



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:23 am GMT

PDB ID : 1SQK  
Title : CRYSTAL STRUCTURE OF CIBOULOT IN COMPLEX WITH SKELETAL ACTIN  
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Deposited on : 2004-03-19  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

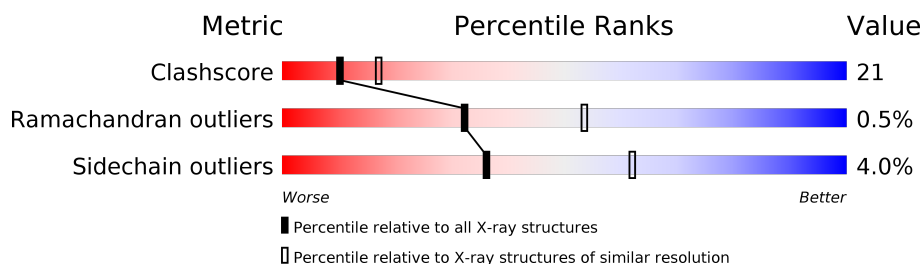
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	377	
2	B	25	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 3159 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called ACTIN, ALPHA SKELETAL MUSCLE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	360	Total	C	N	O	S	0	0	0
			2814	1783	473	540	18			

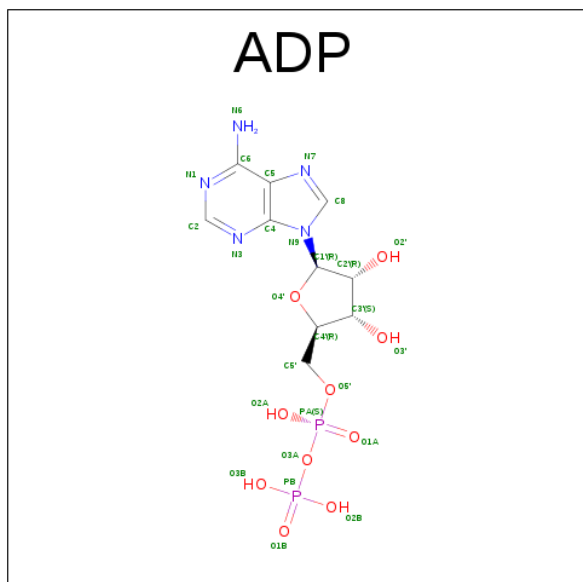
- Molecule 2 is a protein called CIBOULOT.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	25	Total	C	N	O	0	0	0
			196	121	34	41			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

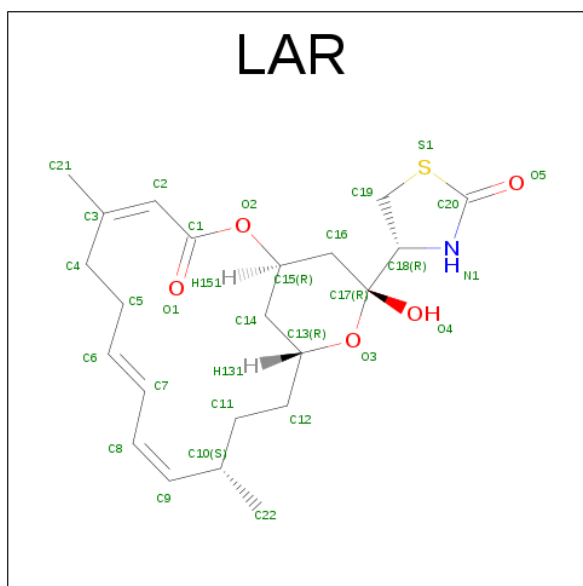
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>).



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

- Molecule 5 is LATRUNCULIN A (three-letter code: LAR) (formula:  $C_{22}H_{31}NO_5S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			29	22	1	5	1		

- Molecule 6 is water.

