



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:09 pm BST

PDB ID : 3Q33  
Title : Structure of the Rtt109-AcCoA/Vps75 Complex and Implications for  
Chaperone-Mediated Histone Acetylation  
Authors : Tang, Y.; Yuan, H.; Meeth, K.; Marmorstein, R.  
Deposited on : 2010-12-21  
Resolution : 2.80 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

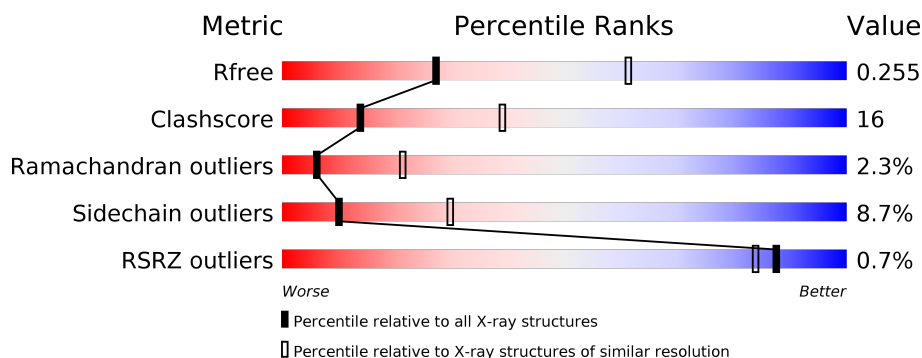
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 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  $\leq 5\%$ . 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.

Mol	Chain	Length	Quality of chain
1	A	438	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0;">%</div> <div style="position: absolute; top: 0; left: 0; width: 58%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 58%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 86%; height: 10px; background-color: orange;"></div> <div style="position: absolute; top: 0; left: 96%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 0; left: 100%; height: 10px; background-color: grey;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>58%</span> <span>28%</span> <span>• • 10%</span> </div> </div>
2	B	232	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: 0; left: 0; width: 59%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 59%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 87%; height: 10px; background-color: orange;"></div> <div style="position: absolute; top: 0; left: 97%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 0; left: 100%; height: 10px; background-color: grey;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>59%</span> <span>28%</span> <span>• 8%</span> </div> </div>
3	D	15	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: 0; left: 0; width: 20%; height: 10px; background-color: green;"></div> <div style="position: absolute; top: 0; left: 20%; height: 10px; background-color: yellow;"></div> <div style="position: absolute; top: 0; left: 27%; height: 10px; background-color: orange;"></div> <div style="position: absolute; top: 0; left: 34%; height: 10px; background-color: red;"></div> <div style="position: absolute; top: 0; left: 67%; height: 10px; background-color: grey;"></div> </div> <div style="display: flex; justify-content: space-between; width: 100%;"> <span>20%</span> <span>7%</span> <span>7%</span> <span>67%</span> </div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5077 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 Histone acetyltransferase RTT109.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3198	2058	539	592	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP A7A0Q4
A	0	SER	-	EXPRESSION TAG	UNP A7A0Q4

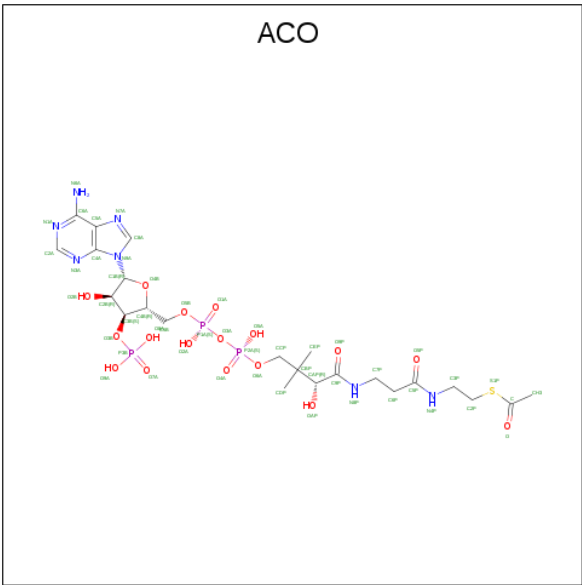
- Molecule 2 is a protein called Vacuolar protein sorting-associated protein 75.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1776	1149	288	334	5			

