



wwPDB EM Map/Model Validation Report ⓘ

Apr 10, 2016 – 01:55 PM BST

PDB ID : 4BT0
EMDB ID: : EMD-2398
Title : MuB is an AAAplus ATPase that forms helical filaments to control target selection for DNA transposition
Authors : Mizuno, N.; Dramicanin, M.; Mizuuchi, M.; Adam, J.; Wang, Y.; Han, Y.W.; Yang, W.; Steven, A.C.; Mizuuchi, K.; Ramon-Maiques, S.
Deposited on : 2013-06-12
Resolution : 17.00 Å(reported)
Based on PDB ID : 1NY6

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 : 1.7.1 (RC1), CSD as537be (2016)
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

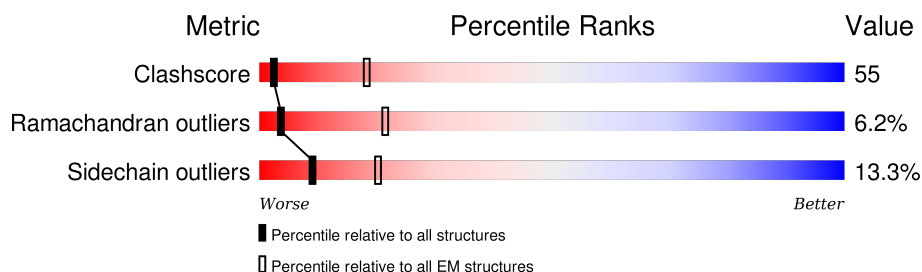
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 17.00 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	73	
2	B	173	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2023 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.

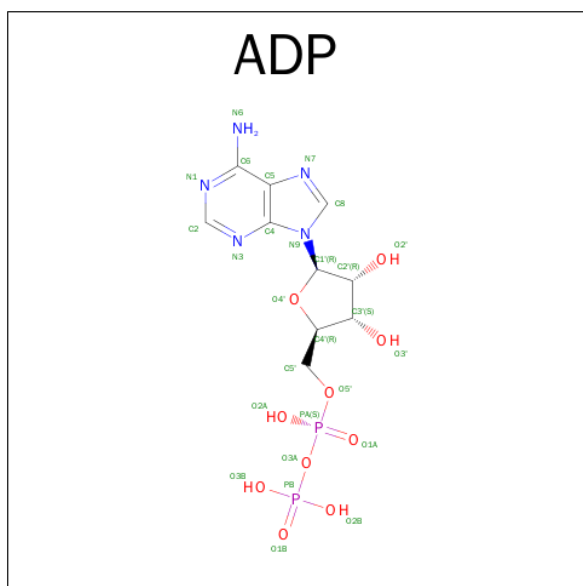
- Molecule 1 is a protein called TRANSCRIPTIONAL REGULATOR.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	73	Total	C	N	O	S	0	0
			604	392	103	108	1		

- Molecule 2 is a protein called TRANSCRIPTIONAL REGULATOR.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	173	Total	C	N	O	S	0	0
			1365	879	226	257	3		

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

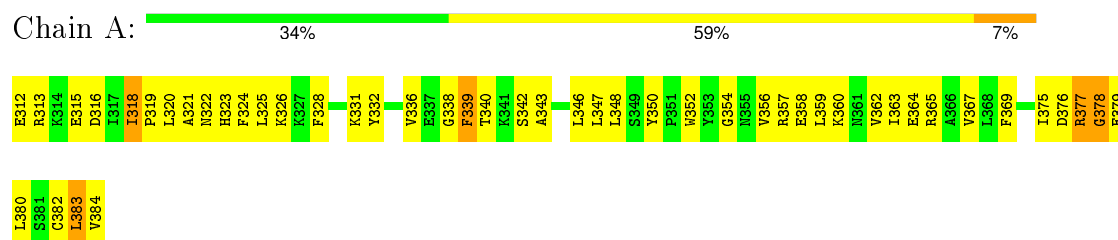


Mol	Chain	Residues	Atoms					AltConf
3	A	1	Total	C	N	O	P	0
			27	10	5	10	2	
3	B	1	Total	C	N	O	P	0
			27	10	5	10	2	

