

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
 Data File : BP017734.D
 Acq On : 19 Oct 2023 19:14
 Operator : MA/JU
 Sample : 04864-03DL 2X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCJV6DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\SFAM-EPA-BP101823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017734.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.210	17	21	27	rVV	31481	48915	1.62%	0.211%
2	3.257	27	29	38	rVB2	15792	27056	0.89%	0.117%
3	3.604	85	88	93	rVB2	10892	14223	0.47%	0.061%
4	3.704	101	105	120	rBV	67185	131891	4.36%	0.569%
5	3.875	130	134	138	rBV	11970	19417	0.64%	0.084%
6	3.999	152	155	160	rVB3	10326	15892	0.53%	0.069%
7	4.393	219	222	232	rVB3	8632	15998	0.53%	0.069%
8	4.928	308	313	317	rBV2	9753	15916	0.53%	0.069%
9	5.340	378	383	390	rVB5	13489	24485	0.81%	0.106%
10	5.698	438	444	448	rBV2	9123	15075	0.50%	0.065%
11	5.734	448	450	459	rVB9	3145	6183	0.20%	0.027%
12	5.934	478	484	491	rVB9	5777	15809	0.52%	0.068%
13	6.481	572	577	582	rVB6	6500	10949	0.36%	0.047%
14	6.893	639	647	651	rBV	74482	145845	4.82%	0.629%
15	7.051	668	674	685	rBV	61076	117766	3.89%	0.508%
16	7.234	699	705	716	rBV	251636	401513	13.27%	1.731%
17	7.328	716	721	728	rVB7	8135	15492	0.51%	0.067%
18	7.410	728	735	744	rBV	242154	392787	12.98%	1.694%
19	7.563	758	761	766	rVB7	4926	7061	0.23%	0.030%
20	7.887	808	816	819	rBV	362982	579428	19.15%	2.498%
21	7.916	819	821	831	rVB	207193	311278	10.29%	1.342%
22	8.169	859	864	872	rBV6	10510	24464	0.81%	0.105%
23	8.434	905	909	910	rBV4	4895	6696	0.22%	0.029%
24	8.463	910	914	919	rVB7	7464	13377	0.44%	0.058%
25	8.616	934	940	956	rBV	151689	275233	9.10%	1.187%
26	8.969	996	1000	1006	rBV9	5870	11660	0.39%	0.050%
27	9.098	1012	1022	1026	rBV	264184	450526	14.89%	1.943%
28	9.145	1026	1030	1039	rVB2	83735	162204	5.36%	0.699%
29	9.522	1092	1094	1103	rVB10	2326	5843	0.19%	0.025%
30	9.810	1137	1143	1154	rBV	208801	355929	11.76%	1.535%
31	9.986	1168	1173	1184	rVB8	8433	26069	0.86%	0.112%
32	10.151	1196	1201	1202	rBV5	4368	5441	0.18%	0.023%
33	10.345	1224	1234	1249	rBV	275422	546780	18.07%	2.358%
34	10.498	1255	1260	1266	rBV10	3959	8073	0.27%	0.035%
35	10.545	1266	1268	1276	rVB8	5982	11122	0.37%	0.048%
36	10.622	1276	1281	1285	rBV8	6898	12217	0.40%	0.053%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
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37	10.722	1288	1298	1305	rBV	465974	780732	25.80%	3.366%
38	10.898	1320	1328	1339	rBV	133602	279698	9.24%	1.206%
39	11.422	1412	1417	1422	rBV2	26610	47036	1.55%	0.203%
40	11.792	1475	1480	1484	rVB8	5342	9020	0.30%	0.039%

41	12.333	1562	1572	1577	rBV9	9042	23661	0.78%	0.102%
42	12.386	1577	1581	1584	rBV5	4687	8874	0.29%	0.038%
43	12.586	1609	1615	1629	rVB	28873	68283	2.26%	0.294%
44	12.798	1646	1651	1658	rBV6	13397	27109	0.90%	0.117%
45	13.145	1706	1710	1715	rVB8	5569	7319	0.24%	0.032%

46	13.327	1735	1741	1745	rBV8	6368	11043	0.36%	0.048%
47	13.410	1748	1755	1762	rBV3	68009	120669	3.99%	0.520%
48	13.539	1771	1777	1781	rBV	58055	101784	3.36%	0.439%
49	13.575	1781	1783	1794	rVB4	28234	54599	1.80%	0.235%
50	13.998	1848	1855	1861	rBV	629066	873156	28.85%	3.765%

51	14.051	1861	1864	1871	rVB5	20367	30950	1.02%	0.133%
52	14.198	1885	1889	1895	rVV8	6114	12820	0.42%	0.055%
53	14.280	1895	1903	1912	rVV	675047	998630	33.00%	4.306%
54	14.586	1948	1955	1961	rVV	688349	1003612	33.17%	4.327%
55	14.645	1961	1965	1969	rVB7	2993	5484	0.18%	0.024%

56	14.863	1998	2002	2008	rBV9	14860	35935	1.19%	0.155%
57	14.992	2019	2024	2027	rBV6	6030	11207	0.37%	0.048%
58	15.045	2029	2033	2037	rVV6	7149	12337	0.41%	0.053%
59	15.092	2037	2041	2047	rVB9	6725	10427	0.34%	0.045%
60	15.333	2078	2082	2087	rBV6	7714	11352	0.38%	0.049%

61	15.416	2092	2096	2098	rVV5	4835	6408	0.21%	0.028%
62	15.457	2101	2103	2111	rVB9	3965	7023	0.23%	0.030%
63	15.586	2117	2125	2133	rBV	888976	1286291	42.51%	5.546%
64	15.733	2145	2150	2162	rBV	190311	299986	9.91%	1.293%
65	15.821	2163	2165	2170	rBV5	4014	6519	0.22%	0.028%

