



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:58 PM BST

PDB ID : 4UOK
EMDB ID: : EMD-2655
Title : Electron Cryo-microscopy of Venezuelan Equine Encephalitis Virus TC-83 in complex with neutralizing antibody Fab 3B4C-4
Authors : Porta, J.; Jose, J.; Roehrig, J.T.; Blair, C.D.; Kuhn, R.J.; G Rossmann, M.
Deposited on : 2014-06-04
Resolution : 18.00 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/EMValidationReportHelp>

MolProbity : 4.02b-467
Mogul : unknown
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : trunk27241

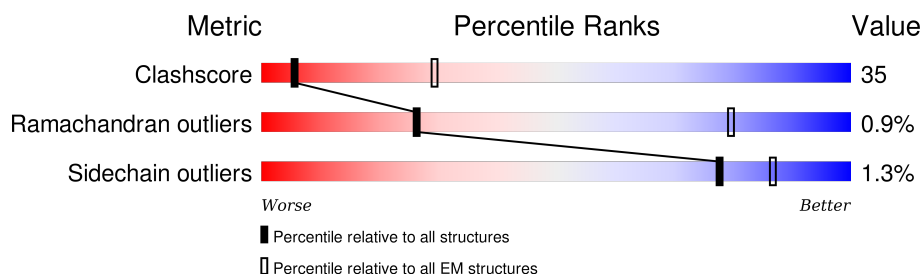
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 18.00 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	218	 74% 24% ..
2	L	215	 68% 29% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3308 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called FAB FRAGMENT HEAVY CHAIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	218	Total	C	N	O	S	0	0
			1645	1041	270	326	8		

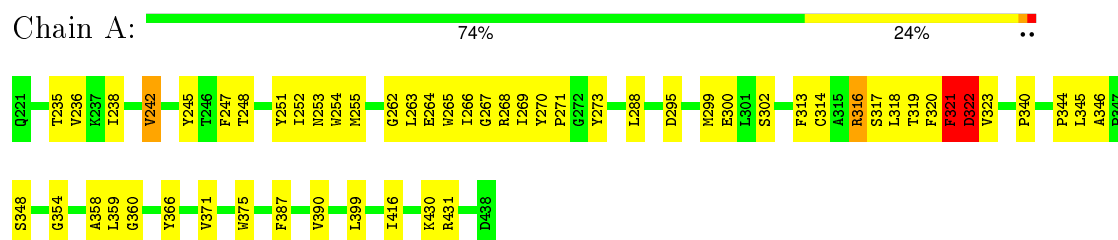
- Molecule 2 is a protein called FAB FRAGMENT LIGHT CHAIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	L	215	Total	C	N	O	S	0	0
			1663	1032	275	350	6		

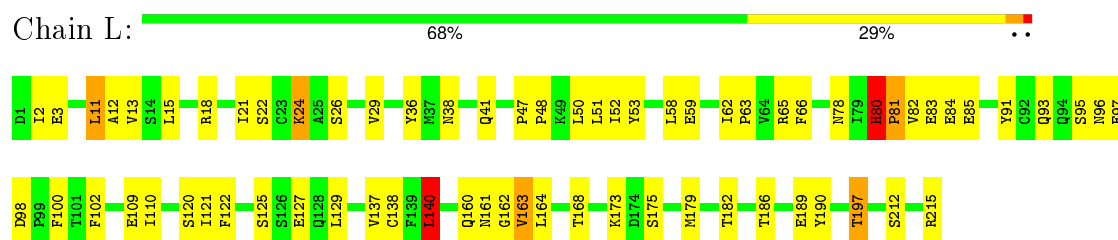
3 Residue-property plots

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. 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. 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: FAB FRAGMENT HEAVY CHAIN



• Molecule 2: FAB FRAGMENT LIGHT CHAIN



4 Experimental information

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	EACH IMAGE, Not provided	Depositor
Microscope	OTHER	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	24	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	59000	Depositor
Image detector	KODAK SO-163 FILM	Depositor

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	$\# Z > 2$	RMSZ	$\# Z > 2$
1	A	0.65	0/1685	1.02	7/2293 (0.3%)
2	L	0.62	0/1701	0.96	2/2312 (0.1%)
All	All	0.63	0/3386	0.99	9/4605 (0.2%)