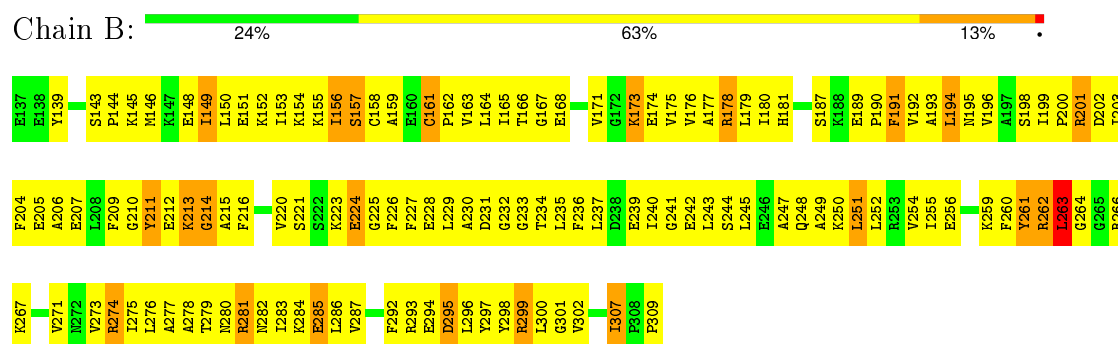
3 Residue-property plots

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: TRANSCRIPTIONAL REGULATOR



• Molecule 2: TRANSCRIPTIONAL REGULATOR



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	PHASE-FLIPPING, Not provided	Depositor
Microscope	FEI/PHILIPS CM200 FEG	Depositor
Voltage (kV)	120	Depositor
Electron dose ($e^-/\text{\AA}^2$)	15	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	38000	Depositor
Image detector	GATAN ULTRASCAN 1000 CCD MODEL 894 (2KX2K)	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ADP

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	0.52	0/616	0.65	0/825
2	B	0.49	0/1386	0.71	0/1859
All	All	0.50	0/2002	0.69	0/2684

There are no bond length outliers.

There are no bond angle outliers.

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	604	0	616	58	0
2	B	1365	0	1405	169	0
3	A	27	0	12	2	0
3	B	27	0	12	0	0
All	All	2023	0	2045	224	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 55.

