

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092622\
 Data File : BP011797.D
 Acq On : 26 Sep 2022 15:15
 Operator : CG/JU
 Sample : PB147897BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147897BL

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-BP092122.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP011797.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.046	6	10	21	rVB	157574	232541	1.22%	0.274%
2	5.011	338	344	357	rBV	542915	825155	4.32%	0.971%
3	5.481	417	424	441	rBV	5447535	8631302	45.19%	10.156%
4	7.081	688	696	718	rBV	5627043	9252014	48.44%	10.887%
5	7.916	831	838	851	rBV	921431	1513913	7.93%	1.781%
6	9.087	1029	1037	1052	rBV	3328095	5475459	28.67%	6.443%
7	10.734	1309	1317	1335	rBV	1304500	2327669	12.19%	2.739%
8	13.193	1727	1735	1752	rBV	8497036	12799052	67.01%	15.060%
9	14.563	1962	1968	1989	rBV	2119649	3212361	16.82%	3.780%
10	16.057	2215	2222	2250	rBV	5966765	8866990	46.42%	10.433%
11	17.322	2430	2437	2451	rBV	2819641	4002292	20.95%	4.709%
12	19.880	2866	2872	2882	rBV	15105851	19101497	100.00%	22.476%
13	21.404	3126	3131	3141	rBV	3393889	4375738	22.91%	5.149%
14	23.857	3542	3548	3562	rVB2	2062720	4370100	22.88%	5.142%

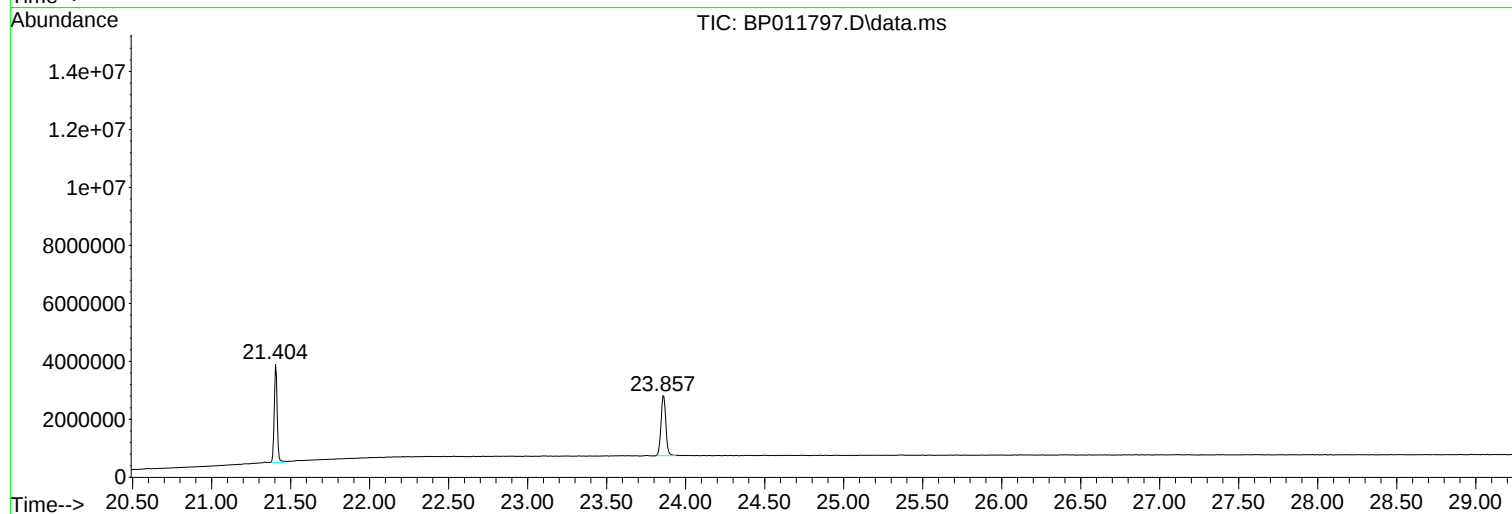
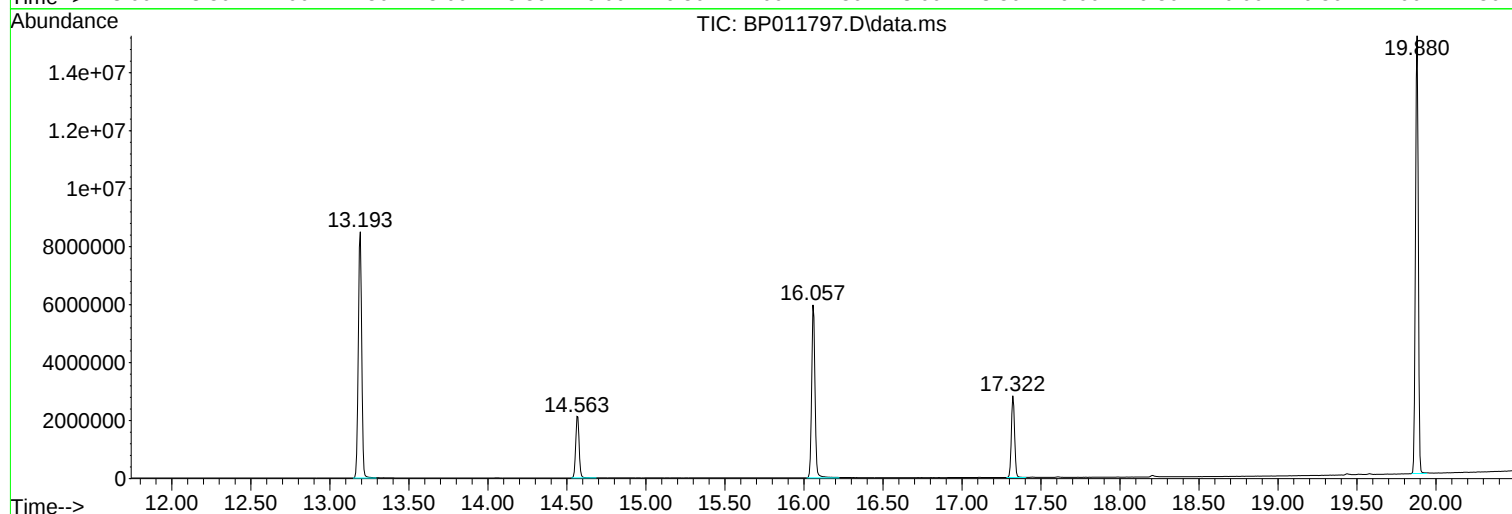
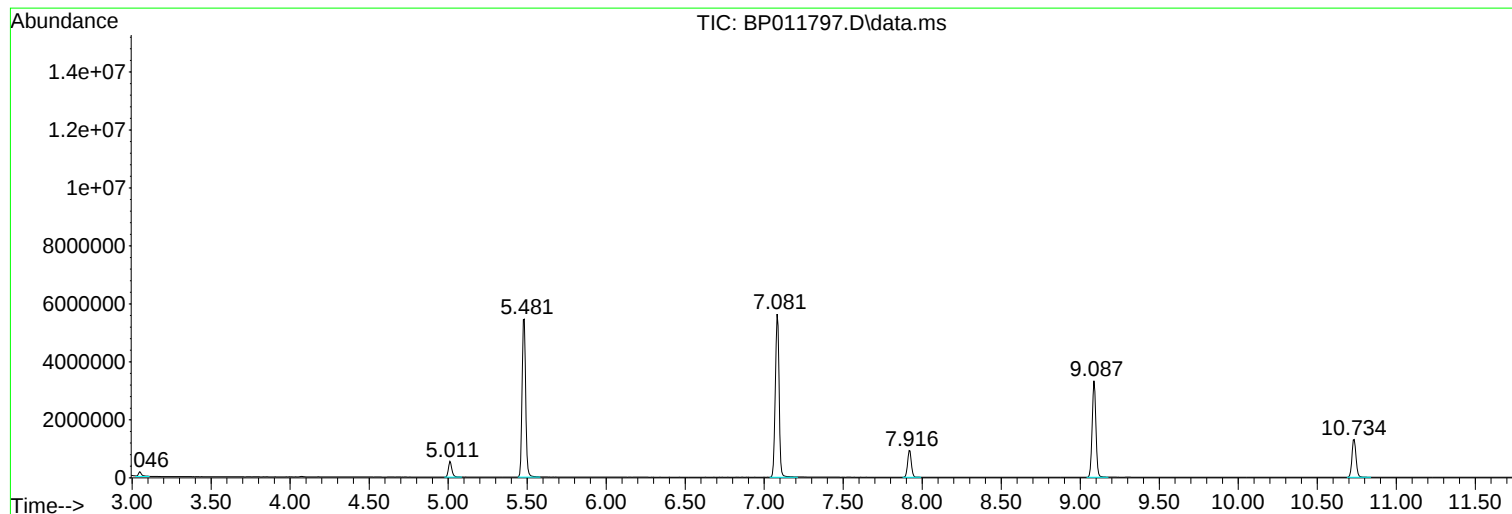
Sum of corrected areas: 84986083

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092122.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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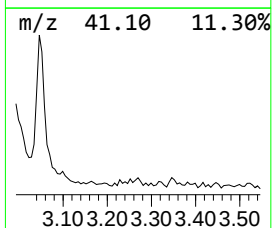
