

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092022\
 Data File : BP011760.D
 Acq On : 20 Sep 2022 13:26
 Operator : CG/JU
 Sample : PB147737BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147737BL

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_P\Methods\8270E-BP091422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP011760.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.075	11	15	19	rVB	428574	465876	1.51%	0.304%
2	5.034	343	348	358	rVB	1208951	1743592	5.66%	1.136%
3	5.493	419	426	442	rBV	12826905	19325274	62.79%	12.592%
4	7.099	690	699	718	rBV	11531050	19230575	62.48%	12.530%
5	7.934	834	841	849	rVB	2060969	3308000	10.75%	2.155%
6	9.099	1031	1039	1056	rBV	7465612	12644128	41.08%	8.239%
7	10.746	1311	1319	1333	rVB	2664294	4610682	14.98%	3.004%
8	13.198	1729	1736	1751	rBV	16498198	24637157	80.05%	16.053%
9	14.575	1961	1970	1983	rBV	3714752	5543506	18.01%	3.612%
10	16.063	2215	2223	2243	rBV	11235400	16310699	52.99%	10.628%
11	17.328	2431	2438	2450	rBV	4188721	5930233	19.27%	3.864%
12	19.880	2866	2872	2877	rBV	25010045	30779117	100.00%	20.055%
13	21.404	3126	3131	3138	rBV	3810988	4972123	16.15%	3.240%
14	23.857	3542	3548	3560	rVB2	1911482	3971101	12.90%	2.588%

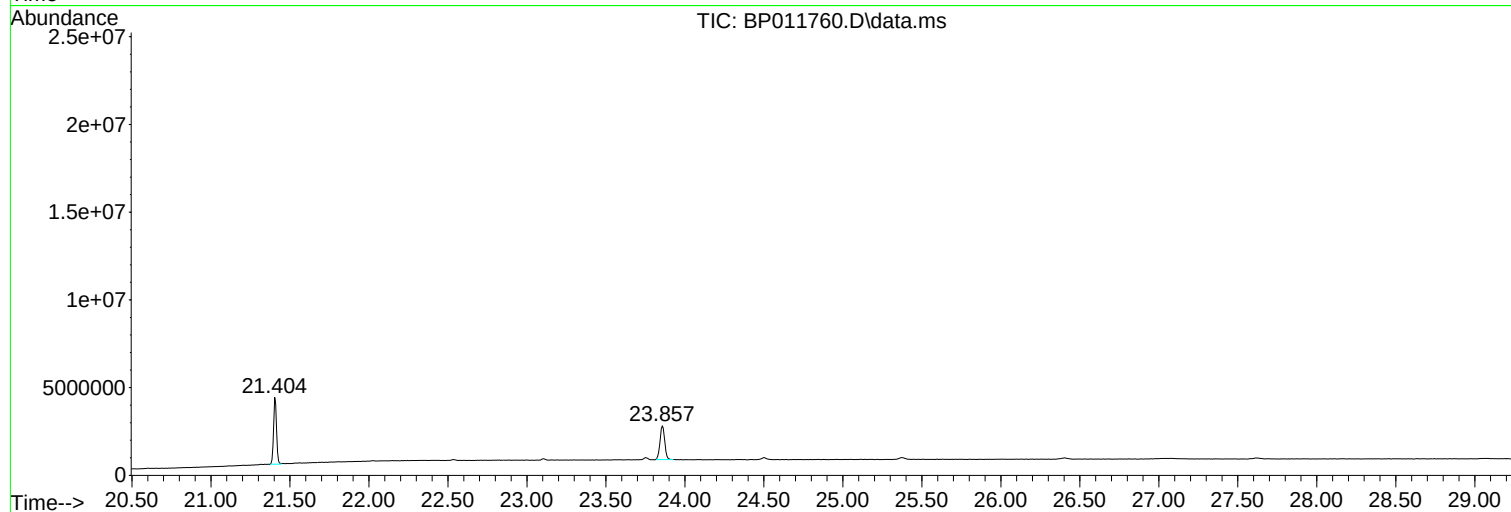
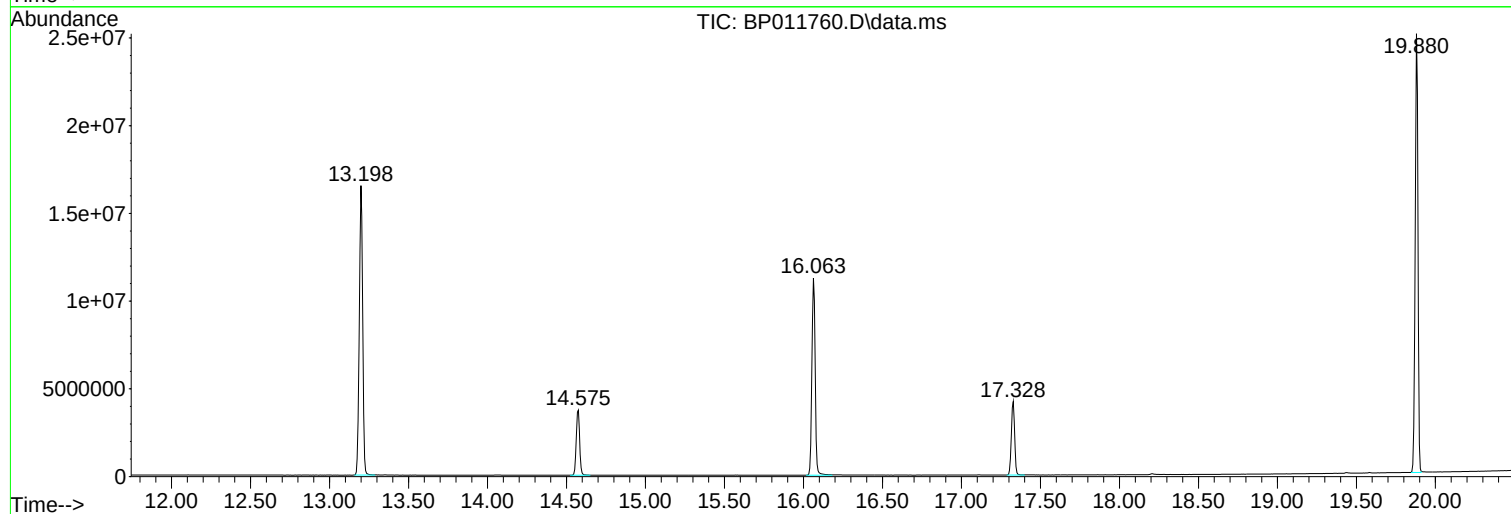
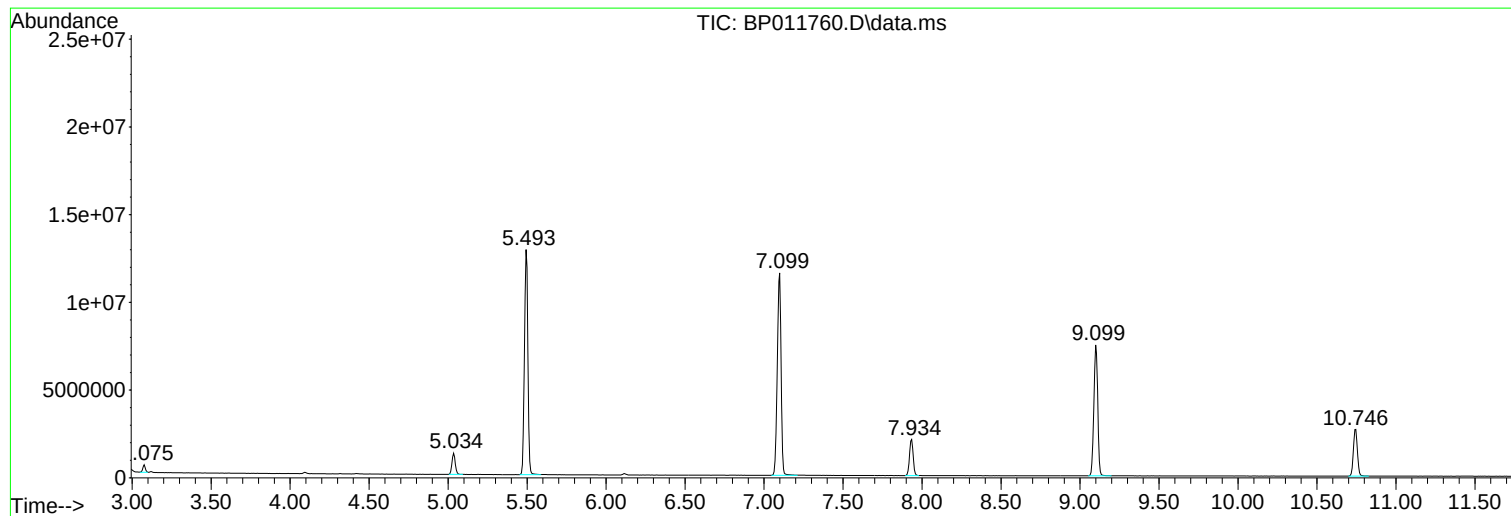
