

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
 Data File : BP015649.D
 Acq On : 22 Jun 2023 16:13
 Operator : MA/JU
 Sample : 03083-15
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFL3

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015649.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	43	rVB	30015	39761	1.56%	0.102%
2	3.846	109	112	126	rBV	206045	338030	13.26%	0.864%
3	3.952	126	130	135	rVB3	21928	34720	1.36%	0.089%
4	4.069	145	150	158	rVB2	21558	36728	1.44%	0.094%
5	4.169	164	167	176	rVB3	10746	18398	0.72%	0.047%
6	4.411	205	208	211	rVB5	5424	6046	0.24%	0.015%
7	4.564	232	234	239	rVB6	5276	6663	0.26%	0.017%
8	5.122	326	329	335	rBV4	5206	10360	0.41%	0.026%
9	5.675	416	423	426	rBV9	3929	8247	0.32%	0.021%
10	5.734	431	433	438	rVB5	2444	3821	0.15%	0.010%
11	6.046	483	486	487	rBV2	4697	4498	0.18%	0.012%
12	6.299	524	529	533	rBV2	9554	17839	0.70%	0.046%
13	6.705	596	598	601	rVV4	4304	5720	0.22%	0.015%
14	7.240	681	689	706	rBV	444936	822746	32.26%	2.104%
15	7.399	710	716	728	rVV	367363	614982	24.12%	1.572%
16	7.605	744	751	760	rBV	501785	843240	33.07%	2.156%
17	7.675	760	763	767	rVB5	9333	14095	0.55%	0.036%
18	8.022	819	822	824	rBV3	3176	4022	0.16%	0.010%
19	8.081	824	832	843	rVV	526001	892962	35.02%	2.283%
20	8.205	851	853	859	rVB7	3081	4340	0.17%	0.011%
21	8.293	863	868	870	rBV5	4892	6003	0.24%	0.015%
22	8.552	906	912	913	rBV5	3058	4450	0.17%	0.011%
23	8.728	933	942	946	rVB5	3146	5650	0.22%	0.014%
24	8.799	946	954	972	rBV	463944	932767	36.58%	2.385%
25	9.257	1025	1032	1045	rBV	486256	888077	34.82%	2.271%
26	9.346	1045	1047	1053	rVV7	4861	9907	0.39%	0.025%
27	9.416	1053	1059	1068	rVB7	10264	24626	0.97%	0.063%
28	9.628	1091	1095	1099	rVV3	7753	11863	0.47%	0.030%
29	9.987	1149	1156	1175	rBV	415772	786580	30.84%	2.011%
30	10.181	1183	1189	1192	rBV8	3592	6965	0.27%	0.018%
31	10.310	1205	1211	1215	rVB9	2383	5018	0.20%	0.013%
32	10.534	1241	1249	1269	rBV	505416	1089195	42.71%	2.785%
33	10.857	1300	1304	1306	rBV5	5809	9504	0.37%	0.024%
34	10.904	1306	1312	1320	rVV	751657	1291247	50.63%	3.301%
35	11.057	1331	1338	1361	rBV	250014	660213	25.89%	1.688%
36	11.746	1450	1455	1458	rBV7	2852	5062	0.20%	0.013%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	11.910	1476	1483	1487	rBV4	7905	12572	0.49%	0.032%
38	12.304	1546	1550	1556	rBV7	8190	15116	0.59%	0.039%
39	12.498	1577	1583	1591	rBV3	7810	17098	0.67%	0.044%
40	12.575	1591	1596	1601	rVB5	8034	12906	0.51%	0.033%

41	12.693	1610	1616	1617	rBV6	2574	4648	0.18%	0.012%
42	12.781	1627	1631	1639	rBV8	5112	10047	0.39%	0.026%
43	12.922	1652	1655	1659	rBV6	3216	4989	0.20%	0.013%
44	13.240	1706	1709	1714	rVB6	8015	10793	0.42%	0.028%
45	13.316	1721	1722	1730	rVB8	3885	5402	0.21%	0.014%

46	13.481	1746	1750	1754	rBV7	3832	5449	0.21%	0.014%
47	13.528	1754	1758	1762	rVV2	14347	19861	0.78%	0.051%
48	13.604	1762	1771	1778	rVB8	16435	41143	1.61%	0.105%
49	13.710	1784	1789	1794	rVB9	6581	10938	0.43%	0.028%
50	13.887	1817	1819	1821	rBV3	3054	3667	0.14%	0.009%

51	13.916	1821	1824	1828	rVB6	6199	9545	0.37%	0.024%
52	13.992	1831	1837	1840	rBV8	3418	5450	0.21%	0.014%
53	14.045	1840	1846	1852	rBV3	32666	55205	2.16%	0.141%
54	14.134	1852	1861	1868	rBV	1150672	1645580	64.53%	4.207%
55	14.234	1876	1878	1883	rVB6	6266	9880	0.39%	0.025%

56	14.422	1903	1910	1925	rBV	1250289	1925137	75.49%	4.922%
57	14.551	1926	1932	1935	rVV8	7126	12670	0.50%	0.032%
58	14.604	1937	1941	1949	rVB3	15076	27078	1.06%	0.069%
59	14.675	1949	1953	1956	rBV4	16586	26697	1.05%	0.068%
60	14.728	1956	1962	1969	rVV	1125462	1672609	65.59%	4.276%

61	14.781	1969	1971	1980	rVB8	15009	30798	1.21%	0.079%
62	14.963	1994	2002	2024	rBV2	184071	608989	23.88%	1.557%
63	15.139	2027	2032	2041	rVB2	17187	53017	2.08%	0.136%
64	15.269	2050	2054	2061	rVB9	17980	31982	1.25%	0.082%
65	15.345	2065	2067	2072	rVB6	5116	7004	0.27%	0.018%

66	15.510	2091	2095	2098	rVV6	7884	10637	0.42%	0.027%
67	15.551	2098	2102	2106	rVB3	20747	24157	0.95%	0.062%
68	15.722	2124	2131	2138	rBV	1547780	2228883	87.40%	5.699%
69	15.863	2149	2155	2165	rBV	276674	520378	20.41%	1.330%
70	16.086	2187	2193	2204	rBV	410407	595927	23.37%	1.524%

71	16.345	2235	2237	2243	rVB6	18039	21427	0.84%	0.055%
72	16.439	2245	2253	2260	rBV5	29218	80187	3.14%	0.205%
73	16.581	2272	2277	2283	rBV9	16887	35523	1.39%	0.091%
74	16.651	2286	2289	2294	rVB5	9901	15879	0.62%	0.041%
75	16.875	2324	2327	2331	rVB6	9449	13556	0.53%	0.035%

