

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090722\
 Data File : BF130159.D
 Acq On : 07 Sep 2022 13:28
 Operator : CG\JU
 Sample : N4446-06
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-06

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF130159.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.887	438	442	456	rVB	278452	338481	4.76%	0.898%
2	5.293	507	511	515	rBV	2726202	2812631	39.57%	7.463%
3	6.316	681	685	689	rBV	1828640	1762324	24.79%	4.676%
4	6.692	746	749	753	rVB	1493124	1241724	17.47%	3.295%
5	7.263	840	846	849	rBV	3611999	3947091	55.53%	10.473%
6	7.975	962	967	970	rBV	2006933	1712356	24.09%	4.544%
7	9.057	1145	1151	1154	rBV	6973394	7108429	100.00%	18.861%
8	9.151	1162	1167	1187	rVB	171588	237832	3.35%	0.631%
9	9.728	1259	1265	1268	rBV	2170266	2087324	29.36%	5.538%
10	10.516	1392	1399	1402	rBV	4980722	5065837	71.27%	13.442%
11	11.204	1512	1516	1519	rBV	2520755	2036214	28.65%	5.403%
12	11.739	1602	1607	1611	rBV	81180	77277	1.09%	0.205%
13	12.798	1781	1787	1790	rBV	6832773	6922961	97.39%	18.369%
14	13.833	1959	1963	1966	rBV	1470666	1312541	18.46%	3.483%
15	15.233	2196	2201	2206	rBV	902423	1024882	14.42%	2.719%

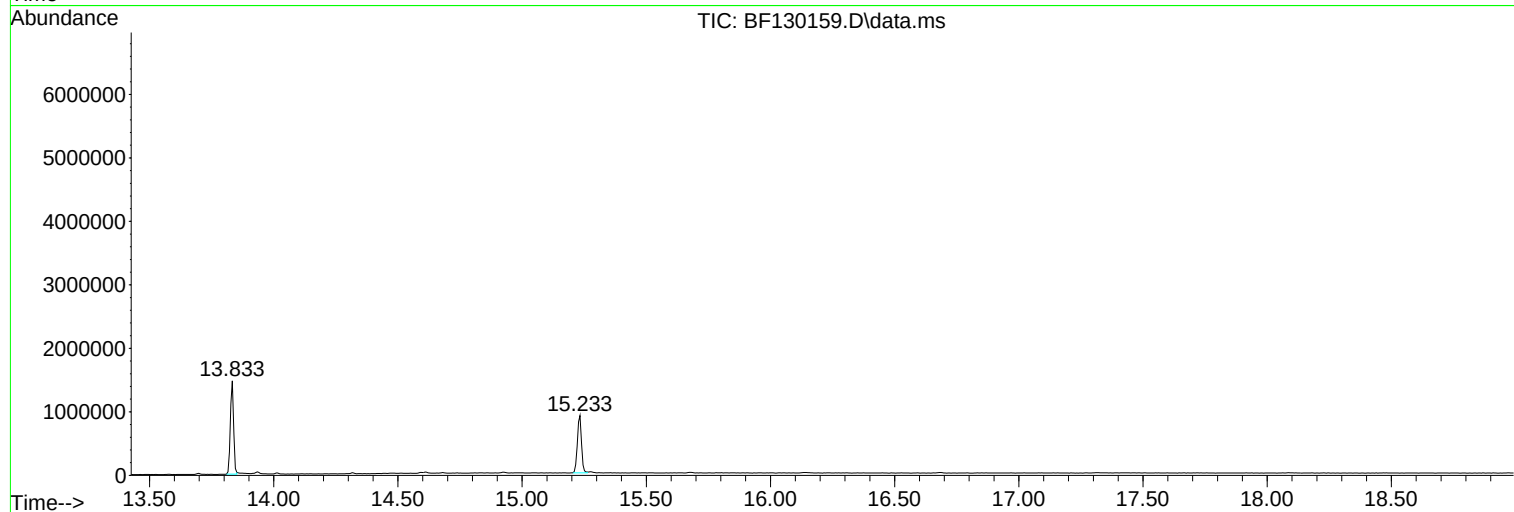
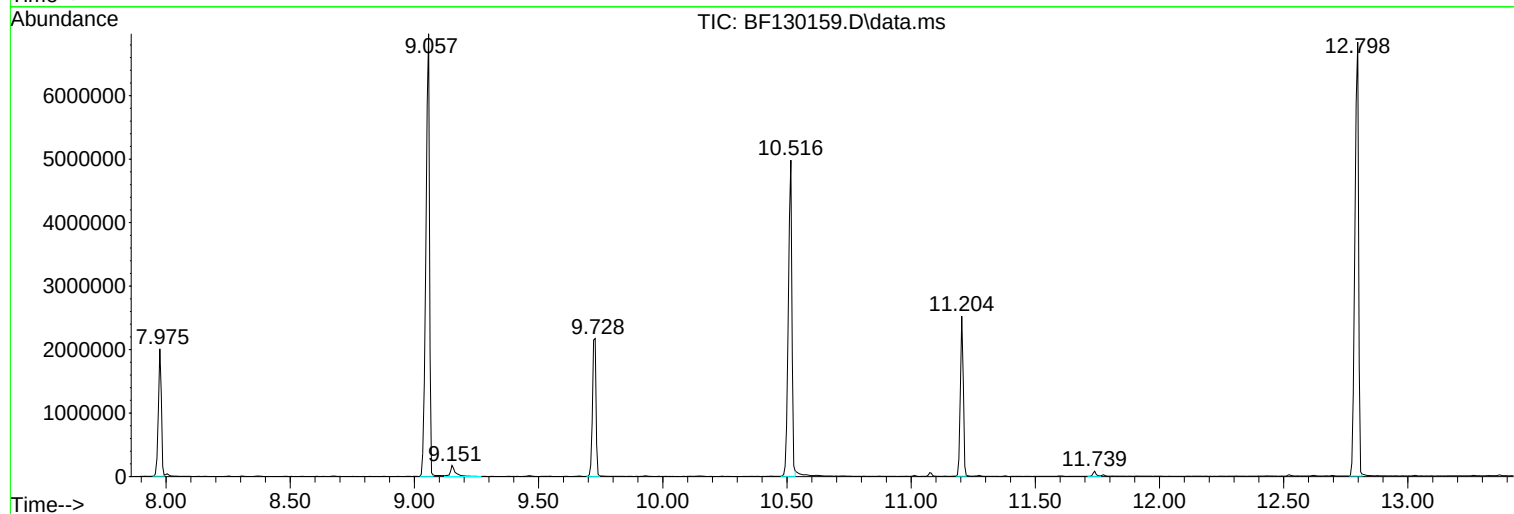
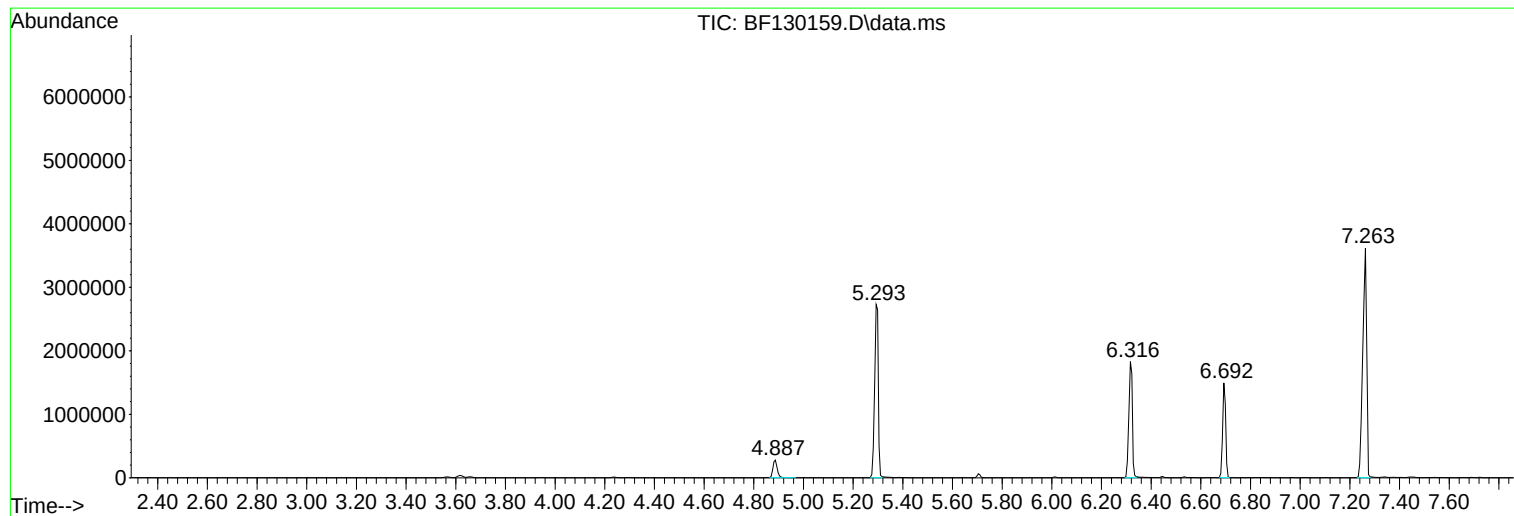
Sum of corrected areas: 37687904

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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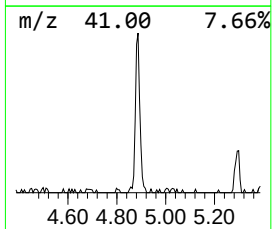
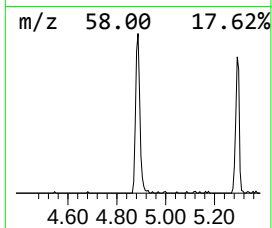
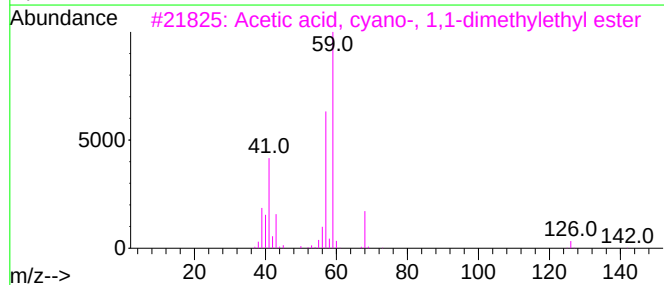
