)	2 (1%)	19	16
1	I	201/252 (80%)	197 (98%)	3 (2%)	1 (0%)	34	34
1	K	200/252 (79%)	194 (97%)	6 (3%)	0	100	100
1	M	200/252 (79%)	196 (98%)	4 (2%)	0	100	100
1	O	201/252 (80%)	196 (98%)	5 (2%)	0	100	100
1	Q	202/252 (80%)	199 (98%)	2 (1%)	1 (0%)	34	34
1	S	200/252 (79%)	196 (98%)	4 (2%)	0	100	100
1	U	201/252 (80%)	197 (98%)	3 (2%)	1 (0%)	34	34
1	X	193/252 (77%)	189 (98%)	2 (1%)	2 (1%)	19	16
2	B	213/220 (97%)	208 (98%)	5 (2%)	0	100	100
2	D	214/220 (97%)	208 (97%)	4 (2%)	2 (1%)	21	18
2	F	213/220 (97%)	208 (98%)	5 (2%)	0	100	100
2	H	214/220 (97%)	212 (99%)	2 (1%)	0	100	100
2	J	213/220 (97%)	208 (98%)	5 (2%)	0	100	100
2	L	213/220 (97%)	210 (99%)	3 (1%)	0	100	100
2	N	213/220 (97%)	208 (98%)	5 (2%)	0	100	100
2	P	213/220 (97%)	206 (97%)	7 (3%)	0	100	100
2	R	213/220 (97%)	206 (97%)	7 (3%)	0	100	100
2	T	213/220 (97%)	207 (97%)	6 (3%)	0	100	100
2	V	208/220 (94%)	203 (98%)	5 (2%)	0	100	100
2	Y	204/220 (93%)	186 (91%)	12 (6%)	6 (3%)	6	2
3	a	10/12 (83%)	10 (100%)	0	0	100	100
3	c	9/12 (75%)	9 (100%)	0	0	100	100
3	e	9/12 (75%)	9 (100%)	0	0	100	100
3	g	10/12 (83%)	10 (100%)	0	0	100	100
3	i	10/12 (83%)	10 (100%)	0	0	100	100
3	k	9/12 (75%)	8 (89%)	1 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	m	10/12 (83%)	10 (100%)	0	0	100	100
3	o	9/12 (75%)	8 (89%)	1 (11%)	0	100	100
3	q	8/12 (67%)	7 (88%)	1 (12%)	0	100	100
3	s	9/12 (75%)	9 (100%)	0	0	100	100
3	u	8/12 (67%)	8 (100%)	0	0	100	100
3	x	10/12 (83%)	10 (100%)	0	0	100	100
All	All	5064/5808 (87%)	4926 (97%)	123 (2%)	15 (0%)	46	50

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	7	THR
2	D	215	ARG
2	Y	159	ARG
1	G	125	PRO
1	X	160	SER
2	Y	156	GLY
1	X	43	LYS
2	Y	130	ALA
2	Y	131	THR
2	Y	184	SER
2	Y	9	PRO
1	I	41	PRO
1	U	41	PRO
1	G	41	PRO
1	Q	41	PRO