- Molecule 3 is a protein called HISTONE H3.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	5	Total	C	N	O	0	0	1
			25	14	6	5			

- Molecule 4 is ACETYL COENZYME \*A (three-letter code: ACO) (formula: C<sub>23</sub>H<sub>38</sub>N<sub>7</sub>O<sub>17</sub>P<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
4	A	1	51	23	7	17	3	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	4	2	2	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	14	Total 14	O 14	0	0
6	B	9	Total 9	O 9	0	0



ALA	ARG	THR	LYS	GLN	THR	ALA	ARG	LYS	SER	T11	G12	G13	K14	T15
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## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.54Å 119.57Å 80.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.16 – 2.80 45.16 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.16-2.80) 96.0 (45.16-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.13 (at 2.81Å)	Xtriage
Refinement program	PHENIX phenix.refine: 1.6.4_486	Depositor
R, $R_{free}$	0.195 , 0.255 0.191 , 0.255	Depositor DCC
$R_{free}$ test set	1167 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	69.2	Xtriage
Anisotropy	0.552	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 59.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5077	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.19% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ALY, ACO, EDO, NH2

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	0.44	0/3255	0.61	0/4402
2	B	0.42	0/1823	0.57	0/2455
3	D	0.60	0/23	0.62	0/28
All	All	0.44	0/5101	0.59	0/6885

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	3198	0	3239	103	0
2	B	1776	0	1721	55	0
3	D	25	0	25	2	0
4	A	51	0	34	0	0
5	A	4	0	6	2	0
6	A	14	0	0	0	0
6	B	9	0	0	1	0
All	All	5077	0	5025	159	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 16.

