

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
 Data File : BM032462.D
 Acq On : 12 Oct 2021 17:44
 Operator : CG/JU
 Sample : M4110-01
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C00T3

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_M\METHODS\SFAM-EPA-BM100521.M
 Title : SVOA CALIBRATION

Signal : TIC: BM032462.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.954	17	20	29	rVB	66780	86587	1.24%	0.116%
2	3.254	69	71	79	rVB	14333	23400	0.33%	0.031%
3	3.390	90	94	110	rBV	272655	431920	6.17%	0.580%
4	3.507	110	114	122	rVV	49241	68581	0.98%	0.092%
5	3.578	124	126	131	rVV3	8886	11770	0.17%	0.016%
6	3.643	132	137	147	rVB2	41031	74576	1.06%	0.100%
7	3.731	148	152	160	rVB	35364	55227	0.79%	0.074%
8	3.831	165	169	174	rVB2	16924	22064	0.32%	0.030%
9	3.984	190	195	202	rBV	179794	229961	3.28%	0.309%
10	4.413	263	268	272	rBV	63730	82566	1.18%	0.111%
11	4.448	272	274	286	rVB3	12200	29478	0.42%	0.040%
12	4.595	294	299	303	rVB5	9056	16260	0.23%	0.022%
13	4.790	326	332	341	rBV	35738	57963	0.83%	0.078%
14	5.066	373	379	385	rBV2	16423	25676	0.37%	0.034%
15	5.689	481	485	491	rVB2	9675	13551	0.19%	0.018%
16	5.984	530	535	539	rBV	15933	25862	0.37%	0.035%
17	6.801	667	674	687	rVB	267653	462138	6.60%	0.621%
18	6.936	690	697	710	rBV	854707	1306700	18.66%	1.755%
19	7.148	725	733	740	rBV	865569	1313454	18.76%	1.764%
20	7.219	740	745	748	rVV	12456	19977	0.29%	0.027%
21	7.607	804	811	821	rBV	907757	1380591	19.71%	1.855%
22	7.689	821	825	836	rVB	13414	25542	0.36%	0.034%
23	8.336	926	935	942	rBV	755344	1151480	16.44%	1.547%
24	8.389	942	944	956	rVB4	18812	38277	0.55%	0.051%
25	8.760	1000	1007	1017	rBV	1022299	1619521	23.13%	2.176%
26	8.913	1028	1033	1035	rBV6	7851	12830	0.18%	0.017%
27	8.936	1035	1037	1043	rVB6	10650	16365	0.23%	0.022%
28	9.236	1085	1088	1093	rVB2	12318	17550	0.25%	0.024%
29	9.483	1123	1130	1137	rBV	806317	1275089	18.21%	1.713%
30	10.030	1215	1223	1234	rBV	1258050	2027581	28.95%	2.724%
31	10.177	1241	1248	1260	rBV	823225	1227655	17.53%	1.649%
32	10.395	1279	1285	1299	rVB	1319328	2151455	30.72%	2.890%
33	10.530	1299	1308	1328	rBV	911388	1571912	22.45%	2.112%
34	11.895	1533	1540	1548	rVB7	9038	16604	0.24%	0.022%
35	11.989	1550	1556	1560	rBV2	23085	38150	0.54%	0.051%
36	12.036	1560	1564	1571	rVV	25224	43156	0.62%	0.058%

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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_M\METHODS\SFAM-EPA-BM100521.M
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37	12.613	1657	1662	1666	rVB4	9351	14256	0.20%	0.019%
38	12.765	1685	1688	1693	rBV2	8692	12866	0.18%	0.017%
39	12.860	1699	1704	1713	rVB4	15052	24878	0.36%	0.033%
40	13.048	1731	1736	1741	rBV3	8228	15583	0.22%	0.021%

41	13.101	1741	1745	1751	rVB7	6877	12543	0.18%	0.017%
42	13.160	1751	1755	1760	rBV5	8009	12169	0.17%	0.016%
43	13.218	1760	1765	1771	rVV2	15130	23341	0.33%	0.031%
44	13.660	1834	1840	1849	rBV	2680017	3573390	51.03%	4.800%
45	13.948	1878	1889	1905	rBV	2976754	4190479	59.84%	5.629%

46	14.165	1921	1926	1930	rBV3	10944	14191	0.20%	0.019%
47	14.259	1935	1942	1953	rVB	2080683	3071326	43.86%	4.126%
48	14.401	1961	1966	1971	rBV3	10580	15731	0.22%	0.021%
49	14.465	1971	1977	1990	rBV	229325	374598	5.35%	0.503%
50	14.595	1994	1999	2002	rBV7	7583	12552	0.18%	0.017%

51	14.671	2007	2012	2019	rBV8	14499	26922	0.38%	0.036%
52	14.830	2035	2039	2046	rVB3	11153	18753	0.27%	0.025%
53	14.995	2062	2067	2073	rBV	32121	41045	0.59%	0.055%
54	15.071	2073	2080	2084	rVV2	61115	87482	1.25%	0.118%
55	15.112	2084	2087	2092	rVB	14620	18995	0.27%	0.026%

56	15.248	2103	2110	2118	rBV	3828229	5444257	77.74%	7.313%
57	15.359	2123	2129	2142	rBV	1185883	1584485	22.63%	2.128%
58	15.483	2145	2150	2154	rBV	48157	73559	1.05%	0.099%
59	15.683	2177	2184	2188	rBV5	17629	29275	0.42%	0.039%
60	15.818	2199	2207	2211	rVB9	7781	16379	0.23%	0.022%

61	15.883	2215	2218	2221	rVV4	15640	21072	0.30%	0.028%
62	15.924	2221	2225	2229	rVV6	12694	27200	0.39%	0.037%
63	15.971	2229	2233	2235	rVV3	28717	42554	0.61%	0.057%
64	15.995	2235	2237	2240	rVV	32601	40415	0.58%	0.054%
65	16.030	2240	2243	2248	rVB5	24306	37537	0.54%	0.050%