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	176/208 (85%)	167 (95%)	9 (5%)	29	33
1	C	175/208 (84%)	165 (94%)	10 (6%)	25	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	176/208 (85%)	166 (94%)	10 (6%)	25	28
1	G	176/208 (85%)	167 (95%)	9 (5%)	29	33
1	I	174/208 (84%)	166 (95%)	8 (5%)	33	39
1	K	174/208 (84%)	161 (92%)	13 (8%)	17	16
1	M	174/208 (84%)	161 (92%)	13 (8%)	17	16
1	O	174/208 (84%)	162 (93%)	12 (7%)	19	19
1	Q	175/208 (84%)	164 (94%)	11 (6%)	22	23
1	S	174/208 (84%)	164 (94%)	10 (6%)	25	28
1	U	175/208 (84%)	163 (93%)	12 (7%)	19	19
1	X	167/208 (80%)	159 (95%)	8 (5%)	31	36
2	B	191/196 (97%)	180 (94%)	11 (6%)	25	27
2	D	192/196 (98%)	182 (95%)	10 (5%)	29	32
2	F	191/196 (97%)	179 (94%)	12 (6%)	22	23
2	H	192/196 (98%)	182 (95%)	10 (5%)	29	32
2	J	191/196 (97%)	182 (95%)	9 (5%)	32	38
2	L	191/196 (97%)	182 (95%)	9 (5%)	32	38
2	N	191/196 (97%)	182 (95%)	9 (5%)	32	38
2	P	191/196 (97%)	181 (95%)	10 (5%)	29	32
2	R	191/196 (97%)	181 (95%)	10 (5%)	29	32
2	T	191/196 (97%)	178 (93%)	13 (7%)	20	20
2	V	189/196 (96%)	181 (96%)	8 (4%)	36	43
2	Y	187/196 (95%)	164 (88%)	23 (12%)	6	4
3	a	11/11 (100%)	10 (91%)	1 (9%)	12	10
3	c	10/11 (91%)	10 (100%)	0	100	100
3	e	10/11 (91%)	9 (90%)	1 (10%)	9	8
3	g	11/11 (100%)	11 (100%)	0	100	100
3	i	11/11 (100%)	10 (91%)	1 (9%)	12	10
3	k	10/11 (91%)	10 (100%)	0	100	100
3	m	11/11 (100%)	9 (82%)	2 (18%)	2	1
3	o	10/11 (91%)	10 (100%)	0	100	100
3	q	9/11 (82%)	9 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	s	10/11 (91%)	10 (100%)	0	100	100
3	u	9/11 (82%)	9 (100%)	0	100	100
3	x	11/11 (100%)	10 (91%)	1 (9%)	12	10
All	All	4501/4980 (90%)	4236 (94%)	265 (6%)	24	26

All (265) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	ARG
1	A	87	ARG
1	A	89	GLU
1	A	93	THR
1	A	107	THR
1	A	114	GLU
1	A	135	VAL
1	A	170	GLN
1	A	189	SER
1	C	42	THR
1	C	61	ARG
1	C	72	ARG
1	C	87	ARG
1	C	89	GLU
1	C	93	THR
1	C	107	THR
1	C	170	GLN
1	C	189	SER
1	C	204	LYS
1	E	7	SER
1	E	61	ARG
1	E	72	ARG
1	E	89	GLU
1	E	93	THR
1	E	104	GLN
1	E	107	THR
1	E	133	SER
1	E	189	SER
1	E	204	LYS
1	G	61	ARG
1	G	87	ARG
1	G	88	SER
1	G	93	THR

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Mol	Chain	Res	Type
1	G	107	THR
1	G	135	VAL
1	G	170	GLN
1	G	189	SER
1	G	204	LYS
1	I	61	ARG
1	I	72	ARG
1	I	89	GLU
1	I	93	THR
1	I	107	THR
1	I	135	VAL
1	I	170	GLN
1	I	189	SER
2	J	1	ASP
2	J	11	LEU
2	J	94	MET
2	J	109	GLU
2	J	140	MET
2	J	146	ARG
2	J	160	ARG
2	J	203	LYS
2	J	207	SER
1	K	13	GLN
1	K	21	SER
1	K	61	ARG
1	K	86	LEU
1	K	89	GLU
1	K	96	CYS
1	K	104	GLN
1	K	106	THR
1	K	107	THR
1	K	135	VAL
1	K	170	GLN
1	K	189	SER
1	K	204	LYS
1	M	13	GLN
1	M	16	ARG
1	M	61	ARG
1	M	72	ARG
1	M	87	ARG
1	M	89	GLU
1	M	93	THR

