

Report # \_\_\_\_\_

**Company:** NuMega Resonance Labs

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**PO#** \_\_\_\_\_      **Date:** 2-25-20      **Results Format:** PDF  and FID

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Sample ID		NMR <sup>①</sup>		Mass <sup>②</sup>		Elemental Analysis <sup>③</sup>			
		H1, P31, C13 etc.	Solvent	Range to scan	M.W.	Single: <input type="checkbox"/> Double: <input checked="" type="checkbox"/> Theoretical %			
						C%	H%	N%	S%
A	A100					50	20	0	0
B	A101					30	5	10	20
C									
D									
E									
F									
G									
H									

- ① **NMR (PDF):** -0.5 to 10 ppm, integration, peak picking for H-1, 0-200 ppm, peak picking for C-13  
Chemical shift (H-1). TMS (0), DMSO (2.50), CDCl<sub>3</sub>(7.26), or D<sub>2</sub>O (4.78).
- ② **Mass:** 1mg in 1mL or 20-50 μM, positive and negative modes. Please specify MW range.
- ③ **Elemental Analysis (3 mg per run):** Duplicate runs are recommended

**Your Instructions/Remarks:** (Hygroscopic, Explosive, Volatile, Toxicity, Storage, EA theoretical %, etc.)

**Confidentiality:** NuMega maintains spectral information in strict confidence

Comment from Analyst: \_\_\_\_\_ **NuMega Use Only**

Routine NMR	__ x \$46 =	
Non-Routine	__ x ____ =	
Others	__ x ____ =	

MASS	__ x \$35 =	

CHN (S/D)	__ x \$29/45 =	
S (S/D)	__ x \$29/45 =	

Analyst: \_\_\_\_\_ Date: \_\_\_\_\_ Total Charge: \$ \_\_\_\_\_