

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091422\
 Data File : BP011704.D
 Acq On : 14 Sep 2022 20:46
 Operator : CG/JU
 Sample : PB147629BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB147629BL

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: BP011704.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	rBV	284171	315068	1.23%	0.269%
2	5.040	342	349	359	rBV	270345	409986	1.60%	0.350%
3	5.499	420	427	440	rBV	9586949	14022756	54.58%	11.957%
4	7.099	691	699	721	rBV	8439737	14081555	54.81%	12.007%
5	7.940	835	842	853	rBV	1431308	2262521	8.81%	1.929%
6	9.110	1029	1041	1056	rBV	5135341	8744016	34.03%	7.456%
7	10.751	1309	1320	1336	rBV	1892054	3216969	12.52%	2.743%
8	13.210	1729	1738	1750	rBV	12353047	19121002	74.42%	16.305%
9	14.581	1964	1971	1985	rBV	2708712	3940244	15.34%	3.360%
10	16.069	2217	2224	2245	rBV	8885878	12651145	49.24%	10.788%
11	17.333	2432	2439	2449	rBV	3093911	4364004	16.99%	3.721%
12	19.886	2867	2873	2878	rBV	20555748	25692967	100.00%	21.909%
13	21.410	3127	3132	3144	rBV	3361494	4366234	16.99%	3.723%
14	23.862	3543	3549	3565	rVB2	1985126	4085023	15.90%	3.483%

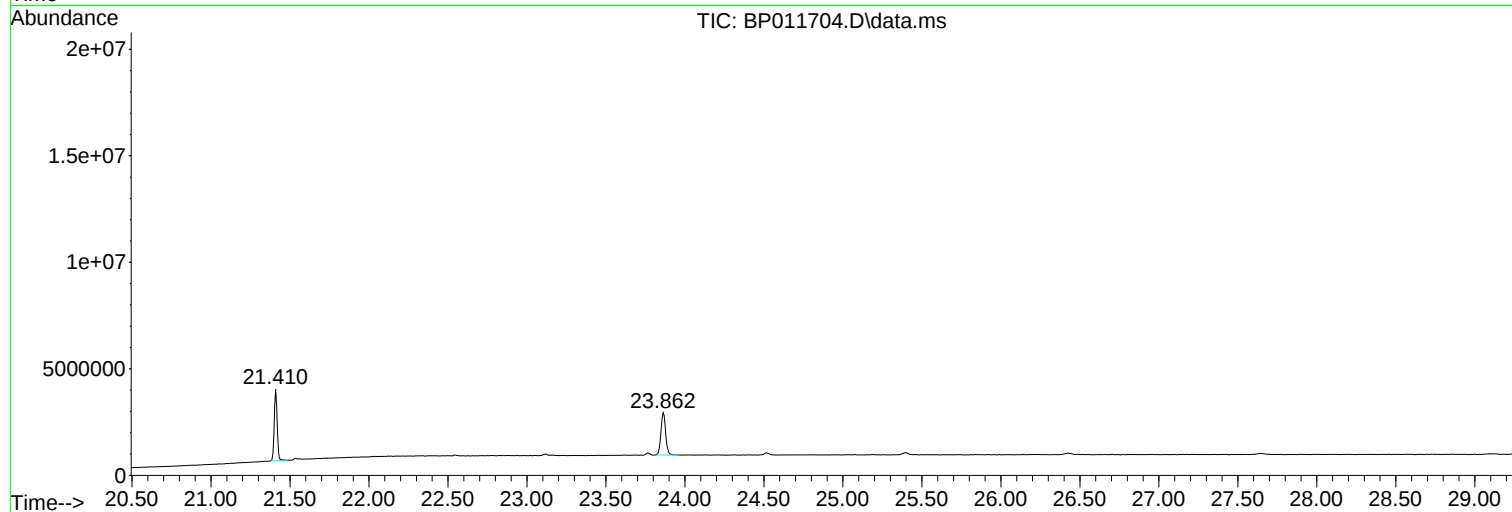
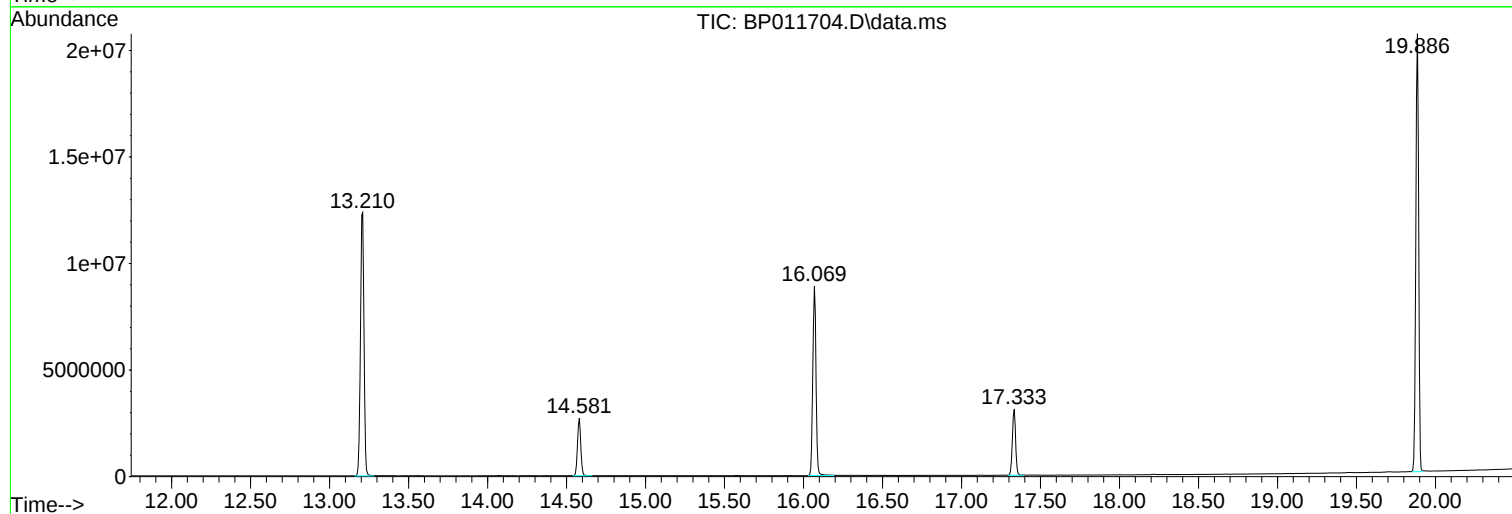
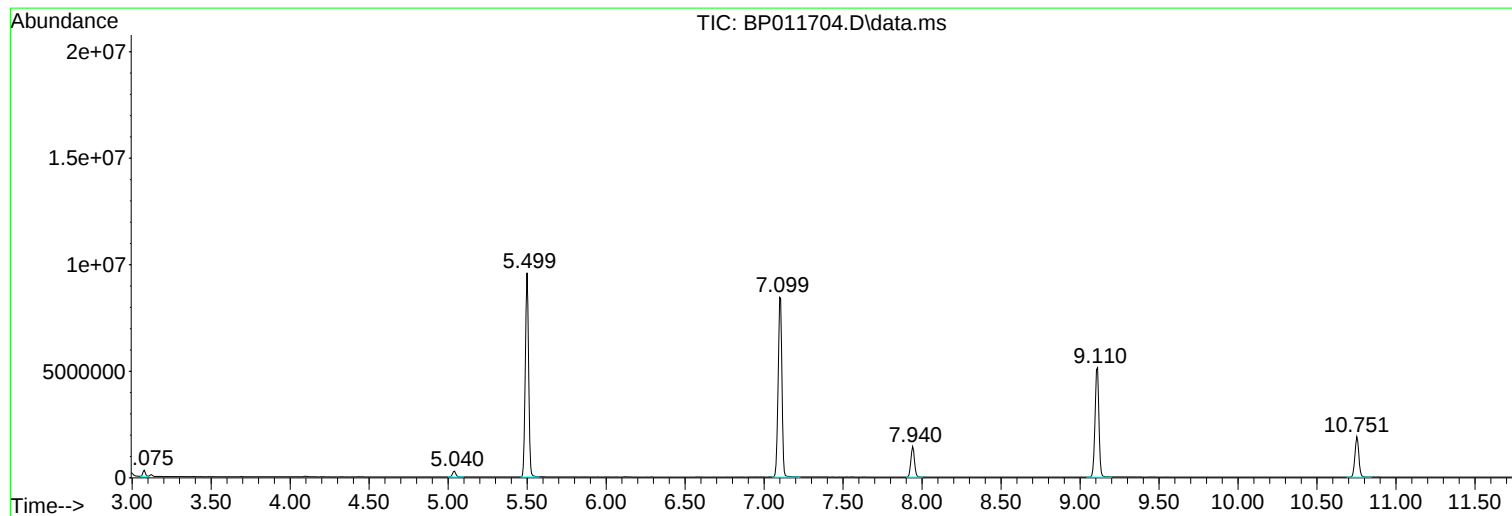
Sum of corrected areas: 117273490

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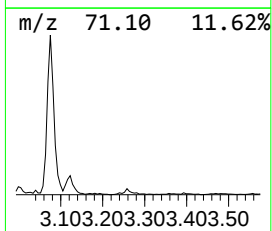