66	16.136	2255	2261	2267	rBV5	10943	24656	0.35%	0.033%
67	16.242	2275	2279	2283	rBV2	13350	18959	0.27%	0.025%
68	16.395	2302	2305	2314	rVB3	26606	39153	0.56%	0.053%
69	16.483	2314	2320	2322	rBV5	11324	20392	0.29%	0.027%
70	16.736	2359	2363	2365	rBV5	10679	13598	0.19%	0.018%

71	16.806	2372	2375	2379	rVB3	26108	35479	0.51%	0.048%
72	16.983	2399	2405	2411	rBV	2727044	3718023	53.09%	4.995%
73	17.036	2411	2414	2417	rVV	379940	459930	6.57%	0.618%
74	17.083	2417	2422	2432	rVB	4385150	6022851	86.00%	8.091%
75	17.212	2440	2444	2451	rVB3	56007	82049	1.17%	0.110%

76	17.353	2461	2468	2473	rBV	120884	160442	2.29%	0.216%
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Integrator: RTE
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 Stop Thrs : 0 Peak Location: TOP

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77	17.524	2493	2497	2501	rBV2	10105	17088	0.24%	0.023%
78	17.653	2514	2519	2523	rBV	144761	178364	2.55%	0.240%
79	17.765	2528	2538	2539	rBV2	109554	178805	2.55%	0.240%
80	17.795	2539	2543	2552	rVV	343981	560370	8.00%	0.753%
81	17.883	2552	2558	2582	rVB	1378563	2136592	30.51%	2.870%
82	18.495	2659	2662	2665	rBV3	41617	47323	0.68%	0.064%
83	18.536	2665	2669	2674	rVV	247285	277932	3.97%	0.373%
84	18.665	2687	2691	2699	rVB2	71090	97972	1.40%	0.132%
85	18.777	2706	2710	2712	rBV3	14875	21473	0.31%	0.029%
86	18.818	2715	2717	2725	rVB7	29101	36868	0.53%	0.050%
87	18.894	2725	2730	2734	rBV	801512	929833	13.28%	1.249%
88	18.930	2734	2736	2742	rVB2	54178	71521	1.02%	0.096%
89	19.000	2745	2748	2752	rVV	52369	70366	1.00%	0.095%
90	19.094	2752	2764	2770	rVV	351208	749827	10.71%	1.007%
91	19.153	2770	2774	2788	rVB	208810	370939	5.30%	0.498%
92	19.371	2805	2811	2817	rBV	5295244	7003183	100.00%	9.408%
93	19.877	2894	2897	2900	rBV4	21600	24540	0.35%	0.033%
94	19.912	2900	2903	2906	rVV	17561	19659	0.28%	0.026%
95	20.853	3059	3063	3069	rBV	582567	681401	9.73%	0.915%
96	21.147	3108	3113	3123	rBV	3183238	4058410	57.95%	5.452%
97	21.288	3135	3137	3141	rVB2	62284	56148	0.80%	0.075%
98	22.612	3359	3362	3366	rVB	184989	230616	3.29%	0.310%
99	23.159	3447	3455	3465	rVB2	3623330	6451046	92.12%	8.666%
100	23.294	3472	3478	3487	rVB2	2128016	3726315	53.21%	5.006%

