Request for LR-MS Sample Analysis UCSD Molecular Mass Spectrometry Facility					
User Name	Submission Date	User Phone	PI Name	PI Phone	Project #/ Task #
User E-mail Address:			Proposed Molecular Structure, Molecular Formula, and Exact Mass: (molecular structure is ESSENTIAL, hand- drawing or cut-and-paste from ChemDraw)		
5-digit/letter Sample ID Code			- drawing or cut-an	a-paste from Chem	Draw)
		CH ₃ CN n-Propanol CH ₂ Cl ₂			
Type of Sample Predissolved samples are preferred. Refer to MS facility website for sample preparation guides: (https://mmsf.ucsd.edu/sample/index.html/)					
Neat Solid	Neat Liquid				
Confidence:TentativeConfidentConfirmed by					
Purity:CrudeFairly PureVery Pure					
Solution: Dissolved in atmg/ml					
Special Handling: (check all that apply)			Proposed Molecular Formula:		
Acid SensitiveLight SensitiveHeat SensitiveAir SensitiveMoisture SensitiveToxicBiohazardsRadioisotopesOther (describe):			Expected Exact Mass:		
Additional Informa	ation: (any information	that is helpful)			
Analyst's Comments: (to be filled out by facility personnel)					
Analyst	Sample Run I	Name	Analysis Date	Hours \$_	Charges

Please acknowledge UCSD Molecular Mass Spectrometry Facility in your publications.