

Supplementary Information

Fluorescence Properties of Systems with Multiple Förster Transfer Pairs

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Table S-1: list of abbreviations

$I_{\text{transmitted}}$	transmitted light
I_{incident}	incident light
I_{detected}	the intensities of (fluorescence) light detected at the detector
d	length of the optical path through the sample
i, j	indices signifying individual dyes, e.g. dye 1, dye 2, etc.;
ϵ_i	molar absorption coefficient of dye i
c_i	molar concentration of the dye i .
λ	wavelength
$f_i(\lambda)$	fluorescence quantum yield of emission within a wavelength range of λ and $\lambda+d\lambda$.
$\lambda_\alpha, \lambda_\beta, \lambda_\gamma$	predefined wavelengths chosen by the experimentalist to characterize the fluorescent properties of his system.
$R_\alpha, R_\beta, R_\gamma$	intensity ratios chosen by the experimentalist
$F_i(\lambda)$	fluorescence efficiency; comprises $f_i(\lambda)$ and all those instrumentation parameters that will cancel out in Eq. (7)-(9).
p_{ij}	probability of Förster transfer of excitation from Donor i to acceptor j
$(1-p_{ij})$	probability of deactivation of the donor via either spontaneous emission or non-radiative decay
τ_D	lifetime of the excited state of the donor in absence of the acceptor,
$n_{D \rightarrow A}$	rate of transfer from donor to acceptor
r_{ij}	Förster radius, the distance between donor and acceptor at which the rate of transfer is equal to the rate of deactivation of the donor
c_{ij}	critical transfer concentration [molecules per volume] characteristic for donor-acceptor pair $i \rightarrow j$
K_i	rate constant for quenching (Stern Vollmer eq.)
p_{ii}	probability for self-quenching of the donor
P_i	probability (in presence of self quenching) of the excitation neither being transferred from an excited donor to an acceptor nor being self-quenched
P_{ij}	probability of energy transfer from the donor i to the acceptor j (in Presence of self quenching or other acceptors than j)
g_{ij}	probability that excitation transferred from donor i to acceptor j is emitted as light.
p_{jj}	probability of self-quenching of the acceptor.
FL	Fluorescein
ES	Eosin
RB	Rhodamine B
AOT	bis(2-ethylhexyl) sulfosuccinate sodium salt

Table S-2a:

R_a as a function of C₂ and C₃

(Datasets graphically represented in figure 11a)

Upper line: exponential representation

Middle line: Stern-Vollmer equation

Bottom line: experimental data

		C ₂ [10 ⁻⁵ mol L ⁻¹]						
		0	1	2	3	4	5	6
C ₃ [10 ⁻⁵ mol L ⁻¹]	0	0.748	0.662	0.593	0.536	0.487	0.446	0.410
		0.749	0.663	0.594	0.538	0.492	0.453	0.419
		0.752	0.663	0.602	0.543	0.490	0.457	0.423
	2	0.688	0.616	0.556	0.505	0.462	0.424	0.392
		0.693	0.621	0.563	0.514	0.473	0.437	0.406
		0.713	0.620	-	-	-	-	-
	4	0.641	0.579	0.526	0.481	0.441	0.407	0.377
		0.648	0.587	0.536	0.493	0.455	0.423	0.394
		0.674	-	0.543	-	-	-	-
	6	0.603	0.547	0.5	0.459	0.423	0.391	0.363
		0.610	0.556	0.512	0.473	0.439	0.409	0.383
		0.635	-	-	0.486	-	-	-
8	0.569	0.520	0.477	0.439	0.406	0.376	0.350	
	0.576	0.529	0.489	0.454	0.423	0.395	0.371	
	0.603	-	-	-	0.435	-	-	
10	0.539	0.494	0.455	0.420	0.390	0.362	0.338	
	0.547	0.505	0.468	0.436	0.407	0.382	0.359	
	0.568	-	-	-	-	0.399	-	
12	0.512	0.471	0.435	0.403	0.374	0.349	0.325	
	0.520	0.482	0.449	0.419	0.392	0.369	0.347	
	0.541	-	-	-	-	-	0.359	

Table S-2b:

R_β as a function of C_2 and C_3

(Datasets graphically represented in figure 11b)