76	17.216	2380	2385	2390	rBV3	55490	78765	3.09%	0.201%
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77	17.363	2406	2410	2418	rBV4	60374	105564	4.14%	0.270%
78	17.433	2418	2422	2425	rVV4	40791	59142	2.32%	0.151%
79	17.486	2425	2431	2436	rVV	1160387	1703668	66.81%	4.356%
80	17.533	2436	2439	2443	rVV	138029	190325	7.46%	0.487%
81	17.586	2443	2448	2460	rVB	1290420	1934585	75.86%	4.946%
82	17.904	2500	2502	2507	rVB3	14204	16747	0.66%	0.043%
83	17.951	2507	2510	2515	rBV2	23830	28215	1.11%	0.072%
84	18.233	2553	2558	2563	rBV3	52510	91628	3.59%	0.234%
85	18.292	2564	2568	2573	rBV	129689	182806	7.17%	0.467%
86	18.345	2573	2577	2584	rBV	648686	977727	38.34%	2.500%
87	18.628	2621	2625	2626	rBV	35063	51579	2.02%	0.132%
88	19.227	2724	2727	2733	rVB2	58173	75550	2.96%	0.193%
89	19.486	2760	2771	2779	rBV	331846	650649	25.51%	1.664%
90	19.575	2782	2786	2794	rVV2	144671	242184	9.50%	0.619%
91	19.816	2822	2827	2831	rBV	1390650	1994573	78.21%	5.100%
92	21.251	3067	3071	3076	rVB2	382028	563554	22.10%	1.441%
93	21.451	3102	3105	3111	rVB	305832	342221	13.42%	0.875%
94	21.574	3121	3126	3131	rBV2	945972	1397651	54.81%	3.573%
95	23.286	3413	3417	3423	rVB	456734	728103	28.55%	1.862%
96	23.963	3526	3532	3539	rBV2	1024614	1987876	77.95%	5.083%
97	24.127	3555	3560	3568	rVB	643353	1300149	50.98%	3.324%
98	24.627	3639	3645	3654	rVB2	1132695	2550117	100.00%	6.520%
99	24.745	3659	3665	3677	rVB	617213	1391618	54.57%	3.558%
100	26.204	3908	3913	3924	rVB2	450994	1159470	45.47%	2.965%

