



Full wwPDB/EMDatabank EM Map/Model Validation Report ⓘ

Apr 8, 2019 – 10:14 PM EDT

PDB ID : 6J0C
EMDB ID: : EMD-9761
Title : Cryo-EM Structure of an Extracellular Contractile Injection System, PVC sheath complex in contracted state
Authors : Jiang, F.; Li, N.; Wang, X.; Cheng, J.; Huang, Y.; Yang, Y.; Yang, J.; Cai, B.; Wang, Y.; Jin, Q.; Gao, N.
Deposited on : 2018-12-23
Resolution : 3.70 Å(reported)

This is a Full wwPDB/EMDatabank EM Map/Model Validation Report
for a publicly released PDB/EMDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

MolProbity : 4.02b-467
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031633

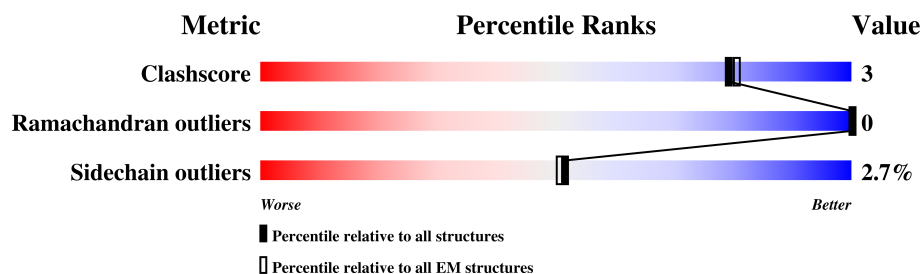
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.70 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	136327	1886
Ramachandran outliers	132723	1663
Sidechain outliers	132532	1531

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	355	83%	17%
1	B	355	82%	18%
1	C	355	83%	17%
1	D	355	81%	19%
1	E	355	82%	17%
1	F	355	82%	17%
1	a	355	97%	.
1	b	355	97%	.
1	c	355	97%	.

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Mol	Chain	Length	Quality of chain
1	d	355	<div><div></div><div>97%</div><div></div></div>
1	e	355	<div><div></div><div>97%</div><div></div></div>
1	f	355	<div><div></div><div>97%</div><div></div></div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 33108 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.

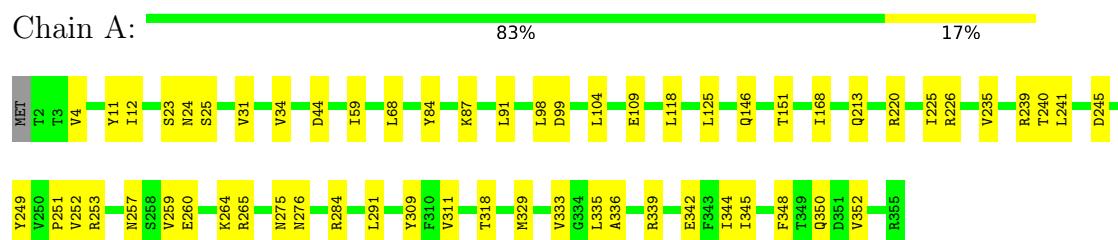
- Molecule 1 is a protein called Pvc2.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	B	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	C	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	D	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	E	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	F	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	a	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	b	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	c	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	d	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	e	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		
1	f	354	Total	C	N	O	S	0	0
			2759	1760	457	533	9		