66	15.898	2175	2178	2184	rVB8	4519	6927	0.23%	0.030%
67	15.980	2187	2192	2198	rBV	62316	91718	3.03%	0.395%
68	16.074	2204	2208	2215	rVB	24779	36855	1.22%	0.159%
69	16.133	2215	2218	2223	rVB7	5664	9217	0.30%	0.040%
70	16.357	2249	2256	2266	rBV3	26697	81598	2.70%	0.352%

71	16.721	2312	2318	2325	rBV3	35934	65312	2.16%	0.282%
72	16.798	2327	2331	2337	rVV9	6624	12966	0.43%	0.056%
73	16.851	2337	2340	2343	rVB5	5760	7728	0.26%	0.033%
74	16.927	2352	2353	2361	rVB8	6294	8426	0.28%	0.036%
75	17.186	2391	2397	2412	rBV2	182659	332006	10.97%	1.432%

76	17.304	2412	2417	2423	rVB7	8600	15469	0.51%	0.067%
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
 Title : SVOA CALIBRATION

77	17.374	2423	2429	2438	rBV	840224	1178595	38.95%	5.082%
78	17.474	2440	2446	2456	rBV	886703	1310065	43.29%	5.649%
79	17.615	2466	2470	2477	rVB5	16239	30393	1.00%	0.131%
80	17.733	2486	2490	2495	rBV8	11296	17496	0.58%	0.075%
81	17.986	2531	2533	2534	rVV2	7164	6540	0.22%	0.028%
82	18.010	2534	2537	2545	rVB10	19837	27086	0.90%	0.117%
83	18.104	2549	2553	2559	rVV6	17439	26991	0.89%	0.116%
84	18.227	2569	2574	2584	rBV	196421	290879	9.61%	1.254%
85	18.321	2584	2590	2595	rVV	2715472	3026042	100.00%	13.048%
86	18.774	2664	2667	2673	rVB8	11379	18533	0.61%	0.080%
87	18.980	2698	2702	2710	rBV3	21029	34561	1.14%	0.149%
88	19.080	2711	2719	2724	rVB3	26504	56237	1.86%	0.242%
89	19.304	2753	2757	2761	rBV7	16076	26010	0.86%	0.112%
90	19.374	2764	2769	2773	rVB4	36936	68108	2.25%	0.294%
91	19.474	2782	2786	2793	rBV4	80338	132743	4.39%	0.572%
92	19.610	2806	2809	2813	rBV	51059	57660	1.91%	0.249%
93	19.733	2825	2830	2837	rBV	1176223	1541619	50.95%	6.647%
94	19.792	2837	2840	2848	rVB2	36594	59052	1.95%	0.255%
95	20.015	2876	2878	2879	rBV2	9334	6715	0.22%	0.029%
96	20.645	2982	2985	2989	rBV2	40764	42308	1.40%	0.182%
97	21.515	3128	3133	3140	rBV2	929818	1216803	40.21%	5.247%
98	22.762	3341	3345	3350	rVB	307429	390580	12.91%	1.684%
99	23.898	3532	3538	3548	rVB2	691142	1422643	47.01%	6.134%
100	24.074	3561	3568	3577	rVB	546215	1135598	37.53%	4.896%