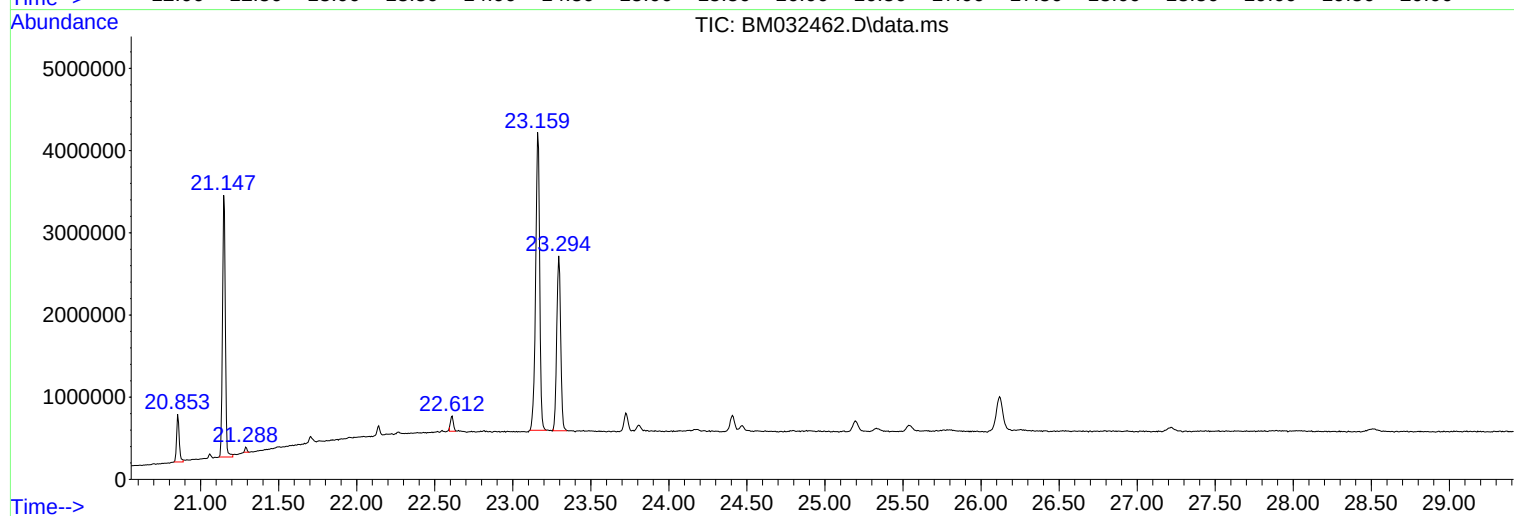
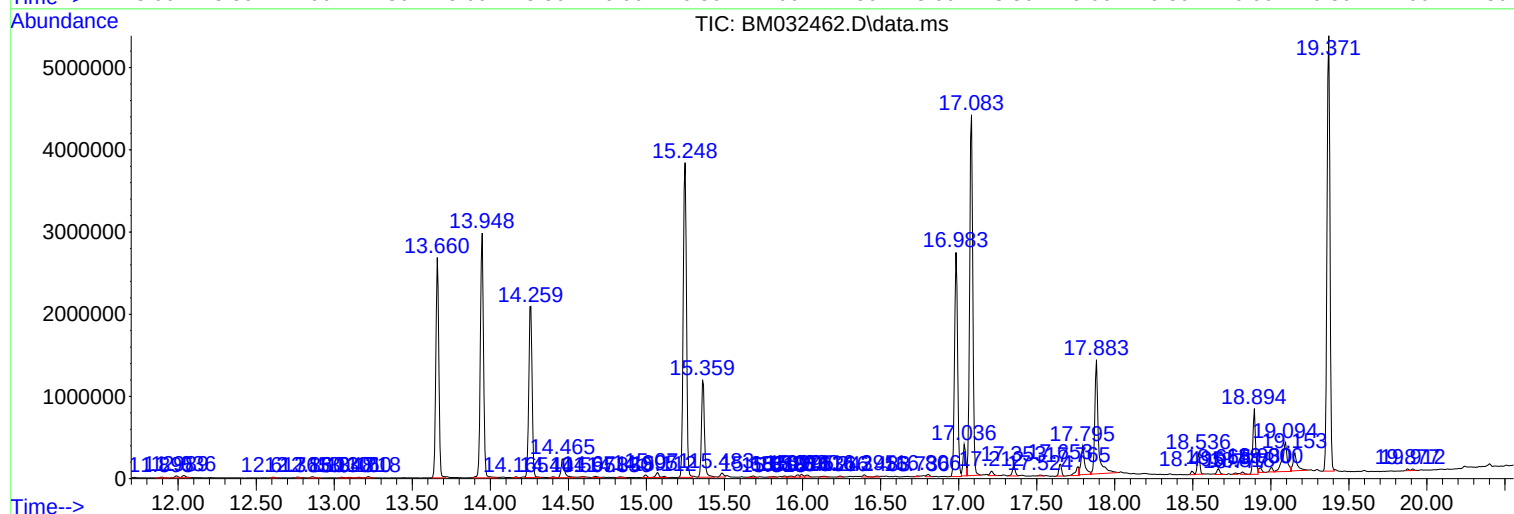
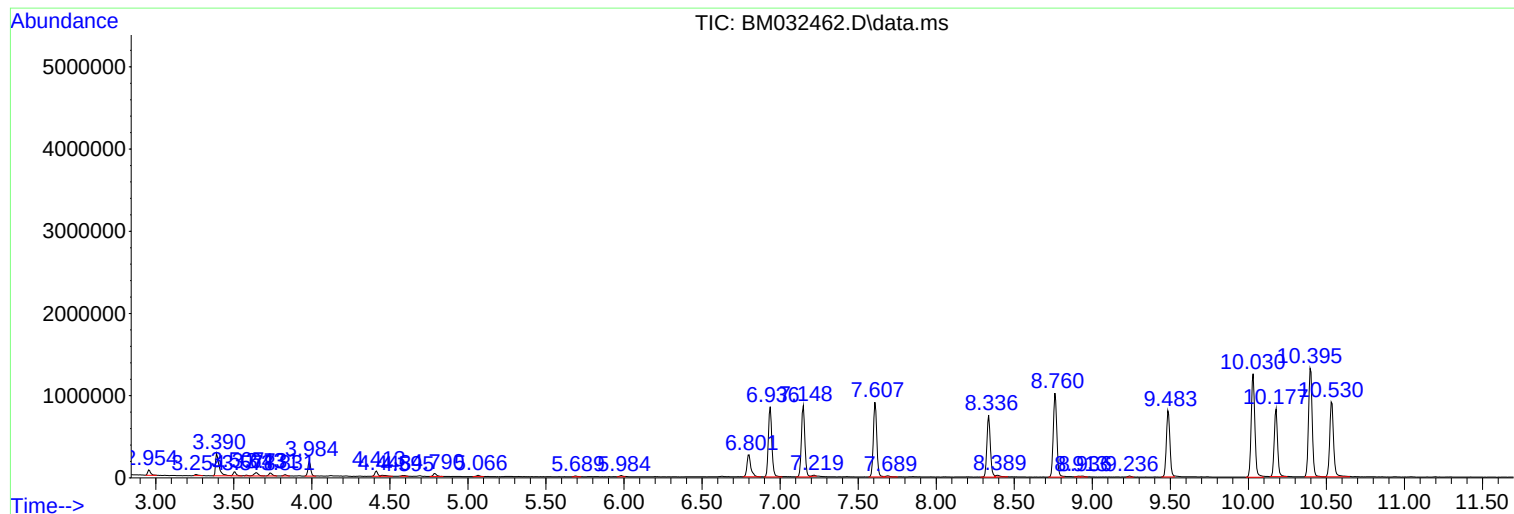
Sum of corrected areas: 74441425

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
Quant Title : SVOA CALIBRATION