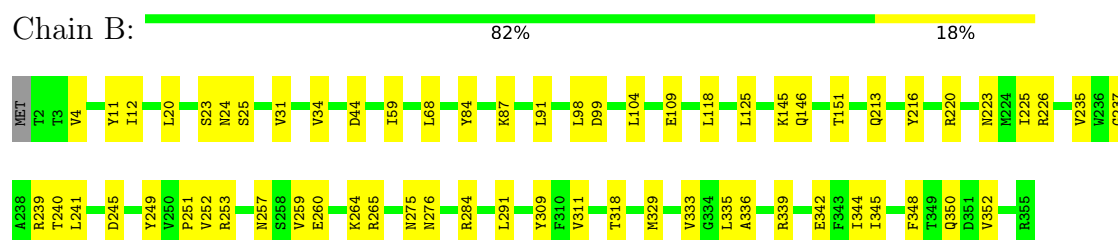
3 Residue-property plots

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

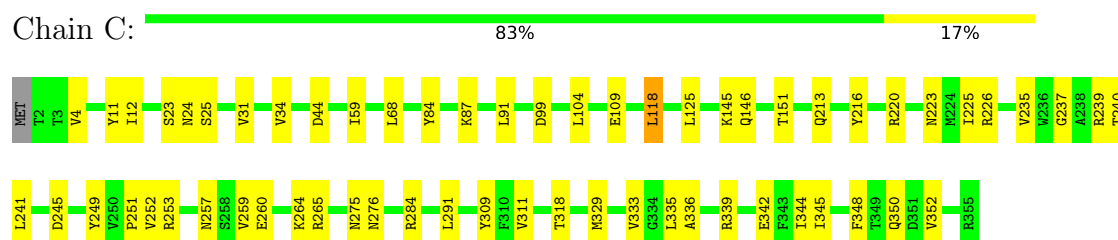
• Molecule 1: Pvc2



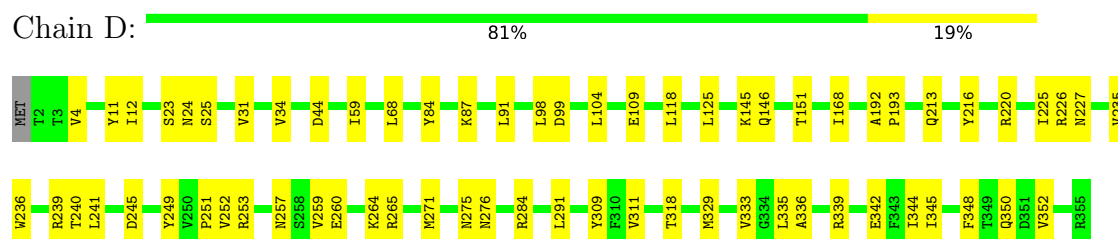
• Molecule 1: Pvc2



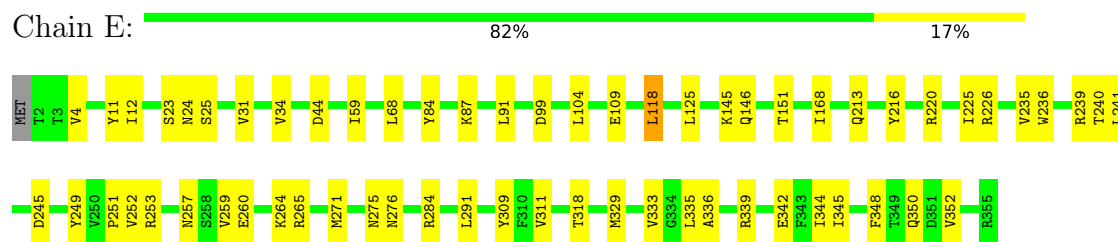
• Molecule 1: Pvc2



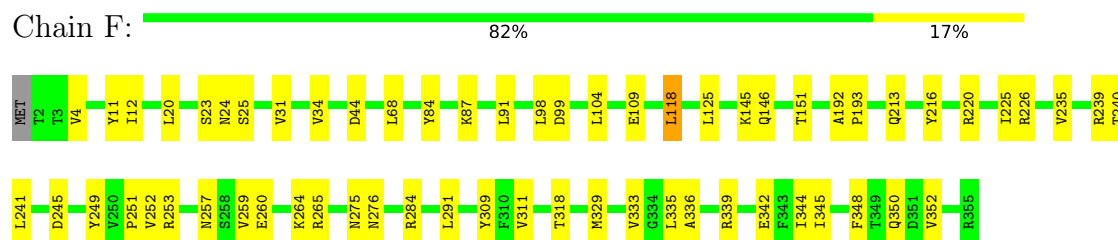
• Molecule 1: Pvc2



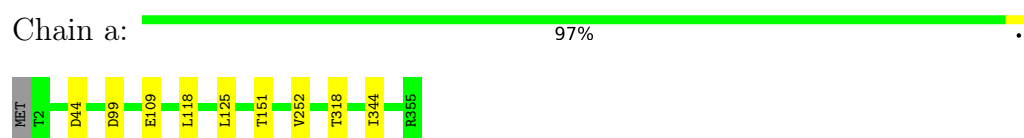
• Molecule 1: Pvc2



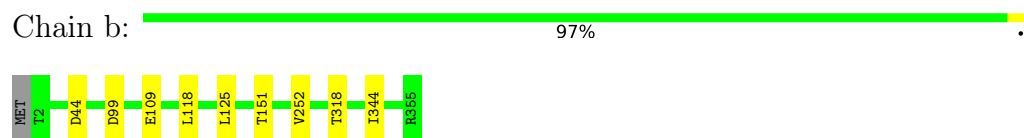
• Molecule 1: Pvc2



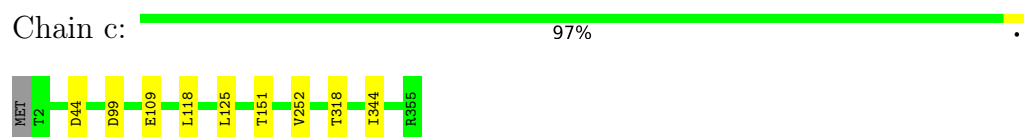
• Molecule 1: Pvc2



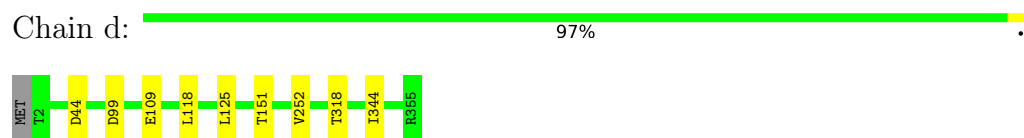
• Molecule 1: Pvc2



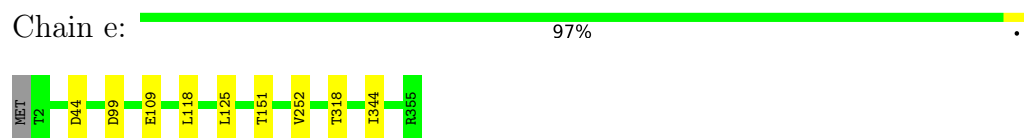
• Molecule 1: Pvc2



• Molecule 1: Pvc2



• Molecule 1: Pvc2



● Molecule 1: Pvc2

Chain f:  97%

4 Experimental information

Property	Value	Source
Reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=31.4°, rise=17 Å, axial sym=C6	Depositor
Number of segments used	36000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	46.4	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	FEI FALCON II (4k x 4k)	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	0.36	0/2819	0.56	1/3842 (0.0%)