All (224) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:259:LYS:HB3	2:B:267:LYS:HG3	1.14	1.12
1:A:343:ALA:HB2	1:A:376:ASP:HA	1.24	1.12
2:B:189:GLU:HG3	2:B:190:PRO:HD2	1.38	1.04
1:A:318:ILE:HG12	1:A:348:LEU:HD21	1.40	1.04
1:A:325:LEU:HD23	1:A:325:LEU:O	1.63	0.97
2:B:168:GLU:O	2:B:171:VAL:HG23	1.63	0.97
2:B:212:GLU:HB2	2:B:262:ARG:O	1.69	0.93
2:B:240:ILE:HG13	2:B:277:ALA:HB1	1.52	0.92
2:B:149:ILE:O	2:B:153:ILE:HG12	1.69	0.92
2:B:224:GLU:HG2	2:B:225:GLY:H	1.39	0.87
1:A:362:VAL:HG13	1:A:383:LEU:HD12	1.59	0.85
2:B:263:LEU:CG	2:B:264:GLY:H	1.92	0.82
1:A:358:GLU:O	1:A:362:VAL:HG23	1.78	0.82
2:B:231:ASP:OD1	2:B:271:VAL:HA	1.81	0.80
2:B:209:PHE:O	2:B:225:GLY:HA3	1.82	0.78
2:B:252:LEU:HB2	2:B:296:LEU:HD23	1.66	0.78
1:A:325:LEU:HD13	1:A:338:GLY:HA2	1.67	0.76
2:B:263:LEU:CD2	2:B:264:GLY:H	1.97	0.76
2:B:263:LEU:HD23	2:B:264:GLY:H	1.51	0.76
2:B:245:LEU:HD22	2:B:248:GLN:NE2	2.01	0.75
1:A:340:THR:OG1	1:A:376:ASP:HB3	1.87	0.74
2:B:299:ARG:HA	2:B:299:ARG:HE	1.54	0.73
2:B:275:ILE:HD12	2:B:275:ILE:N	2.04	0.72
2:B:189:GLU:HG2	2:B:232:GLY:O	1.87	0.72
2:B:194:LEU:HD21	2:B:237:LEU:HD22	1.71	0.72
2:B:189:GLU:CG	2:B:190:PRO:HD2	2.19	0.72
2:B:198:SER:O	2:B:199:ILE:HG13	1.91	0.70
2:B:245:LEU:HD22	2:B:248:GLN:HE22	1.56	0.70
2:B:259:LYS:HB3	2:B:267:LYS:CG	2.08	0.70
2:B:178:ARG:HA	2:B:191:PHE:HE2	1.56	0.69
2:B:263:LEU:HG	2:B:264:GLY:H	1.58	0.69
1:A:369:PHE:HA	2:B:155:LYS:HD3	1.74	0.68
1:A:325:LEU:CD1	1:A:338:GLY:HA2	2.23	0.67
1:A:320:LEU:O	1:A:323:HIS:N	2.27	0.66
1:A:363:ILE:O	1:A:367:VAL:HG23	1.95	0.66
2:B:191:PHE:HE1	2:B:236:PHE:HB3	1.60	0.66
2:B:240:ILE:CG1	2:B:277:ALA:HB1	2.25	0.66
2:B:178:ARG:HG2	2:B:191:PHE:CE2	2.32	0.64
2:B:252:LEU:HD22	2:B:293:ARG:HH12	1.61	0.64
2:B:209:PHE:CE2	2:B:250:LYS:HD3	2.32	0.64
2:B:280:ASN:OD1	2:B:281:ARG:HD3	1.97	0.64
1:A:316:ASP:C	1:A:319:PRO:HD2	2.20	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:261:TYR:N	2:B:261:TYR:CD1	2.68	0.62
2:B:294:GLU:O	2:B:297:TYR:HB3	2.00	0.62
1:A:323:HIS:O	1:A:326:LYS:HB3	2.00	0.61
1:A:339:PHE:CD1	1:A:339:PHE:N	2.68	0.61
1:A:328:PHE:HA	1:A:331:LYS:HB3	1.83	0.61
1:A:316:ASP:O	1:A:319:PRO:HD2	2.00	0.61
2:B:263:LEU:HD23	2:B:264:GLY:N	2.15	0.61
2:B:252:LEU:HD13	2:B:296:LEU:HA	1.83	0.60
2:B:164:LEU:HD12	2:B:165:ILE:N	2.16	0.60
1:A:324:PHE:CE1	1:A:360:LYS:HG3	2.36	0.60
2:B:252:LEU:HD11	2:B:299:ARG:HG3	1.83	0.60
2:B:263:LEU:CG	2:B:264:GLY:N	2.64	0.60
2:B:249:ALA:HA	2:B:293:ARG:NH1	2.16	0.60
2:B:146:MET:CE	2:B:149:ILE:HD12	2.31	0.60
