

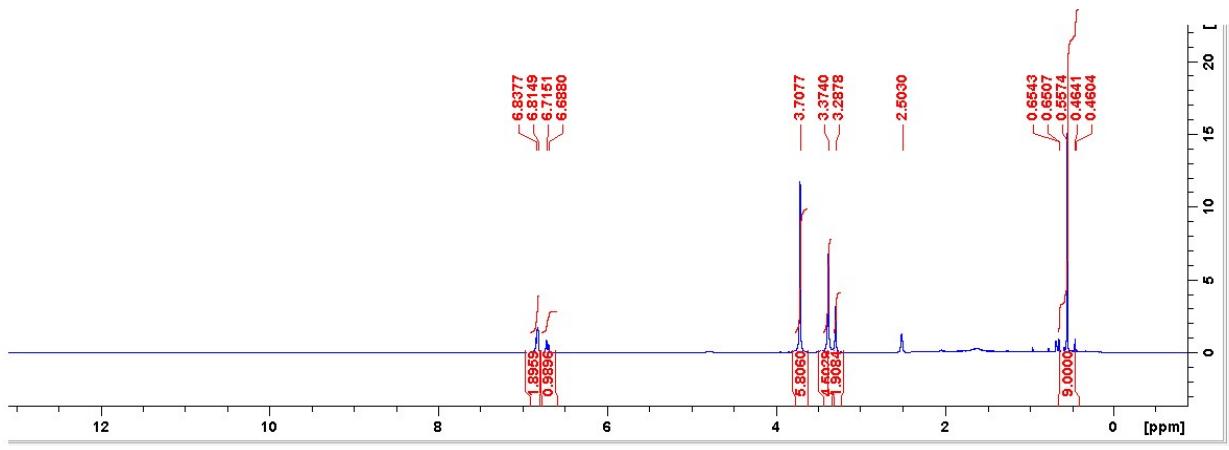
Supporting information

**New Triorganotin(IV) Compounds with Aromatic Carboxylate Ligands: Synthesis and
Evaluation of Pro-apoptotic Mechanism**

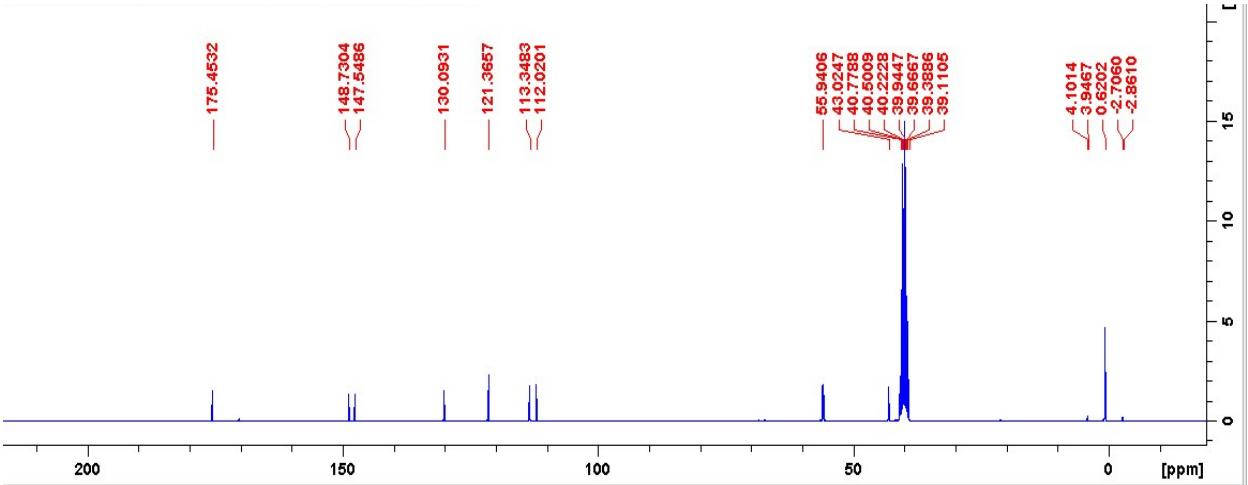
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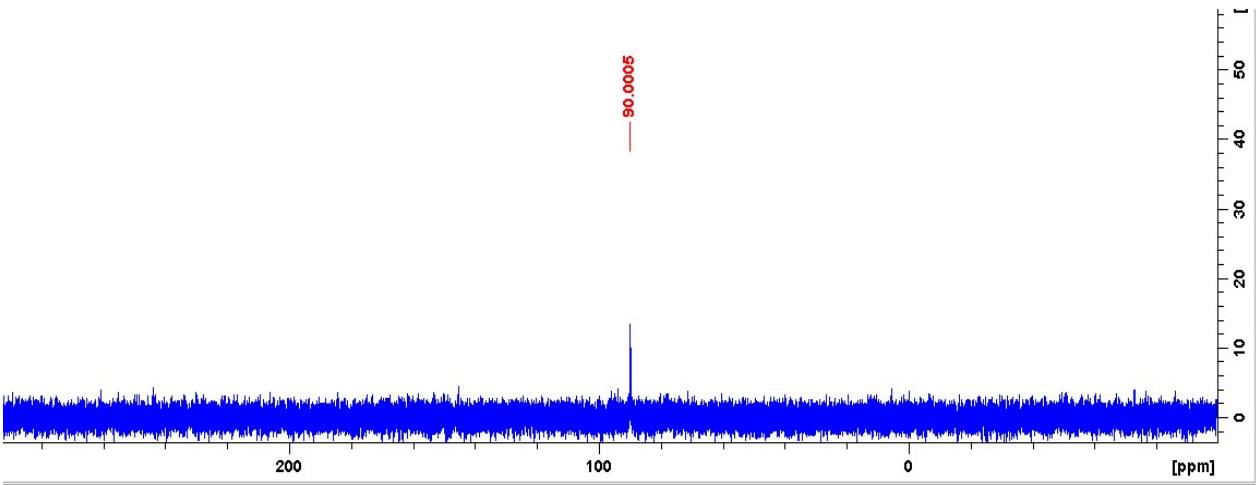
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(a)

