



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 10:44 AM UTC

PDB ID : 3FI1 / pdb_00003fi1
EMDB ID : EMD-5037
Title : NhaA dimer model
Authors : Appel, M.; Hizlan, D.; Vinothkumar, K.R.; Ziegler, C.; Kuehlbrandt, W.
Deposited on : 2008-12-10
Resolution : 7.00 Å(reported)

This is a Full wwPDB EM Validation 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/EMValidationReportHelp>

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:

EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : **NOT EXECUTED**
MapQ : **FAILED**
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

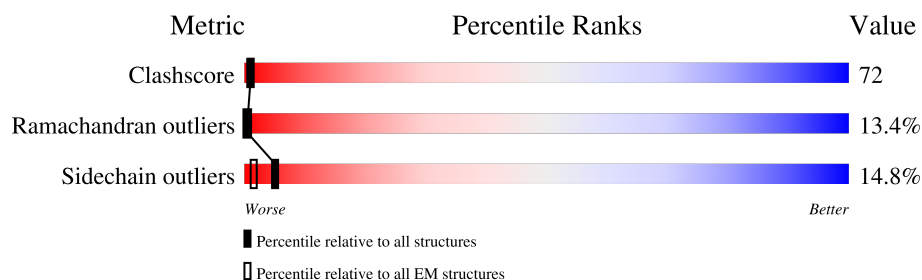
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON CRYSTALLOGRAPHY

The reported resolution of this entry is 7.00 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	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102

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

Mol	Chain	Length	Quality of chain
1	A	376	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2809 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.

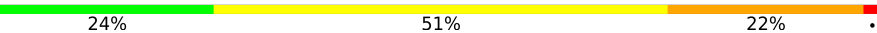
- Molecule 1 is a protein called Na(+)/H(+) antiporter nhaA.

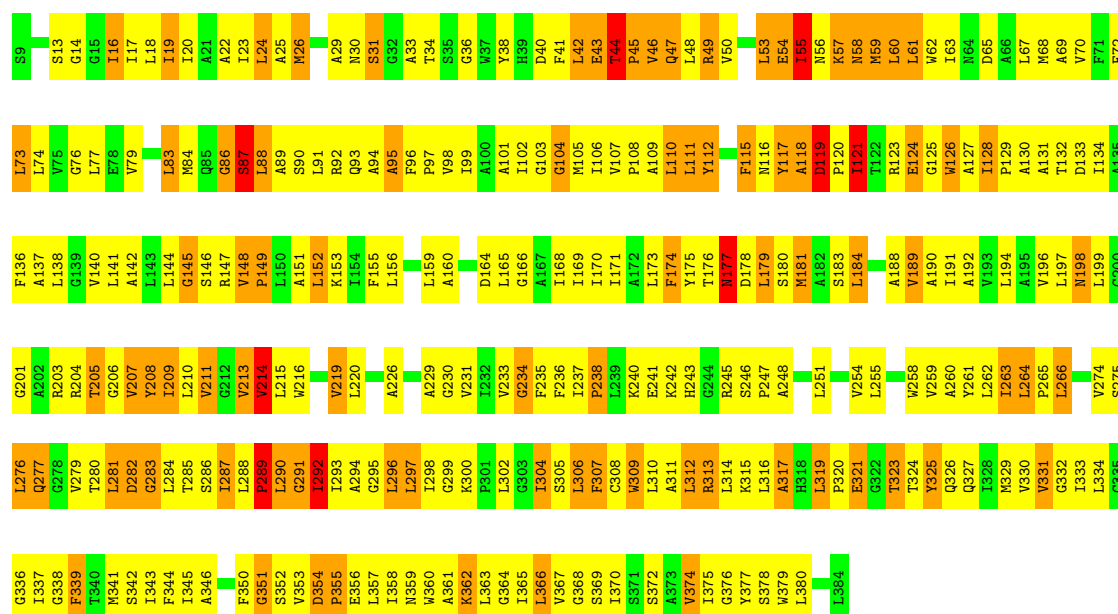
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	376	2809	1865	457	474	13	0	0

3 Residue-property plots

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. 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: Na(+)/H(+) antiporter nhaA

Chain A: 



4 Experimental information

Property	Value	Source
EM reconstruction method	CRYSTALLOGRAPHY	Depositor
Imposed symmetry	2D CRYSTAL, a =Not provided Å, b =Not provided Å, c =Not provided Å, γ =Not provided°, space group=Not provided	Depositor
Number of images used	Not provided	
Resolution determination method	Not provided	
CTF correction method	Not provided	
Microscope	JEOL 3000SFF	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	20	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
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	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.63	1/2868 (0.0%)	1.08	24/3910 (0.6%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	287	ILE	CA-CB	5.37	1.59	1.54

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	356	GLU	N-CA-C	-9.88	102.02	114.56
1	A	307	PHE	N-CA-C	-7.71	102.82	111.07
1	A	111	LEU	N-CA-C	-7.16	103.56	112.72
1	A	115	PHE	N-CA-C	-7.06	104.47	113.23
1	A	73	LEU	N-CA-C	-6.79	103.81	111.14
1	A	319	LEU	CA-C-N	6.73	126.19	118.85
1	A	319	LEU	C-N-CA	6.73	126.19	118.85
1	A	366	LEU	N-CA-C	-6.72	105.08	113.55
1	A	237	ILE	CA-C-N	6.70	128.22	119.84
1	A	237	ILE	C-N-CA	6.70	128.22	119.84
1	A	119	ASP	N-CA-C	6.68	124.57	109.81
1	A	325	TYR	CB-CA-C	-6.52	109.06	116.63
1	A	29	ALA	N-CA-C	6.49	118.01	111.07
1	A	374	VAL	N-CA-C	-6.37	106.77	112.12
1	A	297	LEU	N-CA-C	-6.36	105.52	113.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	121	ILE	N-CA-C	-6.24	107.78	113.71
1	A	87	SER	N-CA-C	6.09	123.77	110.80
1	A	357	LEU	N-CA-C	-5.81	106.05	113.02
1	A	152	LEU	N-CA-C	-5.73	103.79	112.04
1	A	31	SER	N-CA-C	5.58	118.90	111.75
1	A	304	ILE	N-CA-C	-5.49	105.02	111.00
1	A	211	VAL	N-CA-C	-5.28	104.81	110.36
1	A	331	VAL	N-CA-C	-5.25	105.59	110.53
1	A	214	VAL	N-CA-C	-5.00	105.72	110.42