All (159) 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:164:ARG:HG3	2:B:164:ARG:HH11	1.12	1.06
2:B:164:ARG:HH11	2:B:164:ARG:CG	1.90	0.83
2:B:184:ARG:HH11	2:B:184:ARG:HG3	1.44	0.83
2:B:209:TYR:HB3	2:B:210:PRO:HD3	1.62	0.80
2:B:164:ARG:HG3	2:B:164:ARG:NH1	1.82	0.80
1:A:115:ASP:HB2	1:A:341:SER:HA	1.62	0.79
2:B:25:VAL:HA	2:B:28:ILE:HD12	1.70	0.73
2:B:189:LYS:HB3	2:B:192:LYS:HD2	1.70	0.72
1:A:214:ASN:ND2	1:A:217:GLU:HG3	2.04	0.72
1:A:138:ILE:HG21	1:A:147:THR:HG23	1.72	0.71
1:A:316:GLN:HG3	1:A:317:GLU:OE1	1.91	0.71
2:B:165:SER:HB2	2:B:166:PRO:HD2	1.77	0.67
1:A:47:ILE:N	1:A:47:ILE:HD12	2.11	0.66
2:B:129:LYS:HD2	2:B:129:LYS:H	1.59	0.66
1:A:244:ILE:HD12	1:A:252:VAL:HG12	1.78	0.66
1:A:6:PHE:HB3	1:A:106:ILE:HG23	1.78	0.65
2:B:119:GLN:HE22	2:B:147:GLU:HB2	1.61	0.65
2:B:119:GLN:NE2	2:B:147:GLU:HB2	2.11	0.65
1:A:173:GLN:OE1	1:A:173:GLN:HA	1.97	0.65
1:A:325:VAL:HG12	1:A:326:THR:H	1.63	0.64
1:A:319:GLN:O	1:A:321:PHE:N	2.31	0.64
1:A:184:GLU:HB2	1:A:339:THR:HG23	1.80	0.64
2:B:221:ARG:HG3	2:B:221:ARG:HH11	1.63	0.64
2:B:10:GLU:OE2	2:B:13:LYS:HD2	1.98	0.64
2:B:184:ARG:CG	2:B:184:ARG:HH11	2.11	0.63
2:B:156:ASN:HB3	2:B:159:LEU:HB2	1.79	0.63
2:B:94:GLU:O	2:B:95:SER:HB3	1.97	0.63
1:A:384:ARG:HG3	1:A:385:ASP:N	2.13	0.62
1:A:384:ARG:HG3	1:A:385:ASP:H	1.64	0.62
1:A:6:PHE:HB3	1:A:106:ILE:CG2	2.30	0.61
1:A:118:TYR:CE1	1:A:344:PRO:HD2	2.36	0.61
2:B:52:ASP:OD1	2:B:85:LYS:HE2	2.01	0.61
3:D:13:GLY:O	3:D:14:LYS:CB	2.48	0.60
1:A:105:LYS:HB2	1:A:225:ILE:HG23	1.84	0.60
1:A:303:LEU:H	1:A:303:LEU:HD12	1.68	0.59
1:A:368:GLU:HB3	1:A:370:TYR:CE2	2.37	0.58
2:B:65:HIS:CE1	2:B:67:SER:HB3	2.39	0.58
2:B:222:ASP:O	2:B:223:LEU:HB2	2.05	0.57
1:A:154:ARG:CZ	1:A:163:LEU:HD11	2.34	0.57
1:A:191:LEU:C	1:A:191:LEU:HD12	2.25	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:251:ARG:HB3	1:A:251:ARG:NH1	2.20	0.57
2:B:221:ARG:HG3	2:B:221:ARG:NH1	2.18	0.56
1:A:292:ARG:NH1	1:A:320:GLU:HG2	2.19	0.56
2:B:65:HIS:HE1	2:B:67:SER:HB3	1.71	0.56
1:A:325:VAL:HG12	1:A:326:THR:N	2.20	0.56
2:B:189:LYS:N	2:B:189:LYS:HD2	2.21	0.56
1:A:101:ARG:HD3	1:A:224:PHE:CD2	2.41	0.56
1:A:128:ARG:HH22	1:A:176:TYR:HE1	1.55	0.55
1:A:154:ARG:NH2	1:A:163:LEU:HD11	2.22	0.55
2:B:142:GLU:HG3	2:B:143:PRO:HD2	1.89	0.55
2:B:21:CYS:O	2:B:25:VAL:HG23	2.07	0.54
2:B:33:GLU:OE1	2:B:36:ARG:NH1	2.41	0.54
1:A:162:VAL:HG12	1:A:163:LEU:H	1.71	0.54