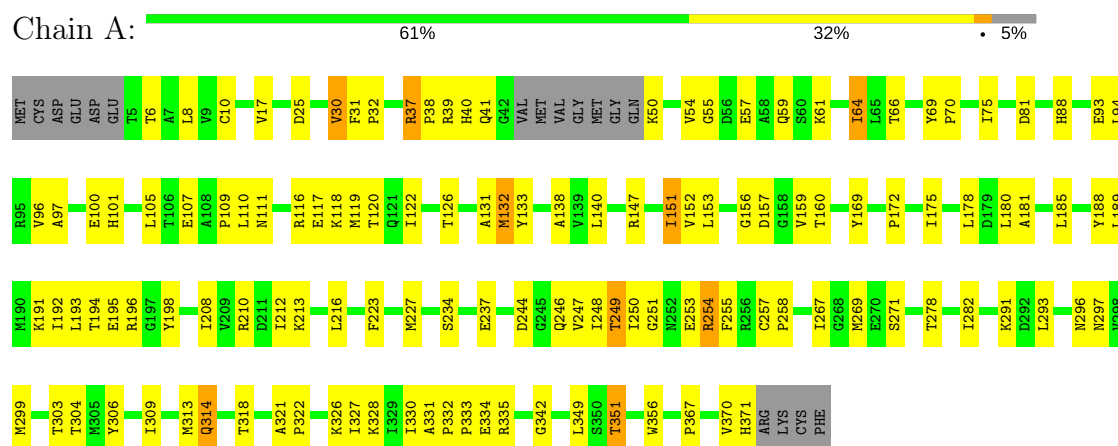
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	90	Total	O	0	0
			90	90		
6	B	2	Total	O	0	0
			2	2		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

#### • Molecule 1: ACTIN, ALPHA SKELETAL MUSCLE



#### • Molecule 2: CIBOULOT



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.49 Å 75.20 Å 85.31 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	99.7 (20.00-2.50)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.09	Depositor
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.219 , 0.278	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3159	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	0.36	0/2875	0.61	0/3898
2	B	0.44	0/197	0.66	0/261
All	All	0.37	0/3072	0.61	0/4159

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	2814	0	2780	113	0
2	B	196	0	197	21	0
3	A	1	0	0	0	0
4	A	27	0	12	1	0
5	A	29	0	31	2	0
6	A	90	0	0	5	0
6	B	2	0	0	0	0
All	All	3159	0	3020	127	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 21.

