

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090222\
 Data File : BF130108.D
 Acq On : 02 Sep 2022 13:31
 Operator : CG\JU
 Sample : N4446-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WELL-01

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: BF130108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.892	439	443	451	rVB	192381	219595	5.49%	0.875%
2	5.298	508	512	516	rBV	2208012	2025861	50.69%	8.074%
3	6.322	682	686	689	rBV	1901801	1577766	39.48%	6.288%
4	6.704	747	751	754	rBV	1422725	1189824	29.77%	4.742%
5	7.263	841	846	849	rBV	2257305	2072564	51.86%	8.260%
6	7.981	964	968	972	rBV	1773680	1585196	39.66%	6.318%
7	9.057	1147	1151	1155	rBV	4290445	3996723	100.00%	15.929%
8	9.733	1261	1266	1269	rBV	2247179	1778125	44.49%	7.087%
9	10.516	1395	1399	1416	rBV	2712104	2500938	62.57%	9.967%
10	11.216	1514	1518	1521	rBV	1913353	1721567	43.07%	6.861%
11	12.798	1783	1787	1791	rBV	3815759	3918141	98.03%	15.615%
12	13.845	1960	1965	1968	rBV	1377983	1288604	32.24%	5.136%
13	13.945	1977	1982	1989	rBV	157917	164884	4.13%	0.657%
14	15.245	2198	2203	2214	rBV	927592	1051746	26.32%	4.192%

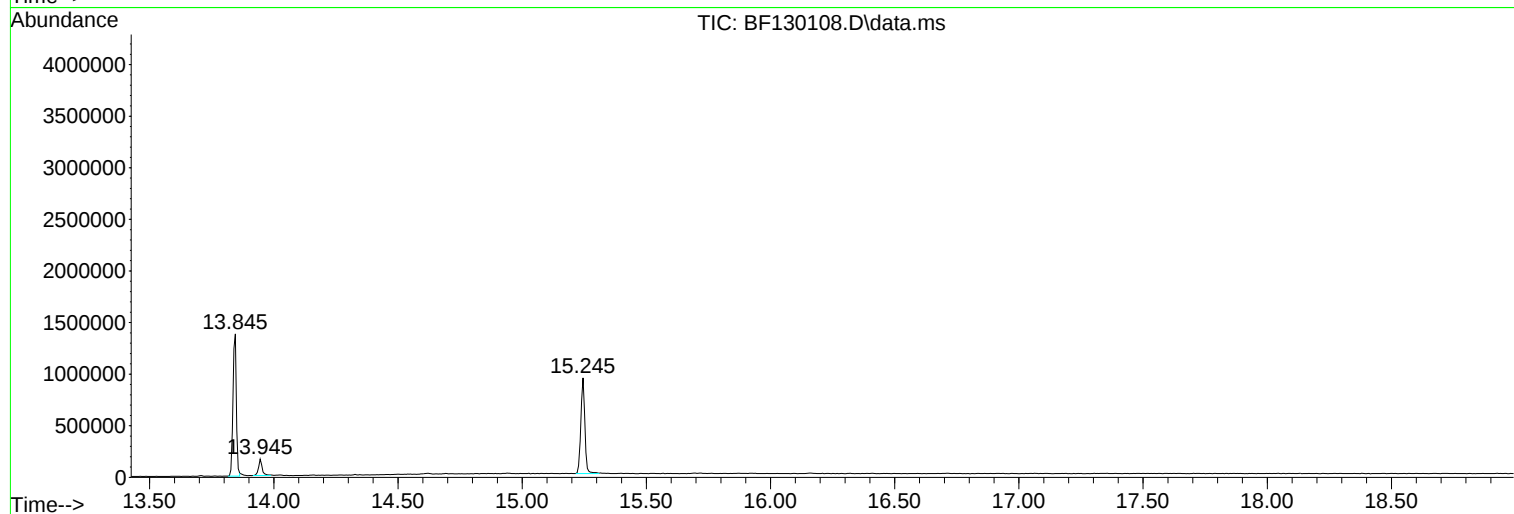
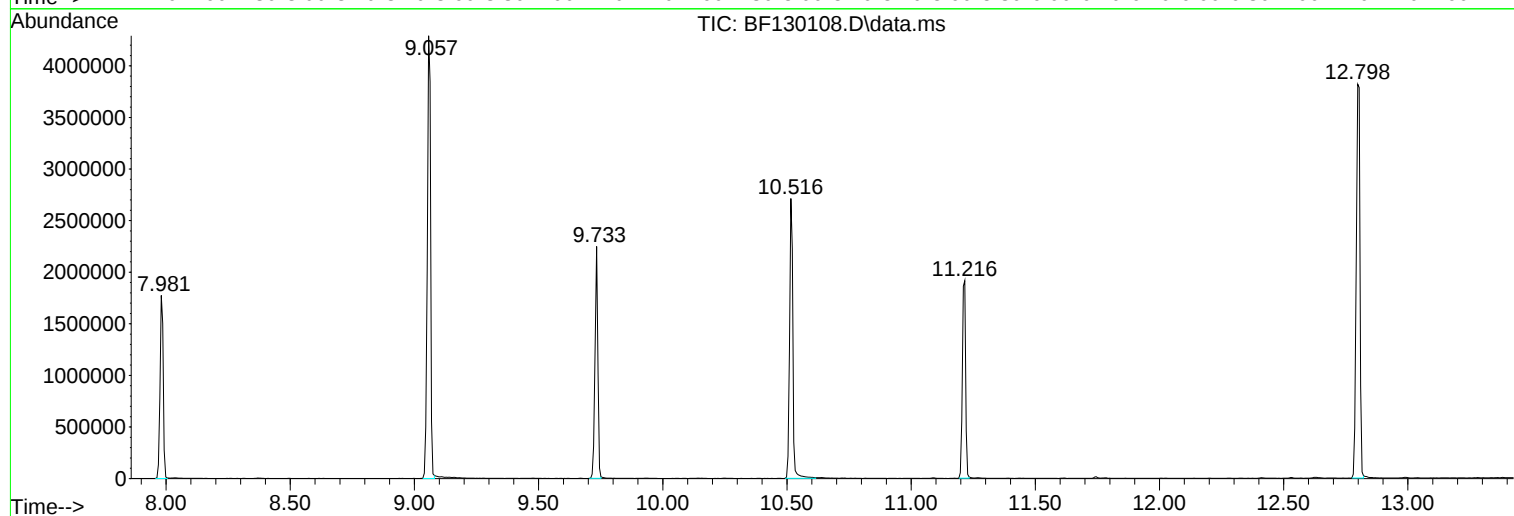
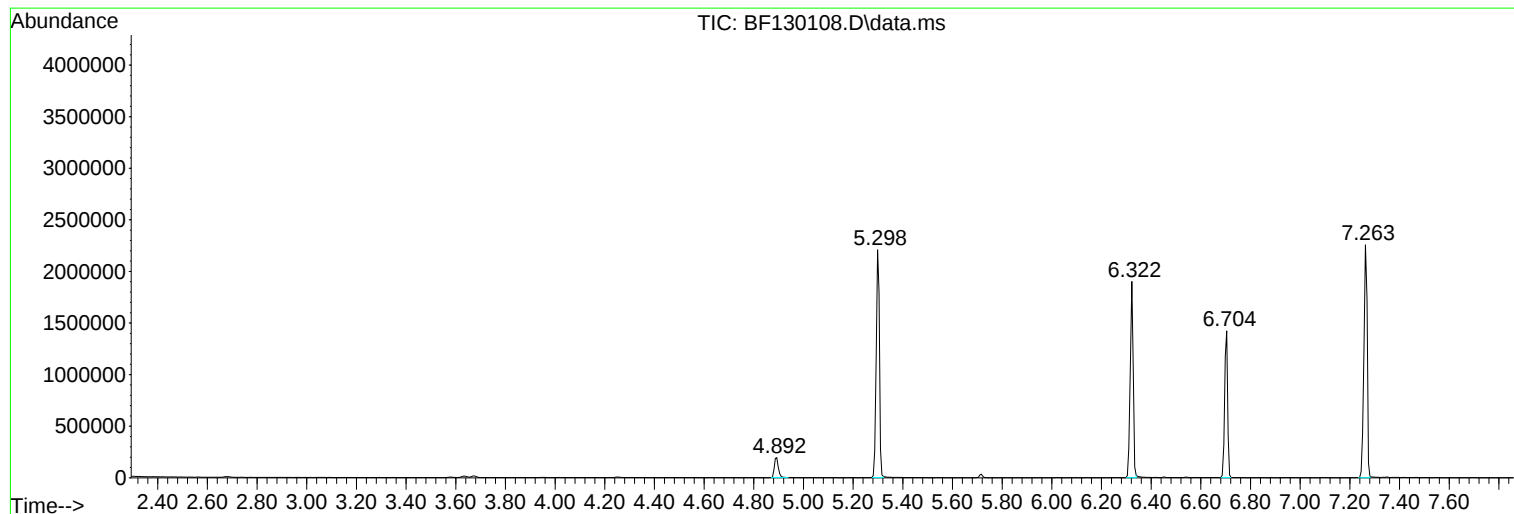
Sum of corrected areas: 25091534

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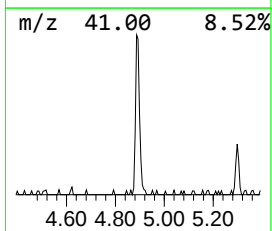
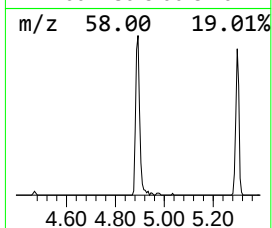
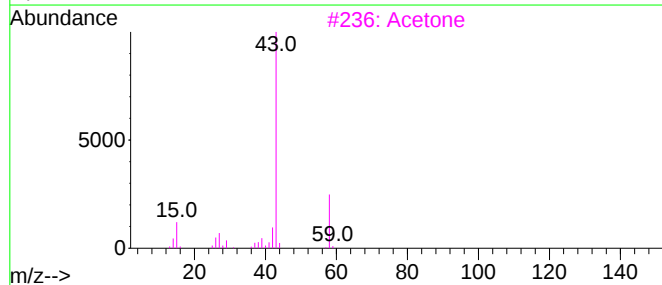
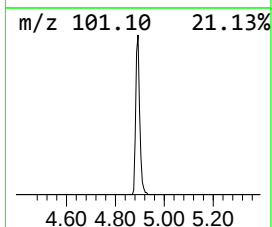
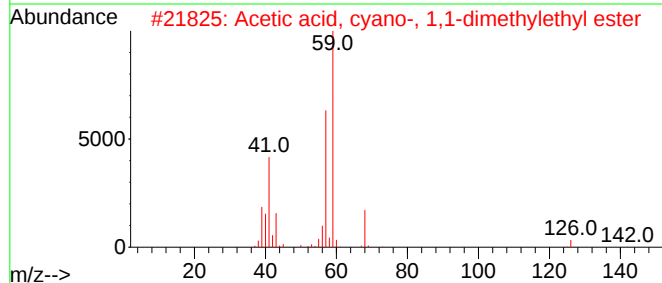