Upper line: exponential representation

Middle line: Stern-Vollmer equation

Bottom line: experimental data

		C_2 [10^{-5} mol L $^{-1}$]						
		0	1	2	3	4	5	6
C_3 [10^{-5} mol L $^{-1}$]	0	0.208	0.275	0.329	0.373	0.411	0.443	0.471
		0.207	0.274	0.328	0.371	0.408	0.438	0.465
		0.206	0.274	0.325	0.368	0.410	0.440	0.466
	2	0.204	0.261	0.308	0.348	0.382	0.411	0.436
		0.200	0.256	0.303	0.341	0.373	0.401	0.426
		0.186	0.253					
	4	0.198	0.248	0.29	0.325	0.355	0.382	0.405
		0.194	0.242	0.282	0.316	0.346	0.371	0.393
		0.172		0.272				
	6	0.191	0.235	0.273	0.305	0.332	0.357	0.378
		0.188	0.23	0.266	0.296	0.322	0.345	0.365
		0.159			0.276			
8	0.185	0.224	0.258	0.287	0.312	0.334	0.354	
	0.184	0.221	0.252	0.279	0.302	0.323	0.341	
	0.147				0.281			
10	0.179	0.214	0.244	0.271	0.294	0.314	0.332	
	0.18	0.212	0.24	0.264	0.285	0.303	0.320	
	0.137					0.279		
12	0.174	0.205	0.232	0.256	0.277	0.296	0.312	
	0.176	0.205	0.229	0.251	0.270	0.286	0.302	
	0.128						0.270	

Table S-2c:

R_γ as a function of C₂ and C₃

(Datasets graphically represented in figure 11c)

Upper line: exponential representation

Middle line: Stern-Vollmer equation

Bottom line: experimental data

		C ₂ [10 ⁻⁵ mol L ⁻¹]						
		0	1	2	3	4	5	6
C ₃ [10 ⁻⁵ mol L ⁻¹]	0	0.044	0.063	0.078	0.09	0.102	0.111	0.120
		0.044	0.063	0.078	0.09	0.101	0.109	0.117
		0.042	0.063	0.073	0.089	0.100	0.103	0.116
	2	0.109	0.123	0.136	0.147	0.156	0.165	0.173
		0.108	0.122	0.134	0.145	0.154	0.161	0.168
		0.101	0.127	-	-	-	-	-
	4	0.161	0.173	0.184	0.194	0.203	0.211	0.218
		0.159	0.171	0.182	0.191	0.199	0.206	0.213
		0.155	-	0.185	-	-	-	-
	6	0.206	0.217	0.227	0.236	0.245	0.252	0.259
		0.202	0.213	0.223	0.231	0.239	0.246	0.252
		0.206	-	-	0.238	-	-	-
8	0.246	0.256	0.266	0.274	0.282	0.29	0.296	
	0.24	0.25	0.259	0.268	0.275	0.282	0.288	
	0.250	-	-	-	0.284	-	-	
10	0.282	0.291	0.3	0.309	0.317	0.324	0.331	
	0.273	0.283	0.292	0.3	0.308	0.315	0.321	
	0.295	-	-	-	-	0.322	-	
12	0.314	0.324	0.333	0.341	0.349	0.356	0.362	
	0.303	0.313	0.322	0.33	0.338	0.345	0.351	
	0.331	-	-	-	-	-	0.372	

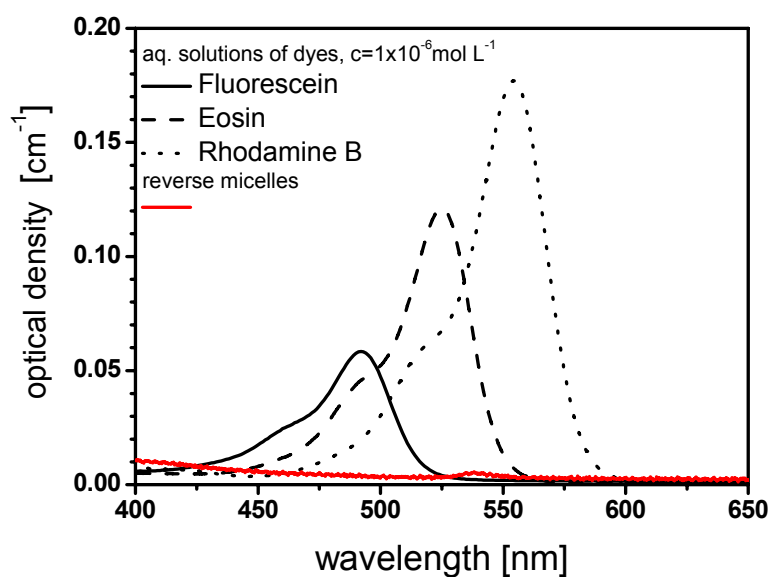


Figure S-1:

Comparison of the optical density [= $\lg(I_0/I)/(\text{path length})$] of diluted aqueous solution of the dyes used in the experiments to the optical density of a dispersion of reverse micelles in isooctane prepared at identical conditions as described in the main paper: in the experiments conducted here, attenuation due to scattering is negligible compared to attenuation due to absorption.