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



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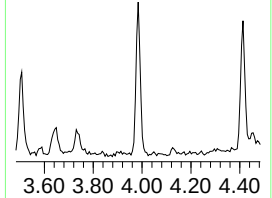
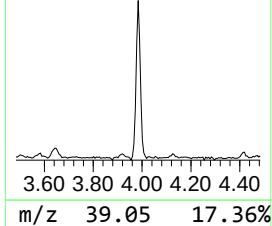
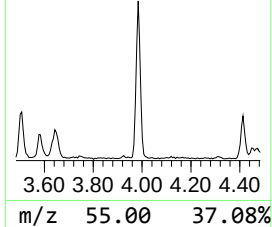
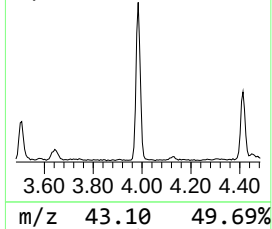
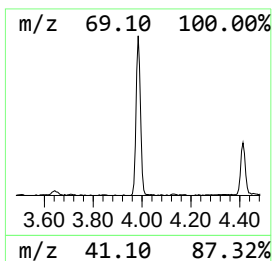
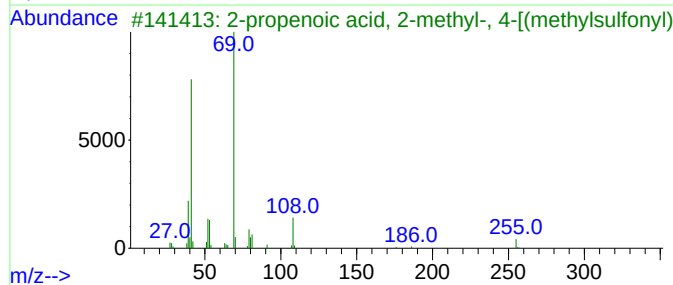
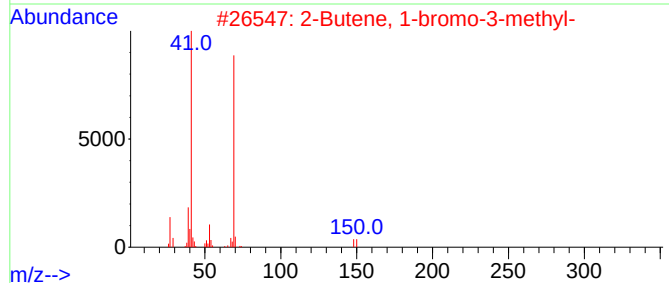
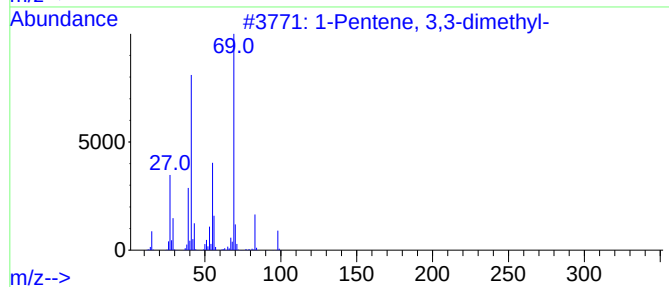
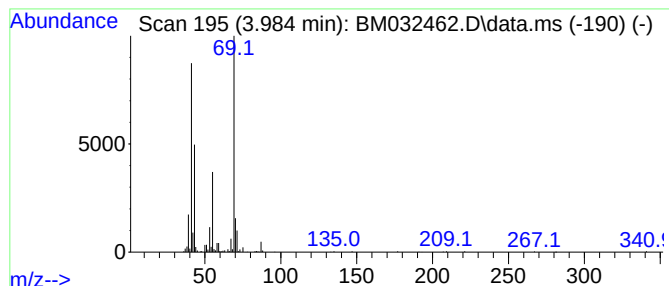
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.984	3.33 ng/ul	229961	1,4-Dichlorobenzene-d4	7.607

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Pentene, 3,3-dimethyl-	98	C7H14	003404-73-7	64
2		2-Butene, 1-bromo-3-methyl-	148	C5H9Br	000870-63-3	45
3		2-propenoic acid, 2-methyl-, 4-[...]	255	C11H13NO4S	1000401-46-9	45
4		1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	45
5		3-Butyn-2-ol, 2-methyl-	84	C5H8O	000115-19-5	43



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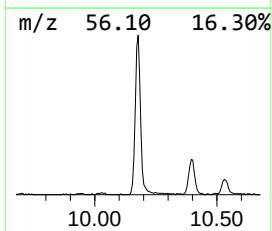
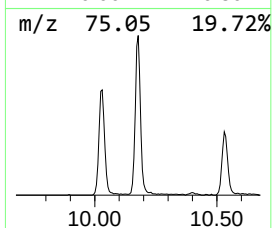
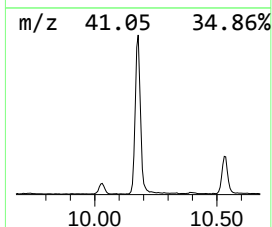
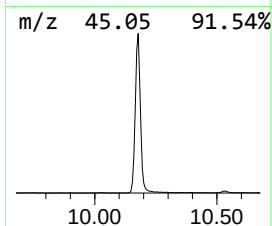
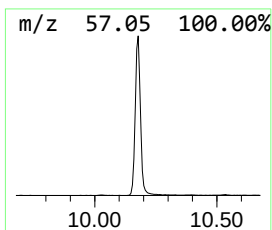
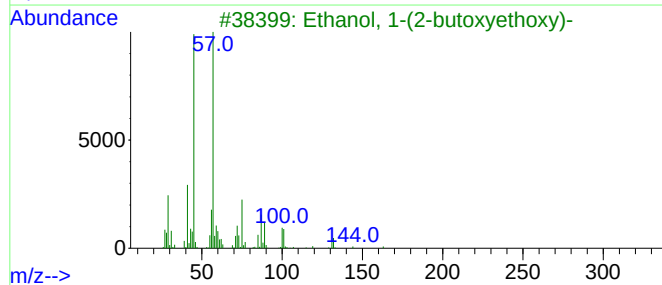
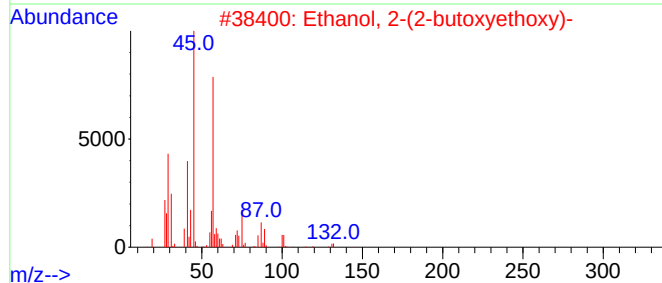
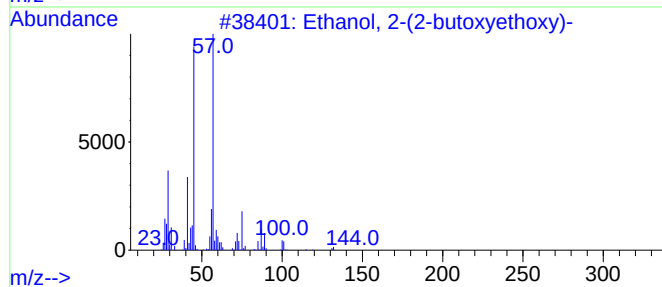
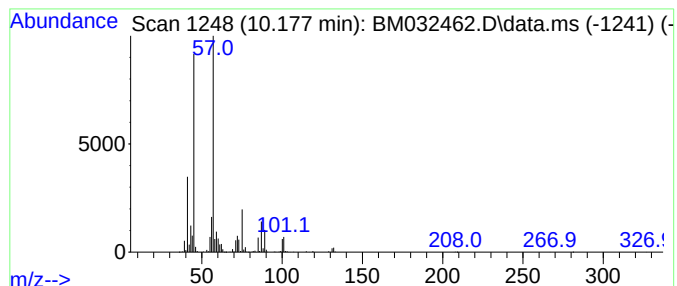
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.177	11.41 ng/ul	1227660	Naphthalene-d8	10.395