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	112	TYR	Sidechain

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	2809	0	2993	415	5
All	All	2809	0	2993	415	5

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 72.

All (415) 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:56:ASN:O	1:A:57:LYS:HG2	1.36	1.20
1:A:119:ASP:HB3	1:A:120:PRO:HD3	1.19	1.18
1:A:48:LEU:HD23	1:A:49:ARG:N	1.63	1.14
1:A:42:LEU:O	1:A:59:MET:HB3	1.50	1.10
1:A:330:VAL:HG11	1:A:380:LEU:HB2	1.33	1.08
1:A:45:PRO:O	1:A:46:VAL:HG12	1.52	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:LEU:HD23	1:A:53:LEU:H	1.26	1.00
1:A:47:GLN:HB2	1:A:56:ASN:ND2	1.77	0.99
1:A:50:VAL:O	1:A:53:LEU:HD21	1.63	0.97
1:A:43:GLU:O	1:A:44:THR:HB	1.63	0.97
1:A:56:ASN:C	1:A:57:LYS:HG2	1.89	0.97
1:A:48:LEU:HD23	1:A:49:ARG:H	1.26	0.95
1:A:47:GLN:CB	1:A:56:ASN:ND2	2.30	0.94
1:A:96:PHE:HB3	1:A:97:PRO:HD3	1.47	0.94
1:A:42:LEU:O	1:A:59:MET:CB	2.16	0.93
1:A:290:LEU:HD22	1:A:291:GLY:H	1.34	0.93
1:A:112:TYR:CD2	1:A:125:GLY:HA2	2.04	0.92
1:A:45:PRO:O	1:A:46:VAL:CG1	2.17	0.92
1:A:119:ASP:HB3	1:A:120:PRO:CD	2.00	0.92
1:A:121:ILE:HB	1:A:124:GLU:OE1	1.71	0.91
1:A:304:ILE:HD13	1:A:333:ILE:HG12	1.55	0.89
1:A:374:VAL:HG23	1:A:375:ILE:HD12	1.55	0.89
1:A:134:ILE:HB	1:A:160:ALA:HB1	1.55	0.88
1:A:281:LEU:HD22	1:A:360:TRP:CH2	2.08	0.87
1:A:125:GLY:O	1:A:127:ALA:N	2.06	0.86
1:A:297:LEU:HD21	1:A:367:VAL:HG12	1.57	0.86
1:A:45:PRO:HA	1:A:57:LYS:O	1.76	0.85
1:A:107:VAL:HB	1:A:306:LEU:HD12	1.58	0.85
1:A:128:ILE:HB	1:A:129:PRO:HD3	1.58	0.85
1:A:47:GLN:HB2	1:A:56:ASN:CG	2.01	0.85
1:A:54:GLU:C	1:A:55:ILE:HG13	2.02	0.84
1:A:54:GLU:O	1:A:55:ILE:HG12	1.76	0.84
1:A:88:LEU:HD23	1:A:88:LEU:O	1.78	0.84
1:A:48:LEU:CD2	1:A:49:ARG:H	1.91	0.84
1:A:148:VAL:HG13	1:A:324:THR:HG21	1.60	0.83
1:A:174:PHE:HD2	1:A:175:TYR:N	1.74	0.83
1:A:112:TYR:CE2	1:A:125:GLY:HA2	2.14	0.82
1:A:53:LEU:HD23	1:A:53:LEU:N	1.95	0.82
1:A:108:PRO:HD3	1:A:306:LEU:HD12	1.62	0.81
1:A:103:GLY:O	1:A:306:LEU:HD11	1.80	0.80
1:A:49:ARG:HA	1:A:53:LEU:O	1.81	0.80
1:A:290:LEU:HD13	1:A:291:GLY:N	1.96	0.80
1:A:312:LEU:C	1:A:314:LEU:H	1.88	0.80
1:A:44:THR:HG23	1:A:44:THR:O	1.80	0.79
1:A:54:GLU:O	1:A:55:ILE:CG1	2.31	0.79
1:A:128:ILE:HG21	1:A:296:LEU:HD12	1.64	0.78
1:A:124:GLU:HB2	1:A:350:PHE:CZ	2.18	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:LEU:O	1:A:288:LEU:HB3	1.84	0.78
1:A:54:GLU:C	1:A:55:ILE:CG1	2.57	0.78
1:A:14:GLY:O	1:A:17:ILE:HG22	1.83	0.77
1:A:50:VAL:HG13	1:A:53:LEU:HD21	1.66	0.77
1:A:99:ILE:HG21	1:A:314:LEU:HD21	1.66	0.77
1:A:104:GLY:N	1:A:306:LEU:HD21	2.00	0.76
1:A:286:SER:C	1:A:289:PRO:HD2	2.10	0.75
1:A:104:GLY:HA2	1:A:306:LEU:CD1	2.17	0.75
1:A:119:ASP:CB	1:A:120:PRO:HD3	2.10	0.75
1:A:45:PRO:O	1:A:46:VAL:CB	2.35	0.74
1:A:104:GLY:CA	1:A:306:LEU:HD11	2.17	0.74
1:A:124:GLU:HB2	1:A:350:PHE:HZ	1.51	0.74
1:A:285:THR:O	1:A:289:PRO:HG2	1.88	0.73
1:A:334:LEU:C	1:A:336:GLY:H	1.96	0.73
1:A:288:LEU:HB3	1:A:289:PRO:HD3	1.69	0.73
1:A:44:THR:O	1:A:44:THR:CG2	2.37	0.72
1:A:48:LEU:CD2	1:A:49:ARG:N	2.46	0.72
1:A:131:ALA:HB1	1:A:341:MET:HB3	1.71	0.72
1:A:50:VAL:O	1:A:53:LEU:CD2	2.38	0.72
1:A:45:PRO:C	1:A:57:LYS:O	2.34	0.71
1:A:194:LEU:HB3	1:A:236:PHE:CD2	2.26	0.71
1:A:334:LEU:O	1:A:337:ILE:HG12	1.89	0.71
1:A:366:LEU:O	1:A:370:ILE:HG13	1.90	0.71
1:A:83:LEU:O	1:A:89:ALA:HA	1.91	0.70
1:A:45:PRO:C	1:A:46:VAL:HG12	2.16	0.70
1:A:129:PRO:HG3	1:A:296:LEU:HB2	1.73	0.70
1:A:198:ASN:C	1:A:198:ASN:ND2	2.49	0.70
1:A:149:PRO:HB3	1:A:153:LYS:HG3	1.74	0.70
1:A:104:GLY:HA2	1:A:306:LEU:HD11	1.74	0.69
1:A:74:LEU:HB2	1:A:255:LEU:HD23	1.75	0.69
1:A:174:PHE:C	1:A:174:PHE:CD2	2.70	0.69