All (127) 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:210:ARG:HB3	1:A:210:ARG:HH11	1.23	0.99
1:A:147:ARG:NE	1:A:296:ASN:HD22	1.63	0.96
1:A:147:ARG:HE	1:A:296:ASN:ND2	1.63	0.94
1:A:210:ARG:HB3	1:A:210:ARG:NH1	1.83	0.93
1:A:160:THR:HG23	1:A:178:LEU:HB3	1.53	0.91
1:A:147:ARG:HE	1:A:296:ASN:HD22	0.91	0.88
1:A:278:THR:HG21	1:A:297:ASN:HD21	1.39	0.86
1:A:169:TYR:HE1	2:B:19:LYS:HZ1	1.22	0.85
1:A:367:PRO:O	1:A:370:VAL:HG23	1.76	0.84
2:B:28:ASP:O	2:B:29:LYS:HG2	1.80	0.81
1:A:64:ILE:HD13	1:A:64:ILE:H	1.50	0.77
1:A:278:THR:O	1:A:282:ILE:HG12	1.85	0.76
1:A:198:TYR:CE1	1:A:248:ILE:HD13	2.21	0.76
1:A:40:HIS:O	1:A:41:GLN:HG2	1.86	0.74
1:A:210:ARG:CB	1:A:210:ARG:HH11	2.01	0.73
1:A:61:LYS:HB3	1:A:64:ILE:HD11	1.72	0.71
1:A:131:ALA:HB1	1:A:356:TRP:HB3	1.70	0.71
2:B:11:LEU:N	2:B:12:PRO:HD2	2.06	0.70
1:A:132:MET:HG2	1:A:133:TYR:N	2.08	0.69
1:A:160:THR:CG2	1:A:178:LEU:HB3	2.23	0.69
1:A:97:ALA:HB3	1:A:100:GLU:HG2	1.75	0.68
1:A:234:SER:HB2	1:A:237:GLU:OE1	1.93	0.67
2:B:10:ASP:N	2:B:12:PRO:HD2	2.09	0.67
1:A:326:LYS:HE3	1:A:328:LYS:HB2	1.78	0.65
1:A:198:TYR:CZ	1:A:248:ILE:HD13	2.31	0.65
1:A:61:LYS:O	1:A:64:ILE:HD13	1.97	0.64
1:A:257:CYS:HB3	1:A:258:PRO:HD3	1.80	0.64
1:A:194:THR:HA	1:A:198:TYR:O	1.97	0.64
2:B:11:LEU:N	2:B:12:PRO:CD	2.61	0.63
1:A:216:LEU:O	1:A:254:ARG:HG2	1.99	0.62
1:A:334:GLU:N	1:A:334:GLU:OE1	2.28	0.61
1:A:185:LEU:HD23	1:A:306:TYR:OH	2.00	0.61
1:A:31:PHE:CE2	1:A:93:GLU:HG3	2.36	0.61
2:B:28:ASP:C	2:B:30:LEU:H	2.02	0.61
1:A:50:LYS:HD2	1:A:50:LYS:N	2.16	0.60
2:B:28:ASP:C	2:B:30:LEU:N	2.55	0.60
1:A:120:THR:OG1	1:A:370:VAL:HG21	2.01	0.59
2:B:28:ASP:OD1	2:B:29:LYS:N	2.33	0.59
2:B:28:ASP:O	2:B:30:LEU:N	2.35	0.59
1:A:39:ARG:NH2	1:A:66:THR:HG22	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:TYR:CG	1:A:267:ILE:HD11	2.38	0.59
1:A:251:GLY:O	1:A:254:ARG:HB2	2.02	0.59
1:A:59:GLN:HA	1:A:59:GLN:HE21	1.67	0.59
1:A:299:MET:HE2	1:A:331:ALA:HB2	1.86	0.58
1:A:117:GLU:OE2	1:A:371:HIS:HE1	1.86	0.58
2:B:11:LEU:HD12	2:B:14:VAL:HG21	1.86	0.58
2:B:10:ASP:C	2:B:12:PRO:HD2	2.24	0.58
1:A:39:ARG:CZ	1:A:66:THR:HG22	2.34	0.57
1:A:88:HIS:HE1	1:A:93:GLU:OE2	1.88	0.57
1:A:303:THR:O	1:A:303:THR:HG22	2.06	0.55
1:A:267:ILE:HG22	1:A:269:MET:HG3	1.88	0.54
1:A:191:LYS:O	1:A:195:GLU:HG3	2.08	0.54
1:A:64:ILE:HD13	1:A:64:ILE:N	2.21	0.54
1:A:97:ALA:HB3	1:A:100:GLU:CG	2.36	0.54
1:A:25:ASP:OD1	2:B:32:ASN:HA	2.08	0.54
1:A:59:GLN:HA	1:A:59:GLN:NE2	2.23	0.54
1:A:189:LEU:HD12	1:A:192:ILE:HD11	1.88	0.54
1:A:250:ILE:HB	1:A:253:GLU:HG2	1.89	0.54
1:A:25:ASP:HA	2:B:30:LEU:HD23	1.90	0.54
1:A:267:ILE:CG2	1:A:267:ILE:O	2.56	0.54
1:A:351:THR:HB	2:B:17:ASN:HD22	1.72	0.53
1:A:248:ILE:HD12	1:A:249:THR:H	1.73	0.53
1:A:321:ALA:HB1	1:A:322:PRO:HD2	1.90	0.53
1:A:160:THR:HG22	1:A:180:LEU:O	2.09	0.53
1:A:351:THR:HB	2:B:17:ASN:ND2	2.24	0.52
1:A:278:THR:CG2	1:A:297:ASN:HD21	2.19	0.52
1:A:37:ARG:NH2	1:A:81:ASP:OD2	2.43	0.52