1:A:66:GLU:HG2	1:A:86:SER:HB3	1.90	0.54
1:A:214:ASN:HD21	1:A:217:GLU:HG3	1.71	0.54
1:A:20:LEU:O	1:A:53:PHE:HA	2.07	0.54
1:A:318:ARG:HH11	1:A:318:ARG:HG3	1.73	0.53
1:A:13:VAL:HG12	1:A:401:LYS:HA	1.91	0.53
2:B:94:GLU:O	2:B:95:SER:CB	2.57	0.53
1:A:193:THR:O	1:A:194:ARG:HG3	2.10	0.52
1:A:243:ARG:NH1	1:A:268:ASP:OD2	2.43	0.52
2:B:27:ALA:HB2	6:B:306:HOH:O	2.10	0.52
1:A:18:GLU:HA	1:A:396:GLN:O	2.10	0.52
2:B:184:ARG:CG	2:B:184:ARG:NH1	2.71	0.51
2:B:24:GLU:O	2:B:28:ILE:HG13	2.10	0.51
1:A:109:GLU:HG2	1:A:233:CYS:SG	2.51	0.51
2:B:109:PHE:O	2:B:118:GLU:HG2	2.11	0.51
2:B:223:LEU:O	2:B:224:GLU:C	2.48	0.51
1:A:290:ALY:HH33	1:A:329:VAL:CG1	2.41	0.51
2:B:78:LYS:HD3	2:B:79:TYR:CE1	2.46	0.51
1:A:298:ALA:HA	1:A:303:LEU:HD21	1.92	0.50
1:A:207:ASN:OD1	1:A:209:LYS:HD2	2.10	0.50
1:A:136:GLU:HG3	1:A:136:GLU:O	2.12	0.50
1:A:393:THR:HG22	1:A:394:ASN:H	1.77	0.50
1:A:13:VAL:HG13	1:A:402:ARG:C	2.32	0.49
1:A:319:GLN:HA	1:A:322:LYS:HB3	1.93	0.49
2:B:179:ILE:HD13	2:B:179:ILE:O	2.12	0.49
1:A:198:GLN:HE21	1:A:198:GLN:C	2.16	0.49
1:A:7:LEU:O	1:A:10:VAL:HG23	2.13	0.48
1:A:160:SER:OG	1:A:161:THR:N	2.43	0.48
1:A:244:ILE:CD1	1:A:252:VAL:HG12	2.42	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:GLN:O	1:A:300:GLU:HG2	2.13	0.48
1:A:136:GLU:HA	1:A:136:GLU:OE2	2.14	0.48
1:A:318:ARG:CZ	1:A:318:ARG:HB2	2.44	0.48
1:A:248:ASP:O	1:A:251:ARG:HB2	2.14	0.48
1:A:4:ASN:HD21	1:A:396:GLN:HB3	1.78	0.48
1:A:245:PRO:HD2	1:A:327:SER:O	2.14	0.47
1:A:11:LEU:HD13	1:A:17:PHE:CD1	2.50	0.47
1:A:165:GLU:O	1:A:172:GLN:HA	2.14	0.47
1:A:65:LEU:HD12	1:A:87:LYS:O	2.14	0.47
2:B:188:LEU:C	2:B:190:PRO:HD3	2.34	0.47
2:B:102:ASP:OD1	2:B:126:GLN:HA	2.15	0.47
3:D:13:GLY:O	3:D:14:LYS:HB3	2.15	0.47
1:A:115:ASP:H	1:A:342:LEU:H	1.62	0.46
1:A:293:PHE:O	1:A:293:PHE:CD1	2.68	0.46
1:A:318:ARG:O	1:A:319:GLN:HG3	2.16	0.46
1:A:378:GLU:OE1	2:B:73:ARG:HD3	2.15	0.46
1:A:57:HIS:NE2	1:A:58:GLN:NE2	2.63	0.46
1:A:128:ARG:NH2	1:A:176:TYR:HE1	2.14	0.46
1:A:16:GLN:HA	1:A:398:LEU:O	2.15	0.46
1:A:298:ALA:CB	1:A:303:LEU:HD21	2.46	0.46
2:B:156:ASN:O	2:B:160:ILE:HG13	2.15	0.46
2:B:30:ARG:NH2	2:B:96:GLU:HA	2.31	0.45
1:A:118:TYR:CD1	1:A:344:PRO:HD2	2.51	0.45
1:A:186:LEU:HD12	1:A:334:GLY:O	2.16	0.45
1:A:59:GLY:HA2	1:A:380:PHE:CD2	2.52	0.45
2:B:141:SER:OG	2:B:142:GLU:N	2.50	0.45
1:A:318:ARG:HH11	1:A:318:ARG:CG	2.29	0.45
1:A:290:ALY:HH33	1:A:329:VAL:HG13	1.97	0.45