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	4
2	L	0	1
All	All	0	5

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	242	VAL	CG1-CB-CG2	-10.94	93.40	110.90
1	A	321	PHE	CB-CG-CD2	9.44	127.41	120.80
1	A	316	ARG	NE-CZ-NH2	6.48	123.54	120.30
1	A	322	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	A	321	PHE	CG-CD2-CE2	5.98	127.37	120.80
2	L	11	LEU	CB-CG-CD1	-5.96	100.86	111.00
2	L	140	LEU	CB-CG-CD2	5.22	119.87	111.00
1	A	322	ASP	CB-CA-C	-5.18	100.04	110.40
1	A	321	PHE	CB-CG-CD1	-5.04	117.27	120.80

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	317	SER	Mainchain
1	A	321	PHE	Mainchain,Peptide
1	A	322	ASP	Mainchain
2	L	80	HIS	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	1645	0	1608	175	0
2	L	1663	0	1564	175	0
All	All	3308	0	3172	228	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 35.

All (228) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:344:PRO:HG2	2:L:125:SER:CB	1.24	1.58
1:A:263:LEU:HD11	2:L:48:PRO:CG	1.19	1.57
1:A:263:LEU:CD1	2:L:48:PRO:HG2	1.37	1.51
1:A:344:PRO:CG	2:L:125:SER:HB2	1.38	1.48
1:A:263:LEU:CD1	2:L:48:PRO:CG	1.88	1.42
1:A:344:PRO:CG	2:L:125:SER:CB	1.97	1.40
1:A:358:ALA:O	2:L:122:PHE:CE2	1.85	1.29
1:A:430:LYS:HE3	2:L:127:GLU:OE2	1.24	1.28
1:A:430:LYS:NZ	2:L:127:GLU:HG2	1.49	1.25
1:A:346:ALA:O	2:L:122:PHE:HD2	1.14	1.23
1:A:255:MET:HE1	2:L:102:PHE:CE1	1.74	1.22
1:A:263:LEU:HD11	2:L:48:PRO:CB	1.71	1.20
1:A:319:THR:OG1	2:L:95:SER:HB2	1.04	1.20
1:A:255:MET:HE3	2:L:102:PHE:HZ	1.12	1.15
1:A:319:THR:OG1	2:L:95:SER:CB	1.94	1.15
1:A:346:ALA:O	2:L:122:PHE:CD2	2.00	1.14
1:A:263:LEU:HD11	2:L:48:PRO:HG3	1.20	1.13
1:A:263:LEU:HD21	2:L:48:PRO:CG	1.80	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:MET:HE3	2:L:102:PHE:CZ	1.85	1.10
1:A:358:ALA:O	2:L:122:PHE:HE2	1.25	1.09
1:A:263:LEU:CD2	2:L:48:PRO:CG	2.31	1.08
1:A:263:LEU:HD13	2:L:48:PRO:HG2	1.34	1.08
1:A:255:MET:CE	2:L:102:PHE:CZ	2.38	1.07
1:A:344:PRO:HG2	2:L:125:SER:HB3	1.09	1.05
1:A:319:THR:HB	2:L:100:PHE:CE1	1.94	1.02
1:A:344:PRO:HD2	2:L:125:SER:OG	1.60	1.02
1:A:263:LEU:CD2	2:L:48:PRO:HG2	1.90	1.02
1:A:387:PHE:CE2	2:L:168:THR:HG23	1.95	1.01
1:A:387:PHE:CE2	2:L:168:THR:CG2	2.44	1.01
1:A:255:MET:CE	2:L:102:PHE:CE1	2.43	1.00
1:A:358:ALA:HB3	2:L:120:SER:CB	1.92	0.99
1:A:263:LEU:HD21	2:L:48:PRO:CD	1.92	0.99