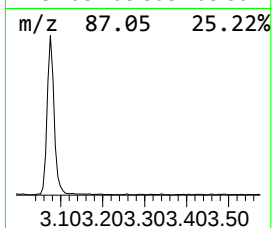
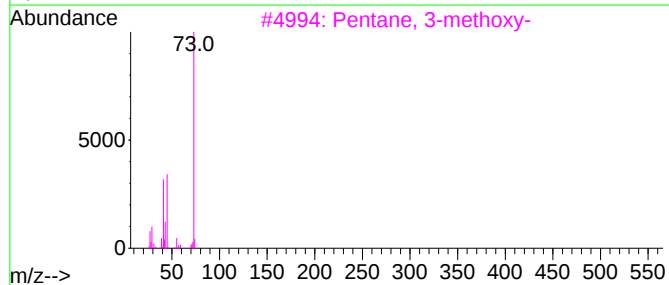
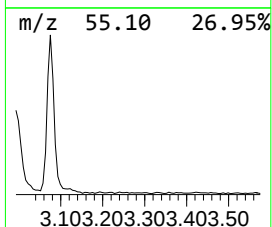
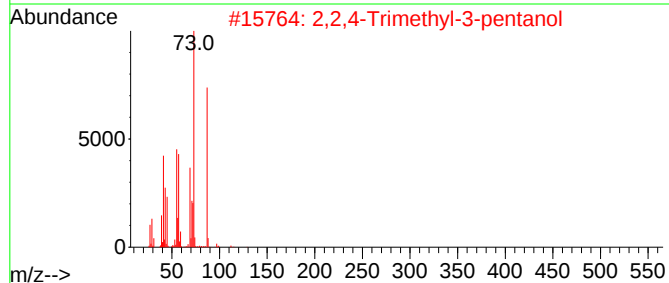
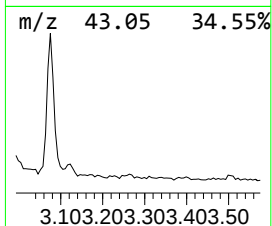
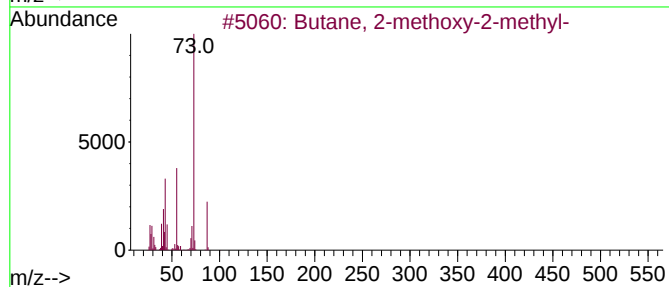
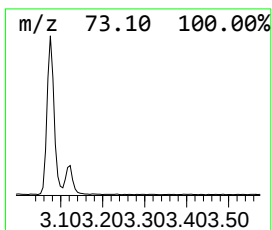
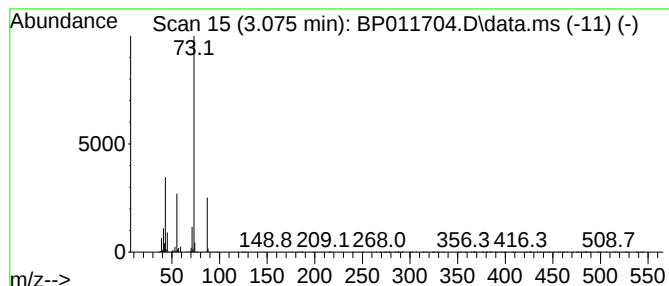
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.075	2.79 ng	315068	1,4-Dichlorobenzene-d4	7.940

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			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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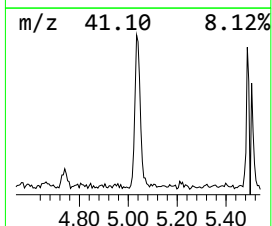
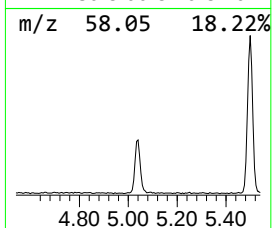