1:A:74:LEU:HD13	1:A:259:VAL:HG21	1.74	0.69
1:A:50:VAL:O	1:A:50:VAL:HG13	1.89	0.69
1:A:286:SER:O	1:A:289:PRO:HD2	1.92	0.69
1:A:296:LEU:C	1:A:296:LEU:HD23	2.18	0.68
1:A:261:TYR:O	1:A:265:PRO:HG2	1.93	0.68
1:A:296:LEU:HD23	1:A:297:LEU:N	2.09	0.68
1:A:47:GLN:HB2	1:A:56:ASN:HD21	1.55	0.68
1:A:36:GLY:O	1:A:40:ASP:HB2	1.94	0.68
1:A:30:ASN:HD21	1:A:276:LEU:H	1.40	0.68
1:A:91:LEU:H	1:A:91:LEU:HD12	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:374:VAL:C	1:A:376:GLY:H	2.01	0.68
1:A:45:PRO:CA	1:A:57:LYS:O	2.42	0.67
1:A:140:VAL:HG11	1:A:334:LEU:HD22	1.77	0.67
1:A:105:MET:HG3	1:A:130:ALA:HB1	1.76	0.67
1:A:60:LEU:O	1:A:60:LEU:HD23	1.94	0.67
1:A:276:LEU:HD12	1:A:277:GLN:H	1.58	0.67
1:A:292:ILE:HD13	1:A:292:ILE:O	1.94	0.66
1:A:337:ILE:HG13	1:A:337:ILE:O	1.95	0.66
1:A:262:LEU:O	1:A:266:LEU:HB2	1.96	0.66
1:A:111:LEU:HD23	1:A:111:LEU:C	2.21	0.66
1:A:210:LEU:O	1:A:213:VAL:HG13	1.95	0.65
1:A:47:GLN:CB	1:A:56:ASN:HD21	2.05	0.65
1:A:291:GLY:HA2	1:A:295:GLY:H	1.61	0.65
1:A:165:LEU:HD23	1:A:165:LEU:O	1.97	0.65
1:A:374:VAL:HA	1:A:377:TYR:HB3	1.79	0.65
1:A:280:THR:C	1:A:281:LEU:HG	2.22	0.65
1:A:374:VAL:HG23	1:A:375:ILE:CD1	2.27	0.65
1:A:188:ALA:O	1:A:191:ILE:HG22	1.96	0.64
1:A:105:MET:O	1:A:105:MET:HG2	1.96	0.64
1:A:116:ASN:HD21	1:A:291:GLY:HA3	1.62	0.64
1:A:44:THR:CG2	1:A:59:MET:SD	2.86	0.64
1:A:96:PHE:HB3	1:A:97:PRO:CD	2.24	0.64
1:A:263:ILE:O	1:A:264:LEU:C	2.40	0.64
1:A:306:LEU:O	1:A:310:LEU:N	2.31	0.64
1:A:56:ASN:O	1:A:57:LYS:CG	2.30	0.64
1:A:68:MET:HG3	1:A:344:PHE:CE1	2.33	0.63
1:A:112:TYR:HD2	1:A:125:GLY:HA2	1.63	0.63
1:A:44:THR:HG22	1:A:59:MET:SD	2.39	0.63
1:A:291:GLY:O	1:A:293:ILE:N	2.31	0.63
1:A:174:PHE:CD2	1:A:175:TYR:N	2.64	0.63
1:A:297:LEU:HD23	1:A:297:LEU:O	1.98	0.63
1:A:101:ALA:HB2	1:A:159:LEU:HD23	1.81	0.63
1:A:141:LEU:HD23	1:A:153:LYS:HG2	1.81	0.63
1:A:44:THR:O	1:A:58:ASN:HA	2.00	0.62
1:A:53:LEU:N	1:A:53:LEU:CD2	2.61	0.62
1:A:277:GLN:O	1:A:279:VAL:N	2.33	0.62
1:A:175:TYR:CE2	1:A:177:ASN:HA	2.34	0.62
1:A:103:GLY:C	1:A:306:LEU:HD11	2.24	0.62
1:A:111:LEU:O	1:A:115:PHE:HB2	1.99	0.62
1:A:293:ILE:O	1:A:297:LEU:HB2	2.00	0.62
1:A:101:ALA:HB2	1:A:159:LEU:CD2	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LEU:HD23	1:A:49:ARG:CA	2.31	0.61
1:A:306:LEU:HD23	1:A:306:LEU:C	2.25	0.61
1:A:127:ALA:O	1:A:129:PRO:N	2.33	0.61
1:A:233:VAL:C	1:A:235:PHE:H	2.07	0.61
1:A:238:PRO:HB2	1:A:247:PRO:HG3	1.83	0.61
1:A:290:LEU:HD22	1:A:291:GLY:N	2.13	0.61
1:A:194:LEU:HB3	1:A:236:PHE:HD2	1.64	0.61
1:A:353:VAL:O	1:A:354:ASP:C	2.44	0.61
1:A:56:ASN:C	1:A:57:LYS:CG	2.70	0.60
1:A:285:THR:O	1:A:289:PRO:CG	2.49	0.60
1:A:296:LEU:HG	1:A:365:ILE:HG22	1.83	0.60
1:A:103:GLY:C	1:A:306:LEU:HD21	2.26	0.60
1:A:55:ILE:C	1:A:56:ASN:OD1	2.44	0.60
1:A:47:GLN:HB3	1:A:56:ASN:ND2	2.14	0.59
1:A:316:LEU:O	1:A:317:ALA:HB2	2.01	0.59
1:A:138:LEU:HD23	1:A:156:LEU:HD23	1.85	0.59
1:A:174:PHE:HD2	1:A:174:PHE:C	2.10	0.59
1:A:188:ALA:HA	1:A:191:ILE:HG22	1.83	0.59
1:A:304:ILE:HG12	1:A:333:ILE:HA	1.84	0.59
1:A:91:LEU:HD12	1:A:91:LEU:N	2.17	0.59
1:A:125:GLY:C	1:A:127:ALA:H	2.11	0.59
1:A:259:VAL:HG23	1:A:260:ALA:N	2.17	0.59
1:A:324:THR:OG1	1:A:327:GLN:HB3	2.02	0.59
1:A:94:ALA:O	1:A:97:PRO:HD2	2.02	0.59
1:A:108:PRO:HD3	1:A:306:LEU:CD1	2.33	0.58
1:A:125:GLY:C	1:A:127:ALA:N	2.61	0.58
1:A:304:ILE:O	1:A:306:LEU:N	2.29	0.58
1:A:330:VAL:CG1	1:A:380:LEU:HB2	2.22	0.58
1:A:115:PHE:C	1:A:116:ASN:HD22	2.12	0.58
1:A:68:MET:HG3	1:A:344:PHE:CD1	2.39	0.58
1:A:304:ILE:HG23	1:A:332:GLY:C	2.29	0.58
1:A:275:SER:H	1:A:359:ASN:HD21	1.51	0.58
1:A:279:VAL:O	1:A:281:LEU:HG	2.04	0.57
1:A:121:ILE:CB	1:A:124:GLU:OE1	2.46	0.57