1:A:271:THR:HG22	1:A:272:SER:N	2.32	0.45
1:A:86:SER:O	5:A:502:EDO:H12	2.16	0.45
1:A:194:ARG:NH1	5:A:502:EDO:O2	2.49	0.44
1:A:17:PHE:C	1:A:18:GLU:HG3	2.36	0.44
1:A:342:LEU:O	1:A:344:PRO:HD3	2.18	0.44
1:A:361:ILE:HG23	1:A:383:ILE:HD13	1.98	0.44
1:A:261:TYR:CD1	1:A:262:PRO:HD2	2.52	0.44
2:B:29:GLU:HA	2:B:32:VAL:CG2	2.48	0.44
1:A:149:ARG:O	1:A:152:ALA:HB3	2.17	0.44
1:A:53:PHE:N	1:A:53:PHE:CD2	2.85	0.44
1:A:350:ILE:O	1:A:352:PRO:HD3	2.18	0.44
2:B:82:THR:O	2:B:109:PHE:HA	2.17	0.44
2:B:77:PHE:HA	2:B:80:ILE:HB	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ILE:N	1:A:47:ILE:CD1	2.81	0.44
1:A:106:ILE:HD13	1:A:106:ILE:HA	1.82	0.44
1:A:242:LEU:HG	1:A:243:ARG:N	2.34	0.43
2:B:146:ILE:HG13	2:B:147:GLU:N	2.33	0.43
1:A:325:VAL:CG1	1:A:326:THR:H	2.30	0.43
1:A:75:ASP:OD1	1:A:76:GLU:N	2.50	0.43
1:A:51:HIS:HB3	1:A:53:PHE:CE2	2.53	0.43
2:B:169:LYS:O	2:B:173:ARG:HG2	2.18	0.43
2:B:29:GLU:HA	2:B:32:VAL:HG22	2.01	0.42
1:A:60:LYS:NZ	1:A:373:GLU:OE1	2.48	0.42
1:A:45:SER:OG	1:A:73:LEU:HD12	2.19	0.42
2:B:193:GLU:O	2:B:194:PHE:C	2.57	0.42
1:A:200:LEU:O	1:A:366:THR:HA	2.19	0.42
2:B:38:ASN:O	2:B:42:PRO:HD2	2.20	0.42
1:A:261:TYR:HA	1:A:262:PRO:HD3	1.69	0.42
1:A:212:ILE:HD12	1:A:212:ILE:N	2.34	0.42
1:A:394:ASN:N	1:A:394:ASN:OD1	2.52	0.42
1:A:398:LEU:HA	1:A:398:LEU:HD23	1.81	0.42
1:A:13:VAL:HG13	1:A:402:ARG:N	2.35	0.41
1:A:36:ASN:ND2	1:A:39:ASP:HB2	2.34	0.41
2:B:128:LYS:O	2:B:137:GLY:HA2	2.20	0.41
2:B:171:LYS:HD3	2:B:171:LYS:HA	1.87	0.41
1:A:163:LEU:HD12	1:A:163:LEU:N	2.36	0.41
1:A:285:PHE:HB3	1:A:286:PRO:HD2	2.01	0.41
2:B:129:LYS:HB3	2:B:129:LYS:HE3	1.87	0.41
2:B:42:PRO:O	2:B:46:LYS:HG2	2.21	0.41
1:A:271:THR:O	1:A:272:SER:O	2.39	0.41
2:B:25:VAL:HA	2:B:28:ILE:CD1	2.44	0.41
2:B:184:ARG:O	2:B:185:TRP:C	2.59	0.41
1:A:318:ARG:O	1:A:319:GLN:C	2.59	0.41
2:B:44:TYR:HE2	2:B:90:TRP:CH2	2.39	0.41
1:A:138:ILE:HD13	1:A:147:THR:HG23	2.03	0.40
2:B:99:ASP:C	2:B:101:ARG:N	2.74	0.40
1:A:135:PRO:O	1:A:136:GLU:HB3	2.21	0.40
1:A:353:LYS:HG2	1:A:357:GLN:OE1	2.21	0.40
1:A:298:ALA:HB2	1:A:303:LEU:HD21	2.02	0.40
1:A:355:ARG:O	1:A:359:ARG:HG3	2.22	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	387/438 (88%)	340 (88%)	37 (10%)	10 (3%)	5	18
2	B	209/232 (90%)	193 (92%)	13 (6%)	3 (1%)	11	34
3	D	3/15 (20%)	2 (67%)	0	1 (33%)	0	0
All	All	599/685 (87%)	535 (89%)	50 (8%)	14 (2%)	6	21