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
3			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



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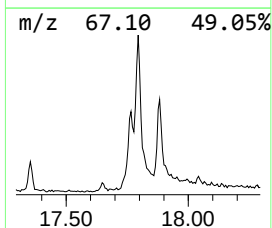
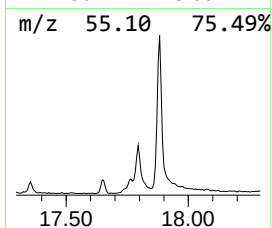
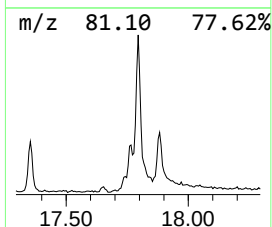
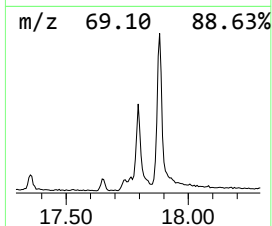
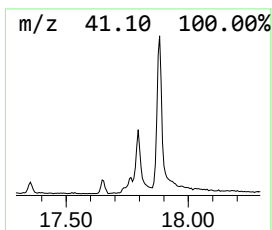
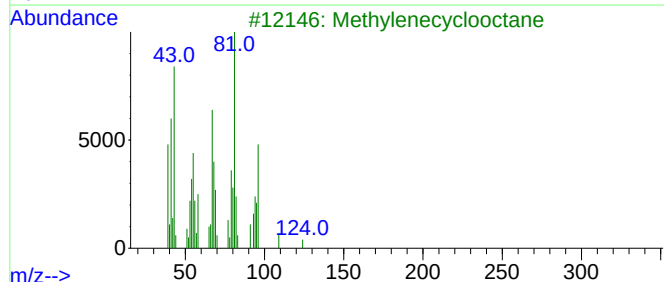
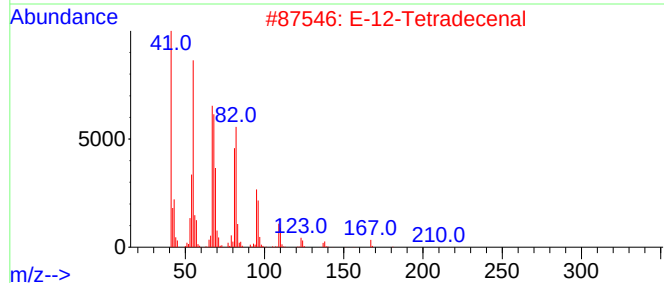
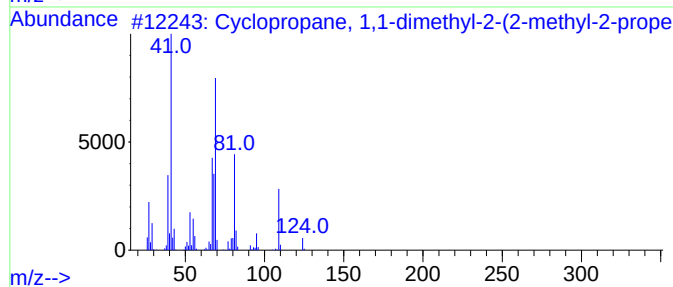
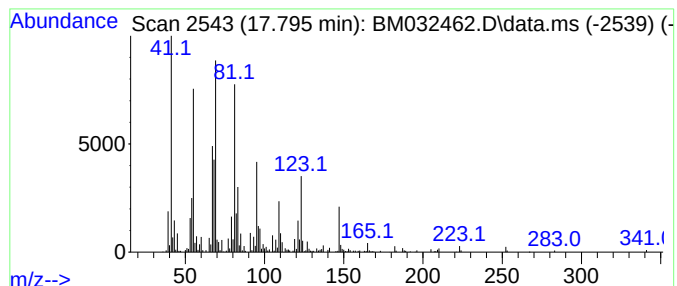
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.795	3.01 ng/ul	560370	Phenanthrene-d10	16.983

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	45
2			E-12-Tetradecenal	210	C14H26O	1000130-96-1	41
3			Methylenecyclooctane	124	C9H16	003618-18-6	38
4			Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	38
5			7,8-Diazabicyclo[4.2.0]octane, 1...	156	C7H12N2O2	169522-32-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
Data File : BM032462.D
Acq On : 12 Oct 2021 17:44
Operator : CG/JU
Sample : M4110-01
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C00T3