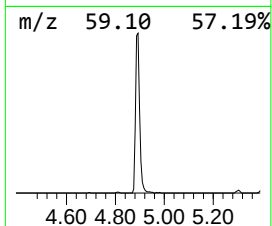
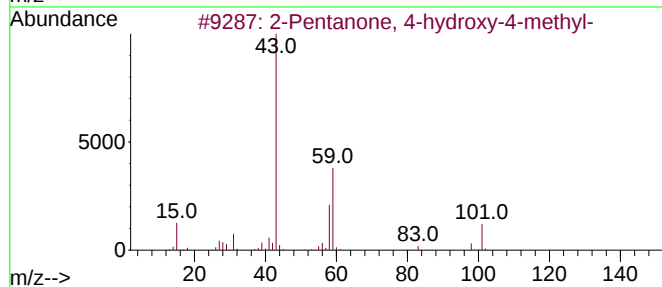
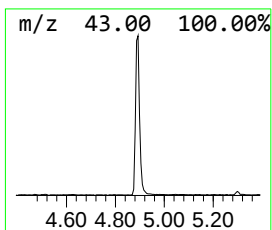
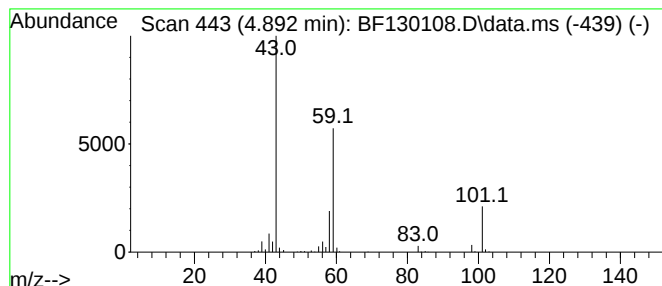
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.892	3.69 ng	219595	1,4-Dichlorobenzene-d4	6.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
3			Acetone	58	C3H6O	000067-64-1	9
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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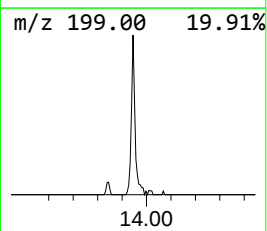
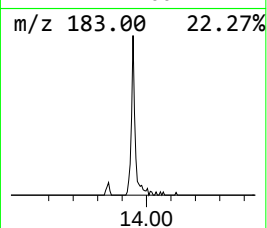
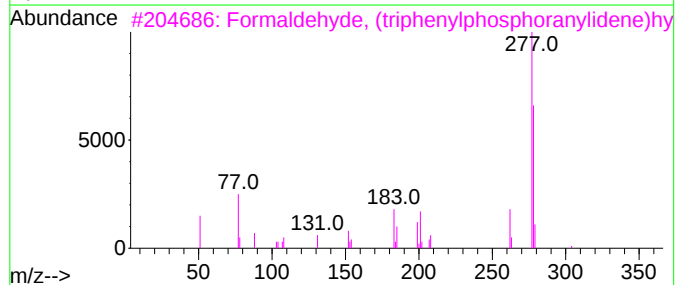
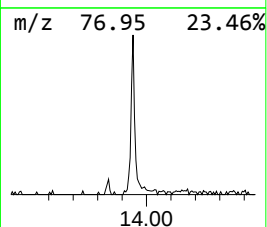
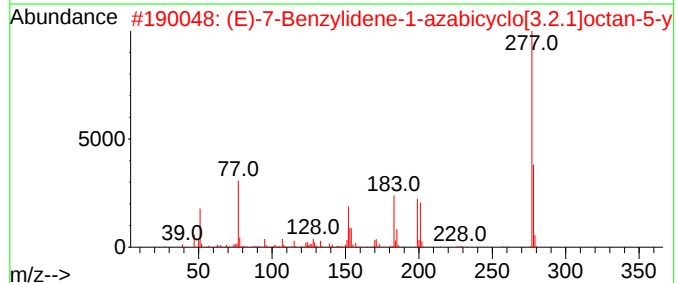
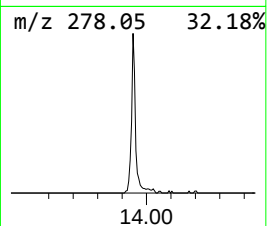
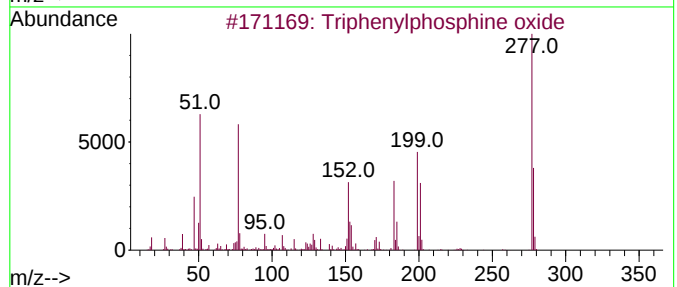
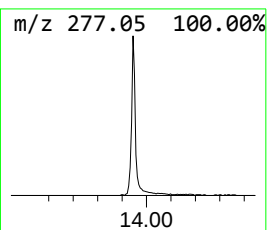