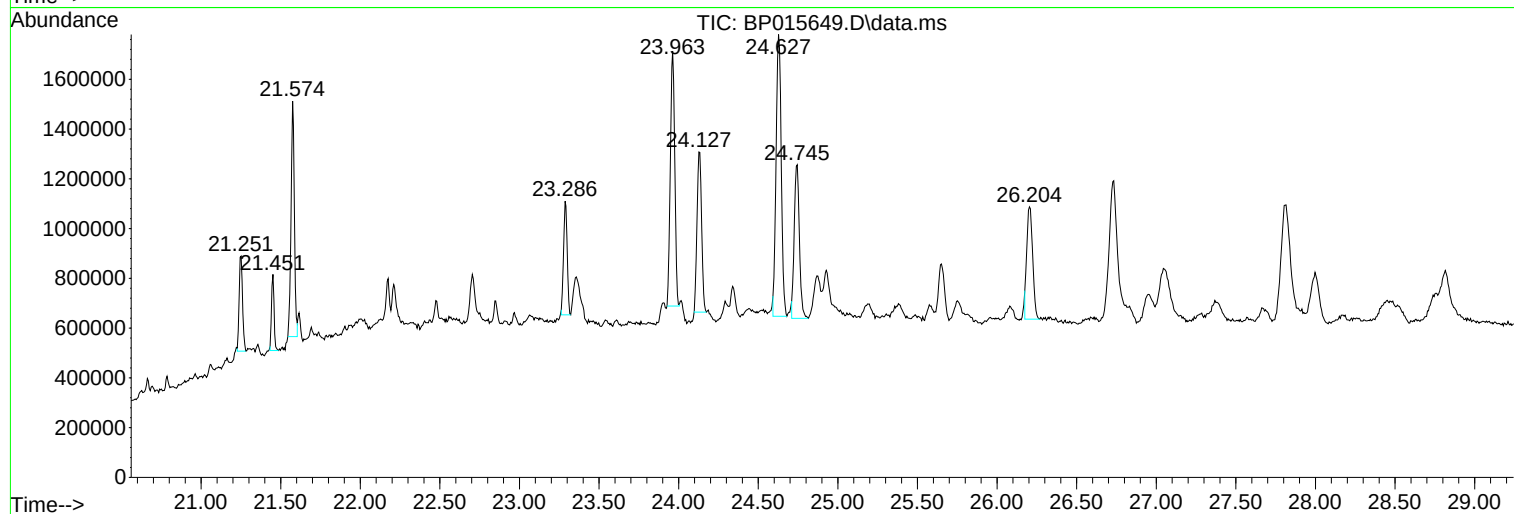
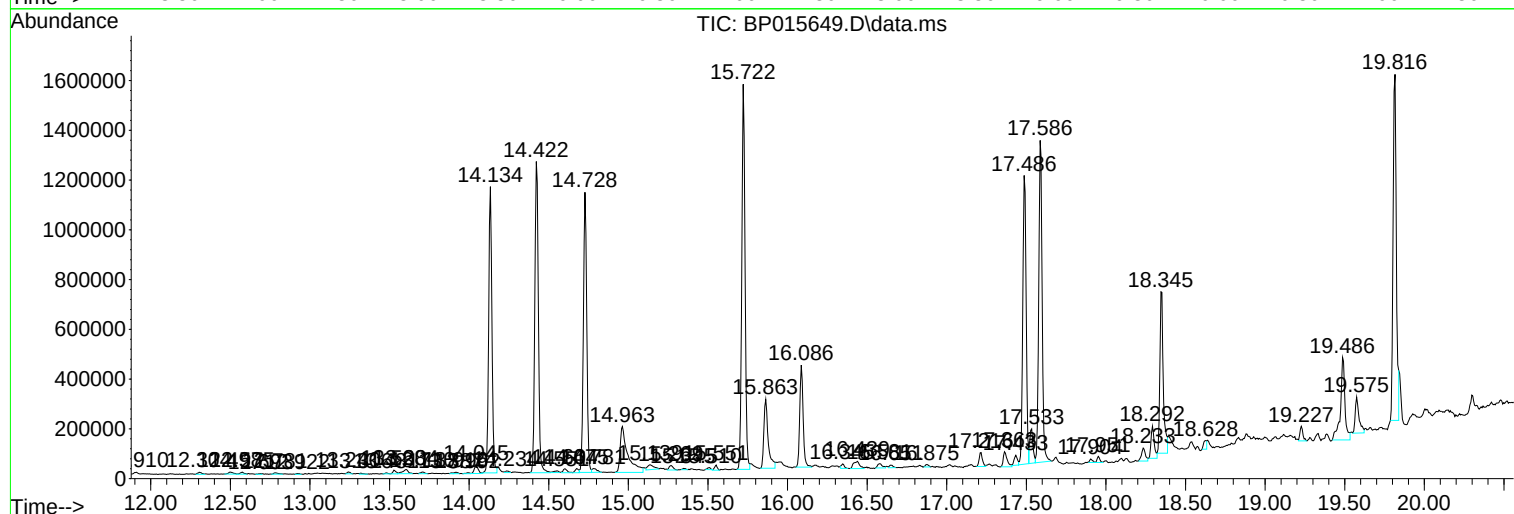
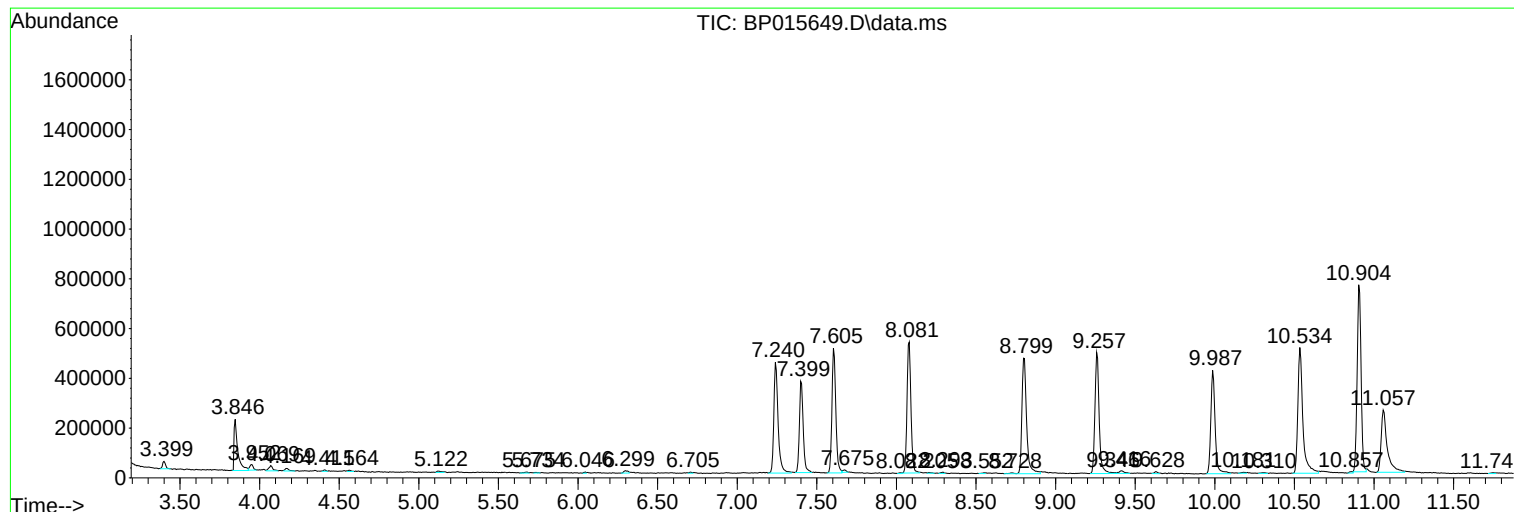
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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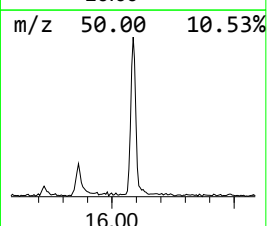
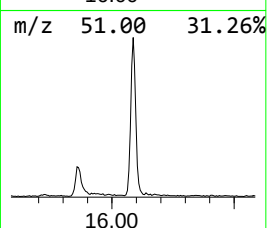
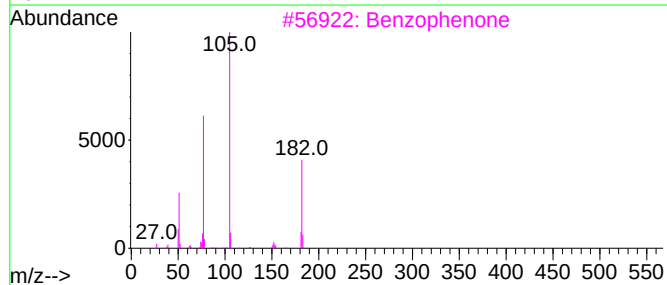
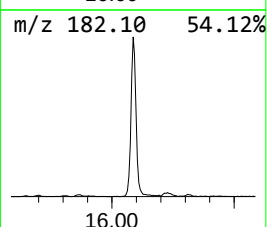
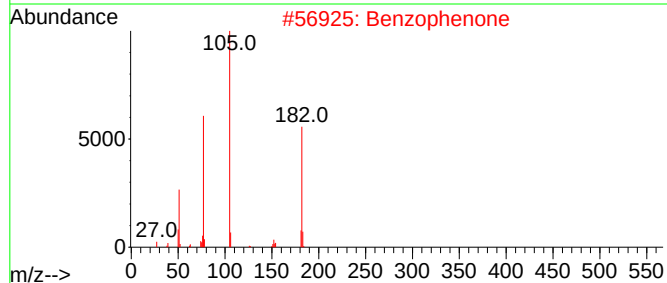
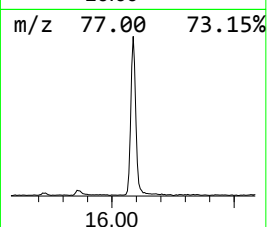
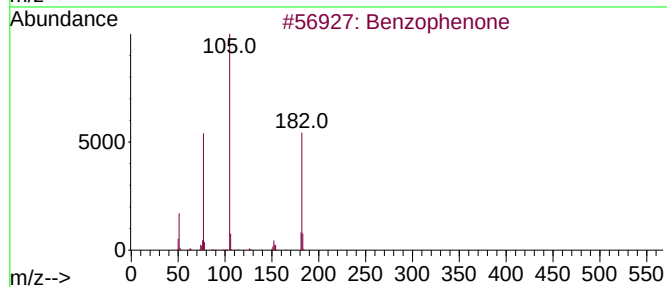
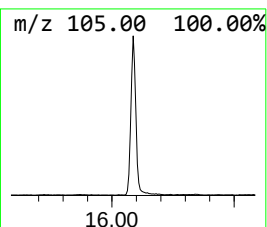
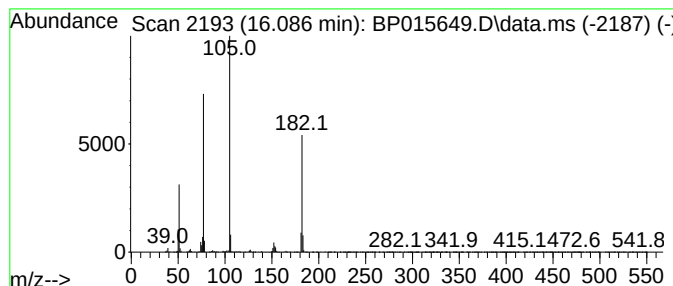
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.087	7.13 ng/ul	595927	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	89