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Mol	Chain	Res	Type
1	M	96	CYS
1	M	107	THR
1	M	135	VAL
1	M	170	GLN
1	M	189	SER
1	M	204	LYS
1	O	13	GLN
1	O	61	ARG
1	O	72	ARG
1	O	89	GLU
1	O	93	THR
1	O	107	THR
1	O	135	VAL
1	O	170	GLN
1	O	171	SER
1	O	189	SER
1	O	195	ASN
1	O	204	LYS
1	Q	43	LYS
1	Q	61	ARG
1	Q	78	THR
1	Q	96	CYS
1	Q	107	THR
1	Q	133	SER
1	Q	135	VAL
1	Q	170	GLN
1	Q	189	SER
1	Q	204	LYS
1	Q	207	LYS
1	S	13	GLN
1	S	61	ARG
1	S	89	GLU
1	S	93	THR
1	S	107	THR
1	S	114	GLU
1	S	135	VAL
1	S	170	GLN
1	S	189	SER
1	S	204	LYS
1	U	7	SER
1	U	61	ARG
1	U	89	GLU

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Mol	Chain	Res	Type
1	U	93	THR
1	U	107	THR
1	U	133	SER
1	U	135	VAL
1	U	170	GLN
1	U	189	SER
1	U	194	CYS
1	U	204	LYS
1	U	207	LYS
1	X	61	ARG
1	X	77	SER
1	X	87	ARG
1	X	93	THR
1	X	107	THR
1	X	170	GLN
1	X	180	VAL
1	X	195	ASN
2	B	1	ASP
2	B	7	THR
2	B	11	LEU
2	B	94	MET
2	B	109	GLU
2	B	140	MET
2	B	146	ARG
2	B	147	ASP
2	B	161	ASP
2	B	203	LYS
2	B	207	SER
2	D	1	ASP
2	D	11	LEU
2	D	94	MET
2	D	109	GLU
2	D	140	MET
2	D	146	ARG
2	D	147	ASP
2	D	203	LYS
2	D	207	SER
2	D	215	ARG
2	F	1	ASP
2	F	11	LEU
2	F	86	GLU
2	F	94	MET

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Mol	Chain	Res	Type
2	F	109	GLU
2	F	140	MET
2	F	146	ARG
2	F	147	ASP
2	F	151	LYS
2	F	203	LYS
2	F	207	SER
2	F	215	ARG
2	H	1	ASP
2	H	11	LEU
2	H	94	MET
2	H	109	GLU
2	H	140	MET
2	H	146	ARG
2	H	147	ASP
2	H	203	LYS
2	H	207	SER
2	H	215	ARG
2	L	11	LEU
2	L	94	MET
2	L	109	GLU
2	L	127	GLU
2	L	140	MET
2	L	146	ARG
2	L	203	LYS
2	L	211	LYS
2	L	214	ASN
2	N	1	ASP
2	N	9	PRO
2	N	11	LEU
2	N	94	MET
2	N	109	GLU
2	N	140	MET
2	N	146	ARG
2	N	211	LYS
2	N	214	ASN
2	P	1	ASP
2	P	11	LEU
2	P	94	MET
2	P	109	GLU
2	P	140	MET
2	P	146	ARG

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Mol	Chain	Res	Type
2	P	171	ASP
2	P	207	SER
2	P	212	SER
2	P	215	ARG
2	R	1	ASP
2	R	11	LEU
2	R	53	ILE
2	R	94	MET
2	R	109	GLU
2	R	140	MET
2	R	146	ARG
2	R	203	LYS
2	R	207	SER
2	R	215	ARG
2	T	1	ASP
2	T	11	LEU
2	T	74	THR
2	T	94	MET
2	T	109	GLU
2	T	140	MET
2	T	146	ARG
2	T	147	ASP
2	T	151	LYS
2	T	203	LYS
2	T	207	SER
2	T	211	LYS
2	T	215	ARG
2	V	1	ASP
2	V	11	LEU
2	V	94	MET
2	V	109	GLU
2	V	140	MET
2	V	146	ARG
2	V	151	LYS
2	V	207	SER
2	Y	1	ASP
2	Y	8	THR
2	Y	11	LEU
2	Y	21	ILE
2	Y	25	SER
2	Y	41	LEU
2	Y	53	ILE