1	B	0.36	0/2819	0.56	1/3842 (0.0%)
1	C	0.36	0/2819	0.56	1/3842 (0.0%)
1	D	0.36	0/2819	0.56	1/3842 (0.0%)
1	E	0.36	0/2819	0.56	1/3842 (0.0%)
1	F	0.36	0/2819	0.56	1/3842 (0.0%)
1	a	0.36	0/2819	0.56	1/3842 (0.0%)
1	b	0.36	0/2819	0.56	1/3842 (0.0%)
1	c	0.36	0/2819	0.56	1/3842 (0.0%)
1	d	0.36	0/2819	0.56	1/3842 (0.0%)
1	e	0.36	0/2819	0.56	1/3842 (0.0%)
1	f	0.36	0/2819	0.56	1/3842 (0.0%)
All	All	0.36	0/33828	0.56	12/46104 (0.0%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	125	LEU	CA-CB-CG	7.26	131.99	115.30
1	e	125	LEU	CA-CB-CG	7.26	131.99	115.30
1	b	125	LEU	CA-CB-CG	7.25	131.98	115.30
1	D	125	LEU	CA-CB-CG	7.25	131.97	115.30
1	A	125	LEU	CA-CB-CG	7.25	131.97	115.30
1	a	125	LEU	CA-CB-CG	7.24	131.95	115.30
1	d	125	LEU	CA-CB-CG	7.24	131.95	115.30
1	C	125	LEU	CA-CB-CG	7.24	131.94	115.30
1	E	125	LEU	CA-CB-CG	7.24	131.94	115.30
1	c	125	LEU	CA-CB-CG	7.23	131.93	115.30
1	B	125	LEU	CA-CB-CG	7.23	131.92	115.30
1	f	125	LEU	CA-CB-CG	7.22	131.91	115.30