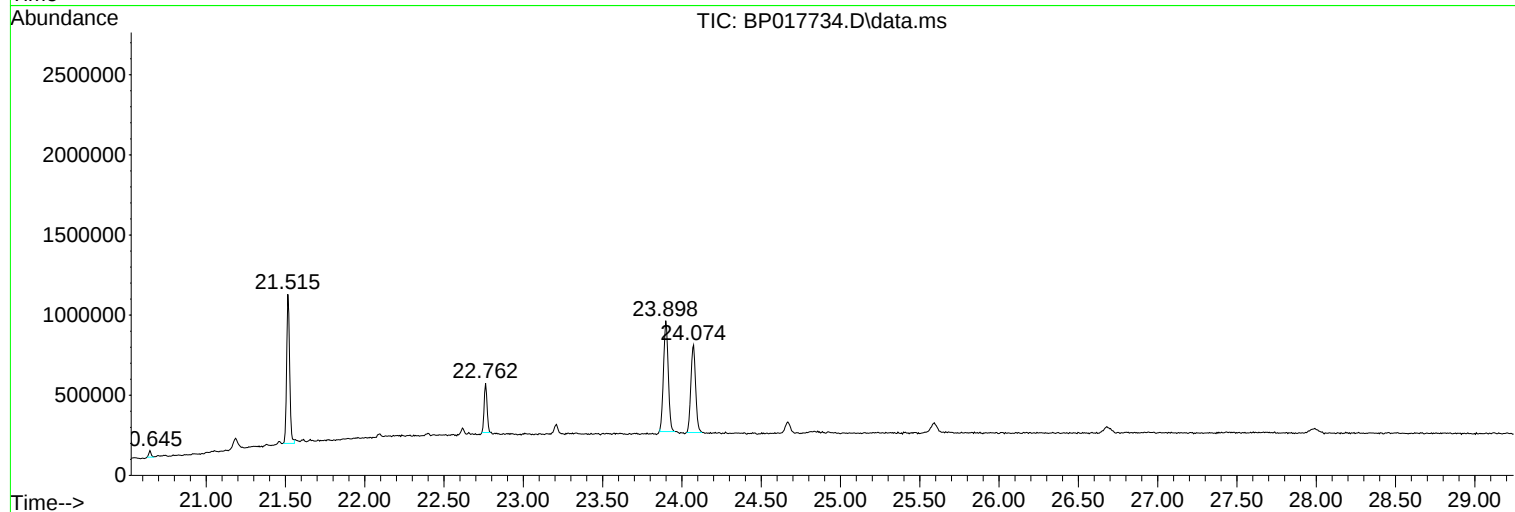
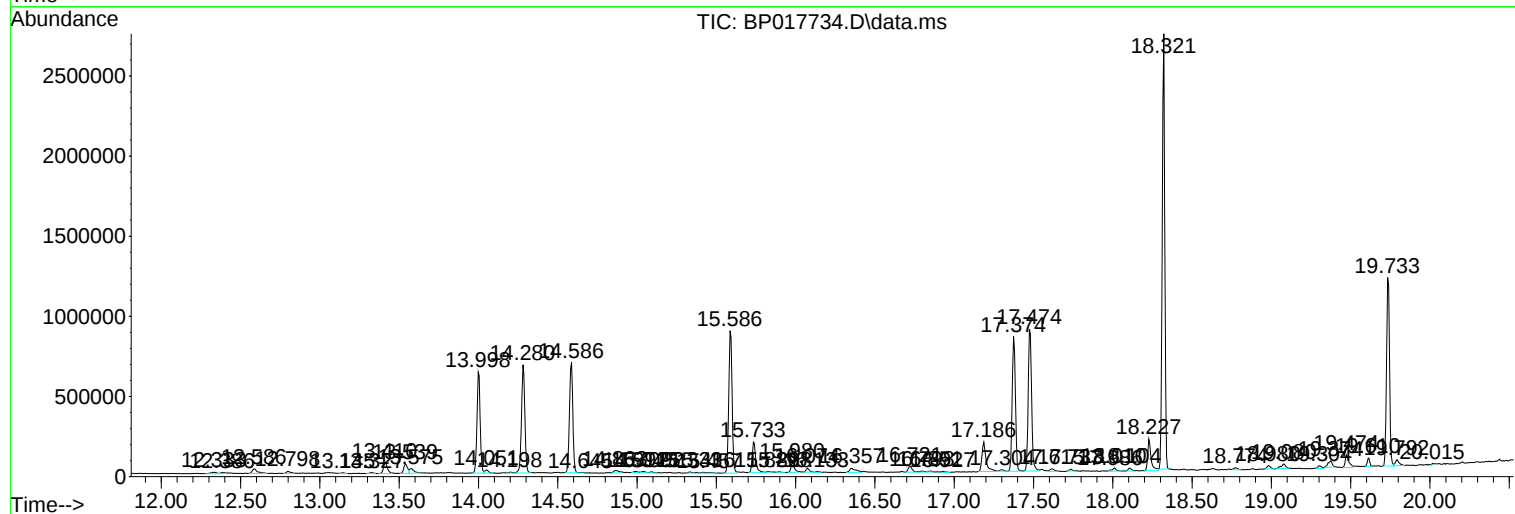
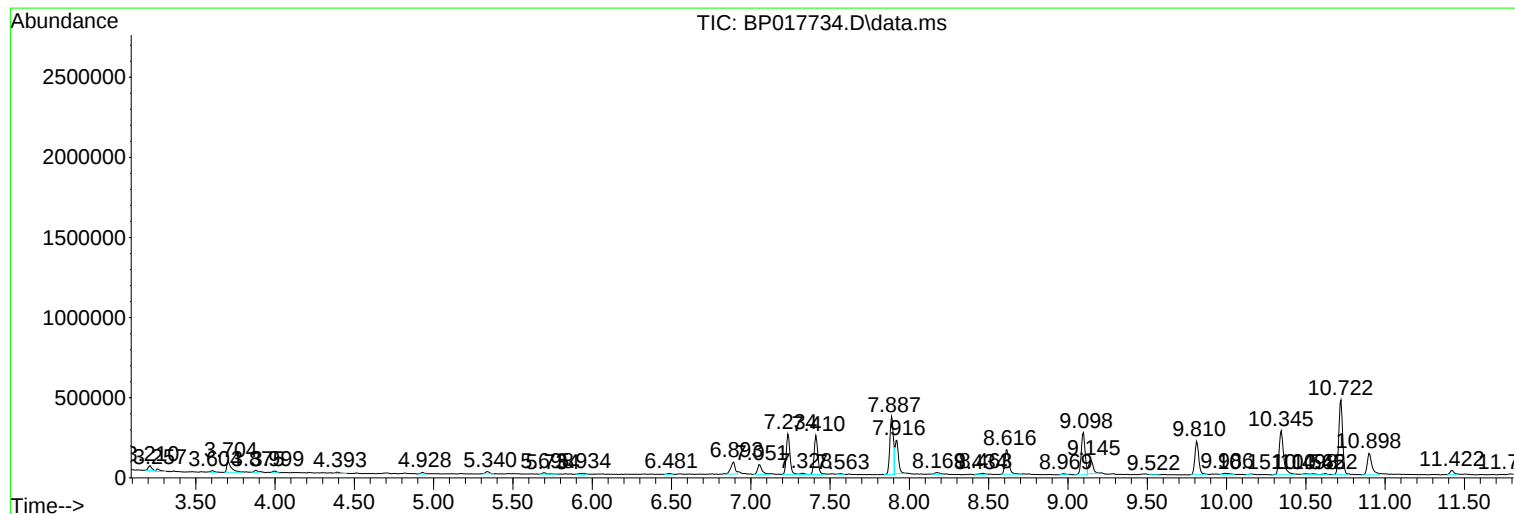
Sum of corrected areas: 23192056

