

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP082720\
 Data File : BP003145.D
 Acq On : 27 Aug 2020 22:39
 Operator : CG/JU
 Sample : L3498-01 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HR-06-082620

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP082420.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.934	3	8	20	rVB	234322	312022	5.11%	0.778%
2	5.269	399	405	417	rBV	393616	585732	9.60%	1.460%
3	6.840	664	672	691	rBV	351910	600850	9.85%	1.498%
4	7.663	805	812	819	rBV	808972	1259722	20.65%	3.140%
5	8.805	993	1006	1020	rBV	237100	444038	7.28%	1.107%
6	10.440	1276	1284	1293	rBV	1081161	1849112	30.31%	4.610%
7	11.899	1524	1532	1544	rBV3	44404	104758	1.72%	0.261%
8	12.375	1608	1613	1618	rBV2	37704	62370	1.02%	0.155%
9	12.922	1700	1706	1717	rBV	787668	1210247	19.84%	3.017%
10	13.151	1738	1745	1751	rBV	63305	116409	1.91%	0.290%
11	13.246	1757	1761	1766	rVB2	40404	66125	1.08%	0.165%
12	13.763	1840	1849	1855	rBV	107346	223589	3.66%	0.557%
13	14.234	1924	1929	1933	rBV	101869	121789	2.00%	0.304%
14	14.304	1934	1941	1953	rBV	1647295	2533681	41.53%	6.316%
15	14.857	2032	2035	2041	rVB7	41641	65201	1.07%	0.163%
16	15.187	2087	2091	2100	rVB	119402	153834	2.52%	0.383%
17	15.598	2154	2161	2166	rBV3	56118	116062	1.90%	0.289%
18	15.798	2189	2195	2205	rVB	498913	753739	12.35%	1.879%
19	16.051	2234	2238	2241	rBV	150517	198433	3.25%	0.495%
20	16.851	2369	2374	2378	rBV2	144816	189588	3.11%	0.473%
21	16.904	2379	2383	2389	rVB2	64880	97770	1.60%	0.244%
22	17.051	2402	2408	2421	rBV	1895865	2854364	46.79%	7.116%
23	17.586	2495	2499	2503	rBV	141253	172543	2.83%	0.430%
24	17.628	2503	2506	2511	rVB2	81354	102028	1.67%	0.254%
25	17.757	2523	2528	2533	rBV	2212177	2636488	43.21%	6.572%
26	17.969	2561	2564	2571	rBV3	74050	116531	1.91%	0.290%
27	18.263	2611	2614	2621	rVB	119774	142752	2.34%	0.356%
28	18.863	2711	2716	2718	rBV	4996026	6100949	100.00%	15.209%