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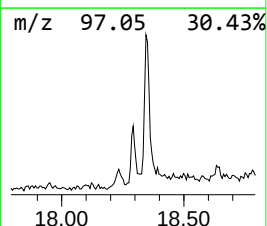
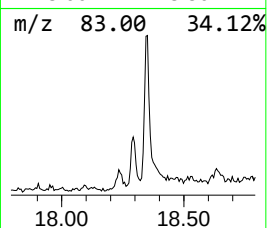
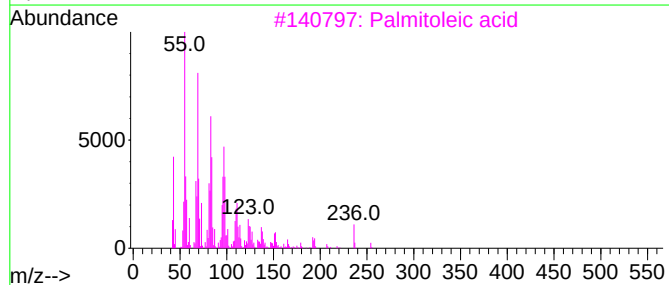
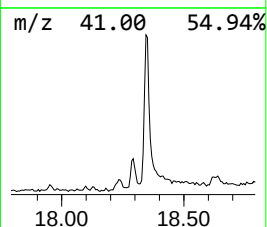
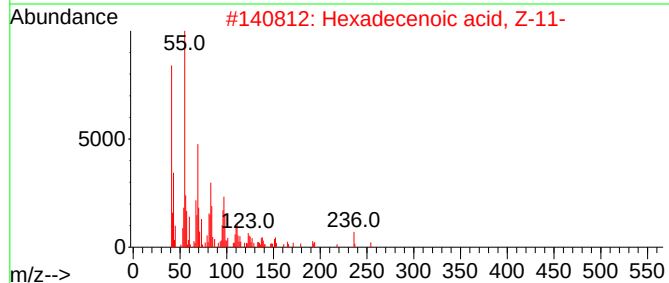
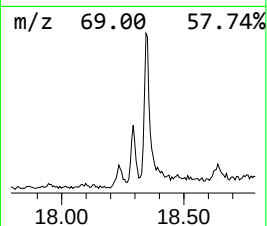
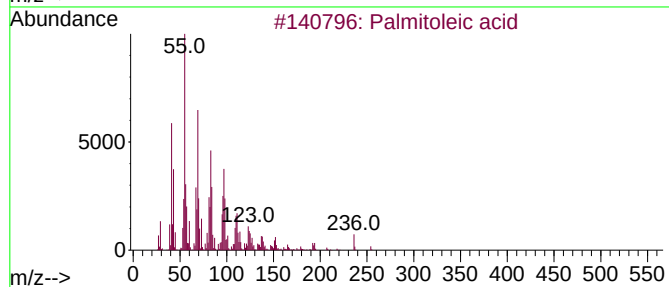
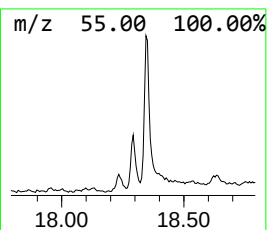
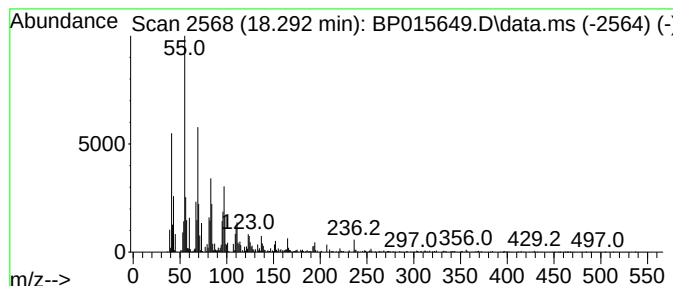
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Palmitoleic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	2.15 ng/ul	182806	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	99
2			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	98
3			Palmitoleic acid	254	C16H30O2	000373-49-9	98
4			11-Eicosenoic acid, methyl ester	324	C21H40O2	003946-08-5	98
5			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	97



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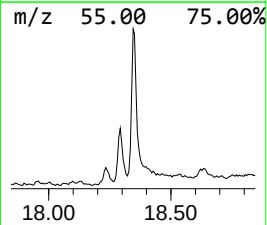
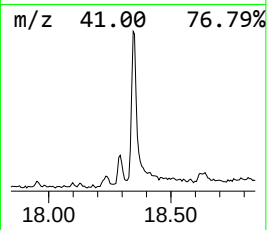
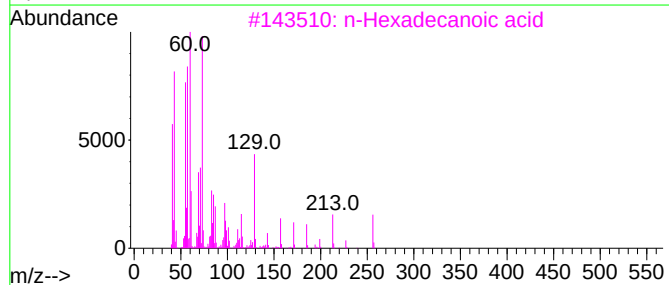
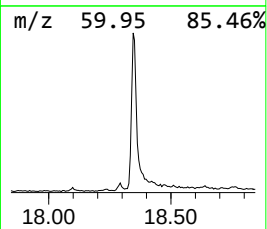
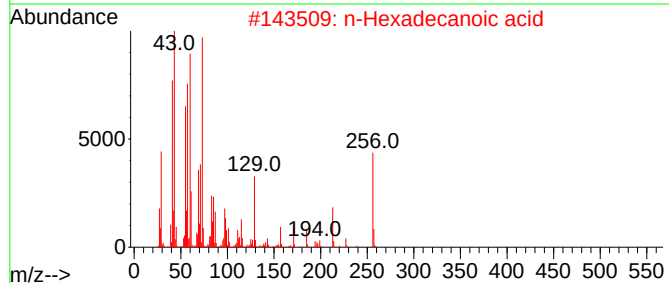
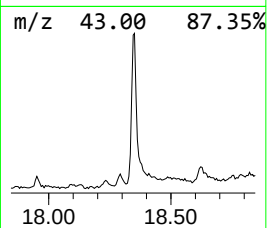
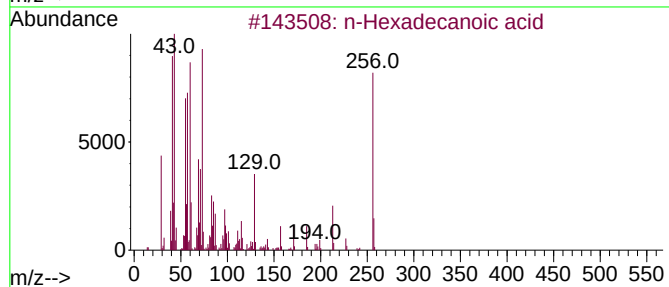
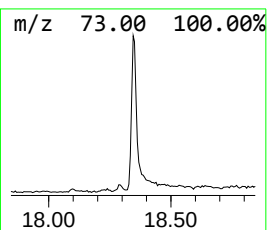
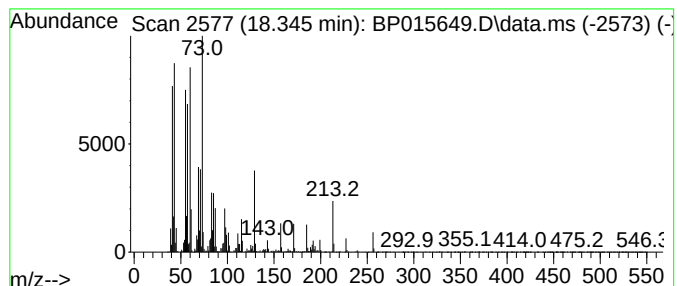
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	11.48 ng/ul	977727	Phenanthrene-d10	17.486