1:A:233:VAL:C	1:A:235:PHE:N	2.59	0.57
1:A:134:ILE:HB	1:A:160:ALA:CB	2.31	0.57
1:A:290:LEU:HD13	1:A:292:ILE:H	1.69	0.57
1:A:297:LEU:HD21	1:A:367:VAL:CG1	2.31	0.57
1:A:45:PRO:O	1:A:46:VAL:HB	2.03	0.57
1:A:90:SER:C	1:A:92:ARG:H	2.12	0.57
1:A:16:ILE:O	1:A:20:ILE:HG13	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:THR:HA	1:A:338:GLY:HA2	1.85	0.57
1:A:375:ILE:HG22	1:A:375:ILE:O	2.05	0.57
1:A:47:GLN:HG3	1:A:48:LEU:N	2.19	0.56
1:A:76:GLY:O	1:A:79:VAL:HG12	2.04	0.56
1:A:210:LEU:HA	1:A:213:VAL:CG1	2.35	0.56
1:A:291:GLY:C	1:A:293:ILE:N	2.61	0.56
1:A:291:GLY:O	1:A:292:ILE:C	2.48	0.56
1:A:109:ALA:C	1:A:111:LEU:H	2.14	0.56
1:A:179:LEU:HA	1:A:184:LEU:HD23	1.86	0.56
1:A:42:LEU:O	1:A:59:MET:HB2	2.05	0.56
1:A:79:VAL:O	1:A:83:LEU:HB2	2.06	0.56
1:A:117:TYR:HD1	1:A:117:TYR:H	1.52	0.56
1:A:300:LYS:HE2	1:A:336:GLY:O	2.04	0.56
1:A:90:SER:OG	1:A:93:GLN:HG3	2.06	0.56
1:A:16:ILE:O	1:A:19:ILE:HG22	2.05	0.56
1:A:67:LEU:HD22	1:A:263:ILE:HG23	1.88	0.56
1:A:101:ALA:C	1:A:103:GLY:N	2.63	0.56
1:A:276:LEU:HD12	1:A:277:GLN:N	2.19	0.55
1:A:43:GLU:O	1:A:44:THR:CB	2.44	0.55
1:A:48:LEU:CG	1:A:49:ARG:H	2.17	0.55
1:A:198:ASN:C	1:A:198:ASN:HD22	2.14	0.55
1:A:233:VAL:O	1:A:235:PHE:N	2.38	0.55
1:A:304:ILE:HG12	1:A:332:GLY:O	2.06	0.55
1:A:159:LEU:O	1:A:160:ALA:C	2.49	0.55
1:A:351:GLY:H	1:A:358:ILE:HD12	1.71	0.55
1:A:49:ARG:CA	1:A:53:LEU:O	2.55	0.54
1:A:374:VAL:C	1:A:376:GLY:N	2.64	0.54
1:A:304:ILE:C	1:A:306:LEU:H	2.15	0.54
1:A:99:ILE:HA	1:A:102:ILE:HG22	1.89	0.54
1:A:116:ASN:ND2	1:A:291:GLY:HA3	2.22	0.54
1:A:128:ILE:CB	1:A:129:PRO:HD3	2.34	0.54
1:A:296:LEU:HD21	1:A:364:GLY:HA3	1.90	0.54
1:A:219:VAL:HG23	1:A:219:VAL:O	2.08	0.54
1:A:307:PHE:O	1:A:311:ALA:N	2.39	0.54
1:A:341:MET:O	1:A:345:ILE:HG12	2.07	0.54
1:A:41:PHE:O	1:A:42:LEU:C	2.51	0.54
1:A:350:PHE:O	1:A:352:SER:N	2.41	0.53
1:A:338:GLY:O	1:A:339:PHE:C	2.52	0.53
1:A:72:PHE:HB3	1:A:231:VAL:HG23	1.90	0.53
1:A:86:GLY:O	1:A:88:LEU:N	2.37	0.53
1:A:307:PHE:C	1:A:309:TRP:H	2.17	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:ILE:O	1:A:288:LEU:C	2.51	0.53
1:A:292:ILE:HG23	1:A:293:ILE:HG12	1.91	0.53
1:A:94:ALA:O	1:A:96:PHE:N	2.42	0.53
1:A:320:PRO:O	1:A:321:GLU:HB2	2.09	0.53
1:A:60:LEU:C	1:A:60:LEU:CD2	2.82	0.52
1:A:312:LEU:C	1:A:314:LEU:N	2.57	0.52
1:A:115:PHE:HD1	1:A:115:PHE:O	1.92	0.52
1:A:326:GLN:O	1:A:329:MET:HG2	2.10	0.52
1:A:98:VAL:O	1:A:102:ILE:HG22	2.10	0.52
1:A:127:ALA:O	1:A:129:PRO:CD	2.58	0.52
1:A:148:VAL:HG11	1:A:327:GLN:OE1	2.10	0.52
1:A:312:LEU:O	1:A:314:LEU:N	2.42	0.52
1:A:147:ARG:O	1:A:149:PRO:HD3	2.10	0.52
1:A:106:ILE:O	1:A:110:LEU:HB2	2.10	0.52
1:A:288:LEU:HB3	1:A:289:PRO:CD	2.40	0.51
1:A:308:CYS:SG	1:A:329:MET:HB2	2.50	0.51
1:A:59:MET:C	1:A:61:LEU:H	2.18	0.51
1:A:304:ILE:C	1:A:306:LEU:N	2.69	0.51
1:A:124:GLU:CG	1:A:125:GLY:N	2.74	0.51
1:A:306:LEU:O	1:A:310:LEU:HB2	2.11	0.51
1:A:104:GLY:HA2	1:A:306:LEU:HD13	1.93	0.51
1:A:197:LEU:HD21	1:A:207:VAL:CG1	2.41	0.51
1:A:324:THR:HB	1:A:327:GLN:NE2	2.26	0.51
1:A:362:LYS:O	1:A:366:LEU:HB2	2.10	0.51
1:A:264:LEU:CB	1:A:265:PRO:HD3	2.41	0.51
1:A:330:VAL:CG1	1:A:380:LEU:HD12	2.41	0.51
1:A:333:ILE:HG21	1:A:376:GLY:HA3	1.93	0.51
1:A:199:LEU:C	1:A:201:GLY:H	2.18	0.51
1:A:47:GLN:HA	1:A:55:ILE:O	2.11	0.50
1:A:112:TYR:CD1	1:A:295:GLY:C	2.90	0.50
1:A:91:LEU:H	1:A:91:LEU:CD1	2.25	0.50
1:A:107:VAL:N	1:A:108:PRO:CD	2.75	0.50
1:A:112:TYR:CE2	1:A:124:GLU:HG2	2.46	0.50
1:A:121:ILE:O	1:A:124:GLU:CD	2.54	0.50
1:A:298:ILE:O	1:A:302:LEU:HB2	2.11	0.50
1:A:203:ARG:HD3	1:A:243:HIS:ND1	2.27	0.50
1:A:178:ASP:O	1:A:180:SER:N	2.43	0.50
1:A:274:VAL:HG21	1:A:363:LEU:HG	1.94	0.50