29	18.892	2718	2721	2725	rVB2	4998887	5846020	95.82%	14.573%
30	19.022	2739	2743	2748	rBV	1102447	1163803	19.08%	2.901%
31	19.433	2810	2813	2823	rVB3	117804	205893	3.37%	0.513%
32	19.628	2842	2846	2851	rBV	1472268	1765019	28.93%	4.400%
33	19.957	2899	2902	2904	rBV	89966	119408	1.96%	0.298%
34	19.980	2904	2906	2916	rVB5	160579	209373	3.43%	0.522%

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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

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35	20.086	2921	2924	2927	rBV	102753	102406	1.68%	0.255%
36	20.375	2971	2973	2978	rVB3	104878	123511	2.02%	0.308%
37	20.692	3022	3027	3030	rVB3	107042	185111	3.03%	0.461%
38	20.886	3056	3060	3067	rBV3	135248	210437	3.45%	0.525%
39	21.145	3099	3104	3110	rBV	2642014	3379398	55.39%	8.424%
40	23.380	3478	3484	3494	rVB	1856132	3613559	59.23%	9.008%

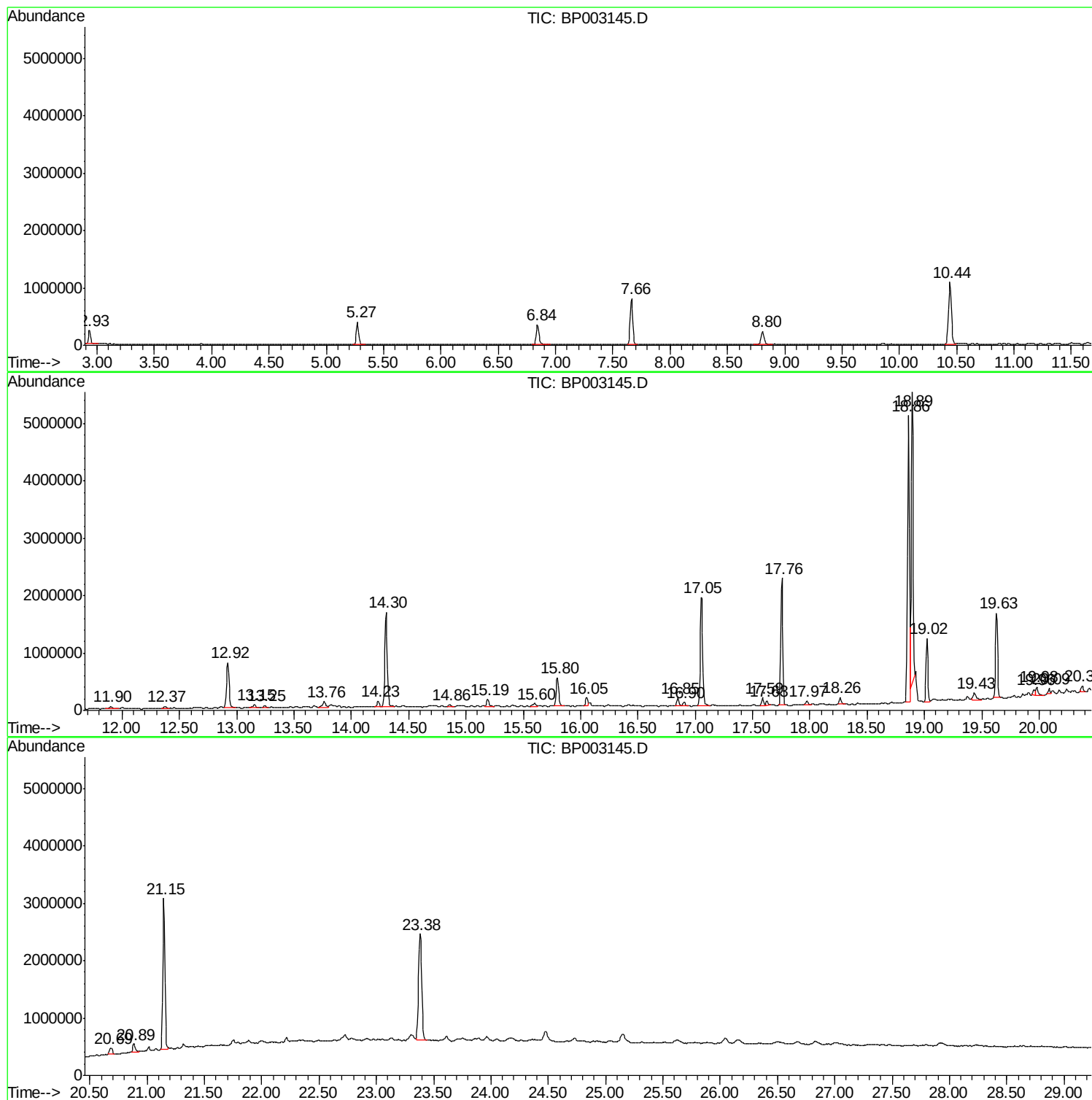