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	2759	0	2697	33	0
1	B	2759	0	2697	34	0
1	C	2759	0	2697	34	0
1	D	2759	0	2697	37	0
1	E	2759	0	2697	36	0
1	F	2759	0	2697	34	0
1	a	2759	0	2697	0	0
1	b	2759	0	2697	0	0
1	c	2759	0	2697	0	0
1	d	2759	0	2697	0	0
1	e	2759	0	2697	0	0
1	f	2759	0	2697	0	0
All	All	33108	0	32364	196	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 (196) 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:B:336:ALA:HB2	1:B:352:VAL:HG21	36.22	0.63
1:C:336:ALA:HB2	1:C:352:VAL:HG21	36.22	0.63
1:E:31:VAL:HG23	1:E:84:TYR:HB2	1.81	0.62
1:D:31:VAL:HG23	1:D:84:TYR:HB2	1.81	0.62
1:A:31:VAL:HG23	1:A:84:TYR:HB2	1.81	0.62
1:B:31:VAL:HG23	1:B:84:TYR:HB2	1.81	0.61
1:F:31:VAL:HG23	1:F:84:TYR:HB2	1.81	0.61
1:E:336:ALA:HB2	1:E:352:VAL:HG21	36.22	0.61
1:C:31:VAL:HG23	1:C:84:TYR:HB2	1.81	0.60
1:A:264:LYS:HG3	1:F:12:ILE:HD11	45.42	0.60
1:D:336:ALA:HB2	1:D:352:VAL:HG21	36.22	0.60
1:F:336:ALA:HB2	1:F:352:VAL:HG21	36.22	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:ILE:HD11	1:E:264:LYS:HG3	45.42	0.60
1:A:12:ILE:HD11	1:B:264:LYS:HG3	45.41	0.59
1:A:336:ALA:HB2	1:A:352:VAL:HG21	36.22	0.58
1:E:12:ILE:HD11	1:F:264:LYS:HG3	45.44	0.57
1:B:12:ILE:HD11	1:C:264:LYS:HG3	45.43	0.56
1:A:345:ILE:HG22	1:F:11:TYR:HB2	1.87	0.56
1:B:275:ASN:HD22	1:B:329:MET:HB3	1.71	0.56
1:F:275:ASN:HD22	1:F:329:MET:HB3	1.71	0.56
1:A:275:ASN:HD22	1:A:329:MET:HB3	1.71	0.55
1:A:11:TYR:HB2	1:B:345:ILE:HG22	1.89	0.55
1:C:311:VAL:HG12	1:C:333:VAL:HG22	1.89	0.55
1:B:311:VAL:HG12	1:B:333:VAL:HG22	1.89	0.55
1:C:12:ILE:HD11	1:D:264:LYS:HG3	45.44	0.55
1:D:311:VAL:HG12	1:D:333:VAL:HG22	1.89	0.55
1:A:311:VAL:HG12	1:A:333:VAL:HG22	1.89	0.55
1:E:311:VAL:HG12	1:E:333:VAL:HG22	1.89	0.55
1:F:311:VAL:HG12	1:F:333:VAL:HG22	1.89	0.55
1:A:253:ARG:NH2	1:A:257:ASN:OD1	2.41	0.54
1:B:253:ARG:NH2	1:B:257:ASN:OD1	2.41	0.54
1:C:253:ARG:NH2	1:C:257:ASN:OD1	2.41	0.54
1:E:253:ARG:NH2	1:E:257:ASN:OD1	2.41	0.54
1:F:253:ARG:NH2	1:F:257:ASN:OD1	2.41	0.54
1:C:275:ASN:HD22	1:C:329:MET:HB3	1.71	0.54
1:D:275:ASN:HD22	1:D:329:MET:HB3	1.71	0.54
1:D:253:ARG:NH2	1:D:257:ASN:OD1	2.41	0.54
1:E:275:ASN:HD22	1:E:329:MET:HB3	1.71	0.54
1:E:11:TYR:HB2	1:F:345:ILE:HG22	1.91	0.53
1:B:11:TYR:HB2	1:C:345:ILE:HG22	1.90	0.53
1:D:11:TYR:HB2	1:E:345:ILE:HG22	1.92	0.52
1:F:225:ILE:HG12	1:F:235:VAL:HG12	1.92	0.52
1:A:225:ILE:HG12	1:A:235:VAL:HG12	1.92	0.52
1:C:4:VAL:HG12	1:C:25:SER:HA	43.15	0.52
1:E:225:ILE:HG12	1:E:235:VAL:HG12	1.92	0.52
1:B:225:ILE:HG12	1:B:235:VAL:HG12	1.92	0.52
1:A:4:VAL:HG12	1:A:25:SER:HA	43.16	0.51
1:B:213:GLN:OE1	1:B:226:ARG:NH2	2.44	0.51
1:A:213:GLN:OE1	1:A:226:ARG:NH2	2.44	0.51
1:C:11:TYR:HB2	1:D:345:ILE:HG22	1.92	0.51
1:C:225:ILE:HG12	1:C:235:VAL:HG12	1.92	0.51
1:D:213:GLN:OE1	1:D:226:ARG:NH2	2.44	0.51
1:D:4:VAL:HG12	1:D:25:SER:HA	43.15	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:213:GLN:OE1	1:E:226:ARG:NH2	2.44	0.51
1:B:4:VAL:HG12	1:B:25:SER:HA	43.15	0.51
1:C:213:GLN:OE1	1:C:226:ARG:NH2	2.44	0.51
1:D:225:ILE:HG12	1:D:235:VAL:HG12	1.92	0.51
1:F:213:GLN:OE1	1:F:226:ARG:NH2	2.44	0.51
1:A:239:ARG:NH1	1:A:342:GLU:OE2	2.45	0.50
1:B:239:ARG:NH1	1:B:342:GLU:OE2	2.45	0.50
1:F:4:VAL:HG12	1:F:25:SER:HA	43.15	0.50
1:E:239:ARG:NH1	1:E:342:GLU:OE2	2.45	0.50
1:E:4:VAL:HG12	1:E:25:SER:HA	43.15	0.50