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Mol	Chain	Res	Type
2	Y	94	MET
2	Y	109	GLU
2	Y	126	THR
2	Y	127	GLU
2	Y	129	LEU
2	Y	157	THR
2	Y	158	GLU
2	Y	159	ARG
2	Y	160	ARG
2	Y	161	ASP
2	Y	181	SER
2	Y	191	GLU
2	Y	195	LEU
2	Y	197	THR
2	Y	203	LYS
2	Y	207	SER
3	a	422	VAL
3	e	422	VAL
3	i	423	ASN
3	m	422	VAL
3	m	423	ASN
3	x	423	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	163	HIS

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

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	207/252 (82%)	-0.08	0 100 100	21, 35, 49, 66	0
1	C	205/252 (81%)	-0.05	1 (0%) 91 91	20, 32, 51, 80	0
1	E	208/252 (82%)	0.09	1 (0%) 91 91	21, 37, 57, 79	0
1	G	207/252 (82%)	0.03	2 (0%) 84 83	22, 41, 62, 84	0
1	I	205/252 (81%)	-0.07	0 100 100	23, 36, 52, 69	0
1	K	204/252 (80%)	0.45	7 (3%) 49 48	28, 48, 71, 93	0
1	M	204/252 (80%)	0.33	2 (0%) 84 83	29, 46, 63, 78	0
1	O	205/252 (81%)	0.23	4 (1%) 68 67	23, 40, 61, 75	0
1	Q	206/252 (81%)	0.50	13 (6%) 23 22	25, 46, 73, 84	0
1	S	204/252 (80%)	0.01	1 (0%) 91 91	18, 34, 55, 78	0
1	U	205/252 (81%)	0.49	13 (6%) 23 22	24, 49, 73, 103	0
1	X	199/252 (78%)	-0.01	1 (0%) 91 91	18, 34, 54, 87	0
2	B	215/220 (97%)	0.07	3 (1%) 78 77	18, 35, 59, 74	0
2	D	216/220 (98%)	0.06	2 (0%) 85 85	18, 31, 55, 87	0
2	F	215/220 (97%)	0.07	2 (0%) 85 85	20, 34, 62, 90	0
2	H	216/220 (98%)	-0.06	1 (0%) 91 91	18, 35, 56, 69	0
2	J	215/220 (97%)	0.00	1 (0%) 91 91	18, 33, 56, 75	0
2	L	215/220 (97%)	0.09	1 (0%) 91 91	22, 39, 57, 68	0
2	N	215/220 (97%)	0.29	3 (1%) 78 77	30, 49, 68, 80	0
2	P	215/220 (97%)	-0.06	0 100 100	23, 37, 61, 75	0
2	R	215/220 (97%)	0.48	12 (5%) 28 27	30, 54, 78, 88	0
2	T	215/220 (97%)	0.45	14 (6%) 22 21	21, 44, 76, 90	0
2	V	212/220 (96%)	0.72	25 (11%) 6 6	27, 52, 88, 106	0
2	Y	210/220 (95%)	0.10	5 (2%) 62 61	15, 33, 72, 82	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
3	a	12/12 (100%)	0.03	0	100	100	23, 31, 50, 79	0
3	c	11/12 (91%)	-0.00	0	100	100	25, 28, 51, 52	0
3	e	11/12 (91%)	0.13	0	100	100	23, 30, 56, 56	0
3	g	12/12 (100%)	-0.29	0	100	100	26, 30, 46, 48	0
3	i	12/12 (100%)	0.01	0	100	100	30, 34, 50, 65	0
3	k	11/12 (91%)	0.43	0	100	100	24, 36, 58, 59	0
3	m	12/12 (100%)	0.31	0	100	100	39, 47, 69, 78	0
3	o	11/12 (91%)	0.28	0	100	100	34, 43, 59, 61	0
3	q	10/12 (83%)	0.41	0	100	100	42, 50, 56, 68	0