1:A:307:PHE:CD2	1:A:332:GLY:HA3	2.47	0.50
1:A:14:GLY:HA2	1:A:17:ILE:HG22	1.94	0.49
1:A:180:SER:O	1:A:181:MET:C	2.54	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:ALA:HA	1:A:164:ASP:OD2	2.12	0.49
1:A:207:VAL:O	1:A:208:TYR:C	2.55	0.49
1:A:304:ILE:CD1	1:A:333:ILE:HA	2.43	0.49
1:A:103:GLY:HA3	1:A:310:LEU:HG	1.95	0.49
1:A:275:SER:N	1:A:359:ASN:HD21	2.10	0.49
1:A:330:VAL:HB	1:A:380:LEU:HD12	1.94	0.49
1:A:47:GLN:HB2	1:A:56:ASN:OD1	2.11	0.49
1:A:290:LEU:HD13	1:A:290:LEU:C	2.38	0.49
1:A:137:ALA:HB1	1:A:156:LEU:HD21	1.94	0.49
1:A:124:GLU:HG2	1:A:125:GLY:N	2.27	0.48
1:A:60:LEU:O	1:A:60:LEU:CD2	2.61	0.48
1:A:46:VAL:O	1:A:56:ASN:HA	2.13	0.48
1:A:152:LEU:O	1:A:155:PHE:HB3	2.12	0.48
1:A:284:LEU:O	1:A:288:LEU:CB	2.58	0.48
1:A:309:TRP:O	1:A:313:ARG:HB2	2.14	0.48
1:A:147:ARG:O	1:A:148:VAL:HB	2.13	0.48
1:A:152:LEU:HD11	1:A:331:VAL:HG22	1.96	0.47
1:A:240:LYS:C	1:A:242:LYS:H	2.22	0.47
1:A:316:LEU:O	1:A:317:ALA:CB	2.61	0.47
1:A:116:ASN:C	1:A:118:ALA:H	2.23	0.47
1:A:99:ILE:HA	1:A:102:ILE:CG2	2.45	0.47
1:A:137:ALA:HA	1:A:337:ILE:HD11	1.96	0.47
1:A:61:LEU:HD13	1:A:61:LEU:O	2.15	0.47
1:A:126:TRP:HD1	1:A:126:TRP:O	1.97	0.47
1:A:214:VAL:HG12	1:A:215:LEU:N	2.28	0.47
1:A:298:ILE:C	1:A:300:LYS:N	2.70	0.47
1:A:375:ILE:O	1:A:375:ILE:CG2	2.61	0.47
1:A:65:ASP:O	1:A:69:ALA:HB2	2.15	0.47
1:A:111:LEU:HD22	1:A:299:GLY:HA2	1.96	0.47
1:A:198:ASN:HD22	1:A:199:LEU:N	2.13	0.47
1:A:229:ALA:O	1:A:233:VAL:HG23	2.15	0.47
1:A:313:ARG:O	1:A:313:ARG:HD3	2.14	0.47
1:A:112:TYR:HE2	1:A:124:GLU:HG2	1.79	0.47
1:A:23:ILE:HG22	1:A:24:LEU:N	2.29	0.46
1:A:199:LEU:C	1:A:201:GLY:N	2.73	0.46
1:A:261:TYR:C	1:A:265:PRO:HG2	2.39	0.46
1:A:77:LEU:HD23	1:A:234:GLY:O	2.15	0.46
1:A:169:ILE:O	1:A:173:LEU:HD22	2.15	0.46
1:A:176:THR:O	1:A:177:ASN:C	2.57	0.46
1:A:216:TRP:CD1	1:A:226:ALA:HB1	2.51	0.46
1:A:17:ILE:HG23	1:A:18:LEU:N	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ASP:C	1:A:121:ILE:H	2.23	0.46
1:A:34:THR:C	1:A:36:GLY:N	2.70	0.46
1:A:101:ALA:O	1:A:104:GLY:N	2.49	0.46
1:A:127:ALA:O	1:A:128:ILE:C	2.58	0.46
1:A:175:TYR:OH	1:A:178:ASP:HB2	2.15	0.46
1:A:60:LEU:HD23	1:A:60:LEU:C	2.40	0.46
1:A:96:PHE:CD2	1:A:314:LEU:HD12	2.51	0.46
1:A:142:ALA:C	1:A:144:LEU:H	2.23	0.46
1:A:149:PRO:HA	1:A:152:LEU:HB3	1.98	0.46
1:A:282:ASP:O	1:A:283:GLY:C	2.58	0.46
1:A:148:VAL:HB	1:A:149:PRO:HD3	1.98	0.46
1:A:264:LEU:HB3	1:A:265:PRO:HD3	1.98	0.46
1:A:169:ILE:O	1:A:170:ILE:C	2.59	0.45
1:A:141:LEU:CD2	1:A:153:LYS:HG2	2.46	0.45
1:A:31:SER:HB3	1:A:34:THR:OG1	2.16	0.45
1:A:320:PRO:O	1:A:323:THR:HG23	2.16	0.45
1:A:112:TYR:OH	1:A:124:GLU:HG3	2.17	0.45
1:A:319:LEU:HA	1:A:320:PRO:HD3	1.80	0.45
1:A:127:ALA:O	1:A:129:PRO:HD2	2.16	0.45
1:A:183:SER:O	1:A:184:LEU:C	2.58	0.45
1:A:343:ILE:HG23	1:A:362:LYS:HD2	1.99	0.45
1:A:337:ILE:C	1:A:339:PHE:H	2.25	0.45
1:A:34:THR:C	1:A:36:GLY:H	2.25	0.45
1:A:147:ARG:C	1:A:149:PRO:HD3	2.42	0.45
1:A:325:TYR:C	1:A:327:GLN:H	2.25	0.45
1:A:94:ALA:O	1:A:95:ALA:C	2.60	0.45
1:A:298:ILE:C	1:A:300:LYS:H	2.24	0.45
1:A:309:TRP:HD1	1:A:310:LEU:HD22	1.82	0.45
1:A:101:ALA:C	1:A:103:GLY:H	2.23	0.45
1:A:147:ARG:C	1:A:149:PRO:CD	2.90	0.44
1:A:275:SER:H	1:A:359:ASN:ND2	2.15	0.44
1:A:211:VAL:O	1:A:214:VAL:HB	2.18	0.44
1:A:47:GLN:HA	1:A:56:ASN:HA	2.00	0.44
1:A:334:LEU:C	1:A:336:GLY:N	2.65	0.44
1:A:274:VAL:HG23	1:A:359:ASN:ND2	2.32	0.44
1:A:287:ILE:O	1:A:294:ALA:HB2	2.16	0.44
1:A:17:ILE:O	1:A:20:ILE:HB	2.18	0.44
1:A:97:PRO:HG3	1:A:155:PHE:CD1	2.52	0.44
1:A:115:PHE:O	1:A:116:ASN:ND2	2.44	0.44
1:A:203:ARG:HB3	1:A:243:HIS:CE1	2.53	0.44
1:A:22:ALA:O	1:A:25:ALA:HB3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:LEU:O	1:A:181:MET:N	2.51	0.44