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Instrument :
BNA_P
ClientSampleId :
DCJV6DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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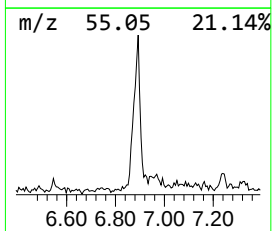
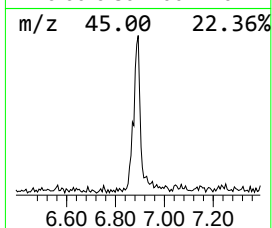
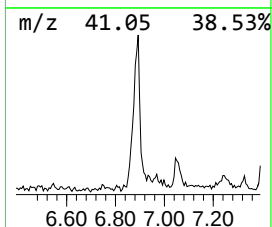
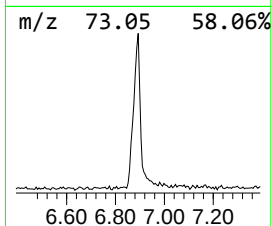
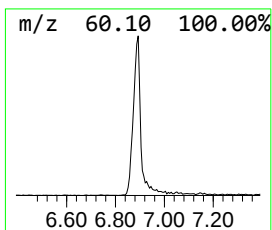
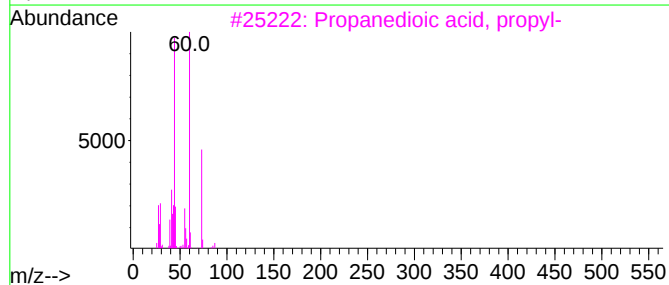
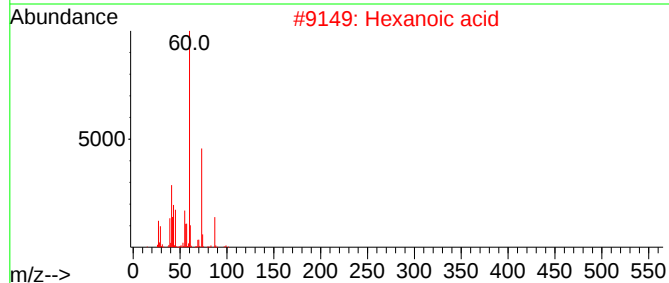
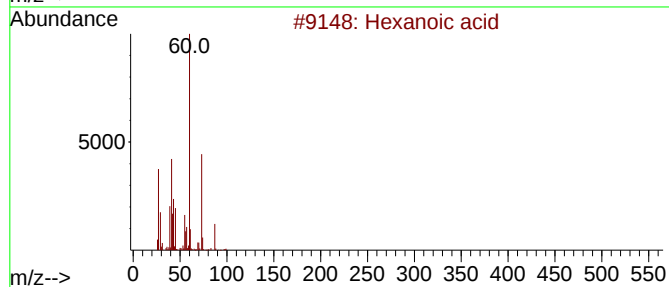
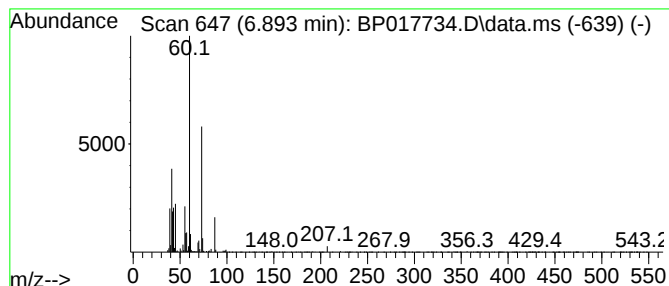
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Hexanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.893	5.03 ng/ul	145845	1,4-Dichlorobenzene-d4	7.887

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid	116	C6H12O2	000142-62-1	83
2			Hexanoic acid	116	C6H12O2	000142-62-1	83
3			Propanedioic acid, propyl-	146	C6H10O4	000616-62-6	78
4			Hexanoic acid	116	C6H12O2	000142-62-1	72
5			Hexanoic acid	116	C6H12O2	000142-62-1	64



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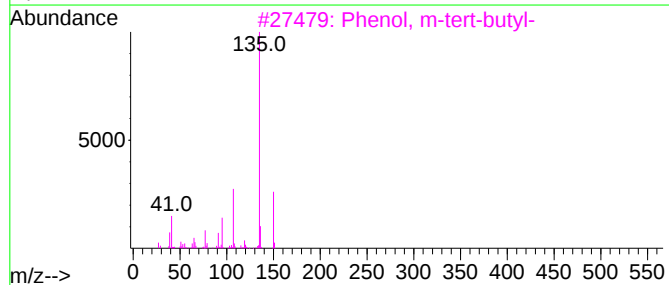
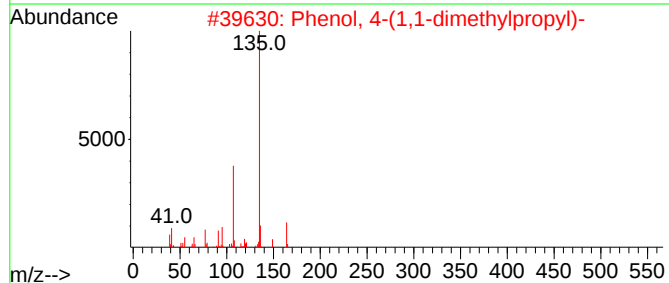
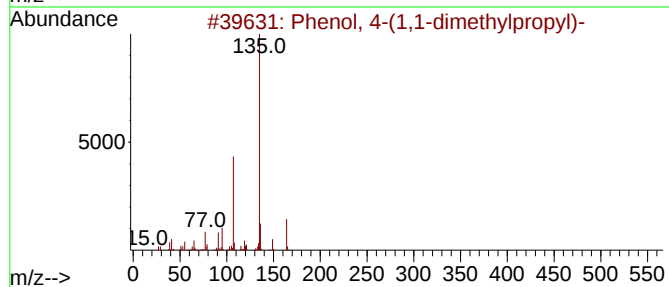
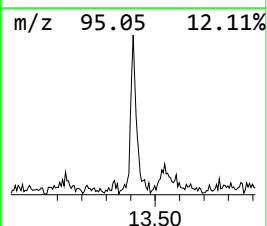
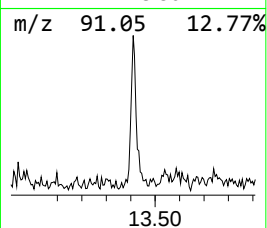
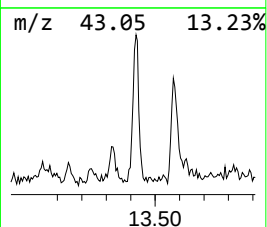
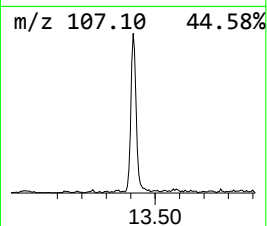
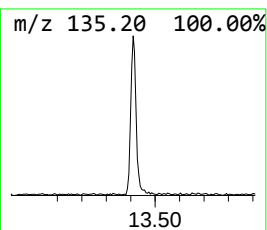
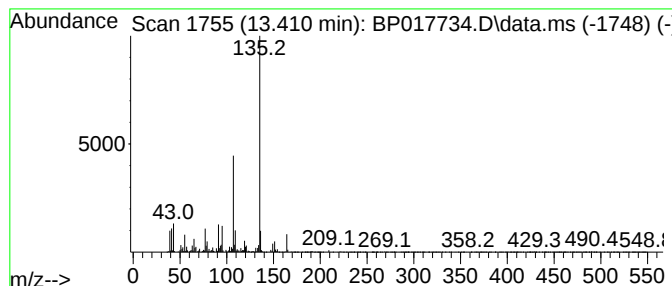
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Phenol, 4-(1,1-dimethylprop... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.410	2.40 ng/ul	120669	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	95
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	91
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	76
4			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	72
5			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64