1:A:344:PRO:CD	2:L:125:SER:CB	2.41	0.99
1:A:263:LEU:CG	2:L:48:PRO:HG2	1.94	0.98
1:A:255:MET:HE1	2:L:102:PHE:HE1	1.09	0.98
2:L:29:VAL:HG23	2:L:96:ASN:HB2	1.44	0.97
1:A:387:PHE:CD2	2:L:168:THR:CG2	2.48	0.96
1:A:430:LYS:HZ1	2:L:127:GLU:HG2	1.24	0.96
1:A:344:PRO:HG3	2:L:125:SER:HB2	1.47	0.96
1:A:358:ALA:C	2:L:122:PHE:CE2	2.39	0.95
1:A:358:ALA:C	2:L:122:PHE:CZ	2.40	0.95
1:A:430:LYS:CE	2:L:127:GLU:OE2	2.13	0.95
1:A:319:THR:CB	2:L:100:PHE:CE1	2.50	0.95
1:A:319:THR:HG1	2:L:100:PHE:HE1	1.03	0.95
1:A:345:LEU:HD21	2:L:137:VAL:HG21	1.48	0.94
1:A:359:LEU:HA	2:L:122:PHE:CE2	2.01	0.94
1:A:263:LEU:CG	2:L:48:PRO:CG	2.47	0.93
1:A:242:VAL:HG11	1:A:247:PHE:CD1	2.04	0.92
1:A:430:LYS:HZ3	2:L:127:GLU:HG2	1.27	0.91
1:A:313:PHE:CZ	2:L:47:PRO:HB3	2.06	0.91
1:A:358:ALA:HB3	2:L:120:SER:HB3	1.53	0.91
1:A:319:THR:HG1	2:L:95:SER:HB2	1.23	0.91
1:A:344:PRO:HD2	2:L:125:SER:CB	2.06	0.85
2:L:3:GLU:HG2	2:L:26:SER:HB3	1.59	0.84
1:A:387:PHE:HE2	2:L:168:THR:HG23	1.41	0.83
1:A:345:LEU:CD2	2:L:137:VAL:HG21	2.09	0.82
1:A:263:LEU:HD21	2:L:48:PRO:HD3	1.57	0.82
1:A:430:LYS:NZ	2:L:127:GLU:CG	2.40	0.81
1:A:242:VAL:HG12	1:A:295:ASP:O	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:GLY:CA	2:L:91:TYR:OH	2.29	0.79
1:A:387:PHE:CD2	2:L:168:THR:HG22	2.15	0.79
2:L:59:GLU:O	2:L:62:ILE:HG12	1.83	0.78
1:A:262:GLY:HA2	2:L:91:TYR:OH	1.84	0.78
1:A:358:ALA:O	2:L:122:PHE:CZ	2.38	0.76
1:A:263:LEU:HD11	2:L:48:PRO:HB3	1.67	0.76
1:A:251:TYR:CD2	1:A:268:ARG:HD2	2.21	0.76
1:A:387:PHE:CD2	2:L:168:THR:HG23	2.19	0.75
2:L:84:GLU:HG3	2:L:175:SER:OG	1.86	0.75
1:A:344:PRO:CG	2:L:125:SER:HB3	1.84	0.75
1:A:430:LYS:HE3	2:L:127:GLU:CD	2.08	0.75
1:A:358:ALA:C	2:L:122:PHE:HZ	1.90	0.75
1:A:319:THR:CB	2:L:95:SER:HB2	2.17	0.75
1:A:263:LEU:HD21	2:L:48:PRO:HG3	1.70	0.73
1:A:358:ALA:CB	2:L:120:SER:CB	2.67	0.73
1:A:319:THR:HB	2:L:100:PHE:CZ	2.24	0.73
1:A:263:LEU:CG	2:L:48:PRO:HG3	2.17	0.72
1:A:318:LEU:HD21	2:L:93:GLN:OE1	1.89	0.72
1:A:319:THR:OG1	2:L:100:PHE:CE1	2.42	0.72
1:A:359:LEU:CA	2:L:122:PHE:CE2	2.73	0.71
1:A:321:PHE:HB2	1:A:322:ASP:HA	1.73	0.70
1:A:344:PRO:HG2	2:L:125:SER:HB2	0.98	0.70
1:A:358:ALA:CB	2:L:122:PHE:HZ	2.03	0.70
1:A:245:TYR:CD1	1:A:316:ARG:NH1	2.59	0.70
1:A:253:ASN:HD21	1:A:265:TRP:HD1	1.37	0.70
1:A:245:TYR:CZ	1:A:316:ARG:HD3	2.27	0.69
1:A:345:LEU:CD2	2:L:137:VAL:CG2	2.71	0.69
