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Mass Spectrometry Request

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NAME:	PHONE:	DATE:
PROFESSOR:	SAMPLE ID:	SAMPLE BOX: Position:
*FREE BASE FORMULA:	*FREE BASE .MONO.MASS:	REQUEST:

PROPOSED STRUCTURE (INCLUDE SALT FORM IF APPROPRIATE)

SAVE STRUCTURE FROM CHEMDRAW AS A .JPEG AND OPEN IT BELOW. TURN ON FORMULA AND EXACT MASS.

SAFETY INFORMATION:

POTENTIAL CARCINOGEN:

LACHRYMATOR:

STENCH:

OTHER (EXPLAIN):

RECOMMENDED SOLVENTS:

	USE	DO NOT USE
METHANOL	<input type="checkbox"/>	<input type="checkbox"/>
ACETONITRILE	<input type="checkbox"/>	<input type="checkbox"/>
WATER	<input type="checkbox"/>	<input type="checkbox"/>
HEXANE	<input type="checkbox"/>	<input type="checkbox"/>
ACETONE	<input type="checkbox"/>	<input type="checkbox"/>

Other:

MS LAB USE ONLY:

INSTRUMENT USED

AGILENT GCMS
WATERS LCMS
BRUKER ULTRAFLEX III
WATERS GCT PREMIER
WATERS LCT PREMIER

OBSERVED MASS:

NOTES:

THEOR. MASS:

ANALYST: _____