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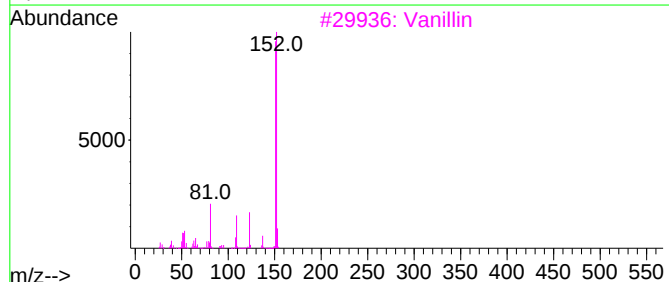
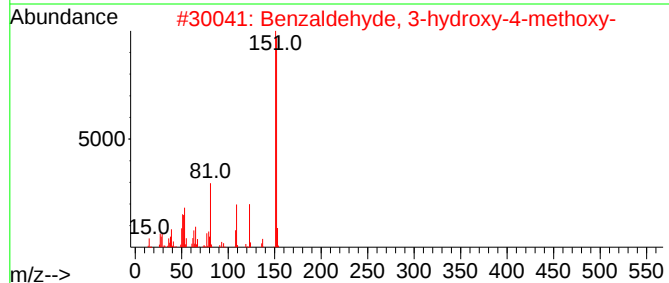
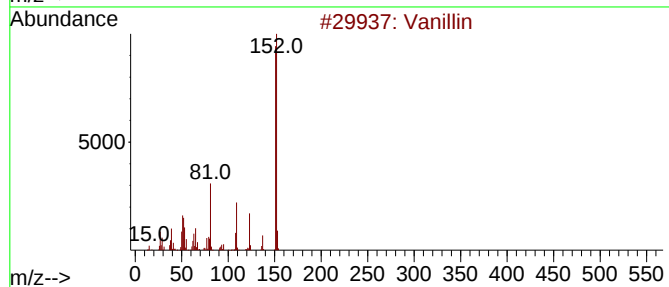
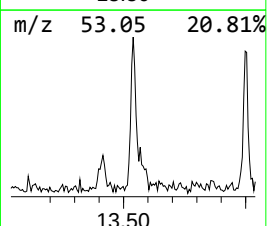
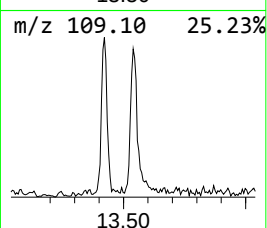
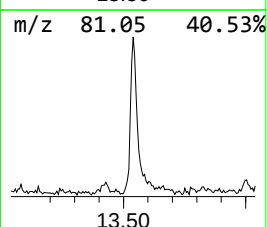
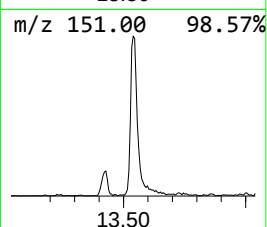
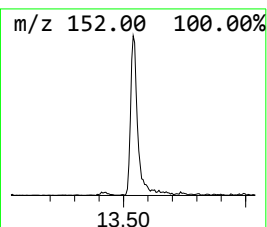
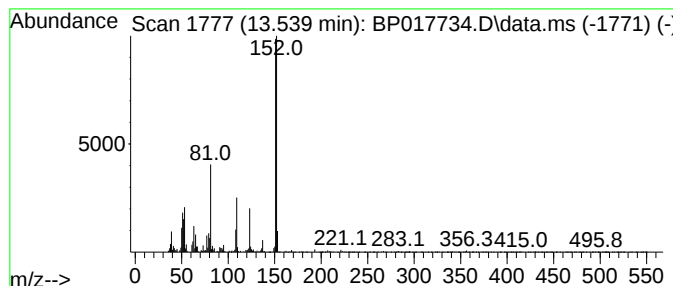
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Vanillin Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.539	2.03 ng/ul	101784	Acenaphthene-d10	14.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Vanillin	152	C8H8O3	000121-33-5	96
2			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	96
3			Vanillin	152	C8H8O3	000121-33-5	95
4			Vanillin	152	C8H8O3	000121-33-5	95
5			Benzaldehyde, 3-hydroxy-4-methoxy-	152	C8H8O3	000621-59-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
Data File : BP017734.D
Acq On : 19 Oct 2023 19:14
Operator : MA/JU
Sample : 04864-03DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCJV6DL