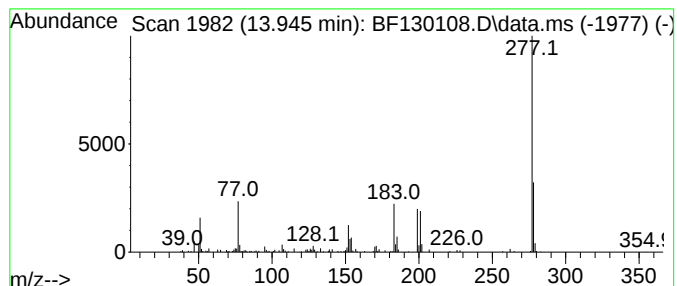
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Peak Number 2 Triphenylphosphine oxide Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.945	2.56 ng	164884	Chrysene-d12	13.845

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triphenylphosphine oxide	278	C18H15OP	000791-28-6	95
2			(E)-7-Benzylidene-1-azabicyclo[3...]	293	C15H19N03S	1000483-83-4	81
3			Formaldehyde, (triphenylphosphor...	304	C19H17N2P	015990-54-2	50
4			2-Phenyl-6-(trifluoromethyl)-1H-...	277	C15H10F3NO	1000387-59-3	43
5			Vanadium, cycloheptatrienyl-(pen...	277	C17H22V	1000155-20-5	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-...	4.892	3.7	ng	219595	1	6.704	1189820	20.0
Triphenylphosph...	13.945	2.6	ng	164884	5	13.845	1288600	20.0