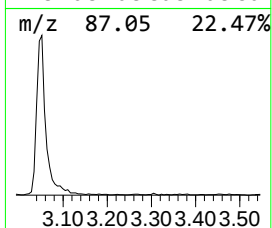
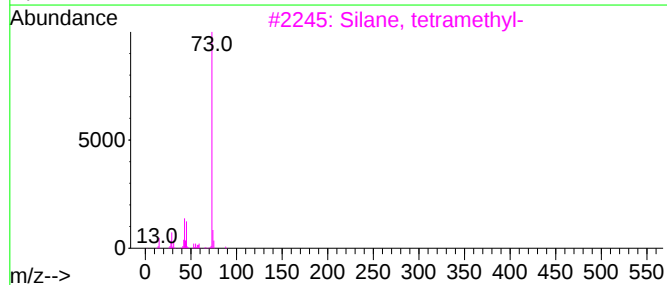
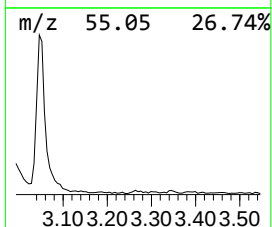
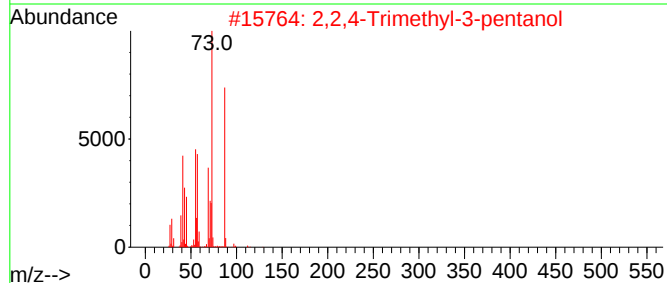
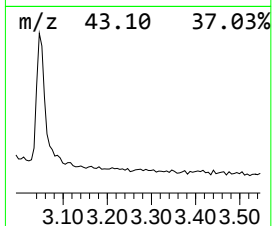
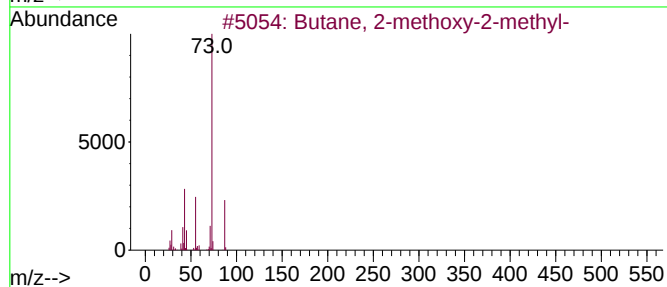
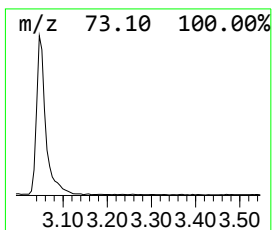
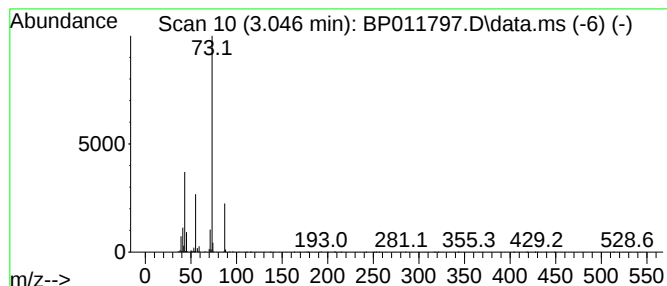
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.046	3.07 ng	232541	1,4-Dichlorobenzene-d4	7.922

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	9
4			N-Ethylformamide	73	C3H7NO	000627-45-2	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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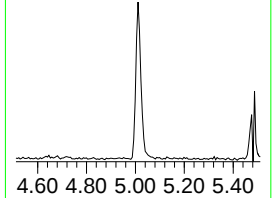