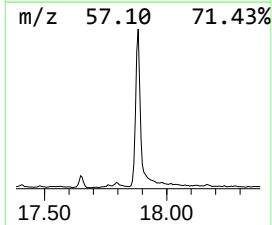
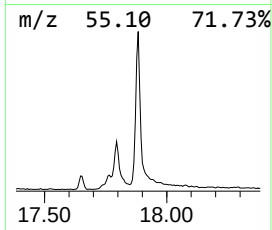
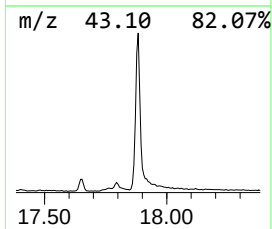
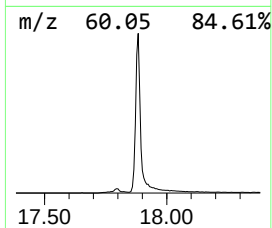
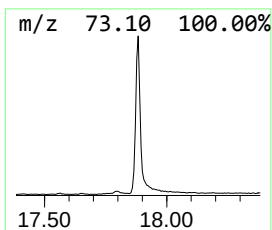
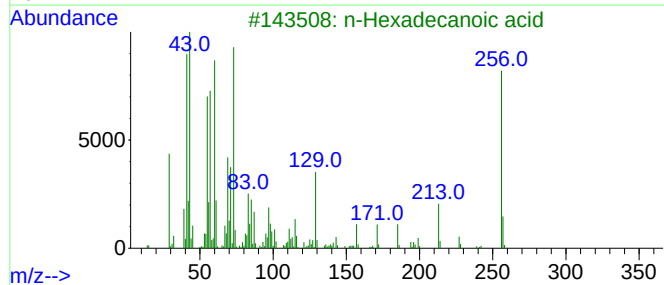
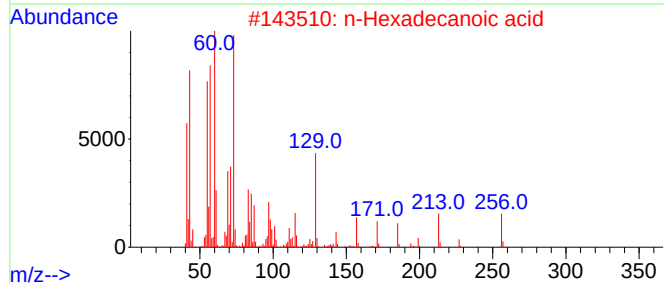
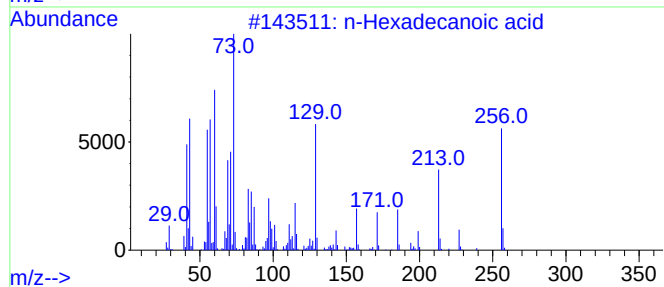
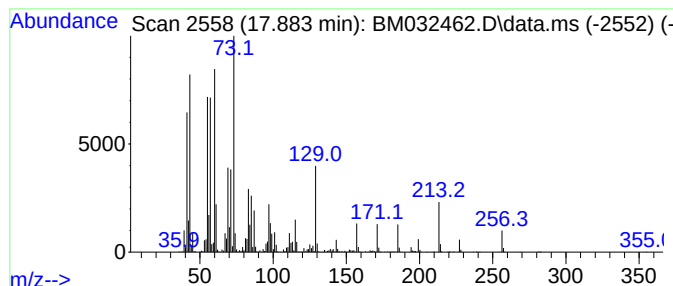
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.883	11.49 ng/ul	2136590	Phenanthrene-d10	16.983

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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
Data File : BM032462.D
Acq On : 12 Oct 2021 17:44
Operator : CG/JU
Sample : M4110-01
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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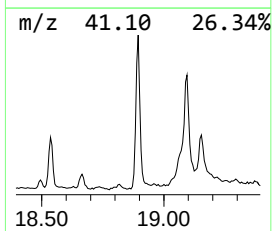
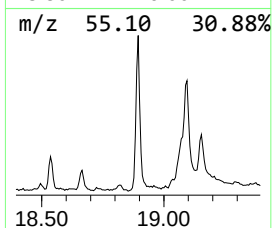
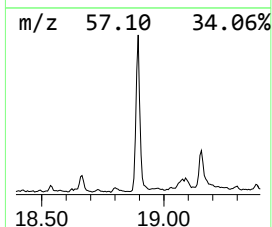
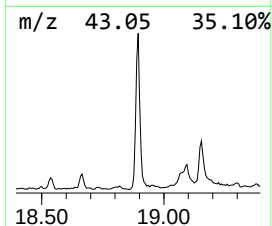
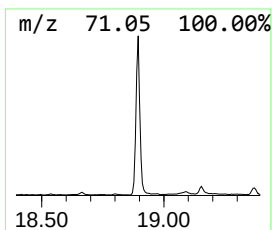
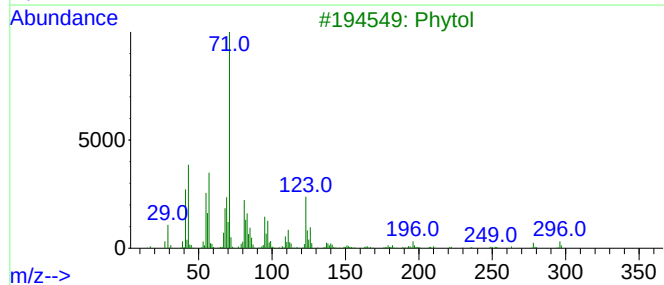
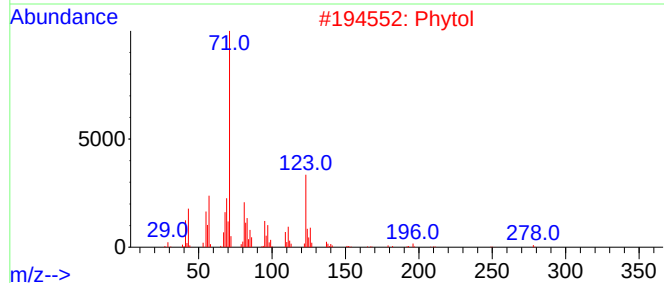
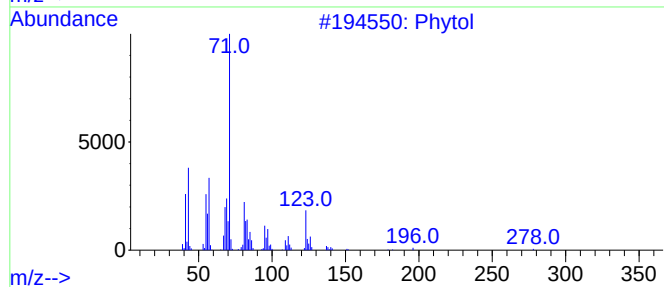
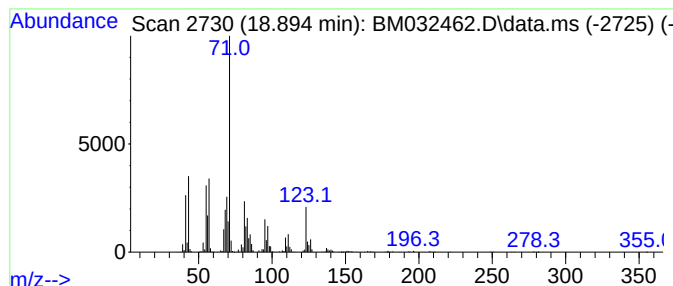
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phytol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.895	5.00 ng/ul	929833	Phenanthrene-d10	16.983