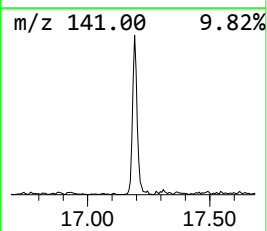
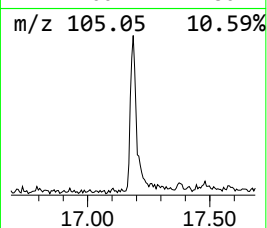
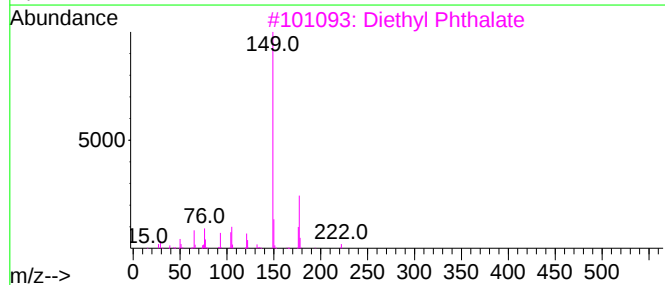
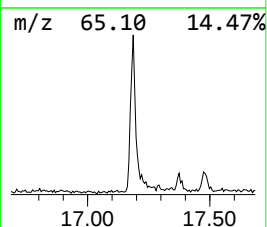
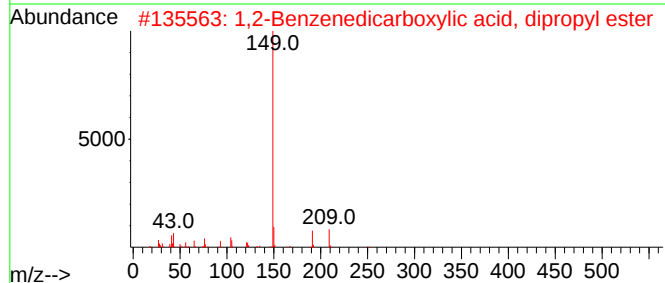
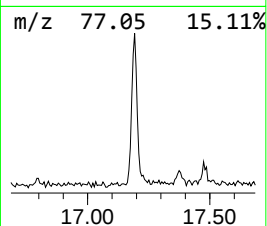
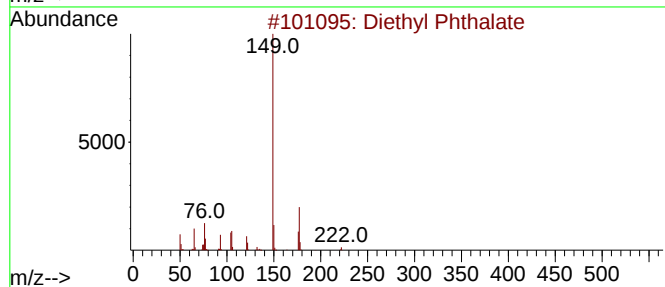
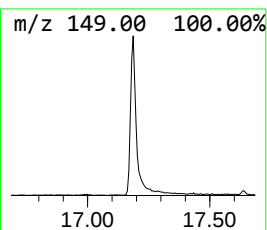
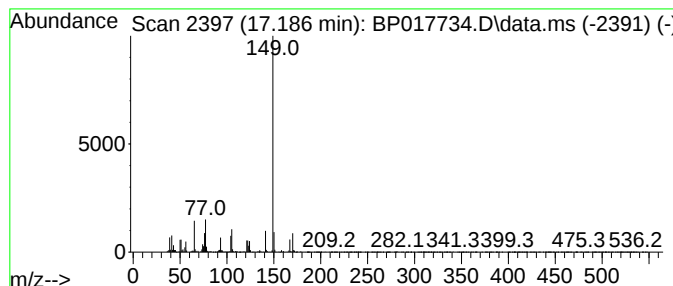
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Diethyl Phthalate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.186	5.63 ng/ul	332006	Phenanthrene-d10	17.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyl Phthalate	222	C12H14O4	000084-66-2	72
2			1,2-Benzenedicarboxylic acid, di...	250	C14H18O4	000131-16-8	64
3			Diethyl Phthalate	222	C12H14O4	000084-66-2	64
4			1,2-Benzenedicarboxylic acid, bi...	250	C14H18O4	000605-45-8	64
5			1,2-Benzenedicarboxylic acid, di...	250	C14H18O4	000131-16-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
Data File : BP017734.D
Acq On : 19 Oct 2023 19:14
Operator : MA/JU
Sample : 04864-03DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCJV6DL