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	TYR
1	A	272	SER
1	A	121	GLN
1	A	175	LEU
1	A	320	GLU
2	B	95	SER
3	D	14	LYS
1	A	135	PRO
1	A	271	THR
2	B	185	TRP
1	A	37	LYS
1	A	127	ILE
1	A	319	GLN
2	B	225	ASP

### 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	359/398 (90%)	327 (91%)	32 (9%)	9	28
2	B	190/207 (92%)	175 (92%)	15 (8%)	12	34
3	D	2/10 (20%)	1 (50%)	1 (50%)	0	0
All	All	551/615 (90%)	503 (91%)	48 (9%)	10	30

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	A	24	SER
1	A	34	THR
1	A	37	LYS
1	A	38	ASP
1	A	122	LYS
1	A	128	ARG
1	A	153	ARG
1	A	164	LYS
1	A	165	GLU
1	A	173	GLN
1	A	176	TYR
1	A	192	PHE
1	A	198	GLN
1	A	205	SER
1	A	208	SER
1	A	209	LYS
1	A	216	GLU
1	A	228	ARG
1	A	243	ARG
1	A	251	ARG
1	A	254	SER
1	A	263	LEU
1	A	273	LYS
1	A	303	LEU
1	A	318	ARG
1	A	319	GLN
1	A	320	GLU
1	A	336	SER
1	A	355	ARG
1	A	384	ARG
1	A	393	THR
2	B	20	LYS
2	B	32	VAL

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Mol	Chain	Res	Type
2	B	45	GLU
2	B	66	VAL
2	B	73	ARG
2	B	95	SER
2	B	117	LYS
2	B	118	GLU
2	B	129	LYS
2	B	146	ILE
2	B	159	LEU
2	B	164	ARG
2	B	179	ILE
2	B	220	GLN
2	B	223	LEU
3	D	14	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	ASN
1	A	198	GLN
1	A	275	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is 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	$\# Z  > 2$
1	ALY	A	290	1	10,11,12	0.34	0	7,12,14	0.63	0



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
1	ALY	A	290	1	-	3/9/10/12	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	290	ALY	OH-CH-NZ-CE
1	A	290	ALY	CH3-CH-NZ-CE
1	A	290	ALY	CG-CD-CE-NZ

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	290	ALY	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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
4	ACO	A	501	-	45,53,53	0.90	3 (6%)	56,79,79	1.68	11 (19%)
5	EDO	A	502	-	3,3,3	0.56	0	2,2,2	0.44	0

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	ACO	A	501	-	-	16/47/67/67	0/3/3/3
5	EDO	A	502	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	501	ACO	C5A-C4A	2.39	1.47	1.40
4	A	501	ACO	C8A-N7A	2.06	1.38	1.34
4	A	501	ACO	C5A-N7A	-2.02	1.32	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	501	ACO	C4A-C5A-N7A	-3.80	105.44	109.40
4	A	501	ACO	C2P-S1P-C	3.77	121.53	101.68
4	A	501	ACO	CAP-C9P-N8P	3.57	123.69	116.58
4	A	501	ACO	C2P-C3P-N4P	-3.52	105.01	112.42
4	A	501	ACO	N3A-C2A-N1A	-3.47	123.25	128.68
4	A	501	ACO	C6P-C7P-N8P	3.08	118.11	111.90
4	A	501	ACO	C1B-N9A-C4A	-2.64	122.00	126.64
4	A	501	ACO	OAP-CAP-CBP	-2.38	104.65	110.25
4	A	501	ACO	C3P-N4P-C5P	2.27	127.06	122.84
4	A	501	ACO	C2A-N1A-C6A	2.02	122.20	118.75
4	A	501	ACO	C2B-C3B-C4B	-2.01	99.66	103.22