1:A:41:GLN:HG3	1:A:41:GLN:O	2.10	0.51
1:A:111:ASN:HD21	1:A:119:MET:CE	2.23	0.51
1:A:118:LYS:O	1:A:122:ILE:HG12	2.12	0.50
1:A:326:LYS:HD3	1:A:326:LYS:C	2.32	0.50
1:A:109:PRO:O	1:A:110:LEU:HB2	2.12	0.49
1:A:17:VAL:O	1:A:30:VAL:HA	2.12	0.49
1:A:244:ASP:OD2	1:A:246:GLN:HG3	2.13	0.49
1:A:8:LEU:HD22	1:A:94:LEU:HD13	1.95	0.49
1:A:208:ILE:O	1:A:212:ILE:HG13	2.12	0.49
1:A:151:ILE:HB	1:A:293:LEU:HD22	1.94	0.48
2:B:33:ALA:O	2:B:34:SER:OG	2.24	0.48
1:A:189:LEU:O	1:A:193:LEU:HG	2.14	0.48
1:A:70:PRO:HG3	1:A:81:ASP:HB3	1.96	0.48
2:B:19:LYS:O	2:B:23:GLU:HG3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:VAL:HG12	1:A:247:VAL:O	2.14	0.47
1:A:159:VAL:HG22	1:A:160:THR:N	2.29	0.47
1:A:132:MET:HG2	1:A:133:TYR:H	1.78	0.47
1:A:223:PHE:O	1:A:227:MET:HB2	2.15	0.47
1:A:248:ILE:HD12	1:A:249:THR:N	2.30	0.47
1:A:172:PRO:HA	1:A:175:ILE:HD12	1.96	0.47
2:B:15:ALA:O	2:B:19:LYS:HG3	2.14	0.46
1:A:138:ALA:HA	6:A:384:HOH:O	2.15	0.46
1:A:153:LEU:HD22	1:A:313:MET:HE1	1.97	0.46
1:A:32:PRO:HB3	5:A:378:LAR:H121	1.98	0.45
1:A:10:CYS:HB3	1:A:105:LEU:HD23	1.98	0.45
1:A:349:LEU:HD13	2:B:21:GLN:HB3	1.99	0.44
1:A:328:LYS:HE2	1:A:330:ILE:CD1	2.47	0.44
1:A:299:MET:HE1	1:A:309:ILE:HG23	1.98	0.44
1:A:151:ILE:HD13	1:A:152:VAL:N	2.33	0.44
1:A:75:ILE:N	1:A:75:ILE:HD12	2.32	0.44
1:A:147:ARG:CD	1:A:296:ASN:HD22	2.27	0.44
1:A:196:ARG:HG3	6:A:425:HOH:O	2.18	0.43
1:A:54:VAL:HG12	1:A:55:GLY:N	2.33	0.43
1:A:88:HIS:CE1	1:A:93:GLU:OE2	2.71	0.43
1:A:304:THR:O	1:A:335:ARG:NH1	2.42	0.43
1:A:314:GLN:O	1:A:318:THR:HG23	2.17	0.43
1:A:271:SER:HB2	6:A:432:HOH:O	2.18	0.43
1:A:111:ASN:ND2	1:A:119:MET:CE	2.82	0.43
1:A:351:THR:CB	2:B:17:ASN:HD22	2.32	0.43
1:A:59:GLN:HB2	5:A:378:LAR:H91	2.01	0.43
1:A:327:ILE:N	1:A:327:ILE:HD12	2.33	0.42
1:A:156:GLY:O	1:A:181:ALA:HB1	2.20	0.42
1:A:31:PHE:CD2	1:A:93:GLU:HG3	2.54	0.42
1:A:216:LEU:HD22	1:A:250:ILE:HD11	2.02	0.42
1:A:59:GLN:NE2	1:A:59:GLN:CA	2.83	0.42
1:A:210:ARG:O	1:A:213:LYS:HB3	2.19	0.41
1:A:37:ARG:HG3	1:A:38:PRO:HD2	2.01	0.41
1:A:96:VAL:HB	1:A:101:HIS:CE1	2.55	0.41
2:B:18:LEU:HD23	2:B:18:LEU:C	2.41	0.41
1:A:107:GLU:OE1	1:A:116:ARG:NH1	2.54	0.41
1:A:140:LEU:O	1:A:342:GLY:HA3	2.21	0.41
1:A:291:LYS:HG3	6:A:423:HOH:O	2.20	0.41
1:A:122:ILE:O	1:A:126:THR:HB	2.21	0.41
1:A:255:PHE:C	1:A:258:PRO:HD2	2.40	0.41
1:A:59:GLN:HE21	1:A:59:GLN:CA	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:TYR:CD1	1:A:169:TYR:N	2.90	0.40
1:A:332:PRO:HA	1:A:333:PRO:HD3	1.86	0.40
1:A:37:ARG:HG3	1:A:38:PRO:CD	2.51	0.40
1:A:248:ILE:HD12	6:A:468:HOH:O	2.21	0.40
1:A:157:ASP:HB2	4:A:377:ADP:H5'1	2.04	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 X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	356/377 (94%)	331 (93%)	25 (7%)	0	100	100
2	B	23/25 (92%)	19 (83%)	2 (9%)	2 (9%)	1	1
All	All	379/402 (94%)	350 (92%)	27 (7%)	2 (0%)	32	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	28	ASP
2	B	29	LYS