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phytol			296	C20H40O	000150-86-7	91
2	Phytol			296	C20H40O	000150-86-7	91
3	Phytol			296	C20H40O	000150-86-7	83
4	1-Hexadecen-3-ol, 3,5,11,15-tetr...			296	C20H40O	1000262-55-5	59
5	Isophytol			296	C20H40O	000505-32-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
Data File : BM032462.D
Acq On : 12 Oct 2021 17:44
Operator : CG/JU
Sample : M4110-01
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C00T3

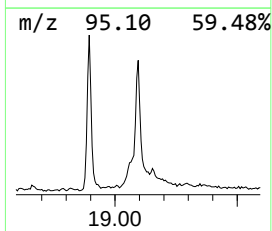
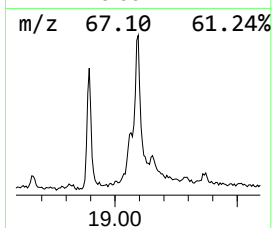
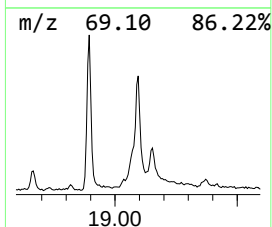
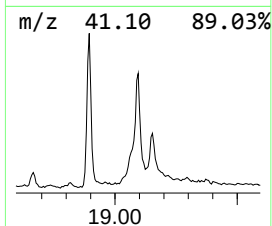
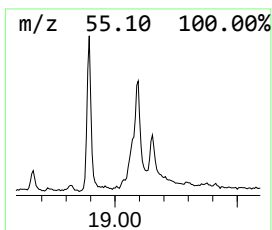
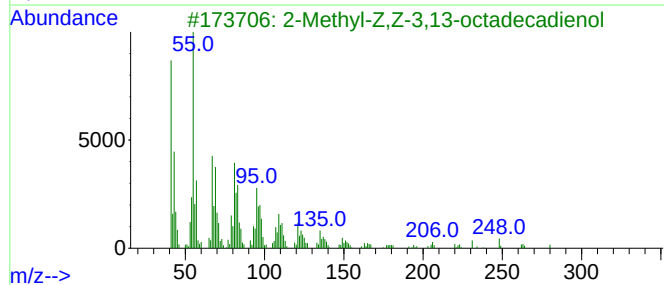
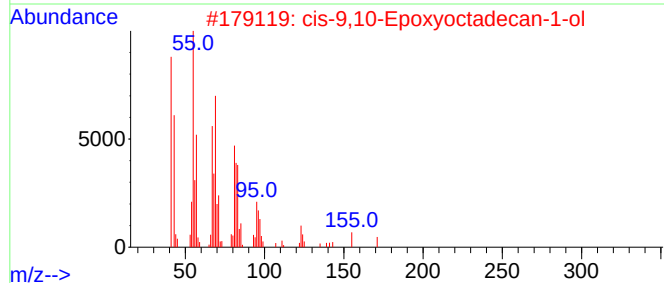
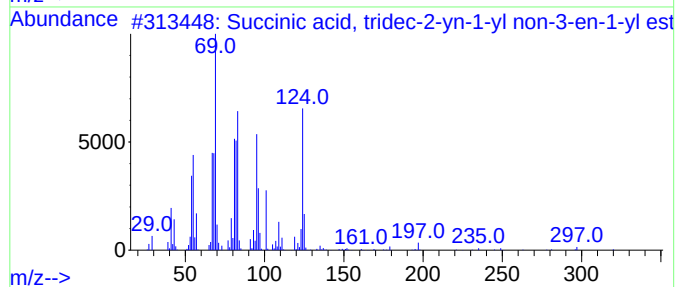
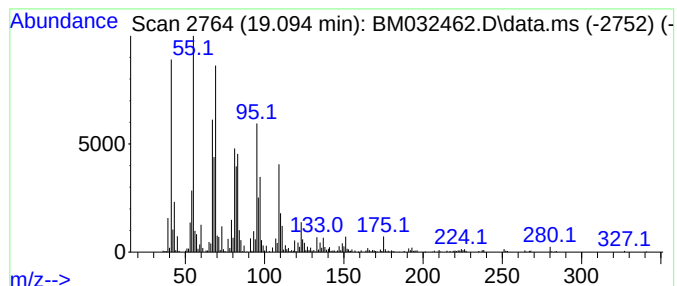
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Succinic acid, tridec-2-yn-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.095	3.70 ng/ul	749827	Chrysene-d12	21.147