1:A:342:SER:O	1:A:346:ALA:N	2.50	0.44
1:A:103:GLY:O	1:A:306:LEU:CD1	2.60	0.44
1:A:77:LEU:HD11	1:A:251:LEU:HD23	2.00	0.44
1:A:171:ILE:O	1:A:171:ILE:HG22	2.17	0.44
1:A:307:PHE:C	1:A:309:TRP:N	2.76	0.44
1:A:229:ALA:O	1:A:230:GLY:C	2.60	0.44
1:A:320:PRO:O	1:A:321:GLU:CB	2.65	0.44
1:A:168:ILE:CD1	1:A:345:ILE:HD11	2.48	0.43
1:A:298:ILE:HG12	1:A:302:LEU:HD23	2.01	0.43
1:A:142:ALA:C	1:A:144:LEU:N	2.77	0.43
1:A:342:SER:O	1:A:343:ILE:C	2.61	0.43
1:A:90:SER:C	1:A:92:ARG:N	2.76	0.43
1:A:245:ARG:HG3	1:A:246:SER:N	2.33	0.43
1:A:367:VAL:O	1:A:368:GLY:C	2.61	0.43
1:A:16:ILE:HD12	1:A:16:ILE:HA	1.76	0.43
1:A:31:SER:C	1:A:33:ALA:H	2.25	0.43
1:A:197:LEU:HD21	1:A:207:VAL:HG11	2.00	0.43
1:A:281:LEU:O	1:A:281:LEU:HD12	2.18	0.43
1:A:109:ALA:O	1:A:111:LEU:N	2.52	0.43
1:A:136:PHE:CD1	1:A:337:ILE:HD12	2.54	0.43
1:A:203:ARG:O	1:A:205:THR:N	2.52	0.43
1:A:277:GLN:C	1:A:279:VAL:N	2.76	0.43
1:A:284:LEU:C	1:A:286:SER:N	2.72	0.43
1:A:144:LEU:HD13	1:A:380:LEU:HD22	2.00	0.43
1:A:153:LYS:C	1:A:155:PHE:H	2.27	0.43
1:A:210:LEU:O	1:A:211:VAL:C	2.59	0.43
1:A:293:ILE:HG22	1:A:294:ALA:N	2.33	0.43
1:A:115:PHE:C	1:A:116:ASN:ND2	2.75	0.43
1:A:209:ILE:HG22	1:A:210:LEU:N	2.33	0.43
1:A:124:GLU:OE2	1:A:125:GLY:N	2.52	0.42
1:A:22:ALA:O	1:A:26:MET:HG2	2.19	0.42
1:A:48:LEU:CG	1:A:49:ARG:N	2.79	0.42
1:A:55:ILE:HG22	1:A:56:ASN:H	1.83	0.42
1:A:138:LEU:HD23	1:A:156:LEU:CD2	2.48	0.42
1:A:241:GLU:O	1:A:241:GLU:HG2	2.18	0.42
1:A:286:SER:O	1:A:293:ILE:HG21	2.18	0.42
1:A:59:MET:O	1:A:61:LEU:N	2.43	0.42
1:A:62:TRP:CZ3	1:A:220:LEU:HD21	2.55	0.42
1:A:96:PHE:CE2	1:A:314:LEU:HD12	2.54	0.42
1:A:125:GLY:O	1:A:126:TRP:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:LEU:O	1:A:291:GLY:C	2.63	0.42
1:A:300:LYS:HE2	1:A:300:LYS:HB3	1.70	0.42
1:A:116:ASN:CG	1:A:290:LEU:HD11	2.43	0.42
1:A:288:LEU:CB	1:A:289:PRO:HD3	2.46	0.42
1:A:63:ILE:O	1:A:67:LEU:HG	2.20	0.42
1:A:380:LEU:O	1:A:380:LEU:CD2	2.68	0.42
1:A:14:GLY:C	1:A:17:ILE:HG22	2.44	0.42
1:A:123:ARG:NH1	1:A:126:TRP:HZ3	2.17	0.41
1:A:188:ALA:C	1:A:191:ILE:HG22	2.45	0.41
1:A:286:SER:O	1:A:289:PRO:CD	2.65	0.41
1:A:153:LYS:C	1:A:155:PHE:N	2.78	0.41
1:A:164:ASP:C	1:A:166:GLY:N	2.76	0.41
1:A:293:ILE:HD11	1:A:360:TRP:C	2.45	0.41
1:A:70:VAL:O	1:A:73:LEU:HB3	2.19	0.41
1:A:144:LEU:O	1:A:145:GLY:C	2.63	0.41
1:A:147:ARG:O	1:A:148:VAL:CB	2.67	0.41
1:A:189:VAL:CG2	1:A:190:ALA:N	2.83	0.41
1:A:48:LEU:N	1:A:55:ILE:O	2.50	0.41
1:A:74:LEU:C	1:A:74:LEU:HD23	2.46	0.41
1:A:97:PRO:HG3	1:A:155:PHE:CE1	2.56	0.41
1:A:238:PRO:O	1:A:248:ALA:HB2	2.20	0.41
1:A:360:TRP:O	1:A:361:ALA:C	2.64	0.41
1:A:378:SER:O	1:A:379:TRP:C	2.62	0.41
1:A:59:MET:C	1:A:61:LEU:N	2.79	0.41
1:A:300:LYS:HE3	1:A:338:GLY:H	1.85	0.41
1:A:50:VAL:O	1:A:50:VAL:CG1	2.61	0.41
1:A:79:VAL:CG1	1:A:235:PHE:CE1	3.04	0.41
1:A:133:ASP:OD2	1:A:339:PHE:HB3	2.21	0.41
1:A:189:VAL:HG22	1:A:190:ALA:N	2.36	0.41
1:A:293:ILE:HD13	1:A:293:ILE:HA	1.98	0.41
1:A:304:ILE:CG1	1:A:333:ILE:HA	2.47	0.41
1:A:367:VAL:C	1:A:369:SER:N	2.78	0.41
1:A:112:TYR:CD2	1:A:125:GLY:CA	2.92	0.41
1:A:189:VAL:O	1:A:192:ALA:HB3	2.21	0.41
1:A:254:VAL:O	1:A:258:TRP:HD1	2.04	0.40
1:A:307:PHE:CE2	1:A:332:GLY:CA	3.04	0.40
1:A:107:VAL:HB	1:A:108:PRO:HD3	2.03	0.40
1:A:380:LEU:O	1:A:380:LEU:HD23	2.21	0.40
1:A:204:ARG:O	1:A:206:GLY:N	2.55	0.40
1:A:343:ILE:HG12	1:A:365:ILE:HD11	2.04	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:LEU:CD1	1:A:48:LEU:CD1[2_555]	1.31	0.89
1:A:198:ASN:ND2	1:A:199:LEU:CD1[2_655]	1.45	0.75
1:A:46:VAL:CG2	1:A:50:VAL:CA[2_555]	1.63	0.57
1:A:282:ASP:OD2	1:A:326:GLN:NE2[4_445]	1.75	0.45
1:A:181:MET:CG	1:A:181:MET:CG[2_655]	1.97	0.23