### 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 X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	305/320 (95%)	293 (96%)	12 (4%)	37	63
2	B	22/22 (100%)	21 (96%)	1 (4%)	32	56
All	All	327/342 (96%)	314 (96%)	13 (4%)	36	62

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	30	VAL
1	A	37	ARG
1	A	57	GLU
1	A	64	ILE
1	A	69	TYR
1	A	132	MET
1	A	151	ILE
1	A	249	THR
1	A	254	ARG
1	A	314	GLN
1	A	351	THR
2	B	21	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	73	HIS
1	A	88	HIS
1	A	92	ASN
1	A	101	HIS
1	A	111	ASN
1	A	121	GLN
1	A	137	GLN
1	A	162	ASN
1	A	296	ASN
1	A	297	ASN
1	A	371	HIS
2	B	17	ASN
2	B	32	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](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

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$
4	ADP	A	377	3	25,29,29	1.58	5 (20%)	24,45,45	2.79	7 (29%)
5	LAR	A	378	-	30,31,31	1.71	7 (23%)	36,43,43	2.00	7 (19%)

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
4	ADP	A	377	3	-	0/12/32/32	0/3/3/3
5	LAR	A	378	-	-	0/23/51/51	0/1/3/3

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	378	LAR	C20-N1	-2.99	1.32	1.36

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	377	ADP	C5-N7	-2.62	1.30	1.39
4	A	377	ADP	C2'-C1'	-2.01	1.50	1.53
5	A	378	LAR	C22-C10	2.07	1.61	1.53
5	A	378	LAR	C21-C3	2.11	1.56	1.50
5	A	378	LAR	O3-C13	2.14	1.49	1.44
5	A	378	LAR	O4-C17	2.40	1.45	1.40
4	A	377	ADP	C4-N3	2.66	1.39	1.35
4	A	377	ADP	O4'-C1'	2.93	1.45	1.41
5	A	378	LAR	C4-C3	3.06	1.58	1.51
4	A	377	ADP	C2-N3	3.51	1.38	1.32
5	A	378	LAR	C18-N1	5.42	1.52	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	377	ADP	N3-C2-N1	-10.97	119.30	128.86
4	A	377	ADP	O3A-PB-O1B	-4.26	85.24	111.44
5	A	378	LAR	S1-C20-N1	-3.31	108.35	111.20
4	A	377	ADP	C4'-O4'-C1'	-2.62	106.98	109.77
4	A	377	ADP	C4-C5-N7	-2.02	107.46	109.41
5	A	378	LAR	C19-C18-N1	2.21	106.68	101.00
4	A	377	ADP	O3B-PB-O2B	2.31	116.93	107.61
5	A	378	LAR	O2-C15-C16	2.50	113.86	107.61
4	A	377	ADP	O4'-C4'-C5'	2.96	119.39	109.40
5	A	378	LAR	C3-C2-C1	3.30	135.99	127.64
4	A	377	ADP	O2B-PB-O1B	3.34	123.57	110.50
5	A	378	LAR	O5-C20-N1	4.34	130.13	125.83
5	A	378	LAR	C19-S1-C20	4.91	94.62	92.00
5	A	378	LAR	C18-N1-C20	5.98	122.08	113.28

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	377	ADP	1	0
5	A	378	LAR	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.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.