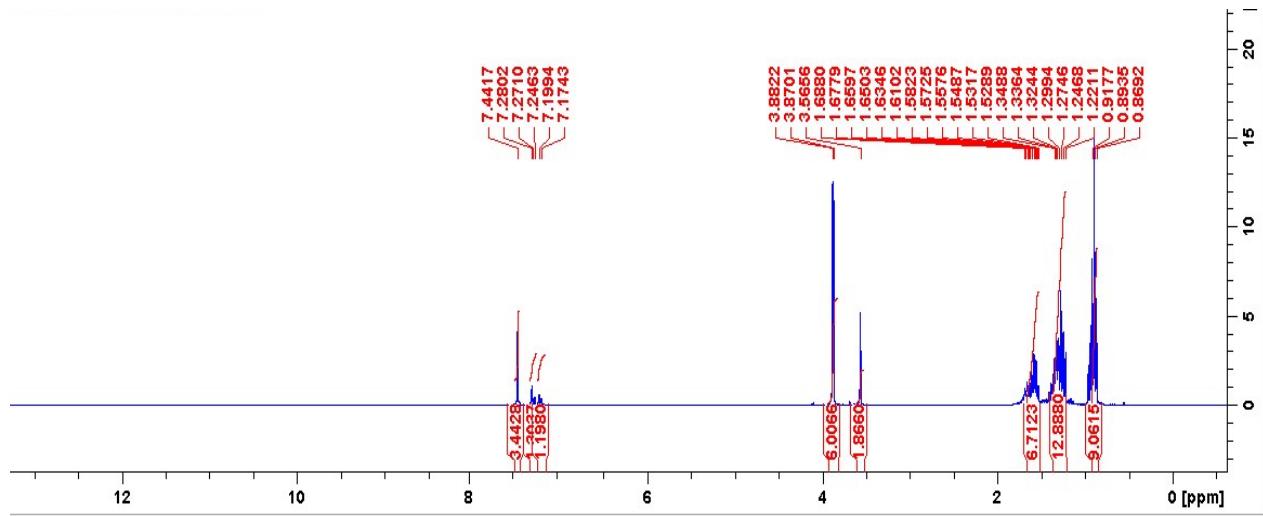


(b)