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	374/376 (100%)	229 (61%)	95 (25%)	50 (13%)	0 4

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	44	THR
1	A	45	PRO
1	A	46	VAL
1	A	86	GLY
1	A	87	SER
1	A	95	ALA
1	A	119	ASP
1	A	126	TRP
1	A	128	ILE
1	A	148	VAL
1	A	263	ILE
1	A	277	GLN
1	A	282	ASP
1	A	283	GLY
1	A	309	TRP
1	A	317	ALA
1	A	321	GLU

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Mol	Chain	Res	Type
1	A	339	PHE
1	A	351	GLY
1	A	372	SER
1	A	38	TYR
1	A	110	LEU
1	A	145	GLY
1	A	146	SER
1	A	149	PRO
1	A	179	LEU
1	A	290	LEU
1	A	292	ILE
1	A	305	SER
1	A	323	THR
1	A	55	ILE
1	A	104	GLY
1	A	151	ALA
1	A	177	ASN
1	A	205	THR
1	A	208	TYR
1	A	313	ARG
1	A	355	PRO
1	A	42	LEU
1	A	88	LEU
1	A	118	ALA
1	A	181	MET
1	A	219	VAL
1	A	49	ARG
1	A	207	VAL
1	A	291	GLY
1	A	354	ASP
1	A	289	PRO
1	A	234	GLY
1	A	238	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 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	291/291 (100%)	248 (85%)	43 (15%)	3 13