1:D:239:ARG:NH1	1:D:342:GLU:OE2	2.45	0.50
1:F:239:ARG:NH1	1:F:342:GLU:OE2	2.45	0.50
1:C:239:ARG:NH1	1:C:342:GLU:OE2	2.45	0.49
1:C:245:ASP:O	1:C:339:ARG:NH1	2.45	0.49
1:D:245:ASP:O	1:D:339:ARG:NH1	2.45	0.48
1:D:24:ASN:OD1	1:D:24:ASN:N	2.47	0.48
1:B:260:GLU:HG3	1:B:348:PHE:CE2	40.15	0.47
1:E:24:ASN:OD1	1:E:24:ASN:N	2.47	0.47
1:F:23:SER:OG	1:F:265:ARG:NH2	2.47	0.47
1:B:23:SER:OG	1:B:265:ARG:NH2	2.47	0.47
1:C:23:SER:OG	1:C:265:ARG:NH2	2.47	0.47
1:E:23:SER:OG	1:E:265:ARG:NH2	2.47	0.47
1:C:260:GLU:HG3	1:C:348:PHE:CE2	40.15	0.47
1:B:291:LEU:HD23	1:B:309:TYR:HB2	1.96	0.47
1:E:245:ASP:O	1:E:339:ARG:NH1	2.45	0.47
1:A:291:LEU:HD23	1:A:309:TYR:HB2	1.96	0.47
1:D:291:LEU:HD23	1:D:309:TYR:HB2	1.96	0.47
1:E:291:LEU:HD23	1:E:309:TYR:HB2	1.96	0.47
1:A:245:ASP:O	1:A:339:ARG:NH1	2.45	0.47
1:B:245:ASP:O	1:B:339:ARG:NH1	2.45	0.47
1:C:24:ASN:OD1	1:C:24:ASN:N	2.47	0.47
1:F:291:LEU:HD23	1:F:309:TYR:HB2	1.96	0.47
1:B:24:ASN:OD1	1:B:24:ASN:N	2.47	0.47
1:C:291:LEU:HD23	1:C:309:TYR:HB2	1.96	0.46
1:F:245:ASP:O	1:F:339:ARG:NH1	2.45	0.46
1:D:260:GLU:HG3	1:D:348:PHE:CE2	40.15	0.46
1:A:23:SER:OG	1:A:265:ARG:NH2	2.47	0.46
1:B:284:ARG:NH1	1:B:311:VAL:O	2.49	0.46
1:C:284:ARG:NH1	1:C:311:VAL:O	2.49	0.46
1:D:23:SER:OG	1:D:265:ARG:NH2	2.47	0.46
1:D:284:ARG:NH1	1:D:311:VAL:O	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:284:ARG:NH1	1:F:311:VAL:O	2.49	0.46
1:A:284:ARG:NH1	1:A:311:VAL:O	2.49	0.46
1:E:260:GLU:HG3	1:E:348:PHE:CE2	40.15	0.45
1:E:284:ARG:NH1	1:E:311:VAL:O	2.49	0.45
1:B:249:TYR:HB3	1:B:251:PRO:HD2	1.98	0.45
1:E:249:TYR:HB3	1:E:251:PRO:HD2	1.98	0.45
1:C:249:TYR:HB3	1:C:251:PRO:HD2	1.98	0.45
1:F:240:THR:OG1	1:F:241:LEU:N	2.50	0.45
1:B:240:THR:OG1	1:B:241:LEU:N	2.50	0.45
1:F:260:GLU:HG3	1:F:348:PHE:CE2	40.15	0.45
1:A:34:VAL:HG22	1:A:87:LYS:HA	1.99	0.45
1:D:249:TYR:HB3	1:D:251:PRO:HD2	1.98	0.45
1:F:249:TYR:HB3	1:F:251:PRO:HD2	1.98	0.45
1:A:240:THR:OG1	1:A:241:LEU:N	2.50	0.45
1:D:240:THR:OG1	1:D:241:LEU:N	2.50	0.45
1:F:34:VAL:HG22	1:F:87:LYS:HA	1.99	0.44
1:A:249:TYR:HB3	1:A:251:PRO:HD2	1.98	0.44
1:C:240:THR:OG1	1:C:241:LEU:N	2.50	0.44
1:C:34:VAL:HB	1:C:68:LEU:HD22	1.99	0.44
1:E:240:THR:OG1	1:E:241:LEU:N	2.50	0.44
1:B:34:VAL:HG22	1:B:87:LYS:HA	1.99	0.44
1:D:145:LYS:HZ3	1:D:216:TYR:HD2	1.69	0.44
1:D:34:VAL:HB	1:D:68:LEU:HD22	1.99	0.44
1:F:146:GLN:HA	1:F:220:ARG:HH21	1.83	0.44
1:A:260:GLU:HG3	1:A:348:PHE:CE2	40.15	0.44
1:C:145:LYS:HZ3	1:C:216:TYR:HD2	1.68	0.44
1:A:24:ASN:OD1	1:A:24:ASN:N	2.47	0.44
1:B:34:VAL:HB	1:B:68:LEU:HD22	1.99	0.43
1:D:339:ARG:HH21	1:D:350:GLN:NE2	38.78	0.43
1:E:34:VAL:HB	1:E:68:LEU:HD22	1.99	0.43
1:F:34:VAL:HB	1:F:68:LEU:HD22	1.99	0.43
1:B:146:GLN:HA	1:B:220:ARG:HH21	1.83	0.43
1:E:168:ILE:HA	1:E:168:ILE:HD12	1.92	0.43
1:A:34:VAL:HB	1:A:68:LEU:HD22	1.99	0.43
1:C:146:GLN:HA	1:C:220:ARG:HH21	1.83	0.43
1:E:59:ILE:HD13	1:E:59:ILE:HA	1.87	0.43
1:F:118:LEU:HA	1:F:118:LEU:HD22	1.91	0.43
1:A:168:ILE:HA	1:A:168:ILE:HD12	1.92	0.43
1:B:145:LYS:HZ3	1:B:216:TYR:HD2	1.66	0.43
1:B:59:ILE:HA	1:B:59:ILE:HD13	1.87	0.43
1:C:34:VAL:HG22	1:C:87:LYS:HA	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:34:VAL:HG22	1:E:87:LYS:HA	1.99	0.43
1:C:59:ILE:HA	1:C:59:ILE:HD13	1.88	0.43
1:E:146:GLN:HA	1:E:220:ARG:HH21	1.83	0.43
1:F:98:LEU:HD23	1:F:98:LEU:HA	1.87	0.43