2:B:154:LYS:O	2:B:157:SER:OG	2.18	0.60
2:B:163:VAL:HB	2:B:276:LEU:CD2	2.32	0.59
2:B:210:GLY:HA2	2:B:224:GLU:O	2.02	0.59
2:B:231:ASP:HA	2:B:273:VAL:HG22	1.85	0.58
1:A:365:ARG:HD2	1:A:383:LEU:CD1	2.34	0.58
2:B:279:THR:HG21	2:B:283:ILE:HD11	1.86	0.57
2:B:275:ILE:CD1	2:B:275:ILE:N	2.67	0.57
2:B:211:TYR:CE1	2:B:263:LEU:HD13	2.39	0.57
2:B:242:GLU:HG2	2:B:281:ARG:NH2	2.19	0.57
1:A:325:LEU:HD23	1:A:325:LEU:C	2.25	0.57
2:B:282:ASN:ND2	2:B:285:GLU:HB2	2.20	0.57
2:B:173:LYS:O	2:B:175:VAL:N	2.38	0.57
1:A:376:ASP:O	1:A:378:GLY:N	2.38	0.56
2:B:207:GLU:HG2	2:B:226:PHE:CE2	2.41	0.56
2:B:211:TYR:HE1	2:B:263:LEU:HD13	1.69	0.56
2:B:211:TYR:HE1	2:B:215:ALA:HB3	1.71	0.56
2:B:168:GLU:OE1	2:B:309:PRO:HB3	2.05	0.56
2:B:150:LEU:O	2:B:153:ILE:HB	2.06	0.56
2:B:224:GLU:HG2	2:B:225:GLY:N	2.16	0.55
2:B:261:TYR:HB3	2:B:266:ARG:C	2.27	0.55
2:B:240:ILE:HG13	2:B:277:ALA:CB	2.33	0.55
2:B:191:PHE:CE1	2:B:193:ALA:HB2	2.42	0.55
2:B:235:LEU:HD21	2:B:237:LEU:HD21	1.89	0.55
2:B:252:LEU:CD1	2:B:299:ARG:HG3	2.37	0.55
2:B:176:VAL:O	2:B:177:ALA:C	2.46	0.54
2:B:252:LEU:CD2	2:B:293:ARG:HH12	2.21	0.54
1:A:318:ILE:CG1	1:A:348:LEU:HD21	2.25	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:178:ARG:HA	2:B:191:PHE:CE2	2.42	0.54
1:A:325:LEU:HD21	1:A:336:VAL:HG12	1.90	0.53
2:B:144:PRO:HG2	2:B:145:LYS:H	1.73	0.53
1:A:325:LEU:HD12	1:A:339:PHE:CE1	2.42	0.53
2:B:296:LEU:HD13	2:B:296:LEU:O	2.08	0.53
1:A:313:ARG:HB3	1:A:316:ASP:OD2	2.08	0.53
2:B:181:HIS:HE1	2:B:187:SER:O	1.91	0.53
1:A:347:LEU:HD21	1:A:380:LEU:HD11	1.91	0.53
2:B:261:TYR:HD2	2:B:263:LEU:O	1.91	0.53
2:B:263:LEU:HG	2:B:264:GLY:N	2.23	0.53
1:A:346:LEU:HD21	1:A:384:VAL:O	2.08	0.53
2:B:229:LEU:HD13	2:B:229:LEU:O	2.09	0.53
2:B:152:LYS:O	2:B:156:ILE:HG12	2.08	0.53
2:B:153:ILE:HG23	2:B:180:ILE:HG12	1.90	0.53
2:B:203:ILE:O	2:B:203:ILE:HG22	2.07	0.53
2:B:260:PHE:HE2	2:B:271:VAL:HG11	1.73	0.52
2:B:166:THR:HG22	2:B:279:THR:HG22	1.92	0.52
2:B:224:GLU:CG	2:B:225:GLY:H	2.18	0.52
2:B:282:ASN:HD22	2:B:285:GLU:HB2	1.75	0.52
2:B:228:GLU:HG3	2:B:260:PHE:HZ	1.75	0.51
1:A:365:ARG:HH11	1:A:383:LEU:HD22	1.75	0.51
1:A:339:PHE:CZ	1:A:363:ILE:HD13	2.46	0.51
1:A:365:ARG:HD2	1:A:383:LEU:HD11	1.92	0.51
1:A:375:ILE:HG22	1:A:380:LEU:HD23	1.93	0.51
2:B:173:LYS:O	2:B:176:VAL:N	2.41	0.50
2:B:153:ILE:HG23	2:B:180:ILE:CG1	2.42	0.50
2:B:148:GLU:O	2:B:150:LEU:N	2.45	0.50
2:B:256:GLU:HG2	2:B:299:ARG:HH11	1.76	0.50
2:B:234:THR:HG21	2:B:276:LEU:HD12	1.95	0.49
2:B:191:PHE:HD1	2:B:191:PHE:C	2.15	0.49
