

# Researcher / Enterprise Profile form FP7 - Health



Date	2009	09	30	Valid until:	2009	12	31	

#### **CONTACT DETAILS**

Research organisation / Enterprise	Research Organization
Organisation Name	University of Malaya
Department	Chemistry
Address	Dept of Chemistry, University of Malaya,
City Country	Kuala Lumpur, Malaysia
www address	http://www.kimia.um.edu.my/

Researcher / Contact person	Noorsaadah Abd. Rahman		
Name / Surname	Noorsaadah <u>Abd. Rahman</u>		
Gender	□M		
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#### **ORGANISATION TYPE**

Research organisation type	☐ Resea☐ Comp	•	sation	Is your compa and Medium S Enterprise ( S Number of em	Sized ME* )?	□ YES	⊠ NO
Description of research activity:							
Former participan FP Europea project?		YES	□NO				
Project title / A	•	INTEGRATING NANOMATERIALS IN FORMULATIONS (InForm)  Research Activities					

- \* Your enterprise is an SME if:
- it is engaged in economic activity
- it has less than 250 employees
- it has either an annual turnover not exceeding €50M, or an balance sheet total not exceeding €43 M
- it is autonomous

For the definition of SMEs, look at:

http://ec.europa.eu/enterprise/enterprise\_policy/sme\_definition/index\_en.htm

#### **SMEs go Health**

Activities supporting SME participation in the health programme of the 7<sup>th</sup> Framework Programme www.smesgohealth.org



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### **EXPERTISE/COMMITMENT OFFERED**

Keywords specifying the expertise:		Drug discovery, molecular modelling, bioassay, organic synthesis					
Description of the expertise:		Our research involved searching for potential inhibitors for RNA virus, in particular the dengue virus (a single-stranded RNA virus). Different approaches ranging from bio-assay guided screening of natural product extracts and compounds from combinatorial library to molecular modeling and synthesis were used in searching and developing a potential lead inhibitor for the virus. In searching for a potential inhibitor for the viruses, in vitro antiviral activities were assessed using protease assay as well as virus inhibition assay employing HepG2 cells. We also perform maximum non-toxic dose experiments on the extracts and compounds to determine the relative toxicity of these compounds to the cells. The molecular modeling approaches for the designed of novel inhibitors in our group involve in silico screen of virtual library, docking as well as QSAR, often based on natural product or compounds that have shown some activities as inhibitor. Following the computational design of the compounds with good theoretical activities, the compounds were synthesised in the lab, either individually or combinatorially for further testing.					
Commitment offered							
		☐ Technology	☐ Dissemination ☐ Othe	r:			
				T			
Interested in		e-scale	Small or medium-     Small or medium-	☐ Targeted	Other (Marie		
participation in Project types		egrating aborative project	scale focused research collaborative project	to SMEs	Curie Actions, ERA-NET):		
	☐ Cod	ordination and	Network of	☐ Research for	1		
	Support Action		Excellence	the benefit of SMEs			
Call references	□ 4 <sup>th</sup>	call HEALTH					
2 10.0.0							

**SMEs go Health**Activities supporting SME participation in the health programme of the 7<sup>th</sup> Framework Programme <a href="https://www.smesgohealth.org">www.smesgohealth.org</a>



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**Main Research Topics :** For details and description of research topics, please visit: <a href="https://www.smesqohealth.org">www.smesqohealth.org</a> and <a href="http://cordis.europa.eu/fp7/">http://cordis.europa.eu/fp7/</a>

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FP7-HEALTH-2010-single-stage. HEALTH.2010.2.3.3-2 and -3 (mostly interested in Dengue fever), or 2.3.3-4.

<b>EXPECTATIONS</b>			
Term commitment	X Short (< 1 year)	Medium (1 to 3 years)	X Long (more than 3 years)
Expected results for your organisation:	Some compounds the State-of-the-art tech Human capacity dev		RNA virus inhibitor

I agree with the publication of my contact data:

#### PLEASE FILL IN THE PROFILE FORM AND RETURN IT TO:

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tegas@apre.it

bergonzi@apre.it

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