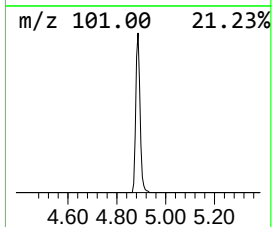
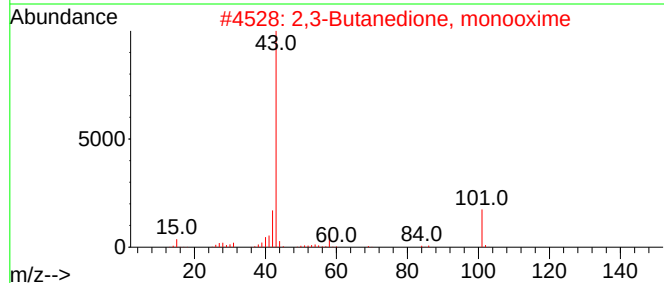
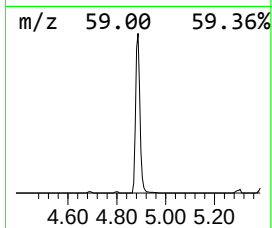
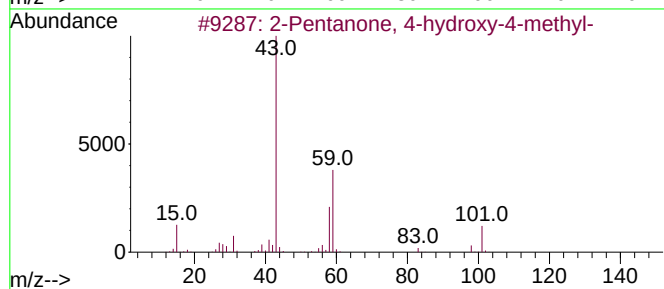
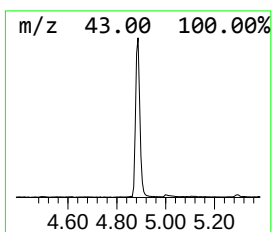
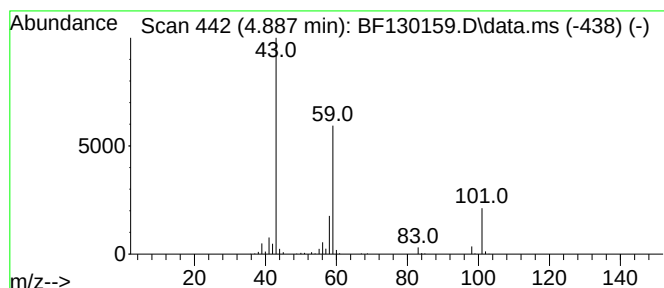
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.887	5.45 ng	338481	1,4-Dichlorobenzene-d4	6.692

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50		
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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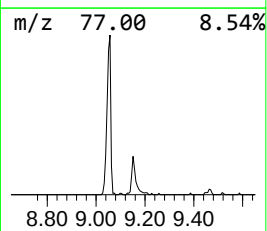
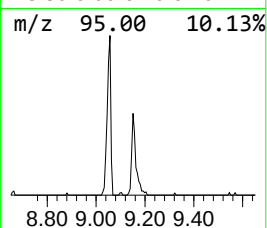
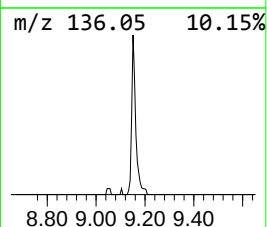
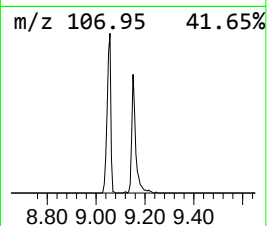
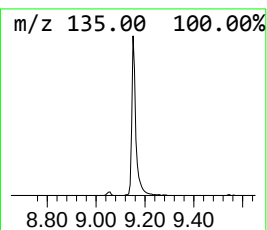
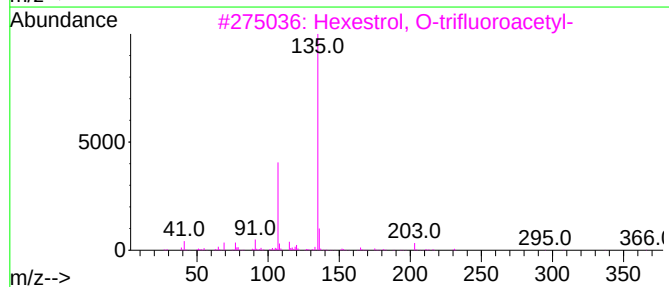