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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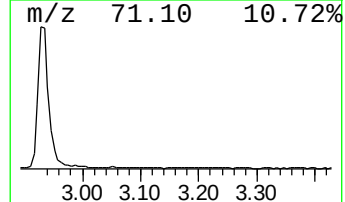
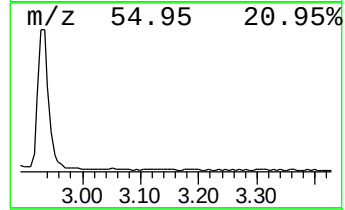
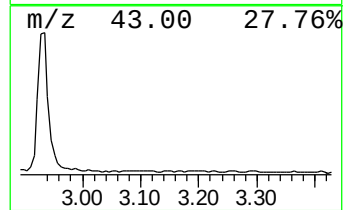
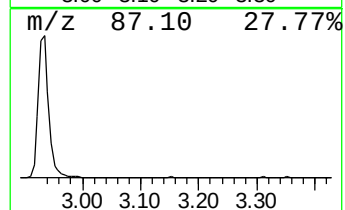
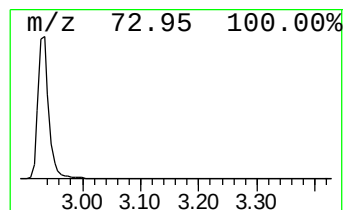
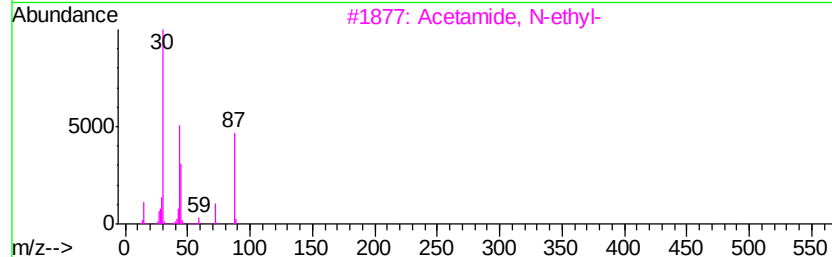
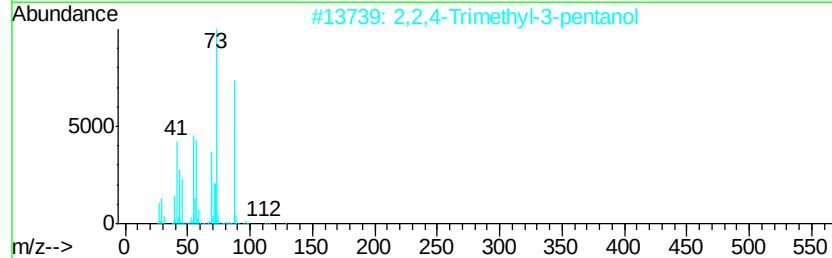
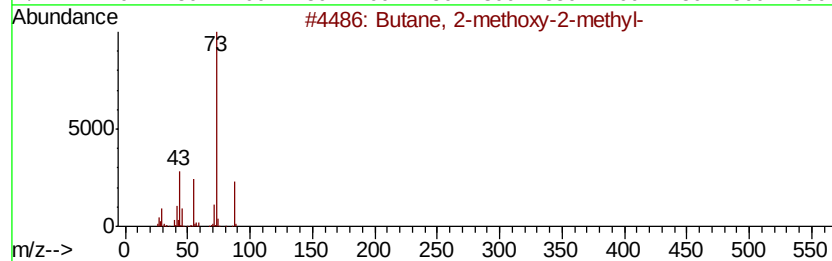
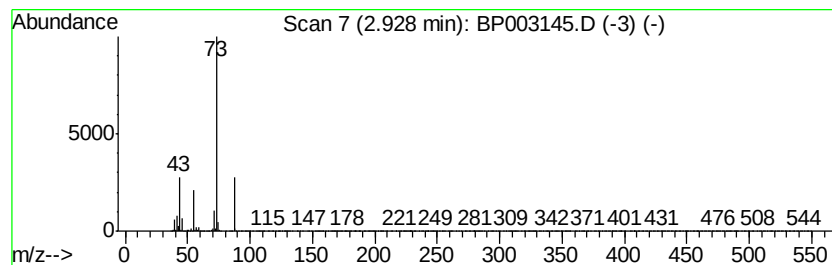
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.93	4.95 ng	312022	1,4-Dichlorobenzene-d4	7.66	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	9



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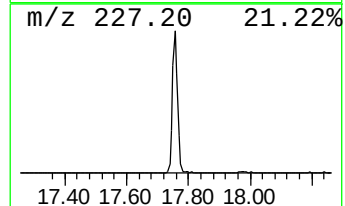
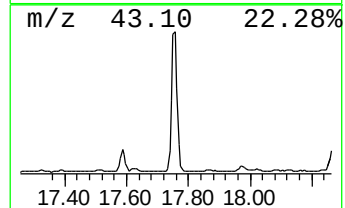
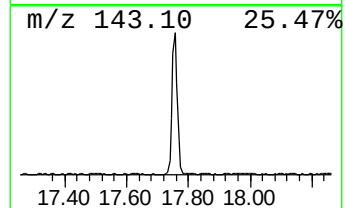
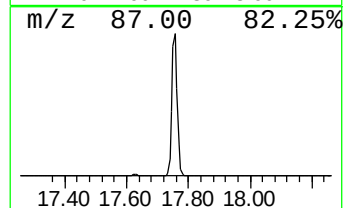
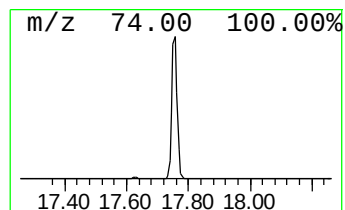
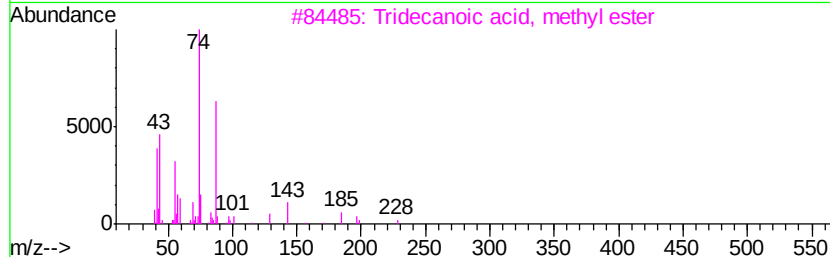
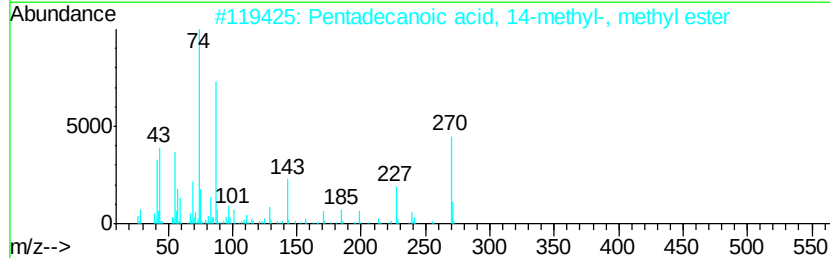
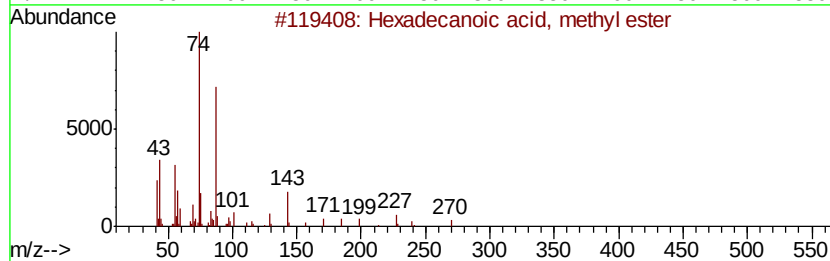