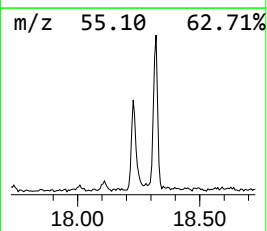
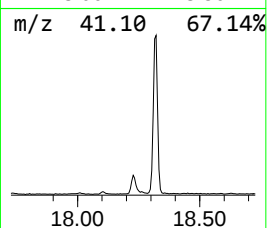
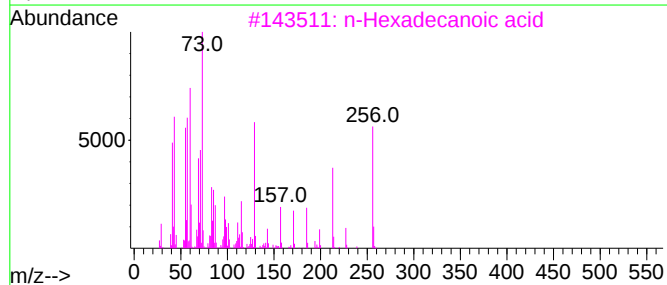
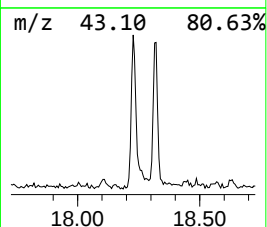
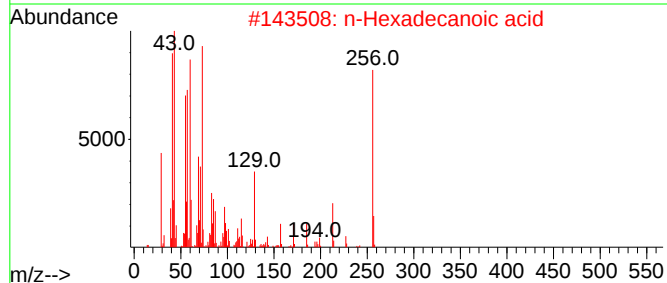
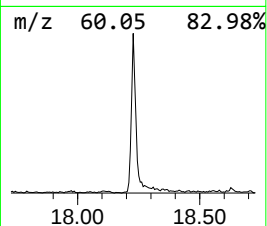
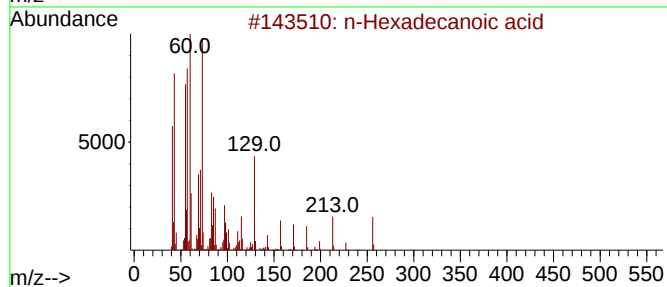
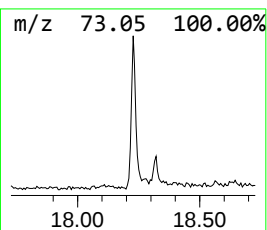
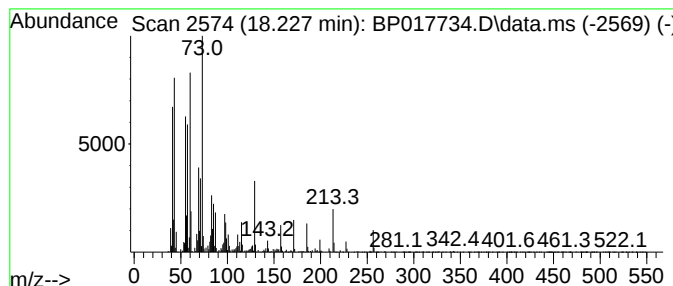
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	4.94 ng/ul	290879	Phenanthrene-d10	17.374

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
Data File : BP017734.D
Acq On : 19 Oct 2023 19:14
Operator : MA/JU
Sample : 04864-03DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCJV6DL