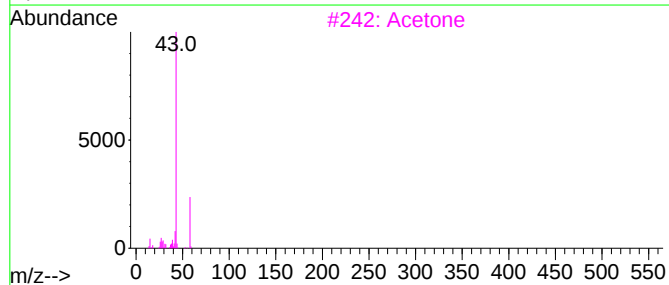
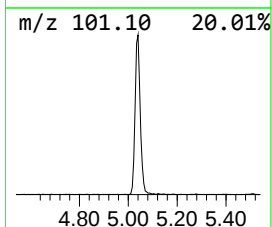
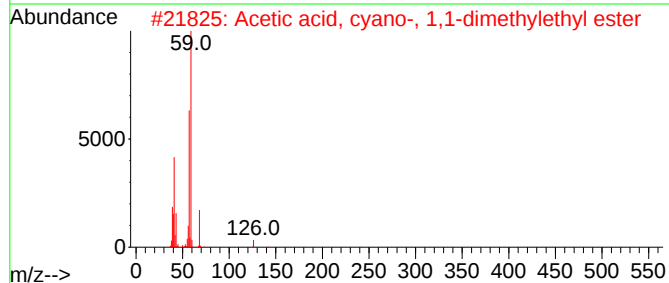
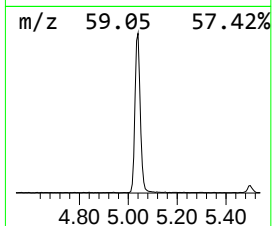
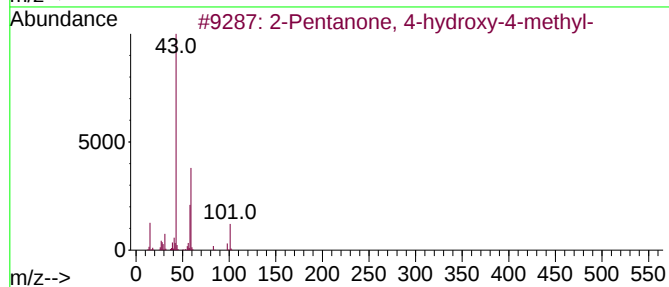
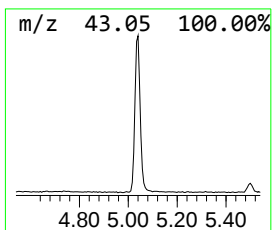
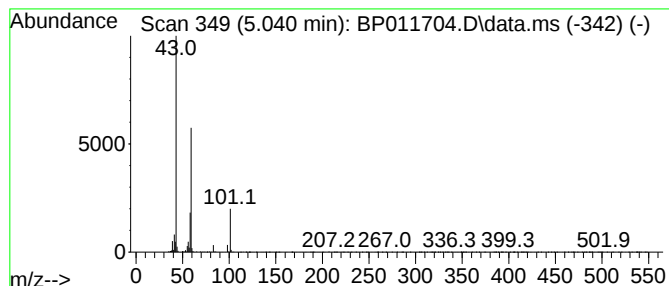
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.040	3.62 ng	409986	1,4-Dichlorobenzene-d4	7.940

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			Hexanamide	115	C6H13NO	000628-02-4	9
5			Tetraglyme	222	C10H22O5	000143-24-8	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	315068	1	7.940	2262520	20.0
2-Pentanone, 4-...	5.040	3.6	ng	409986	1	7.940	2262520	20.0