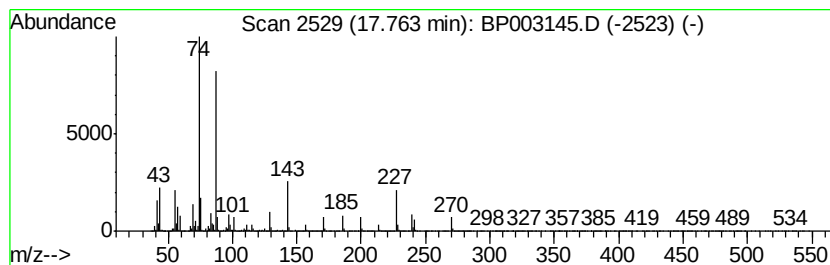
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.76	18.47 ng	2636490	Phenanthrene-d10	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	97
2		Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	90
3		Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4		Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
5		Methyl 9-methyltetradecanoate	256	C16H32O2	213617-69-7	83



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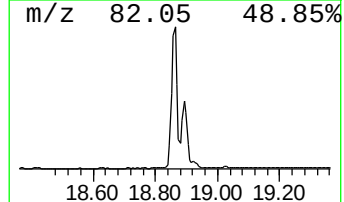
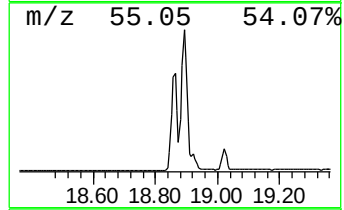
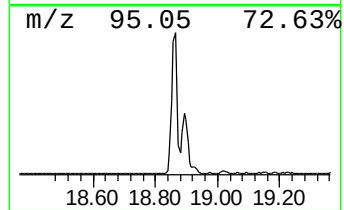
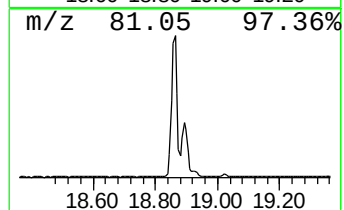
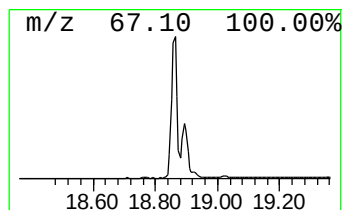
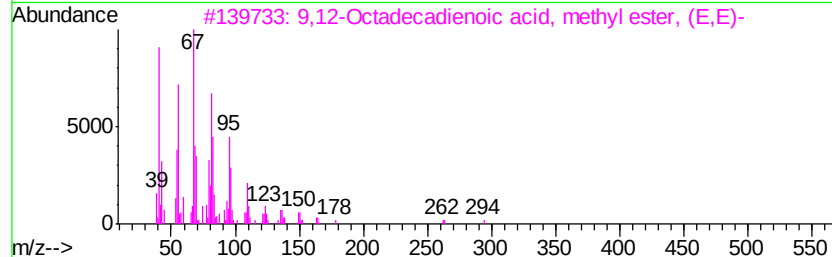
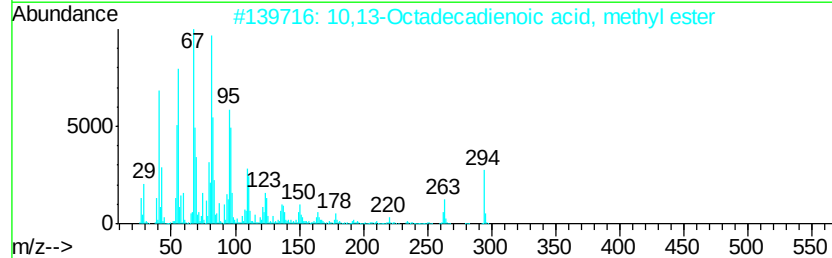
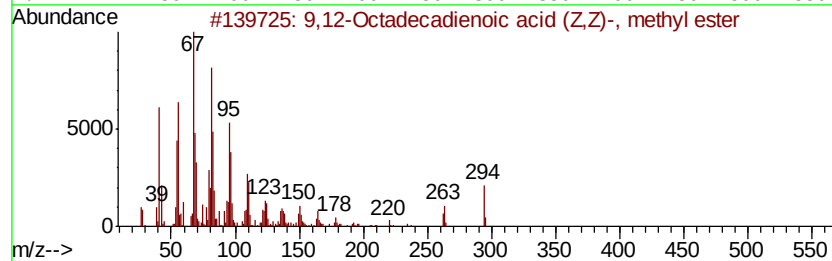
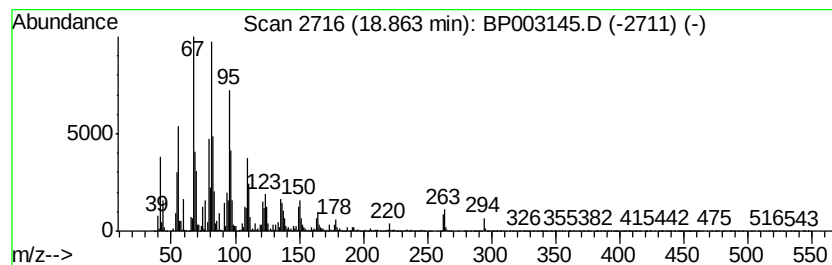