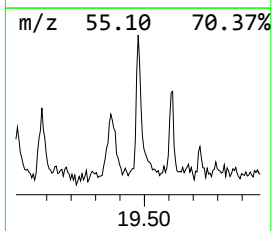
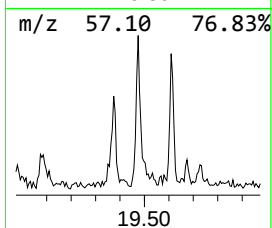
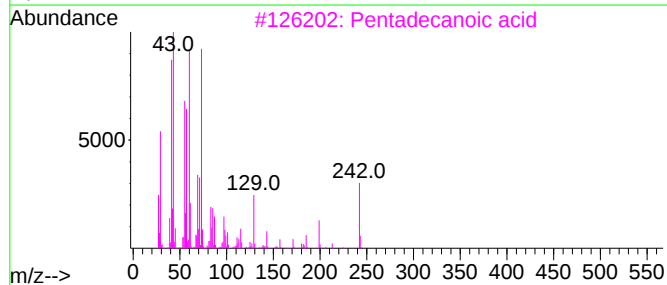
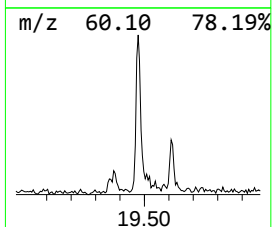
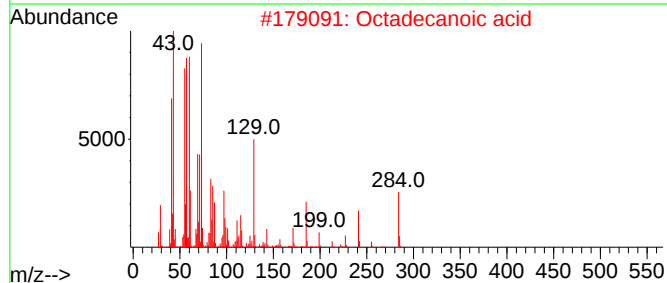
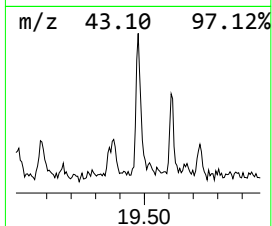
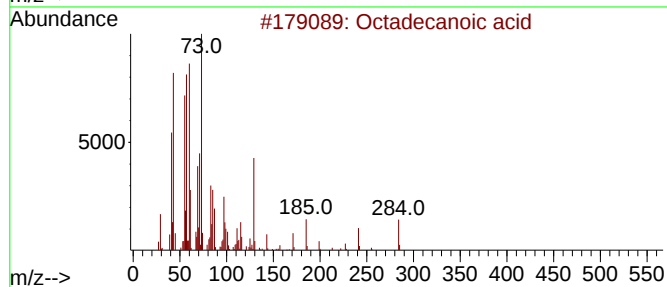
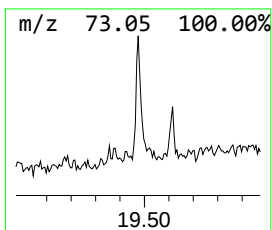
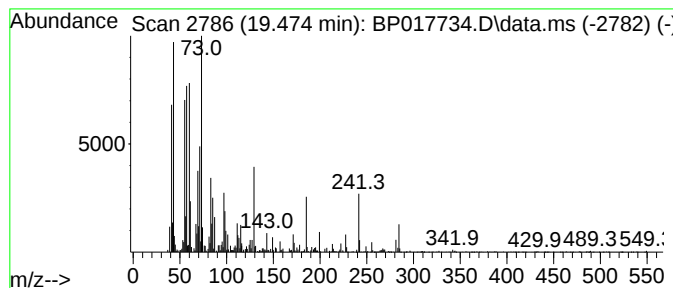
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.474	2.18 ng/ul	132743	Chrysene-d12	21.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	95
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
Data File : BP017734.D
Acq On : 19 Oct 2023 19:14
Operator : MA/JU
Sample : 04864-03DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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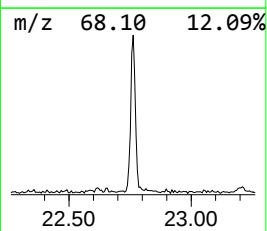
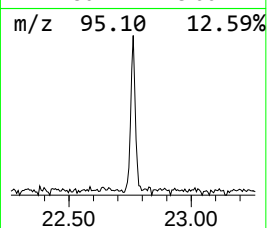
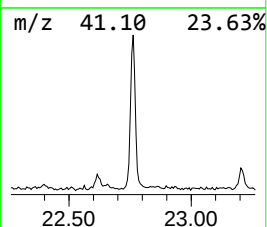
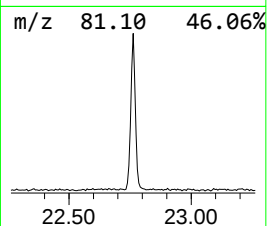
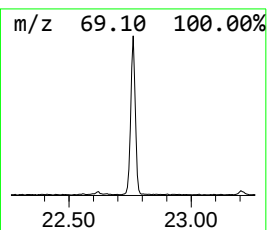
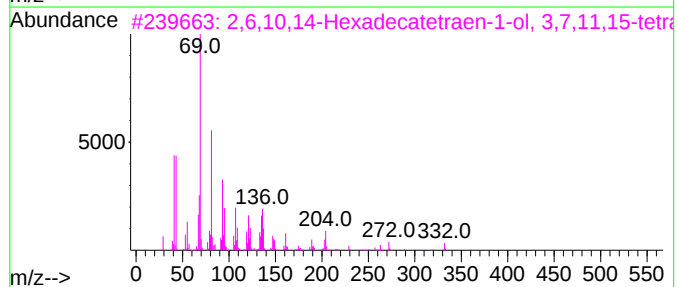
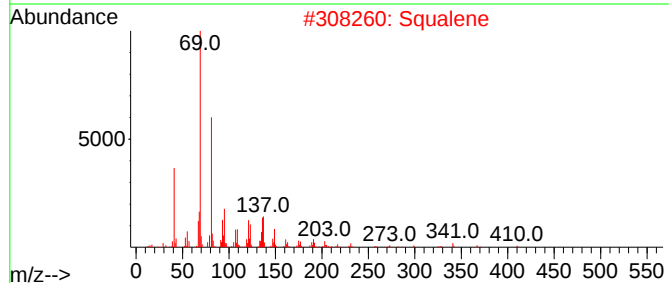
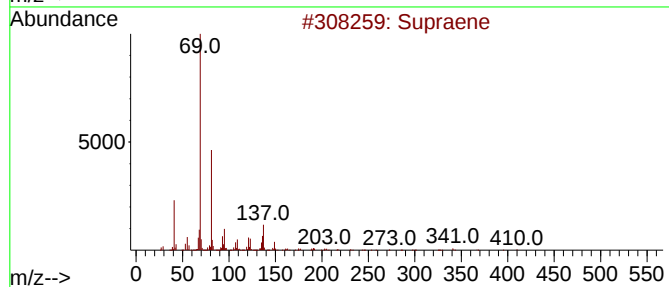
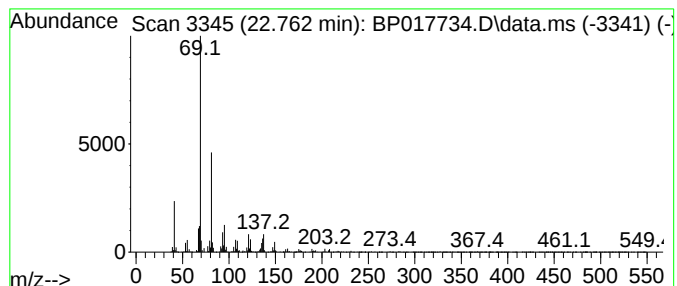
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Supraene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.762	6.42 ng/ul	390580	Chrysene-d12	21.515

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	90
3			2,6,10,14-Hexadecatetraen-1-ol, ...	332	C22H36O2	061691-98-3	78
4			2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	72
5			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101923\
Data File : BP017734.D
Acq On : 19 Oct 2023 19:14
Operator : MA/JU
Sample : 04864-03DL 2X
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101823.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexanoic acid	6.893	5.0	ng/ul	145845	1	7.887	579428	20.0
Phenol, 4-(1,1-...	13.410	2.4	ng/ul	120669	3	14.586	1003610	20.0
Vanillin	13.539	2.0	ng/ul	101784	3	14.586	1003610	20.0
Diethyl Phthalate	17.186	5.6	ng/ul	332006	4	17.374	1178600	20.0
n-Hexadecanoic ...	18.227	4.9	ng/ul	290879	4	17.374	1178600	20.0
Octadecanoic acid	19.474	2.2	ng/ul	132743	5	21.515	1216800	20.0
Supraene	22.762	6.4	ng/ul	390580	5	21.515	1216800	20.0