1:A:318:LEU:HD22	2:L:93:GLN:HE22	1.58	0.68
2:L:58:LEU:HD21	2:L:66:PHE:O	1.93	0.67
2:L:140:LEU:HB2	2:L:179:MET:HG3	1.76	0.67
2:L:80:HIS:CD2	2:L:81:PRO:HD3	2.30	0.67
2:L:95:SER:HA	2:L:100:PHE:CD1	2.29	0.66
1:A:345:LEU:HD22	2:L:137:VAL:HB	1.78	0.66
2:L:11:LEU:HD11	2:L:21:ILE:HG22	1.78	0.66
1:A:263:LEU:CD1	2:L:48:PRO:CB	2.52	0.65
2:L:52:ILE:CD1	2:L:58:LEU:HD23	2.26	0.65
1:A:288:LEU:HD23	1:A:299:MET:HA	1.80	0.64
1:A:242:VAL:CG1	1:A:295:ASP:HB3	2.27	0.64
1:A:262:GLY:HA3	2:L:91:TYR:OH	1.97	0.64
1:A:242:VAL:CG2	1:A:245:TYR:CE1	2.81	0.64
2:L:22:SER:HB3	2:L:24:LYS:HE2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LEU:CD1	2:L:48:PRO:HG3	1.90	0.63
2:L:58:LEU:HA	2:L:62:ILE:HD11	1.80	0.63
1:A:319:THR:OG1	2:L:100:PHE:HE1	1.73	0.62
1:A:359:LEU:HA	2:L:122:PHE:HE2	1.64	0.61
1:A:344:PRO:HG3	2:L:127:GLU:OE1	2.00	0.61
1:A:242:VAL:HG21	1:A:245:TYR:CZ	2.37	0.60
1:A:263:LEU:CD2	2:L:48:PRO:CD	2.70	0.59
2:L:58:LEU:CA	2:L:62:ILE:HD11	2.32	0.58
1:A:430:LYS:CE	2:L:127:GLU:CD	2.69	0.58
2:L:13:VAL:O	2:L:110:ILE:HD12	2.03	0.58
2:L:83:GLU:HG3	2:L:85:GLU:HB3	1.85	0.58
1:A:345:LEU:HD21	2:L:137:VAL:CG2	2.29	0.58
1:A:387:PHE:CE2	2:L:168:THR:HG21	2.36	0.58
1:A:345:LEU:HD22	2:L:137:VAL:CB	2.34	0.57
2:L:15:LEU:N	2:L:110:ILE:HD11	2.20	0.57
1:A:320:PHE:CE2	1:A:321:PHE:CZ	2.93	0.56
1:A:251:TYR:CE1	1:A:270:TYR:HD1	2.23	0.56
1:A:387:PHE:HD2	2:L:168:THR:HG22	1.67	0.56
1:A:358:ALA:CB	2:L:122:PHE:CZ	2.87	0.56
2:L:95:SER:HA	2:L:100:PHE:HD1	1.71	0.56
2:L:84:GLU:HG2	2:L:173:LYS:O	2.04	0.56
1:A:318:LEU:CD2	2:L:93:GLN:HE22	2.19	0.56
1:A:313:PHE:CE1	2:L:47:PRO:HB3	2.40	0.56
2:L:197:THR:HB	2:L:212:SER:HB2	1.87	0.56
1:A:345:LEU:HD22	2:L:137:VAL:CG2	2.36	0.55
1:A:242:VAL:HG23	1:A:245:TYR:CE1	2.42	0.55
1:A:344:PRO:CD	2:L:125:SER:HB2	2.16	0.55
1:A:387:PHE:HD2	2:L:168:THR:CG2	2.15	0.55
1:A:245:TYR:CG	1:A:316:ARG:NH1	2.75	0.55
2:L:11:LEU:HD11	2:L:21:ILE:CG2	2.37	0.54
1:A:319:THR:CB	2:L:100:PHE:CZ	2.86	0.54
1:A:316:ARG:HB3	1:A:323:VAL:CG1	2.36	0.54
1:A:319:THR:HG21	2:L:100:PHE:CZ	2.42	0.54
1:A:358:ALA:C	2:L:122:PHE:HE2	1.88	0.54
2:L:95:SER:HB2	2:L:100:PHE:CE1	2.43	0.54
2:L:41:GLN:HB2	2:L:51:LEU:HD11	1.90	0.53
1:A:235:THR:HG22	1:A:302:SER:HA	1.90	0.53
2:L:51:LEU:O	2:L:62:ILE:HD13	2.08	0.53
1:A:430:LYS:HZ3	2:L:127:GLU:CG	2.11	0.53
1:A:358:ALA:HB2	2:L:120:SER:OG	2.08	0.53