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	SER
1	A	16	ILE
1	A	19	ILE
1	A	24	LEU
1	A	26	MET
1	A	43	GLU
1	A	44	THR
1	A	47	GLN
1	A	53	LEU
1	A	54	GLU
1	A	55	ILE
1	A	57	LYS
1	A	58	ASN
1	A	59	MET
1	A	60	LEU
1	A	61	LEU
1	A	83	LEU
1	A	84	MET
1	A	87	SER
1	A	117	TYR
1	A	121	ILE
1	A	124	GLU
1	A	174	PHE
1	A	177	ASN
1	A	184	LEU
1	A	189	VAL
1	A	196	VAL
1	A	198	ASN
1	A	209	ILE
1	A	213	VAL
1	A	214	VAL
1	A	264	LEU
1	A	266	LEU
1	A	276	LEU
1	A	281	LEU
1	A	289	PRO
1	A	292	ILE
1	A	296	LEU

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Mol	Chain	Res	Type
1	A	306	LEU
1	A	312	LEU
1	A	315	LYS
1	A	355	PRO
1	A	362	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	64	ASN
1	A	198	ASN
1	A	271	ASN
1	A	318	HIS
1	A	359	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 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 ⓘ

There are no chain breaks in this entry.

6 Map visualisation

This section contains visualisations of the EMDB entry EMD-5037. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections

This section was not generated.

6.2 Central slices

This section was not generated.

6.3 Largest variance slices

This section was not generated.

6.4 Orthogonal standard-deviation projections (False-color)

This section was not generated.

6.5 Orthogonal surface views

This section was not generated.

6.6 Mask visualisation

This section was not generated. No masks/segmentation were deposited.

7 Map analysis ⓘ

This section contains the results of statistical analysis of the map.

7.1 Map-value distribution ⓘ

This section was not generated.

7.2 Volume estimate versus contour level ⓘ

This section was not generated.

7.3 Rotationally averaged power spectrum ⓘ

This section was not generated. The rotationally averaged power spectrum had issues being displayed.

8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit

This section was not generated.