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 9,12-Octadecadienoic acid (... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.86	42.75 ng	6100950	Phenanthrene-d10	17.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99	
2	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99	
3	9,12-Octadecadienoic acid, methv...	294	C19H34O2	002566-97-4	99	
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99	
5	Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	1000336-44-0	99	



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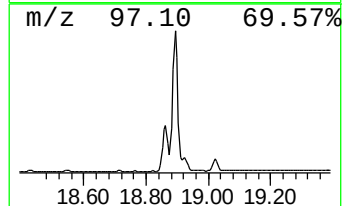
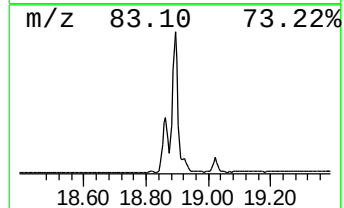
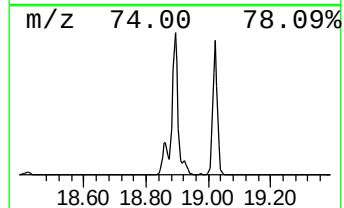
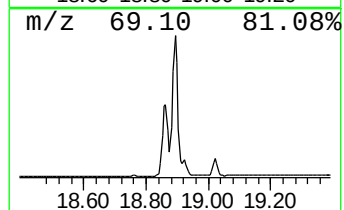
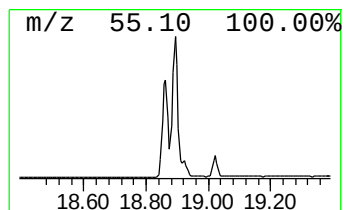
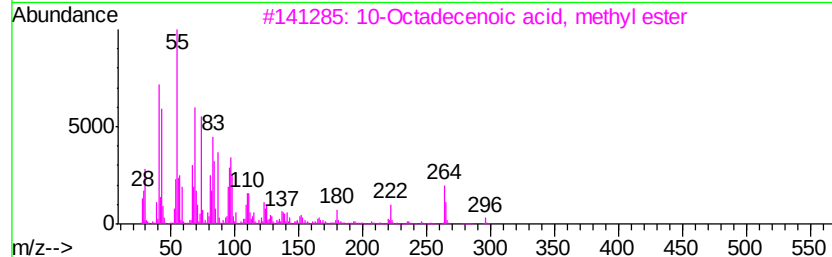
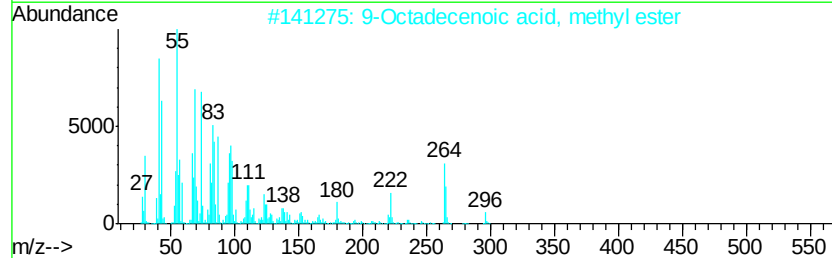
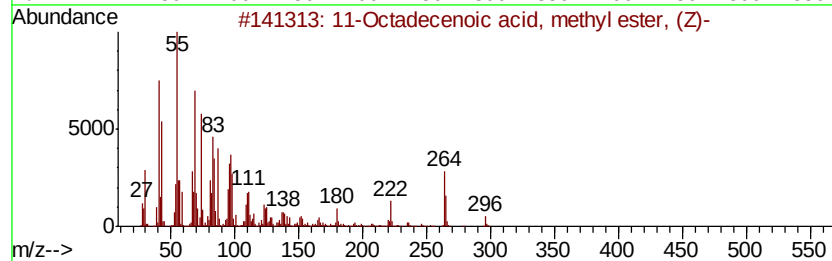
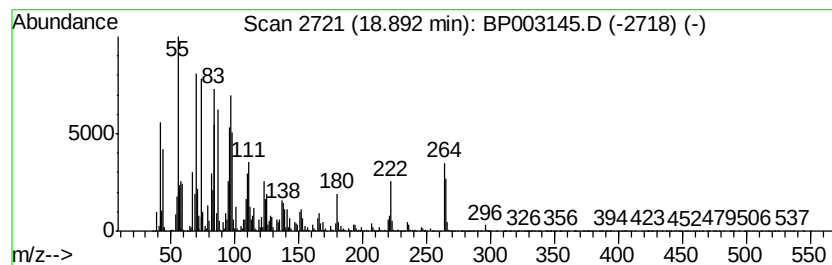
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 11-Octadecenoic acid, methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.89	40.96 ng	5846020	Phenanthrene-d10	17.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99