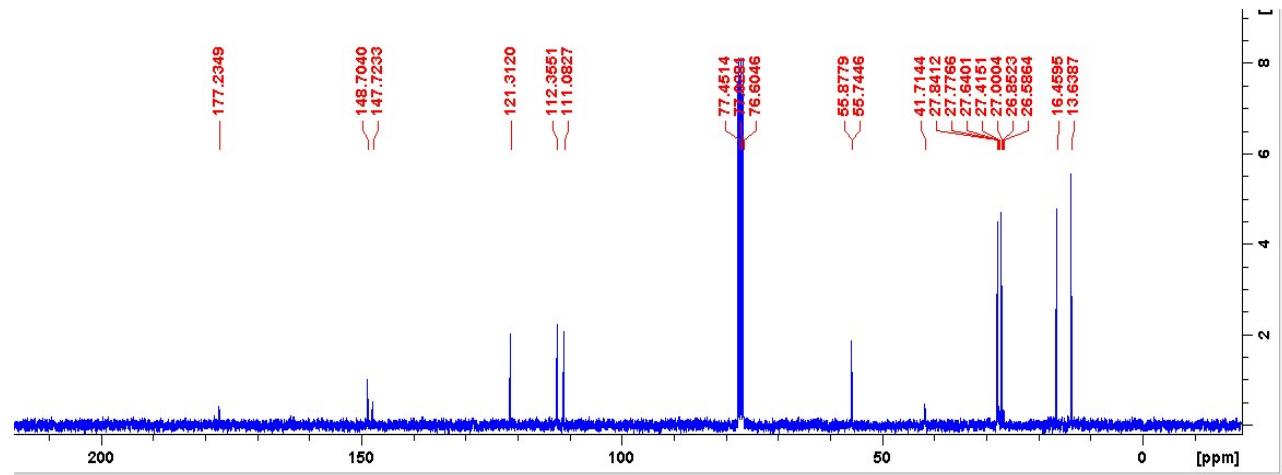


(c)

Fig. S1. Multinuclear (a) ^1H - , (b) ^{13}C - and (c) ^{119}Sn -NMR spectra of C1



(a)



(b)

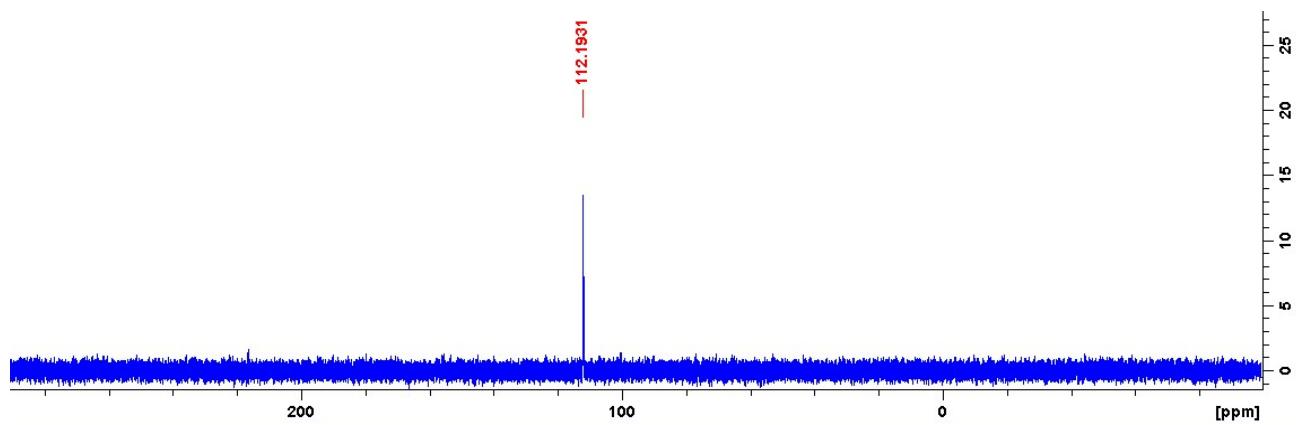


Fig. S2. Multinuclear (a) ^1H - , (b) ^{13}C - and (c) ^{119}Sn -NMR spectra of C2