1:A:146:GLN:HA	1:A:220:ARG:HH21	1.83	0.43
1:D:146:GLN:HA	1:D:220:ARG:HH21	1.83	0.43
1:D:34:VAL:HG22	1:D:87:LYS:HA	1.99	0.43
1:A:98:LEU:HA	1:A:98:LEU:HD23	1.88	0.43
1:C:339:ARG:HH21	1:C:350:GLN:NE2	38.77	0.43
1:E:339:ARG:HH21	1:E:350:GLN:NE2	38.77	0.43
1:D:168:ILE:HA	1:D:168:ILE:HD12	1.92	0.42
1:E:31:VAL:HG12	1:E:104:LEU:HA	2.02	0.42
1:D:31:VAL:HG12	1:D:104:LEU:HA	2.02	0.42
1:D:236:TRP:CH2	1:D:271:MET:HG2	45.69	0.42
1:E:118:LEU:HA	1:E:118:LEU:HD22	1.91	0.42
1:B:259:VAL:HG11	1:B:335:LEU:HD11	2.02	0.42
1:C:31:VAL:HG12	1:C:104:LEU:HA	2.02	0.42
1:F:91:LEU:HD21	1:F:104:LEU:HD11	2.01	0.42
1:F:31:VAL:HG12	1:F:104:LEU:HA	2.02	0.42
1:A:259:VAL:HG11	1:A:335:LEU:HD11	2.02	0.42
1:B:339:ARG:HH21	1:B:350:GLN:NE2	38.77	0.42
1:C:91:LEU:HD21	1:C:104:LEU:HD11	2.01	0.42
1:E:91:LEU:HD21	1:E:104:LEU:HD11	2.01	0.42
1:A:31:VAL:HG12	1:A:104:LEU:HA	2.02	0.42
1:B:91:LEU:HD21	1:B:104:LEU:HD11	2.01	0.42
1:E:276:ASN:OD1	1:E:276:ASN:N	2.53	0.42
1:C:118:LEU:HD22	1:C:118:LEU:HA	1.91	0.42
1:D:227:ASN:N	1:D:227:ASN:OD1	2.53	0.42
1:E:145:LYS:HZ3	1:E:216:TYR:HD2	1.68	0.42
1:F:24:ASN:N	1:F:24:ASN:OD1	2.47	0.42
1:B:31:VAL:HG12	1:B:104:LEU:HA	2.02	0.42
1:B:276:ASN:OD1	1:B:276:ASN:N	2.53	0.42
1:C:259:VAL:HG11	1:C:335:LEU:HD11	2.02	0.42
1:D:339:ARG:HH21	1:D:350:GLN:HE22	39.43	0.42
1:D:91:LEU:HD21	1:D:104:LEU:HD11	2.01	0.42
1:D:98:LEU:HA	1:D:98:LEU:HD23	1.87	0.42
1:F:20:LEU:HA	1:F:20:LEU:HD12	1.92	0.42
1:A:276:ASN:OD1	1:A:276:ASN:N	2.53	0.41
1:F:145:LYS:HZ3	1:F:216:TYR:HD2	1.67	0.41
1:A:91:LEU:HD21	1:A:104:LEU:HD11	2.01	0.41
1:C:276:ASN:OD1	1:C:276:ASN:N	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:276:ASN:N	1:D:276:ASN:OD1	2.53	0.41
1:D:59:ILE:HD13	1:D:59:ILE:HA	1.87	0.41
1:F:259:VAL:HG11	1:F:335:LEU:HD11	2.02	0.41
1:A:59:ILE:HA	1:A:59:ILE:HD13	1.87	0.41
1:D:192:ALA:HA	1:D:193:PRO:HD3	1.93	0.41
1:C:339:ARG:HH21	1:C:350:GLN:HE22	39.43	0.41
1:F:192:ALA:HA	1:F:193:PRO:HD3	1.93	0.41
1:E:241:LEU:HD23	1:E:241:LEU:HA	1.94	0.41
1:D:259:VAL:HG11	1:D:335:LEU:HD11	2.02	0.41
1:F:339:ARG:HH21	1:F:350:GLN:NE2	38.77	0.41
1:E:259:VAL:HG11	1:E:335:LEU:HD11	2.02	0.41
1:A:339:ARG:HH21	1:A:350:GLN:NE2	38.77	0.41
1:B:98:LEU:HA	1:B:98:LEU:HD23	1.88	0.40
1:B:223:ASN:HB3	1:B:237:GLY:HA3	2.03	0.40
1:C:223:ASN:HB3	1:C:237:GLY:HA3	2.03	0.40
1:A:339:ARG:HH21	1:A:350:GLN:HE22	39.43	0.40
1:B:20:LEU:HD12	1:B:20:LEU:HA	1.92	0.40
1:E:236:TRP:CH2	1:E:271:MET:HG2	45.69	0.40
1:E:339:ARG:HH21	1:E:350:GLN:HE22	39.43	0.40
1:F:276:ASN:OD1	1:F:276:ASN:N	2.53	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 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	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	B	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	C	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	D	352/355 (99%)	329 (94%)	23 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	F	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	a	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	b	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	c	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	d	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	e	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
1	f	352/355 (99%)	329 (94%)	23 (6%)	0	100	100
All	All	4224/4260 (99%)	3948 (94%)	276 (6%)	0	100	100