2		9-Octadecenoic acid, methyl ester	296	C19H36O2	002462-84-2	99
3		10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4		13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
5		12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99



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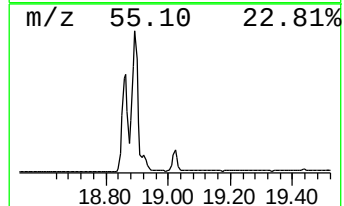
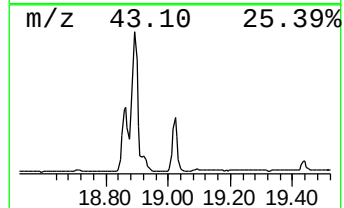
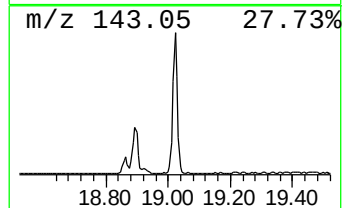
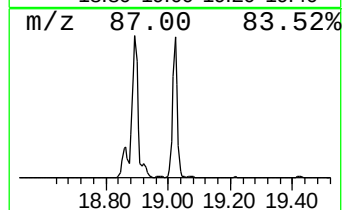
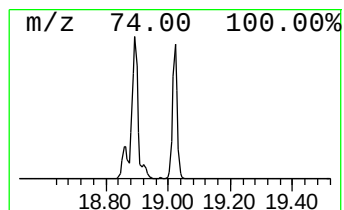
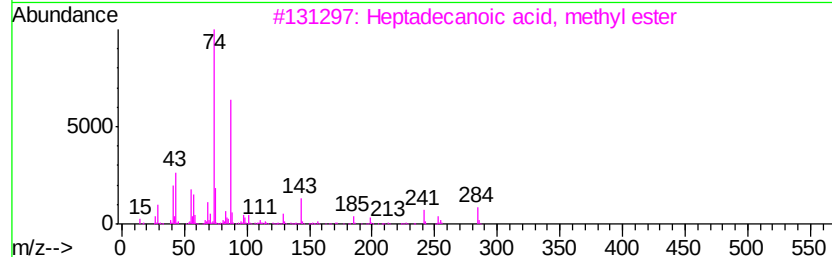
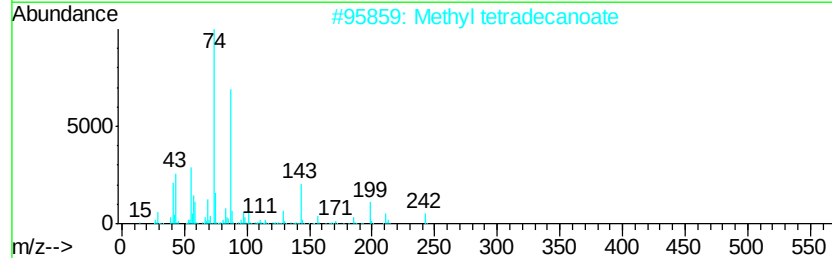
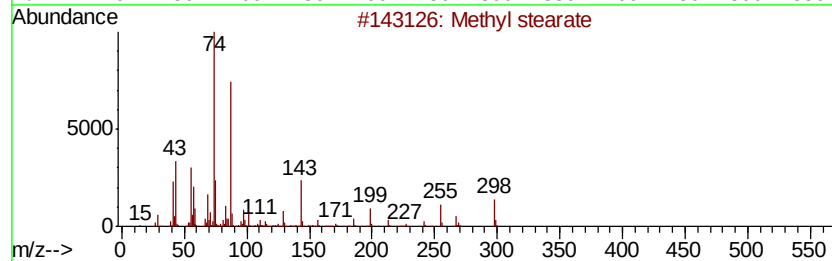
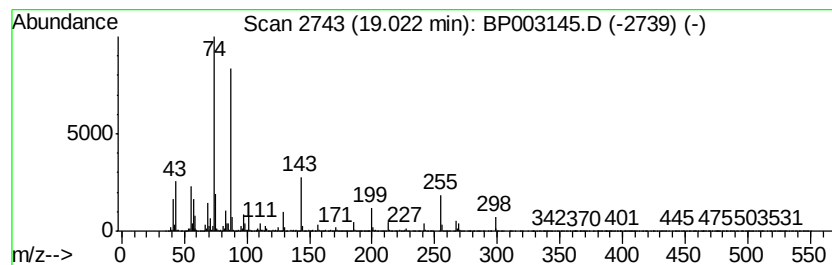
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Peak Number 5 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.02	8.15 ng	1163800	Phenanthrene-d10	17.05

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Methyl stearate	298	C19H38O2	000112-61-8	99
2		Methyl tetradecanoate	242	C15H30O2	000124-10-7	93
3		Heptadecanoic acid, methyl ester	284	C18H36O2	001731-92-6	80
4		Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	80
5		Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	78



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP082720\
Data File : BP003145.D
Acq On : 27 Aug 2020 22:39
Operator : CG/JU
Sample : L3498-01 5X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HR-06-082620

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP082420.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.93	5.0	ng	312022	1	7.66	1259720	20.0
Hexadecanoic acid...	17.76	18.5	ng	2636490	4	17.05	2854360	20.0
9,12-Octadecadien...	18.86	42.8	ng	6100950	4	17.05	2854360	20.0
11-Octadecenoic a...	18.89	41.0	ng	5846020	4	17.05	2854360	20.0
Methyl stearate	19.02	8.2	ng	1163800	4	17.05	2854360	20.0