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	501	ACO	C5B-O5B-P1A-O3A
4	A	501	ACO	O9P-C9P-CAP-CBP
4	A	501	ACO	N8P-C9P-CAP-CBP

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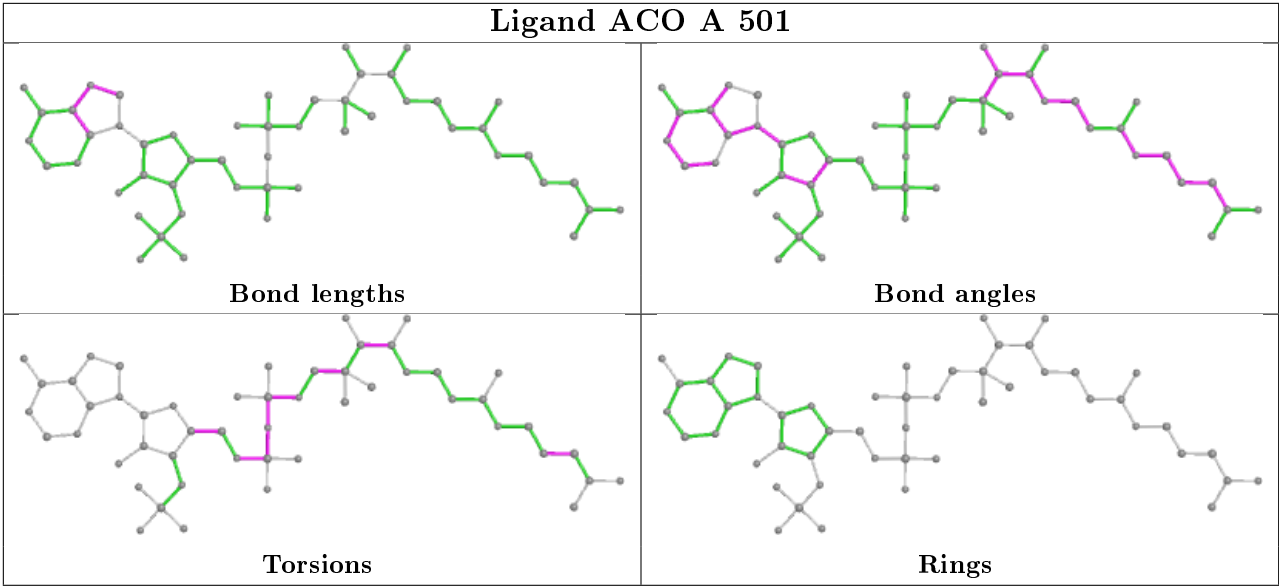
Mol	Chain	Res	Type	Atoms
4	A	501	ACO	N8P-C9P-CAP-OAP
4	A	501	ACO	C3P-C2P-S1P-C
4	A	501	ACO	O9P-C9P-CAP-OAP
5	A	502	EDO	O1-C1-C2-O2
4	A	501	ACO	O4B-C4B-C5B-O5B
4	A	501	ACO	C5B-O5B-P1A-O1A
4	A	501	ACO	C5B-O5B-P1A-O2A
4	A	501	ACO	CAP-CBP-CCP-O6A
4	A	501	ACO	CDP-CBP-CCP-O6A
4	A	501	ACO	CEP-CBP-CCP-O6A
4	A	501	ACO	P2A-O3A-P1A-O2A
4	A	501	ACO	C3B-C4B-C5B-O5B
4	A	501	ACO	P1A-O3A-P2A-O5A
4	A	501	ACO	CCP-O6A-P2A-O4A

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	502	EDO	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	14:LYS	C	15:NH2	N	1.06

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	393/438 (89%)	-0.24	3 (0%) 86 81	42, 78, 132, 162	0
2	B	213/232 (91%)	-0.21	1 (0%) 91 88	57, 84, 118, 158	0
3	D	4/15 (26%)	-0.01	0 100 100	62, 69, 103, 116	0
All	All	610/685 (89%)	-0.23	4 (0%) 87 84	42, 81, 130, 162	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	135	PRO	3.2
2	B	129	LYS	3.0
1	A	40	LYS	2.3
1	A	141	ALA	2.1