1:A:254:TRP:HB2	1:A:266:ILE:HD11	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:ASN:ND2	1:A:265:TRP:CD1	2.77	0.53
2:L:29:VAL:CG2	2:L:96:ASN:HB2	2.30	0.53
1:A:430:LYS:HZ1	2:L:127:GLU:CG	2.09	0.53
1:A:252:ILE:HD11	1:A:314:CYS:HB2	1.91	0.52
1:A:263:LEU:HD22	2:L:48:PRO:HG2	1.87	0.52
2:L:38:ASN:ND2	2:L:53:TYR:HB3	2.24	0.52
1:A:358:ALA:CB	2:L:120:SER:OG	2.57	0.52
2:L:95:SER:HB2	2:L:100:PHE:HE1	1.75	0.52
2:L:140:LEU:HG	2:L:179:MET:SD	2.50	0.52
1:A:390:VAL:HG21	2:L:164:LEU:HD21	1.91	0.52
1:A:242:VAL:HG13	1:A:295:ASP:HB3	1.92	0.51
1:A:316:ARG:HB3	1:A:323:VAL:HG12	1.93	0.50
1:A:253:ASN:ND2	1:A:318:LEU:HD13	2.26	0.50
1:A:270:TYR:HD2	1:A:273:TYR:H	1.58	0.50
2:L:163:VAL:HA	2:L:182:THR:O	2.12	0.50
1:A:242:VAL:HG11	1:A:247:PHE:HD1	1.67	0.50
1:A:358:ALA:HB3	2:L:122:PHE:CZ	2.47	0.49
2:L:52:ILE:HD11	2:L:58:LEU:HD23	1.92	0.49
1:A:321:PHE:HB2	1:A:322:ASP:CA	2.41	0.49
1:A:318:LEU:HD22	2:L:93:GLN:NE2	2.25	0.49
1:A:359:LEU:N	2:L:122:PHE:CE2	2.80	0.49
2:L:80:HIS:CG	2:L:81:PRO:HD3	2.47	0.49
2:L:15:LEU:HG	2:L:82:VAL:O	2.13	0.49
1:A:430:LYS:CE	2:L:127:GLU:HG2	2.36	0.48
1:A:254:TRP:O	1:A:266:ILE:HG12	2.13	0.48
1:A:254:TRP:HB2	1:A:266:ILE:CG1	2.43	0.48
1:A:344:PRO:O	2:L:125:SER:HB3	2.13	0.48
1:A:242:VAL:HG12	1:A:295:ASP:HB3	1.95	0.48
1:A:319:THR:CG2	2:L:100:PHE:CZ	2.97	0.48
1:A:251:TYR:CE2	1:A:268:ARG:HD2	2.49	0.47
1:A:318:LEU:CD2	2:L:93:GLN:NE2	2.77	0.47
1:A:248:THR:HA	1:A:271:PRO:HB2	1.95	0.47
2:L:36:TYR:HB3	2:L:95:SER:OG	2.14	0.47
1:A:245:TYR:CE1	1:A:316:ARG:NH1	2.83	0.47
1:A:251:TYR:CE1	1:A:270:TYR:CD1	3.03	0.47
2:L:63:PRO:HB2	2:L:65:ARG:HG2	1.98	0.46
1:A:348:SER:HA	2:L:121:ILE:O	2.15	0.46
1:A:346:ALA:C	2:L:122:PHE:HD2	2.07	0.46
1:A:265:TRP:CZ2	1:A:267:GLY:HA2	2.50	0.46
1:A:344:PRO:CD	2:L:125:SER:HB3	2.31	0.46
2:L:98:ASP:OD1	2:L:100:PHE:CE2	2.69	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LEU:CD2	2:L:48:PRO:HG3	2.26	0.46
2:L:38:ASN:HD21	2:L:50:LEU:HD11	1.81	0.46
1:A:430:LYS:CE	2:L:127:GLU:CG	2.93	0.45
1:A:340:PRO:HB3	1:A:366:TYR:HB3	1.98	0.45
1:A:371:VAL:CG2	1:A:399:LEU:HD21	2.46	0.45
1:A:321:PHE:HD2	1:A:322:ASP:CG	2.20	0.45
1:A:359:LEU:CA	2:L:122:PHE:HE2	2.21	0.45
1:A:319:THR:HG21	2:L:100:PHE:HZ	1.81	0.45
1:A:320:PHE:O	1:A:321:PHE:CG	2.70	0.45
1:A:254:TRP:CB	1:A:266:ILE:HD11	2.47	0.45
2:L:190:TYR:CE2	2:L:215:ARG:HD2	2.52	0.45