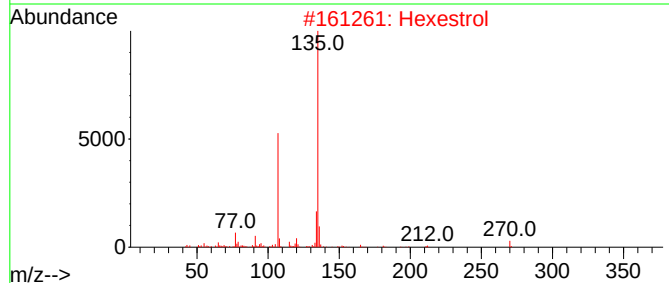
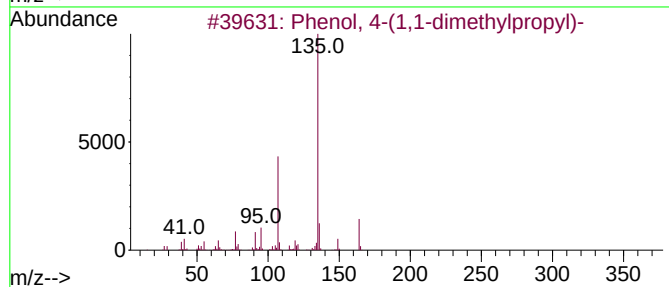
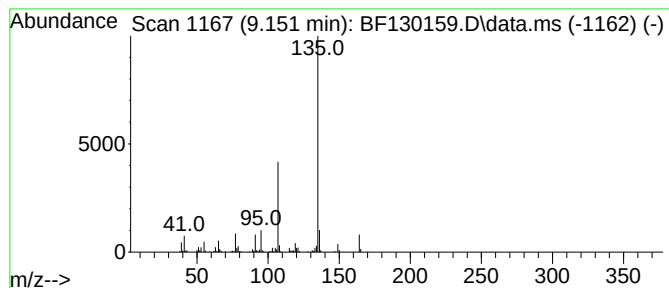
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Peak Number 2 Phenol, 4-(1,1-dimethylprop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.151	2.28 ng	237832	Acenaphthene-d10	9.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	94
2			Hexestrol	270	C18H22O2	000084-16-2	72
3			Hexestrol, O-trifluoroacetyl-	366	C20H21F3O3	1000365-31-7	72
4			Imidazo[1,2-c]pyrimidin-5-ol	135	C6H5N3O	055662-66-3	64
5			4,6-Bis(4-ethoxybenzylthio)-5-ni...	457	C22H23N3O4S2	325957-76-4	64



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.887	5.5	ng	338481	1	6.692	1241720	20.0
Phenol, 4-(1,1-...	9.151	2.3	ng	237832	3	9.728	2087320	20.0