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
1	ALY	A	290	12/13	0.97	0.17	40,52,64,64	0

### 6.3 Carbohydrates

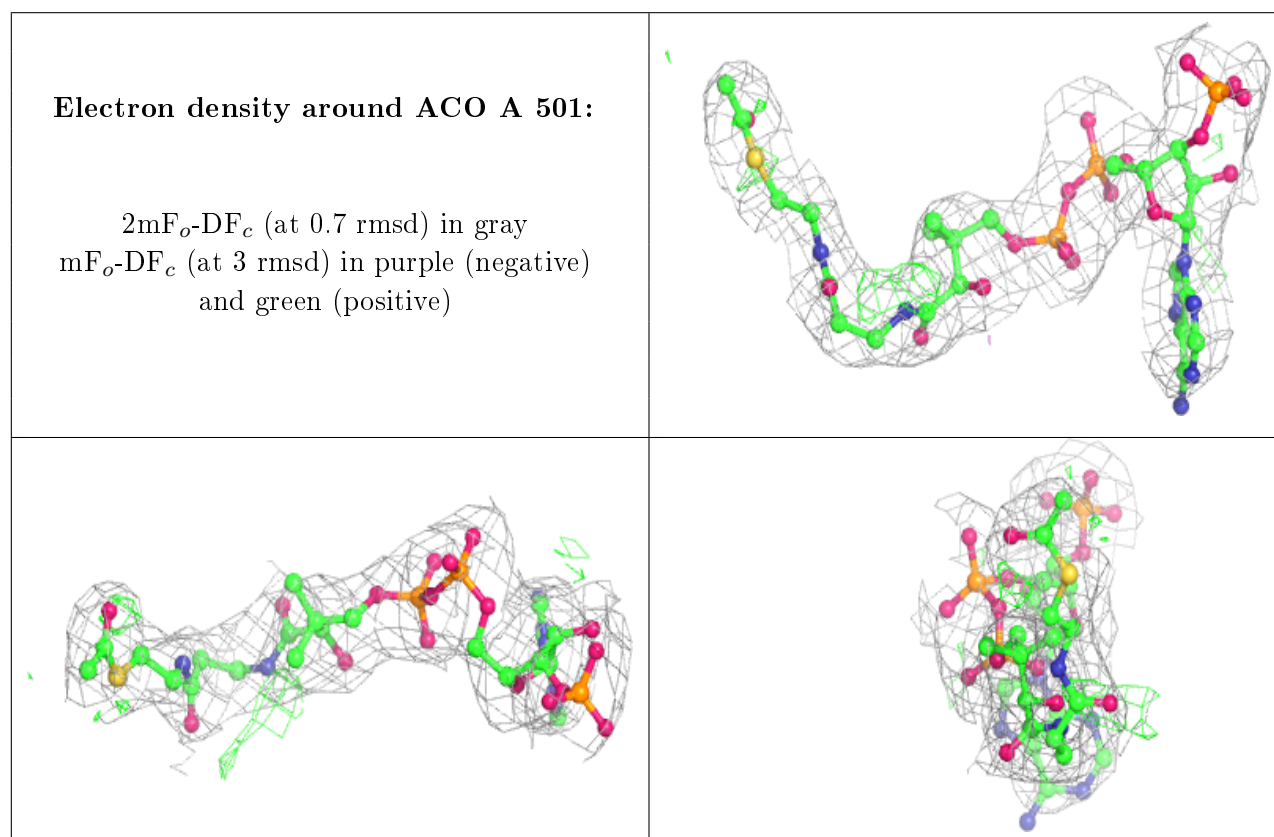
There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	A	502	4/4	0.91	0.16	72,76,77,78	0
4	ACO	A	501	51/51	0.96	0.16	47,79,119,121	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.