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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Data File : BP015649.D
Acq On : 22 Jun 2023 16:13
Operator : MA/JU
Sample : 03083-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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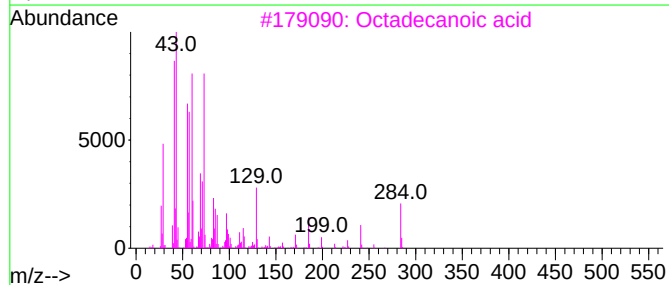
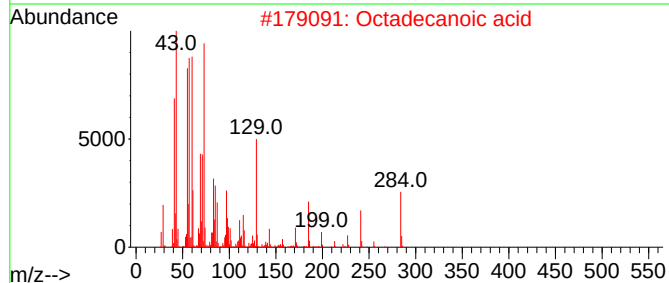
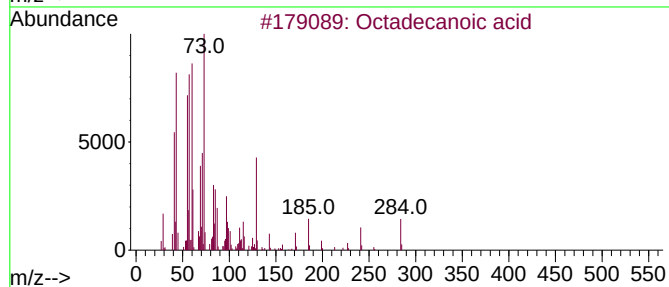
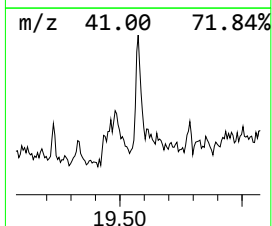
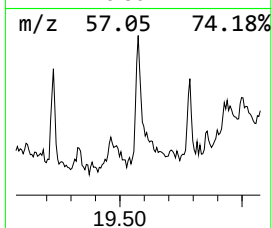
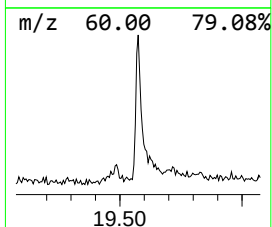
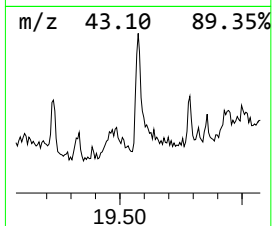
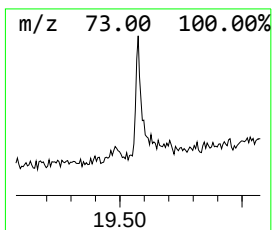
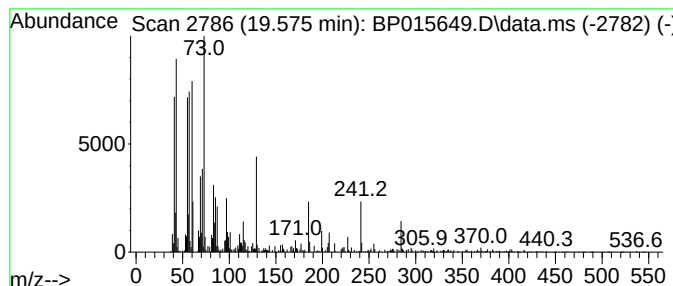
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.575	3.47 ng/ul	242184	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	97
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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Sample : 03083-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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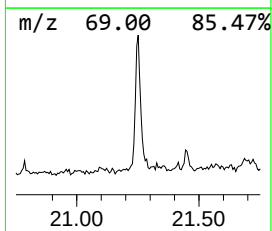
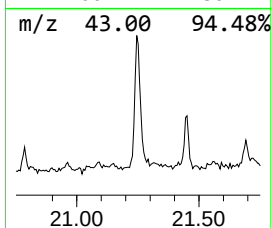
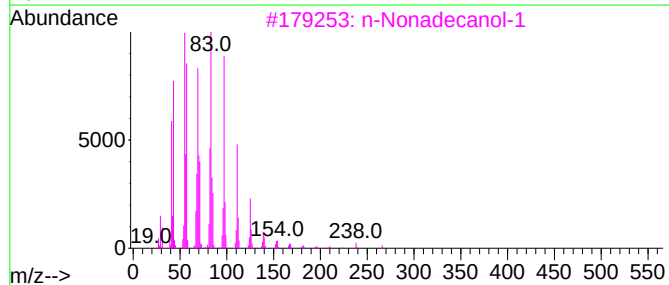
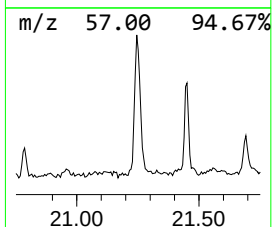
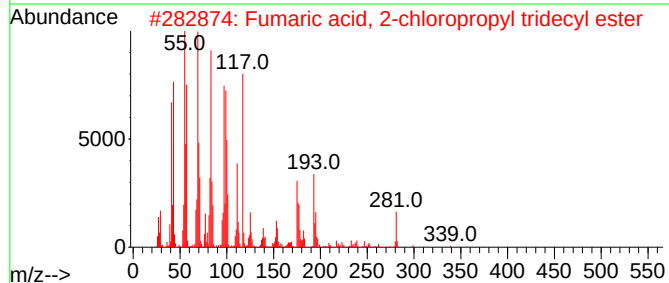
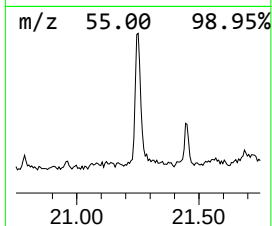
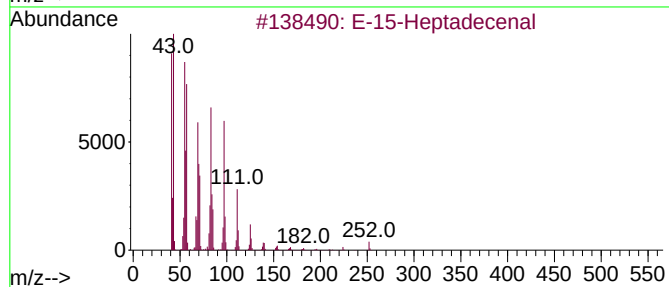
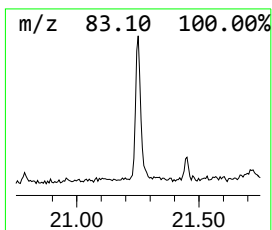
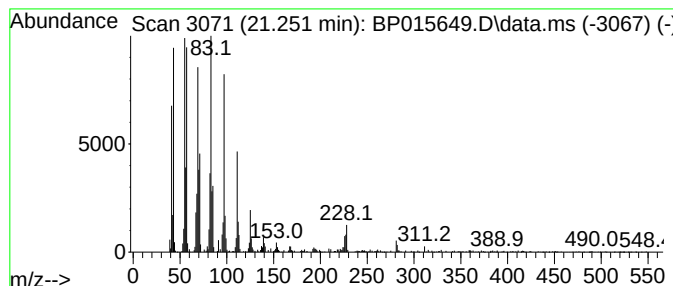
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 E-15-Heptadecenal Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	8.06 ng/ul	563554	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			E-15-Heptadecenal	252	C17H32O	1000130-97-9	99
2			Fumaric acid, 2-chloropropyl tri...	374	C20H35ClO4	1010348-57-1	96
3			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
4			1-Tetracosene	336	C24H48	010192-32-2	94
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015649.D
Acq On : 22 Jun 2023 16:13
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Sample : 03083-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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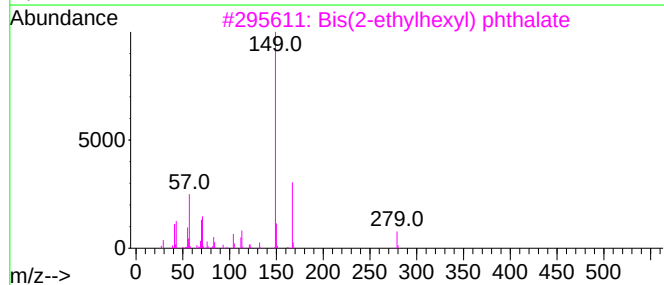
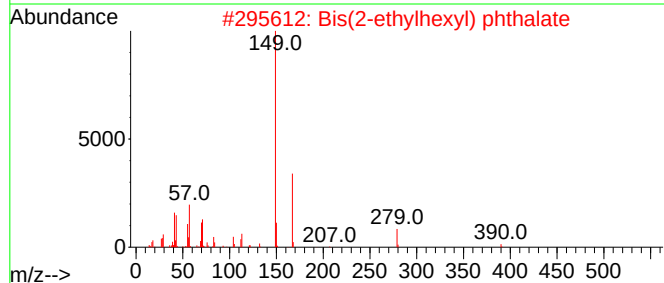
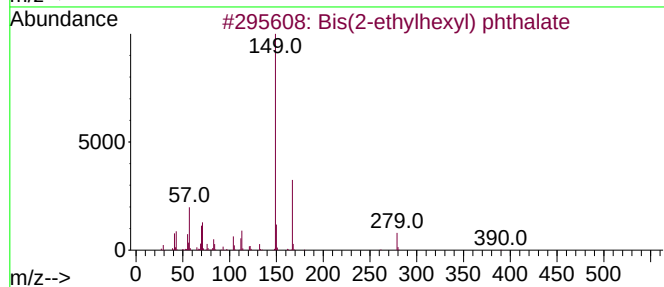
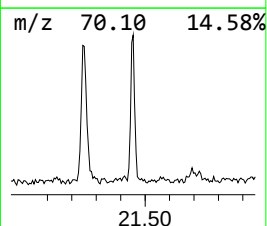
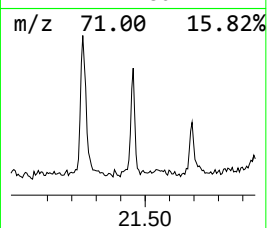
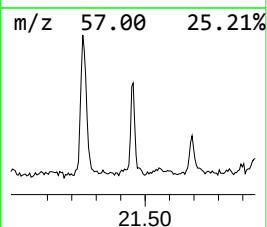
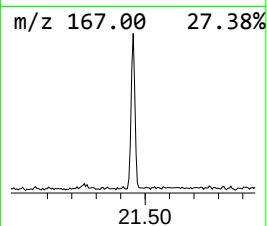
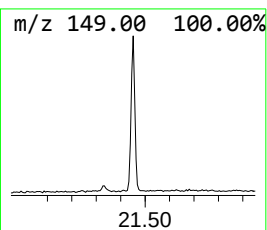
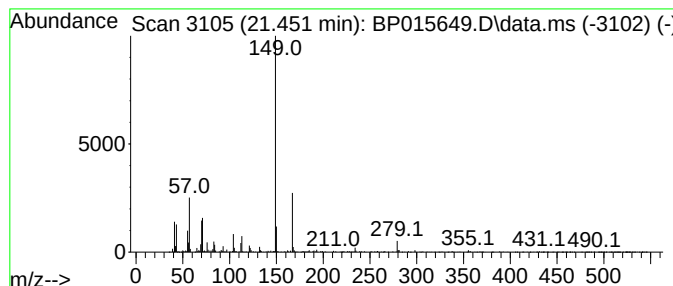
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Bis(2-ethylhexyl) phthalate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.451	4.90 ng/ul	342221	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	93
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	90
4			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	86
5			Phthalic acid, decyl 2-ethylhexy...	418	C26H42O4	1000309-00-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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Sample : 03083-15
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ALS Vial : 12 Sample Multiplier: 1