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Succinic acid, tridec-2-yn-1-yl ...	420	C26H44O4	1010391-10-5	64
2			cis-9,10-Epoxyoctadecan-1-ol	284	C18H36O2	013980-12-6	53
3			2-Methyl-Z,Z-3,13-octadecadienol	280	C19H36O	1000130-90-5	53
4			11-Dodecen-1-ol trifluoroacetate	280	C14H23F3O2	128792-46-1	49
5			Pentanoic acid, 10-undecenyl ester	254	C16H30O2	1000159-93-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
Data File : BM032462.D
Acq On : 12 Oct 2021 17:44
Operator : CG/JU
Sample : M4110-01
Misc :
ALS Vial : 48 Sample Multiplier: 1

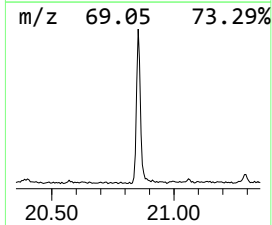
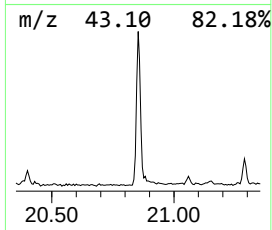
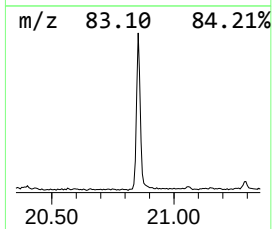
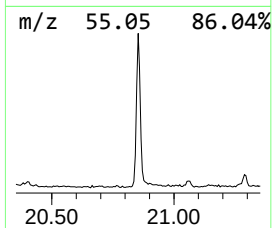
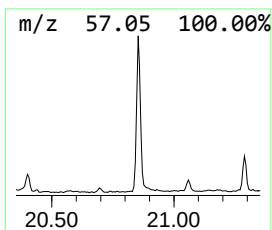
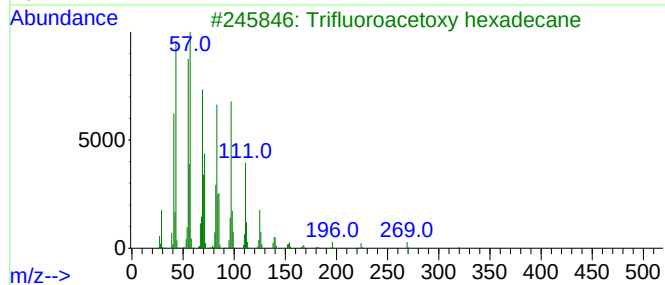
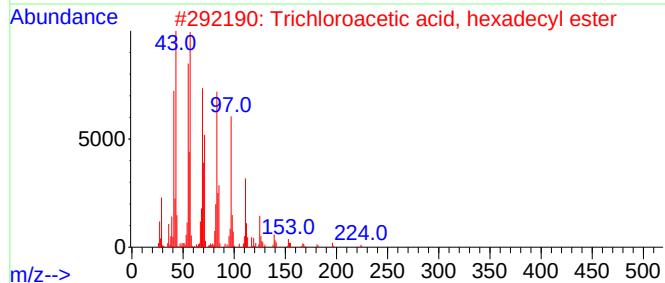
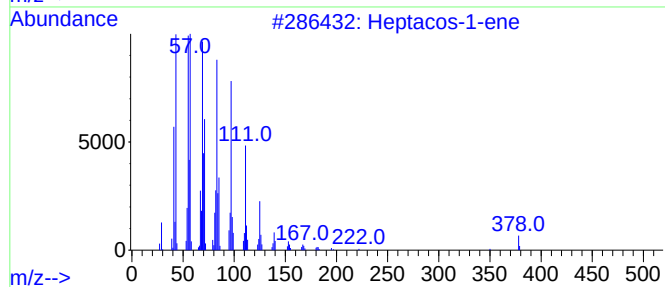
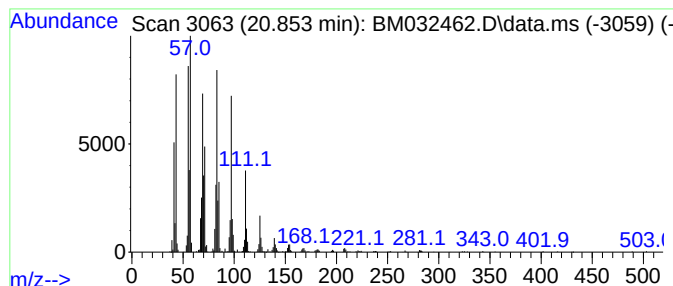
Instrument :
BNA_M
ClientSampleId :
C00T3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.853	3.36 ng/ul	681401	Chrysene-d12	21.147	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacos-1-ene	378	C27H54	015306-27-1	94
2	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
4	Pentacos-1-ene	350	C25H50	016980-85-1	94
5	n-Tetracosanol-1	354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101121\
Data File : BM032462.D
Acq On : 12 Oct 2021 17:44
Operator : CG/JU
Sample : M4110-01
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C00T3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM100521.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
(DEL) Alkane: S...	3.984	3.3	ng/ul	229961	1	7.607	1380590	20.0
Ethanol, 2-(2-b...	10.177	11.4	ng/ul	1227660	2	10.395	2151460	20.0
unknown-01	17.795	3.0	ng/ul	560370	4	16.983	3718020	20.0
n-Hexadecanoic ...	17.883	11.5	ng/ul	2136590	4	16.983	3718020	20.0
Phytol	18.895	5.0	ng/ul	929833	4	16.983	3718020	20.0
Succinic acid, ...	19.095	3.7	ng/ul	749827	5	21.147	4058410	20.0
(DEL) Alkane: S...	20.853	3.4	ng/ul	681401	5	21.147	4058410	20.0