2:B:248:GLN:OE1	2:B:292:PHE:HA	2.12	0.49
2:B:198:SER:C	2:B:199:ILE:HG13	2.31	0.49
2:B:167:GLY:H	2:B:173:LYS:HD3	1.76	0.49
1:A:331:LYS:HD3	1:A:332:TYR:CZ	2.48	0.49
2:B:233:GLY:O	2:B:273:VAL:HG13	2.11	0.49
2:B:192:VAL:O	2:B:235:LEU:HD12	2.13	0.49
2:B:267:LYS:O	2:B:267:LYS:HD3	2.12	0.49
2:B:163:VAL:HB	2:B:276:LEU:HD22	1.94	0.49
1:A:365:ARG:NH2	2:B:302:VAL:O	2.44	0.49
2:B:191:PHE:C	2:B:191:PHE:CD1	2.86	0.49
2:B:145:LYS:O	2:B:149:ILE:HG13	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:161:CYS:HB2	2:B:302:VAL:HG11	1.95	0.48
2:B:148:GLU:O	2:B:151:GLU:N	2.46	0.48
1:A:339:PHE:H	1:A:339:PHE:HD1	1.55	0.48
2:B:235:LEU:HB2	2:B:273:VAL:HG11	1.96	0.48
2:B:299:ARG:O	2:B:302:VAL:HG23	2.12	0.48
2:B:234:THR:HG22	2:B:235:LEU:N	2.29	0.48
2:B:196:VAL:HG12	2:B:196:VAL:O	2.14	0.48
1:A:376:ASP:O	1:A:377:ARG:C	2.51	0.48
2:B:191:PHE:CE1	2:B:236:PHE:HB3	2.45	0.48
2:B:164:LEU:HD12	2:B:164:LEU:C	2.34	0.48
2:B:181:HIS:HD2	2:B:234:THR:OG1	1.97	0.47
2:B:181:HIS:C	2:B:181:HIS:ND1	2.67	0.47
2:B:251:LEU:HD22	2:B:255:ILE:HG13	1.96	0.47
2:B:204:PHE:C	2:B:206:ALA:H	2.17	0.47
2:B:209:PHE:CZ	2:B:250:LYS:HB3	2.50	0.47
2:B:211:TYR:HB2	2:B:223:LYS:HB2	1.97	0.47
2:B:194:LEU:CD2	2:B:237:LEU:HD22	2.42	0.47
2:B:242:GLU:HG2	2:B:281:ARG:HH21	1.80	0.47
2:B:151:GLU:O	2:B:152:LYS:C	2.52	0.47
1:A:350:TYR:HD2	1:A:352:TRP:CD2	2.33	0.47
1:A:379:GLU:H	1:A:379:GLU:CD	2.17	0.47
2:B:167:GLY:N	2:B:173:LYS:HD3	2.29	0.47
2:B:251:LEU:HD22	2:B:255:ILE:HD11	1.95	0.47
2:B:224:GLU:HA	2:B:262:ARG:NH1	2.30	0.47
2:B:262:ARG:O	2:B:263:LEU:HB3	2.14	0.47
2:B:155:LYS:O	2:B:157:SER:N	2.47	0.47
2:B:261:TYR:CD2	2:B:263:LEU:O	2.67	0.46
2:B:240:ILE:O	2:B:242:GLU:N	2.48	0.46
2:B:181:HIS:O	2:B:181:HIS:ND1	2.47	0.46
2:B:299:ARG:HE	2:B:299:ARG:CA	2.24	0.46
1:A:312:GLU:O	1:A:313:ARG:HG3	2.16	0.46
2:B:252:LEU:HD22	2:B:293:ARG:NH1	2.29	0.46
1:A:356:VAL:HG11	3:A:1385:ADP:C8	2.50	0.46
2:B:251:LEU:HD22	2:B:255:ILE:CG1	2.46	0.46
2:B:275:ILE:H	2:B:275:ILE:CD1	2.29	0.46
1:A:328:PHE:HB2	1:A:367:VAL:HG21	1.97	0.46
2:B:205:GLU:HG2	2:B:247:ALA:HB2	1.97	0.46
2:B:156:ILE:O	2:B:156:ILE:HG22	2.16	0.46
2:B:244:SER:O	2:B:248:GLN:HG3	2.15	0.46
1:A:325:LEU:CD1	1:A:339:PHE:CE1	2.99	0.46
2:B:250:LYS:O	2:B:254:VAL:HG23	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:LEU:HB2	1:A:384:VAL:H	1.66	0.46
1:A:313:ARG:O	1:A:316:ASP:N	2.39	0.45
2:B:259:LYS:HD3	2:B:267:LYS:HD2	1.97	0.45
2:B:181:HIS:CD2	2:B:234:THR:OG1	2.70	0.45
1:A:365:ARG:NH2	2:B:301:GLY:O	2.50	0.45
1:A:356:VAL:HG11	3:A:1385:ADP:C5	2.52	0.45