3	s	11/12 (91%)	0.11	0	100	100	27, 32, 45, 48	0
3	u	10/12 (83%)	0.20	0	100	100	32, 37, 40, 54	0
3	x	12/12 (100%)	-0.04	0	100	100	23, 29, 57, 59	0
All	All	5168/5808 (88%)	0.17	114 (2%)	65	64	15, 40, 67, 106	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	K	75	ALA	5.9
1	Q	31	ASP	5.1
2	Y	157	THR	4.6
2	H	132	GLY	4.5
2	Y	156	GLY	4.4
1	U	157	ALA	4.2
2	V	204	THR	4.2
2	V	201	VAL	4.1
2	R	85	ALA	3.9
2	V	154	ILE	3.9
2	T	131	THR	3.8
2	V	209	VAL	3.8
2	V	210	VAL	3.7
1	U	186	THR	3.6
2	R	115	ALA	3.6
2	R	173	LYS	3.5
2	V	194	ASN	3.4
2	V	8	THR	3.3
1	E	75	ALA	3.2
1	Q	76	LYS	3.2
2	T	191	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
2	V	214	ASN	3.1
2	V	162	GLY	3.1
1	U	160	SER	3.1
1	Q	56	GLY	3.1
1	U	170	GLN	3.1
2	T	195	LEU	3.0
2	D	216	ASN	2.9
1	U	213	ASN	2.9
2	R	119	VAL	2.9
1	Q	26	GLY	2.9
2	V	195	LEU	2.9
1	Q	28	THR	2.8
2	V	183	LEU	2.8
1	K	80	TYR	2.8
2	V	207	SER	2.7
1	Q	2	VAL	2.7
2	V	173	LYS	2.7
2	R	134	ALA	2.7
2	R	206	SER	2.7
1	S	170	GLN	2.6
1	Q	66	GLY	2.6
2	V	7	THR	2.6
2	T	196	TYR	2.6
2	V	196	TYR	2.6
2	B	210	VAL	2.6
2	V	132	GLY	2.6
2	T	130	ALA	2.6
1	O	76	LYS	2.6
1	K	14	PRO	2.6
2	N	1	ASP	2.6
1	K	190	GLN	2.5
1	G	213	ASN	2.5
2	T	155	ASP	2.5
2	V	160	ARG	2.5
2	Y	7	THR	2.5
1	Q	75	ALA	2.4
2	T	188	ALA	2.4
2	T	156	GLY	2.4
2	V	158	GLU	2.4
2	T	204	THR	2.4
2	Y	9	PRO	2.4
1	K	20	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
2	L	195	LEU	2.4
2	T	185	LEU	2.4
2	J	133	GLY	2.4
1	M	32	SER	2.4
1	U	155	SER	2.4
2	R	213	PHE	2.4
2	Y	193	HIS	2.4
2	V	134	ALA	2.4
1	U	210	VAL	2.4
1	Q	25	SER	2.3
1	X	171	SER	2.3
2	D	207	SER	2.3
1	Q	3	LYS	2.3
1	Q	185	SER	2.3
2	F	134	ALA	2.2
1	O	5	GLN	2.2
2	N	27	GLN	2.2
2	N	161	ASP	2.2
2	B	192	SER	2.2
2	V	159	ARG	2.2
2	R	7	THR	2.2
1	C	170	GLN	2.2
1	Q	4	LEU	2.2
1	K	67	ARG	2.2
1	U	209	ILE	2.1
2	T	154	ILE	2.1
2	V	153	LYS	2.1
1	O	2	VAL	2.1
2	T	88	LEU	2.1
2	R	132	GLY	2.1
1	U	189	SER	2.1
2	R	205	SER	2.1
1	O	12	VAL	2.1
1	U	18	LEU	2.1
2	F	129	LEU	2.1
1	M	3	LYS	2.1
2	B	132	GLY	2.1
1	Q	13	GLN	2.1
1	K	21	SER	2.1
2	V	205	SER	2.1
1	U	150	THR	2.0
2	V	213	PHE	2.0

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Mol	Chain	Res	Type	RSRZ
2	R	144	TYR	2.0
2	V	190	TYR	2.0
1	G	87	ARG	2.0
2	R	16	GLY	2.0
2	T	194	ASN	2.0
2	V	119	VAL	2.0
1	U	145	PHE	2.0
1	U	185	SER	2.0
2	T	144	TYR	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.