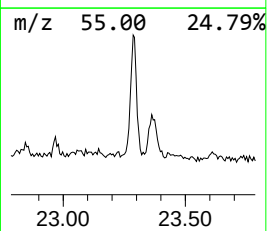
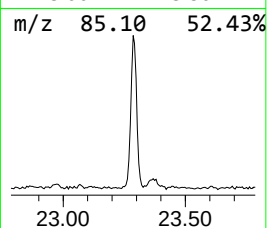
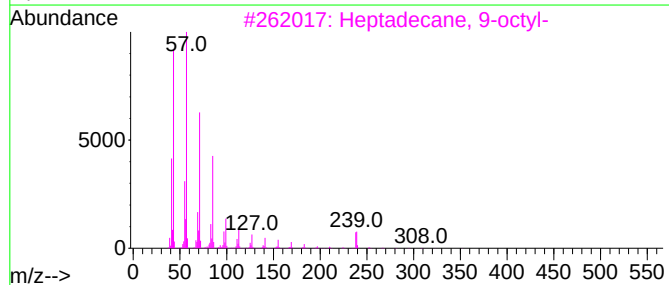
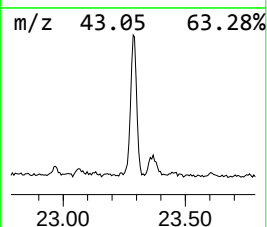
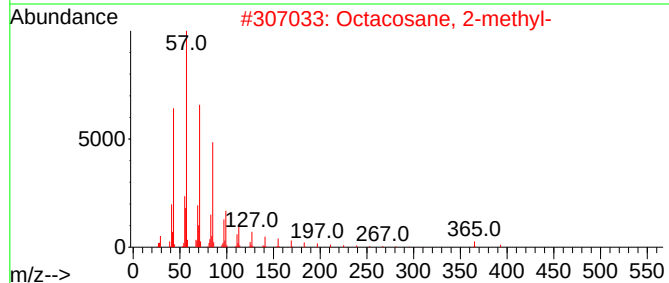
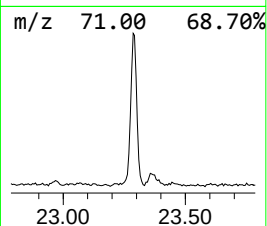
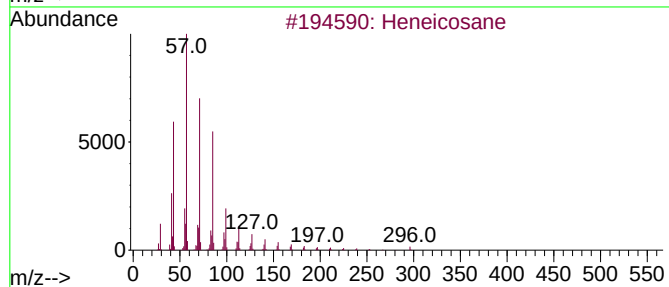
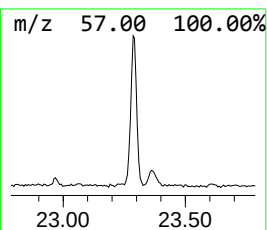
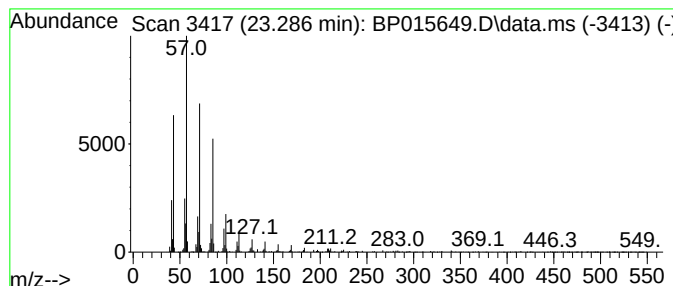
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Quant Title : SVOA CALIBRATION