There are no Ramachandran outliers to report.

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 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	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	B	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	C	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	D	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	E	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	F	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	a	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	b	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	c	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	d	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	e	291/294 (99%)	283 (97%)	8 (3%)	48	77
1	f	291/294 (99%)	283 (97%)	8 (3%)	48	77
All	All	3492/3528 (99%)	3396 (97%)	96 (3%)	51	77

All (96) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	ASP
1	A	99	ASP
1	A	109	GLU
1	A	118	LEU
1	A	151	THR
1	A	252	VAL
1	A	318	THR
1	A	344	ILE
1	B	44	ASP
1	B	99	ASP
1	B	109	GLU
1	B	118	LEU
1	B	151	THR
1	B	252	VAL
1	B	318	THR
1	B	344	ILE
1	C	44	ASP
1	C	99	ASP
1	C	109	GLU
1	C	118	LEU
1	C	151	THR
1	C	252	VAL
1	C	318	THR
1	C	344	ILE
1	D	44	ASP
1	D	99	ASP
1	D	109	GLU
1	D	118	LEU
1	D	151	THR
1	D	252	VAL
1	D	318	THR
1	D	344	ILE
1	E	44	ASP
1	E	99	ASP
1	E	109	GLU
1	E	118	LEU
1	E	151	THR
1	E	252	VAL
1	E	318	THR
1	E	344	ILE
1	F	44	ASP
1	F	99	ASP