Sum of corrected areas: 153472063

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.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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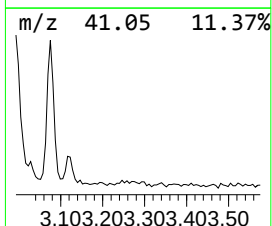
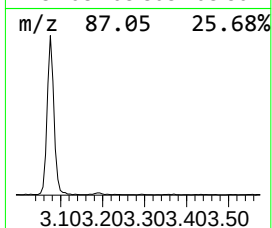
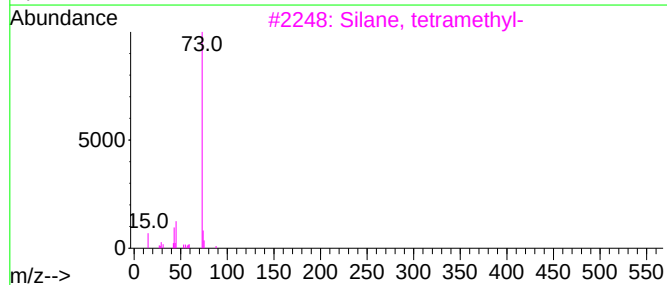
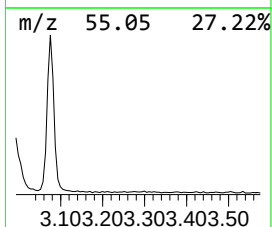
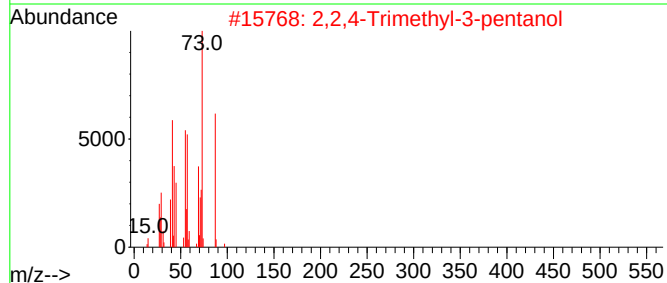
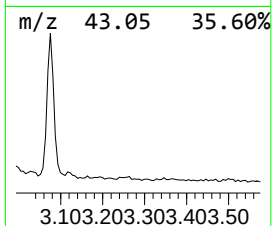
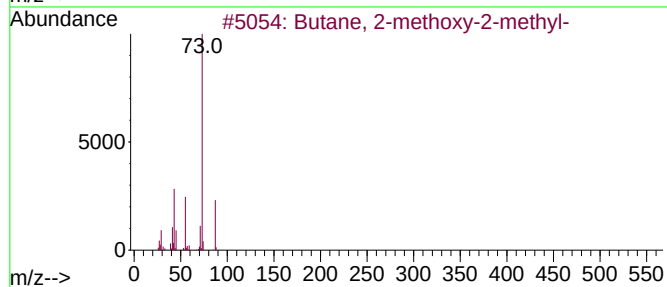
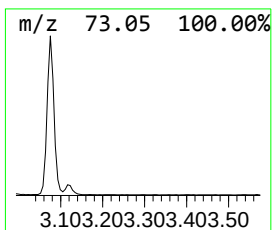
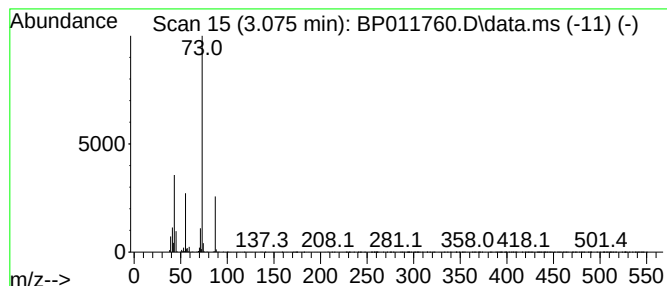
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	2.82 ng	465876	