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.	
23.286	11.20 ng/ul	728103	Perylene-d12	24.133	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	97
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
4	Docosane, 11-butyl-	366	C26H54	013475-76-8	87
5	Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015649.D
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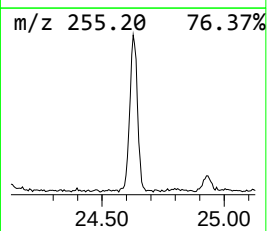
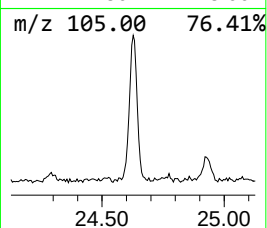
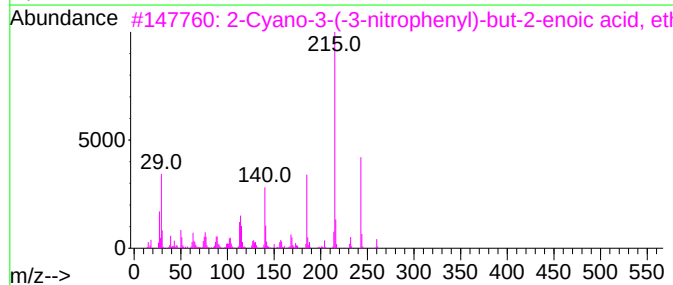
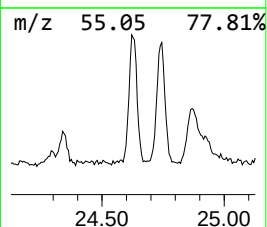
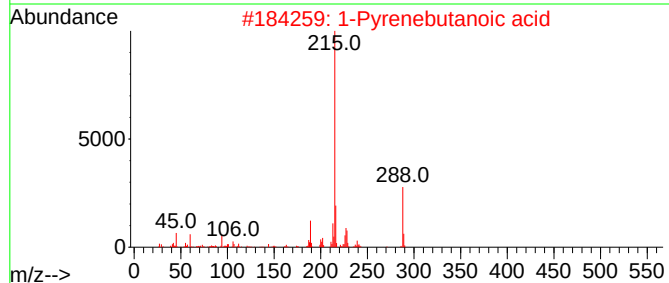
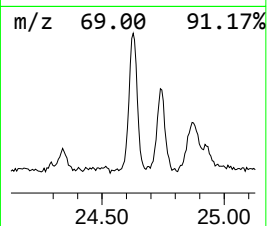
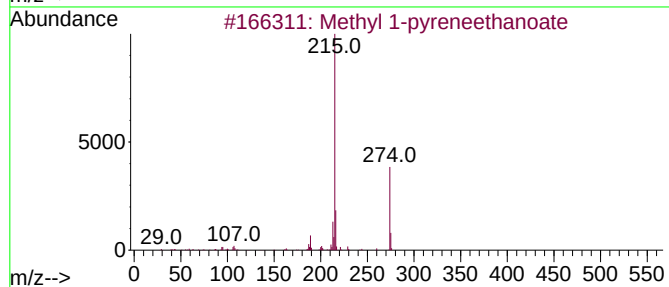
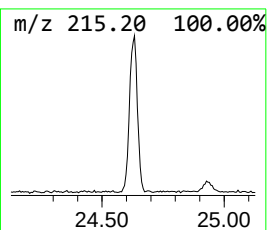
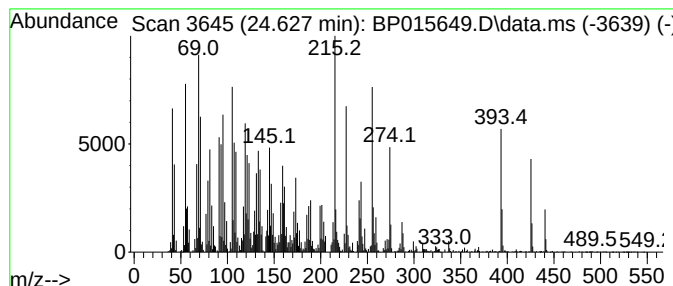
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	39.23 ng/ul	2550120	Perylene-d12	24.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	38
2			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	25
3			2-Cyano-3-(-3-nitrophenyl)-but-2...	260	C13H12N2O4	014442-35-4	25
4			Metharbital, N-trimethylsilyl-	270	C12H22N2O3Si	1000442-60-0	15
5			3-Imidazo[2,1-b][1,3]thiazol-6-y...	215	C11H9N3S	861206-26-0	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
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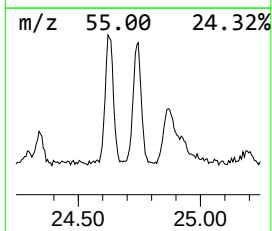
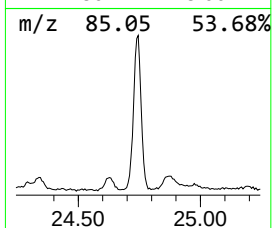
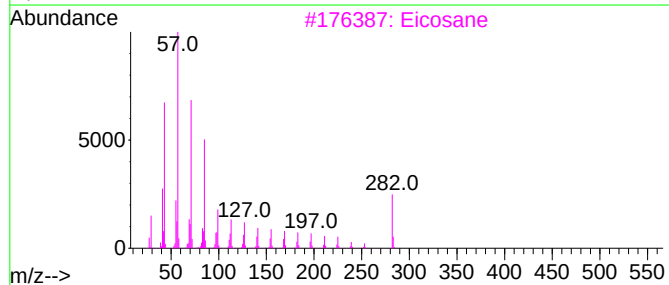
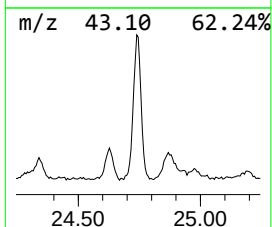
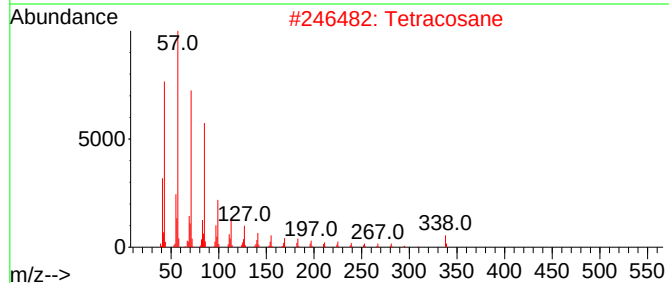
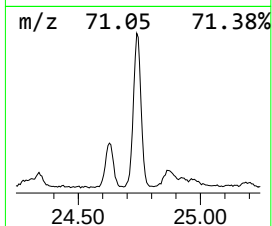
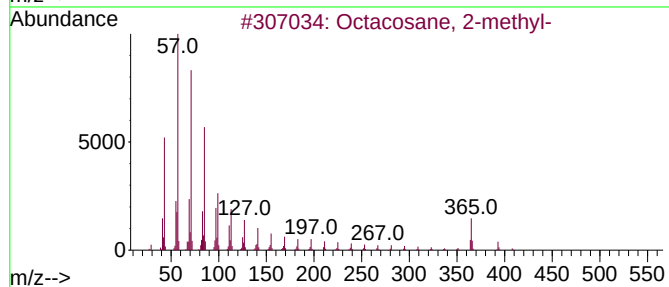
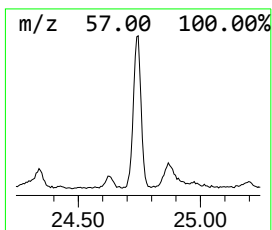
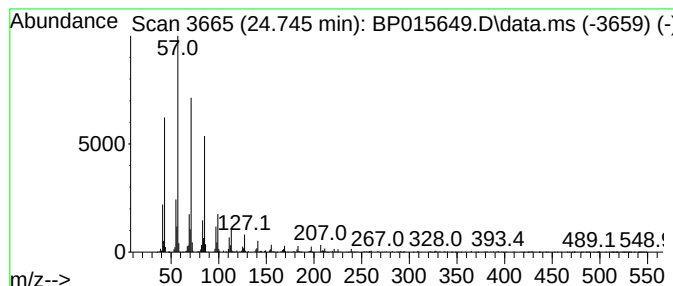
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.745	21.41 ng/ul	1391620	Perylene-d12	24.133	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane, 2-methyl-	408	C29H60	001560-98-1	95
2	Tetracosane	338	C24H50	000646-31-1	94
3	Eicosane	282	C20H42	000112-95-8	93
4	Heptacosane	380	C27H56	000593-49-7	93