2:B:192:VAL:HB	2:B:230:ALA:HB2	1.98	0.44
2:B:239:GLU:N	2:B:278:ALA:O	2.50	0.44
2:B:207:GLU:HG2	2:B:226:PHE:HE2	1.80	0.44
2:B:204:PHE:CE2	2:B:243:LEU:HD21	2.52	0.44
2:B:155:LYS:C	2:B:157:SER:N	2.70	0.44
1:A:380:LEU:HD12	1:A:384:VAL:HG21	1.98	0.43
2:B:173:LYS:C	2:B:175:VAL:N	2.72	0.43
1:A:339:PHE:HZ	1:A:363:ILE:HD13	1.81	0.43
1:A:382:CYS:C	1:A:383:LEU:HD23	2.38	0.43
2:B:294:GLU:OE2	2:B:298:TYR:HE2	2.01	0.43
2:B:236:PHE:HE1	2:B:278:ALA:HB2	1.83	0.43
2:B:237:LEU:HB2	2:B:240:ILE:HD11	2.01	0.43
1:A:325:LEU:HB2	1:A:339:PHE:CZ	2.54	0.43
2:B:284:LYS:O	2:B:285:GLU:C	2.57	0.42
1:A:352:TRP:HZ3	1:A:362:VAL:HG21	1.82	0.42
2:B:284:LYS:O	2:B:287:VAL:HB	2.19	0.42
2:B:252:LEU:HD22	2:B:295:ASP:OD1	2.18	0.42
2:B:157:SER:O	2:B:159:ALA:N	2.51	0.42
2:B:216:PHE:CG	2:B:216:PHE:O	2.71	0.42
2:B:143:SER:HB2	2:B:144:PRO:HD2	2.00	0.42
2:B:161:CYS:O	2:B:274:ARG:NH1	2.53	0.42
2:B:259:LYS:HD3	2:B:267:LYS:HE3	2.01	0.42
2:B:155:LYS:C	2:B:157:SER:H	2.22	0.42
1:A:318:ILE:HG13	1:A:318:ILE:H	1.49	0.42
2:B:173:LYS:C	2:B:175:VAL:H	2.23	0.42
2:B:209:PHE:O	2:B:225:GLY:CA	2.63	0.41
1:A:350:TYR:CD1	1:A:384:VAL:HG12	2.55	0.41
2:B:161:CYS:CB	2:B:302:VAL:HG11	2.50	0.41
2:B:235:LEU:HG	2:B:237:LEU:HD23	2.02	0.41
1:A:313:ARG:C	1:A:315:GLU:N	2.72	0.41
2:B:200:PRO:O	2:B:202:ASP:N	2.53	0.41
2:B:191:PHE:CE1	2:B:236:PHE:CB	3.04	0.41
2:B:220:VAL:O	2:B:220:VAL:HG12	2.19	0.41
1:A:354:GLY:HA3	1:A:358:GLU:HB2	2.03	0.41
2:B:162:PRO:HB3	2:B:275:ILE:HB	2.00	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:ASP:C	1:A:378:GLY:N	2.74	0.41
2:B:213:LYS:O	2:B:214:GLY:C	2.59	0.41
2:B:211:TYR:CE1	2:B:215:ALA:HB3	2.54	0.41
2:B:259:LYS:HD3	2:B:267:LYS:CE	2.51	0.41
2:B:254:VAL:HG12	2:B:254:VAL:O	2.21	0.41
1:A:379:GLU:O	1:A:382:CYS:HB2	2.20	0.41
1:A:383:LEU:N	1:A:383:LEU:HD23	2.36	0.41
2:B:162:PRO:HA	2:B:275:ILE:O	2.21	0.41
1:A:320:LEU:O	1:A:321:ALA:C	2.59	0.41
2:B:166:THR:CG2	2:B:279:THR:HG22	2.51	0.41
2:B:146:MET:HE3	2:B:307:ILE:HG23	2.03	0.41
2:B:215:ALA:CB	2:B:263:LEU:HD13	2.51	0.40
2:B:178:ARG:CG	2:B:178:ARG:HH11	2.34	0.40
1:A:332:TYR:OH	1:A:364:GLU:OE2	2.37	0.40
2:B:251:LEU:HD22	2:B:255:ILE:CD1	2.51	0.40
2:B:243:LEU:HA	2:B:243:LEU:HD23	1.79	0.40
2:B:252:LEU:CB	2:B:296:LEU:HD23	2.44	0.40
2:B:275:ILE:HG21	2:B:300:LEU:HD21	2.03	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	71/73 (97%)	51 (72%)	18 (25%)	2 (3%)	6	44
2	B	171/173 (99%)	127 (74%)	31 (18%)	13 (8%)	1	20
All	All	242/246 (98%)	178 (74%)	49 (20%)	15 (6%)	4	26