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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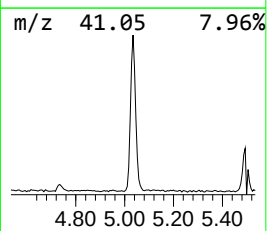
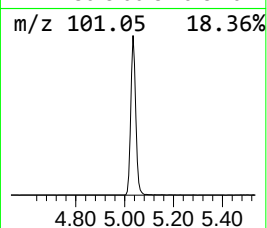
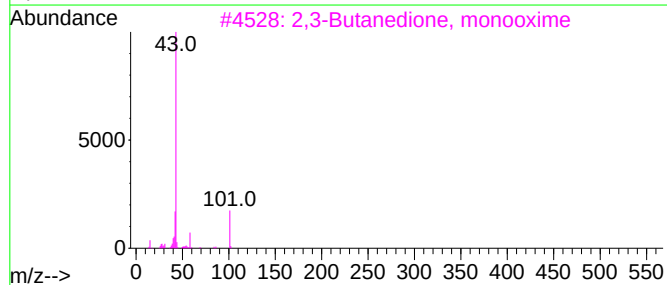
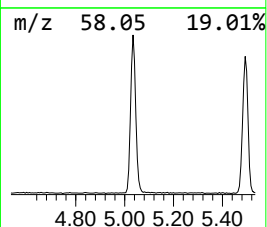
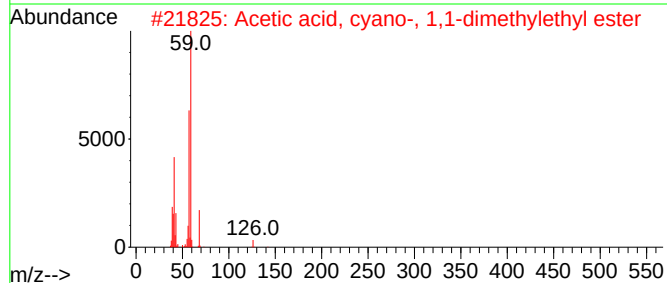
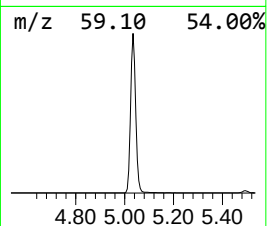
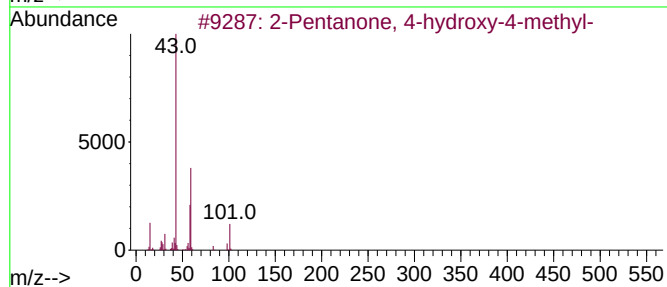
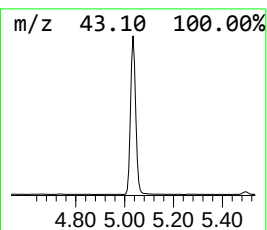
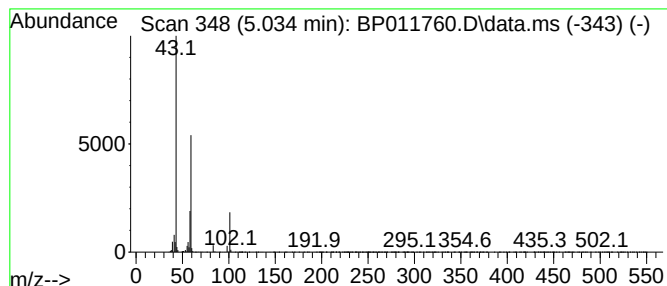
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.034	10.54 ng	1743590	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.075	2.8	ng	465876	1	7.934	3308000	20.0
2-Pentanone, 4-...	5.034	10.5	ng	1743590	1	7.934	3308000	20.0