



## **Supplementary Materials**

## Synergistical Use of Electrostatic and Hydrophobic Interactions for the Synthesis of a New Class of Multifunctional Nanohybrids: Plasmonic Magneto-Liposomes

Gabriela Fabiola Știufiuc <sup>1</sup>, Ștefan Nițică <sup>2</sup>, Valentin Toma <sup>2</sup>, Cristian Iacoviță <sup>3</sup>, Dietrich Zahn <sup>4</sup>, Romulus Tetean <sup>1</sup>, Emil Burzo <sup>1</sup>, Constantin Mihai Lucaciu <sup>3</sup> and Rareș Ionuț Știufiuc <sup>2,3,\*</sup>

- <sup>1</sup> Faculty of Physics, "Babeș-Bolyai" University, M. Kogălniceanu 1, 400084, Cluj-Napoca, Romania; gabi.stiufiuc@phys.ubbcluj.ro (G.Ș.); romulus.tetean@phys.ubbcluj.ro (R.T.); emil.burzo@phys.ubbcluj.ro (E.B.)
- <sup>2</sup> MedFuture Research Center for Advance Medicine, "Iuliu Hațieganu" University of Medicine and Pharmacy, L. Pasteur 4-6, 400349, Cluj-Napoca, Romania; stefan\_nitica@yahoo.com (Ş.N.); valentin.toma@umfcluj.ro (V.T.)
- <sup>3</sup> Faculty of Pharmacy, "Iuliu Hațieganu" University of Medicine and Pharmacy, L. Pasteur 4-6, 400349, Cluj-Napoca, Romania; cristian.iacovita@umfcluj.ro (C.I.); clucaciu@umfcluj.ro (C.L.)
- <sup>4</sup> Semiconductor Physics, Chemnitz University of Technology, D-09107 Chemnitz, Germany; zahn@physik.tu-chemnitz.de (D.Z.)
- \* Correspondence: rares.stiufiuc@umfcluj.ro



Figure S1. X-ray diffractogram acquired on SPIONs powder.



Figure S2. Zero Field Cooling and Field Cooling curves acquired on Fe<sub>3</sub>O<sub>4</sub> spherical magnetic nanoparticles.



Figure S3. Chemical structures of DOTAP and of soybean phosphatidylcholine (SPC) lipids.



**Figure S4.** Optical image of 5 batches of liposomal dispersions ordered from left to right with respect to the volume of SPIONs dispersion used in the synthesis: 1. 0  $\mu$ L (bare liposomes) 2. 10  $\mu$ L 3. 50  $\mu$ L 4. 250  $\mu$ L 5. 1000  $\mu$ L.



Figure S5. Energy Dispersive Spectroscopy (EDS) analysis of MLP250 sample.

Sample	Dilution	pН	ζ (mV)
AuPEG	1:16	10.1	-29.6
Lipo 0 µL	1:4	3.56	+59.7
Lipo 10 µL	1:4	3.45	+43.9
Lipo 50 µL	1:4	3.53	+43.4
Lipo 250 μL	1:4	3.99	+47.9
Lipo1000 µL	1:4	3.24	+51.1

Table S1. Zeta potential of nanoobjects.

Table S2. PCS measurements results for the liposomal dispersions.

Lincome hetek	Dilution	PDI	Hydrodynamic diameter (nm)			
Liposome batch			Peak <sub>1</sub>	Peak <sub>2</sub>		
Simple (0 µL)	-	0.295	180.9	31.27		
10 µL	-	0.431	179.8	36.31		
50 μL	-	0.342	227.5	33.55		
250 μL	1:4	0.460	393.9	71.60		
1000 μL	1:16	0.302	389.2	93.35		
PDI = polydispersity index						

Wavenumber (cm <sup>-1</sup> )	Assignment		
718	symmetric stretch vibration of the C-N bonds from the N <sup>+</sup> (CH <sub>3</sub> ) <sub>3</sub> choline group		
765	stretching vibration of the quaternary ammonium group & stretching vibration of the esterified P-O bonds from the phosphate groups		
850	Asymmetric stretching vibration of the esterified O-P-O bonds from the phosphate groups		
875	asymmetric stretch vibration of the C-N bonds from the N <sup>+</sup> (CH <sub>3</sub> ) <sub>3</sub> choline group		
972	bending vibration of the CH groups		
1044	symmetric stretching vibration of $P0_4^{3-}$ group		
1085	stretching vibration of C-C bond		
1265	deformation vibration of the unsaturated =CH group		
1302	twisting vibration of CH <sub>2</sub> groups		
1444	CH <sub>2</sub> /CH <sub>3</sub> scissoring vibrations		
1658	C=C bond stretching vibrations		
1741	stretching vibration of the C=O bond from the ester groups		
2728	in-plane scissoring vibrations of CH2		
2854	symmetric stretching vibration of the CH2 groups (fatty acids)		
2895	asymmetric stretching vibration of the CH2 groups (fatty acids)		
2929	symmetric stretching vibration of the CH <sub>3</sub> groups		
3010	stretching vibrations of unsaturated =C-H groups (fatty acids)		

Table S3. Assignment of the major vibrational bands recorded on cationic liposomes.

Table S4. Hyperthermia properties of PEGylated AuNPs and their liposomal complexes.

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Sample	<b>m</b> Au (g)	ΔΤ (Κ)	SAR (W/g)		
PEGylated AuNPs	4.115.10-5	2.61	327		
MMPNHs	4.113.10-5	3.08	450		