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	213	LYS
2	B	224	GLU
1	A	377	ARG
1	A	378	GLY
2	B	149	ILE
2	B	158	CYS
2	B	173	LYS
2	B	174	GLU
2	B	201	ARG
2	B	263	LEU
2	B	214	GLY
2	B	156	ILE
2	B	157	SER
2	B	286	LEU
2	B	241	GLY

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	65/65 (100%)	58 (89%)	7 (11%)	8	35
2	B	146/146 (100%)	125 (86%)	21 (14%)	4	25
All	All	211/211 (100%)	183 (87%)	28 (13%)	9	28

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	318	ILE
1	A	322	ASN
1	A	339	PHE
1	A	342	SER
1	A	357	ARG
1	A	359	LEU
1	A	383	LEU
2	B	139	TYR
2	B	161	CYS

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Mol	Chain	Res	Type
2	B	178	ARG
2	B	179	LEU
2	B	191	PHE
2	B	194	LEU
2	B	195	ASN
2	B	201	ARG
2	B	211	TYR
2	B	221	SER
2	B	227	PHE
2	B	251	LEU
2	B	261	TYR
2	B	262	ARG
2	B	263	LEU
2	B	274	ARG
2	B	281	ARG
2	B	285	GLU
2	B	295	ASP
2	B	299	ARG
2	B	307	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	355	ASN
1	A	361	ASN
2	B	181	HIS
2	B	282	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

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ADP	A	1385	-	24,29,29	1.55	6 (25%)	23,45,45	3.09	3 (13%)
3	ADP	B	1310	-	24,29,29	1.46	4 (16%)	23,45,45	3.03	3 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	A	1385	-	-	0/12/32/32	0/3/3/3
3	ADP	B	1310	-	-	0/12/32/32	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1385	ADP	C5-N7	-2.82	1.29	1.39
3	B	1310	ADP	C5-N7	-2.40	1.30	1.39
3	A	1385	ADP	C8-N7	-2.15	1.30	1.34
3	B	1310	ADP	PA-O2A	-2.06	1.46	1.55
3	A	1385	ADP	PA-O2A	-2.06	1.46	1.55
3	A	1385	ADP	C2-N1	2.09	1.37	1.33
3	B	1310	ADP	O4'-C1'	2.11	1.44	1.41
3	A	1385	ADP	C2-N3	2.56	1.36	1.32
3	B	1310	ADP	C2-N3	2.58	1.36	1.32
3	A	1385	ADP	O4'-C1'	2.79	1.45	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1310	ADP	N3-C2-N1	-13.65	118.15	128.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1385	ADP	N3-C2-N1	-13.57	118.22	128.87
3	B	1310	ADP	C2'-C3'-C4'	2.49	107.73	102.64
3	A	1385	ADP	C2'-C3'-C4'	2.55	107.86	102.64
3	B	1310	ADP	C4'-O4'-C1'	2.68	112.49	109.64
3	A	1385	ADP	C4'-O4'-C1'	3.16	112.99	109.64

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1385	ADP	2	0

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.