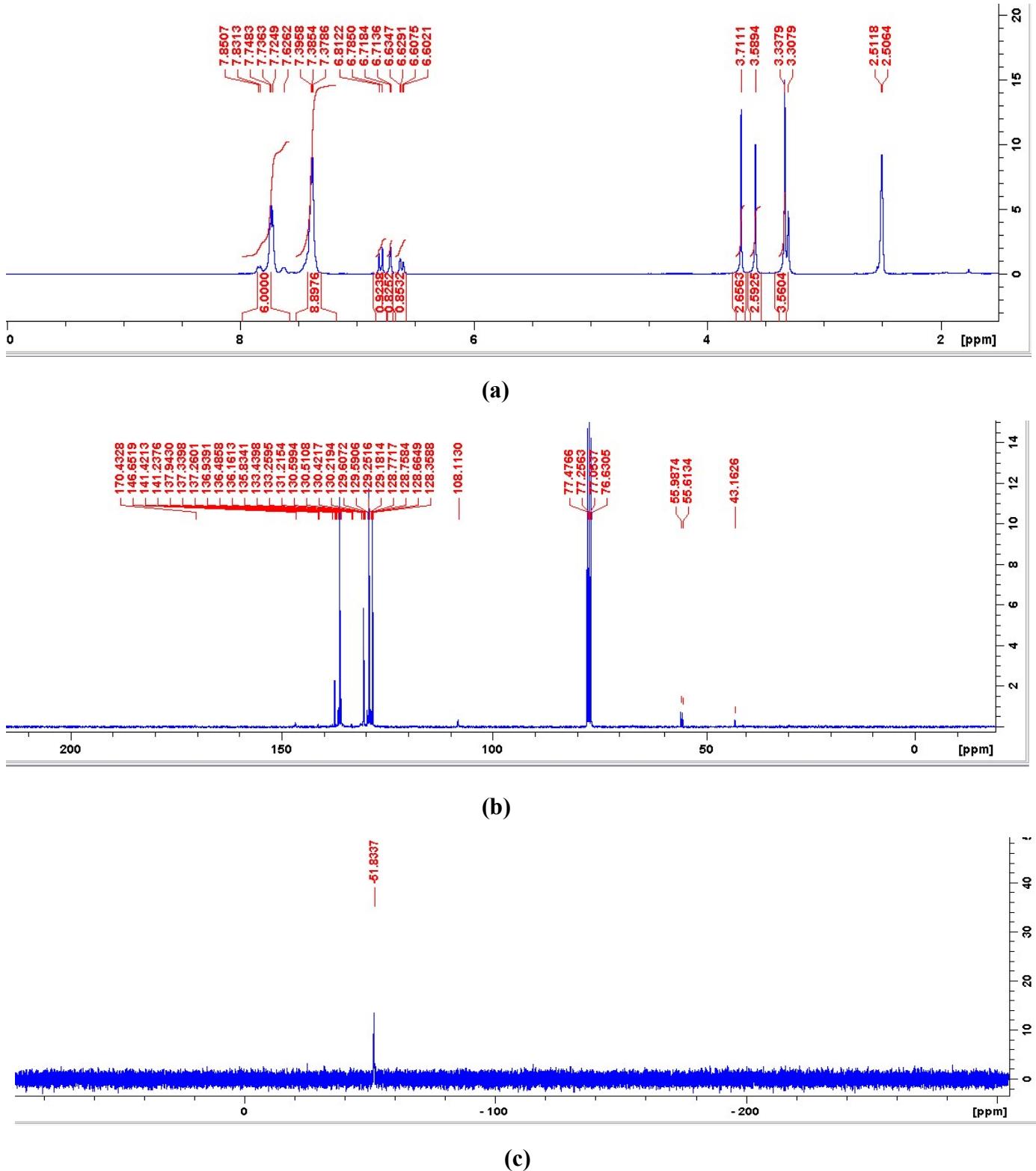


Fig. S3. Multinuclear (a) ^1H - , (b) ^{13}C - and (c) ^{119}Sn -NMR spectra of C3

Table S1. Crystal data and refinement parameter for complex **1**.

Crystal data	1
Empirical formula	C ₂₂ H ₃₈ O ₄ Sn
Formula weight	485.21
Temperature/K	100(2)
Crystal system	Monoclinic
Space group	P2 ₁ /c
a/Å	26.639(3)
b/Å	10.7495(11)
c/Å	16.3176(16)
α/°	90
β/°	98.1189(16)
γ/°	90
Volume/Å ³	4625.8(8)
Z	8
ρ _{calc} g/cm ³	1.393
μ/mm ⁻¹	1.127
F(000)	2016.0
Crystal size/mm ³	.3 × .2 × .1
Radiation	MoKα ($\lambda = 0.71073$)
2Θ range for data collection/°	3.088 to 61.814
Index ranges	-37 ≤ h ≤ 38, -15 ≤ k ≤ 15, -23 ≤ l ≤ 22
Reflections collected	60460
Independent reflections	13595 [R _{int} = 0.0280, R _{sigma} = 0.0257]
Data/restraints/parameters	13595/0/497
Goodness-of-fit on F ²	1.214
Final R indexes [I>=2σ (I)]	R ₁ = 0.0574, wR ₂ = 0.1271
Final R indexes [all data]	R ₁ = 0.0759, wR ₂ = 0.1396
Largest diff. peak/hole / e Å ⁻³	5.34/-1.90