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Mol	Chain	Res	Type
1	F	109	GLU
1	F	118	LEU
1	F	151	THR
1	F	252	VAL
1	F	318	THR
1	F	344	ILE
1	a	44	ASP
1	a	99	ASP
1	a	109	GLU
1	a	118	LEU
1	a	151	THR
1	a	252	VAL
1	a	318	THR
1	a	344	ILE
1	b	44	ASP
1	b	99	ASP
1	b	109	GLU
1	b	118	LEU
1	b	151	THR
1	b	252	VAL
1	b	318	THR
1	b	344	ILE
1	c	44	ASP
1	c	99	ASP
1	c	109	GLU
1	c	118	LEU
1	c	151	THR
1	c	252	VAL
1	c	318	THR
1	c	344	ILE
1	d	44	ASP
1	d	99	ASP
1	d	109	GLU
1	d	118	LEU
1	d	151	THR
1	d	252	VAL
1	d	318	THR
1	d	344	ILE
1	e	44	ASP
1	e	99	ASP
1	e	109	GLU
1	e	118	LEU

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Mol	Chain	Res	Type
1	e	151	THR
1	e	252	VAL
1	e	318	THR
1	e	344	ILE
1	f	44	ASP
1	f	99	ASP
1	f	109	GLU
1	f	118	LEU
1	f	151	THR
1	f	252	VAL
1	f	318	THR
1	f	344	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (56) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	78	ASN
1	A	195	ASN
1	A	246	ASN
1	B	57	ASN
1	B	63	ASN
1	B	78	ASN
1	B	195	ASN
1	B	246	ASN
1	C	57	ASN
1	C	78	ASN
1	C	195	ASN
1	C	246	ASN
1	D	57	ASN
1	D	78	ASN
1	D	195	ASN
1	D	246	ASN
1	E	57	ASN
1	E	78	ASN
1	E	195	ASN
1	E	246	ASN
1	F	57	ASN
1	F	78	ASN
1	F	195	ASN
1	F	246	ASN
1	a	57	ASN

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Mol	Chain	Res	Type
1	a	63	ASN
1	a	78	ASN
1	a	195	ASN
1	a	246	ASN
1	a	350	GLN
1	b	57	ASN
1	b	78	ASN
1	b	195	ASN
1	b	246	ASN
1	b	350	GLN
1	c	57	ASN
1	c	78	ASN
1	c	195	ASN
1	c	246	ASN
1	c	350	GLN
1	d	57	ASN
1	d	78	ASN
1	d	195	ASN
1	d	246	ASN
1	d	350	GLN
1	e	57	ASN
1	e	78	ASN
1	e	195	ASN
1	e	246	ASN
1	e	350	GLN
1	f	57	ASN
1	f	78	ASN
1	f	195	ASN
1	f	246	ASN
1	f	350	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](#)

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.