5	Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015649.D
Acq On : 22 Jun 2023 16:13
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Sample : 03083-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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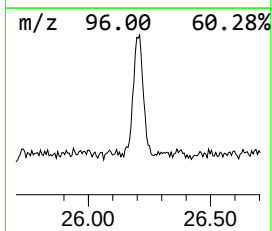
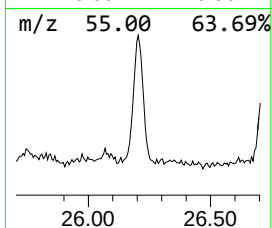
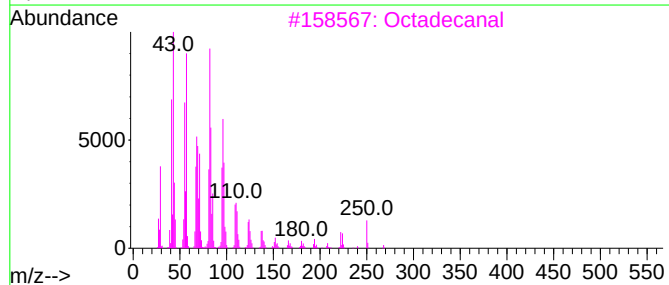
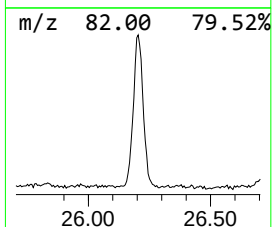
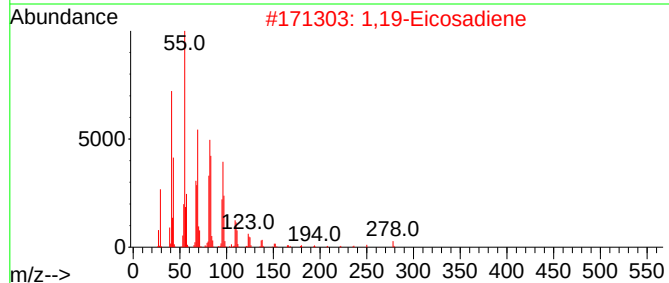
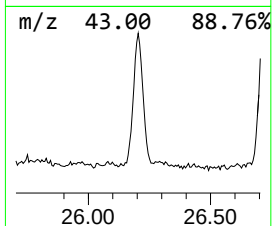
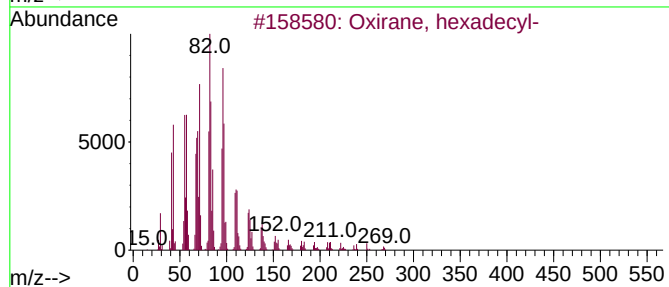
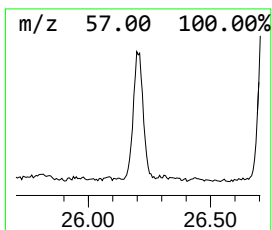
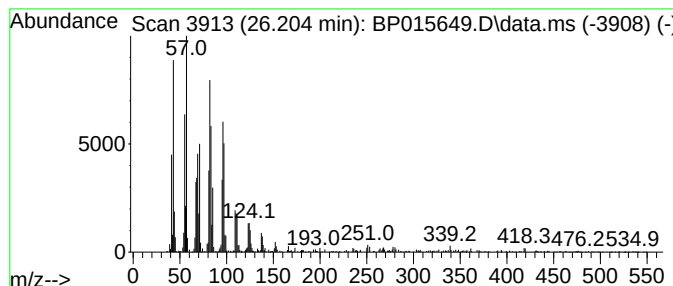
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Oxirane, hexadecyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.204	17.84 ng/ul	1159470	Perylene-d12	24.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	99
2			1,19-Eicosadiene	278	C20H38	014811-95-1	95
3			Octadecanal	268	C18H36O	000638-66-4	94
4			Hexacosanal	380	C26H52O	026627-85-0	93
5			Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015649.D
Acq On : 22 Jun 2023 16:13
Operator : MA/JU
Sample : 03083-15
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFL3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.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
Benzophenone	16.087	7.1	ng/ul	595927	3	14.728	1672610	20.0
Palmitoleic acid	18.292	2.1	ng/ul	182806	4	17.486	1703670	20.0
n-Hexadecanoic ...	18.345	11.5	ng/ul	977727	4	17.486	1703670	20.0
Octadecanoic acid	19.575	3.5	ng/ul	242184	5	21.574	1397650	20.0
E-15-Heptadecenal	21.251	8.1	ng/ul	563554	5	21.574	1397650	20.0
Bis(2-ethylhexy...	21.451	4.9	ng/ul	342221	5	21.574	1397650	20.0
(DEL) Alkane: S...	23.286	11.2	ng/ul	728103	6	24.133	1300150	20.0
unknown-01	24.627	39.2	ng/ul	2550120	6	24.133	1300150	20.0
(DEL) Alkane: S...	24.745	21.4	ng/ul	1391620	6	24.133	1300150	20.0
Oxirane, hexade...	26.204	17.8	ng/ul	1159470	6	24.133	1300150	20.0