1:A:360:GLY:HA2	1:A:375:TRP:CH2	2.52	0.45
1:A:242:VAL:HG21	1:A:245:TYR:CE1	2.52	0.44
1:A:321:PHE:HD2	1:A:322:ASP:HB2	1.82	0.44
2:L:3:GLU:HG2	2:L:26:SER:CB	2.40	0.44
2:L:18:ARG:HD3	2:L:78:ASN:HD21	1.82	0.44
2:L:161:ASN:ND2	2:L:162:GLY:H	2.16	0.44
2:L:38:ASN:HD22	2:L:53:TYR:HB3	1.82	0.43
1:A:238:ILE:CD1	1:A:254:TRP:CZ3	3.02	0.43
2:L:190:TYR:CZ	2:L:215:ARG:HD2	2.53	0.43
1:A:359:LEU:N	2:L:122:PHE:CZ	2.87	0.43
1:A:251:TYR:HE1	1:A:270:TYR:CD1	2.35	0.43
2:L:62:ILE:HA	2:L:63:PRO:HD3	1.82	0.43
2:L:50:LEU:HD22	2:L:59:GLU:HB2	2.00	0.43
1:A:269:ILE:O	1:A:271:PRO:HD3	2.18	0.43
2:L:52:ILE:HD13	2:L:58:LEU:HD23	1.98	0.42
1:A:358:ALA:HB1	2:L:122:PHE:HZ	1.81	0.42
1:A:255:MET:SD	1:A:264:GLU:C	2.98	0.42
2:L:186:THR:OG1	2:L:189:GLU:HG3	2.19	0.42
2:L:58:LEU:CB	2:L:62:ILE:HD11	2.50	0.41
1:A:320:PHE:C	1:A:321:PHE:CD1	2.94	0.41
1:A:266:ILE:HD11	1:A:299:MET:HG3	2.02	0.41
1:A:238:ILE:CG1	1:A:299:MET:HB3	2.51	0.41
1:A:358:ALA:HB3	2:L:120:SER:HB2	1.92	0.41
2:L:120:SER:O	2:L:138:CYS:HA	2.20	0.41
1:A:268:ARG:NH1	2:L:98:ASP:CG	2.74	0.41
2:L:2:ILE:CD1	2:L:97:GLU:HG2	2.50	0.41
2:L:12:ALA:HA	2:L:109:GLU:O	2.21	0.41
1:A:238:ILE:HG12	1:A:299:MET:HB3	2.04	0.40
1:A:236:VAL:O	1:A:300:GLU:HA	2.21	0.40
2:L:160:GLN:HE21	2:L:160:GLN:HB3	1.77	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:95:SER:HA	2:L:100:PHE:CE1	2.57	0.40
1:A:318:LEU:HD23	1:A:318:LEU:C	2.42	0.40
1:A:416:ILE:HG12	1:A:431:ARG:HG2	2.03	0.40
2:L:51:LEU:O	2:L:62:ILE:CD1	2.70	0.40
2:L:161:ASN:HD22	2:L:162:GLY:H	1.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	216/218 (99%)	211 (98%)	3 (1%)	2 (1%)	21	67
2	L	213/215 (99%)	203 (95%)	8 (4%)	2 (1%)	21	67
All	All	429/433 (99%)	414 (96%)	11 (3%)	4 (1%)	26	67

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	322	ASP
2	L	81	PRO
1	A	354	GLY
2	L	80	HIS

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	186/186 (100%)	186 (100%)	0	100	100
2	L	189/189 (100%)	184 (97%)	5 (3%)	54	80
All	All	375/375 (100%)	370 (99%)	5 (1%)	78	89

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	L	24	LYS
2	L	129	LEU
2	L	140	LEU
2	L	163	VAL
2	L	197	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	253	ASN
2	L	38	ASN
2	L	160	GLN
2	L	161	ASN
2	L	214	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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.