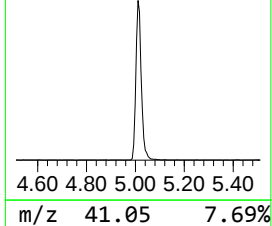
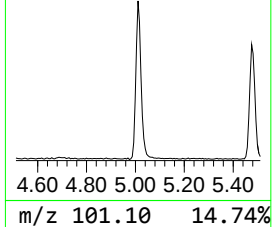
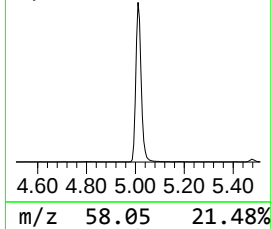
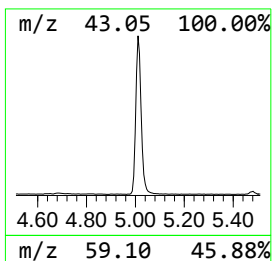
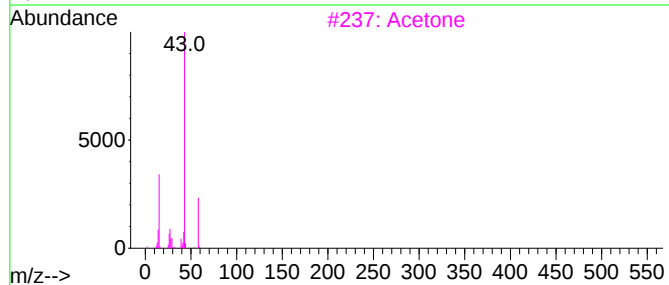
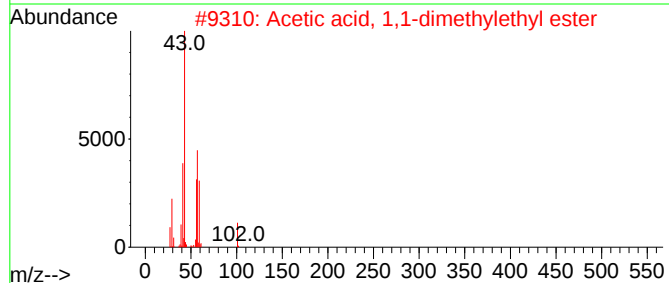
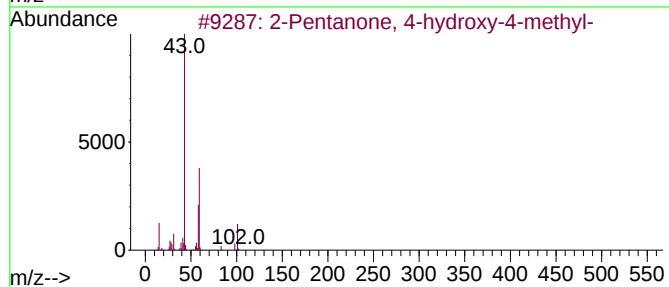
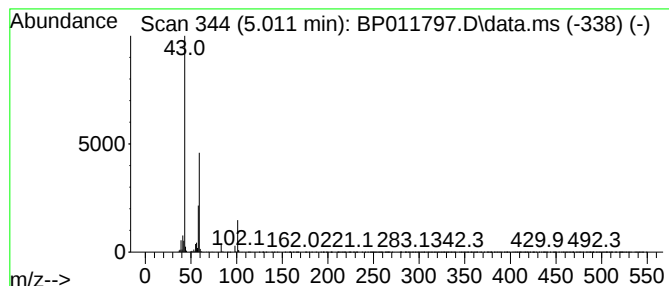
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.011	10.90 ng	825155	1,4-Dichlorobenzene-d4	7.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	83
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	42
3			Acetone	58	C3H6O	000067-64-1	16
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.046	3.1	ng	232541	1	7.922	1513910	20.0
2-Pentanone, 4-...	5.011	10.9	ng	825155	1	7.922	1513910	20.0