

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
 Data File : BP015366.D
 Acq On : 31 May 2023 18:49
 Operator : CG/JU
 Sample : 02851-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0FK2

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-BP051123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015366.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.417	35	39	48	rVB	29427	47930	1.10%	0.094%
2	3.887	118	119	124	rBV3	11422	17919	0.41%	0.035%
3	4.093	149	154	159	rBV2	17682	26900	0.62%	0.053%
4	4.193	168	171	179	rVB2	8541	12937	0.30%	0.025%
5	4.587	233	238	241	rBV7	5000	8185	0.19%	0.016%
6	4.652	244	249	255	rBV	25318	39454	0.90%	0.078%
7	4.905	287	292	296	rBV8	6097	11186	0.26%	0.022%
8	5.146	328	333	341	rVB3	12165	23237	0.53%	0.046%
9	5.634	412	416	421	rVB2	5821	11753	0.27%	0.023%
10	5.764	433	438	443	rBV	10853	17696	0.40%	0.035%
11	7.205	677	683	686	rBV	24550	37112	0.85%	0.073%
12	7.258	686	692	706	rVV	506520	926468	21.20%	1.821%
13	7.422	710	720	734	rVB	407622	673739	15.41%	1.324%
14	7.628	748	755	766	rBV	530529	870407	19.91%	1.711%
15	7.911	798	803	810	rBV8	4438	8379	0.19%	0.016%
16	8.105	828	836	847	rBV	555070	895070	20.48%	1.759%
17	8.234	854	858	863	rBV3	9029	13980	0.32%	0.027%
18	8.816	950	957	963	rBV	664486	1097600	25.11%	2.157%
19	8.875	963	967	973	rVV	333022	547462	12.53%	1.076%
20	8.940	973	978	986	rVB	148590	246963	5.65%	0.485%
21	9.281	1027	1036	1047	rBV	547634	913469	20.90%	1.796%
22	9.669	1095	1102	1107	rBV2	67563	101405	2.32%	0.199%
23	10.011	1152	1160	1168	rBV	450230	762945	17.46%	1.500%
24	10.187	1183	1190	1192	rBV3	12292	21383	0.49%	0.042%
25	10.240	1195	1199	1215	rVV4	25773	67974	1.56%	0.134%
26	10.369	1216	1221	1227	rVB4	15205	28173	0.64%	0.055%
27	10.552	1244	1252	1263	rBV	662079	1161321	26.57%	2.283%
28	10.934	1309	1317	1326	rBV	826545	1362205	31.17%	2.678%
29	11.075	1333	1341	1355	rBV	203310	414862	9.49%	0.815%
30	11.475	1400	1409	1414	rBV	8800	21689	0.50%	0.043%
31	11.740	1447	1454	1462	rBV6	10389	28361	0.65%	0.056%
32	12.163	1521	1526	1532	rVB2	47764	69993	1.60%	0.138%
33	12.328	1545	1554	1557	rBV10	7263	17627	0.40%	0.035%
34	12.428	1565	1571	1577	rBV	28783	55340	1.27%	0.109%
35	12.487	1577	1581	1583	rVV4	14385	22908	0.52%	0.045%
36	12.510	1583	1585	1594	rVB3	16170	23945	0.55%	0.047%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
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37	12.728	1617	1622	1626	rBV7	5878	10051	0.23%	0.020%
38	13.028	1665	1673	1684	rBV	237410	392166	8.97%	0.771%
39	13.622	1767	1774	1781	rBV3	27083	55804	1.28%	0.110%
40	14.069	1843	1850	1854	rVB9	7078	12279	0.28%	0.024%
41	14.151	1854	1864	1872	rBV	1428902	2048856	46.88%	4.027%
42	14.210	1872	1874	1878	rVV5	7056	8920	0.20%	0.018%
43	14.328	1887	1894	1899	rBV	13961	20035	0.46%	0.039%
44	14.446	1906	1914	1927	rBV	1517756	2332843	53.37%	4.585%
45	14.563	1928	1934	1941	rBV	49090	89209	2.04%	0.175%
46	14.746	1956	1965	1974	rBV	1350686	2001219	45.79%	3.934%
47	14.946	1993	1999	2018	rBV	384142	795681	18.20%	1.564%
48	15.104	2022	2026	2029	rVV6	11671	23270	0.53%	0.046%
49	15.163	2029	2036	2044	rVB6	26601	70844	1.62%	0.139%
50	15.487	2086	2091	2092	rBV5	5123	8005	0.18%	0.016%
51	15.687	2121	2125	2128	rBV2	40168	48761	1.12%	0.096%
52	15.740	2128	2134	2142	rVB	2040184	2917824	66.76%	5.735%
53	15.863	2149	2155	2167	rBV	385595	559219	12.79%	1.099%
54	15.981	2171	2175	2182	rVB5	10948	18434	0.42%	0.036%
55	16.098	2189	2195	2205	rBV	812742	1143067	26.15%	2.247%
56	16.181	2207	2209	2216	rVB8	5643	8277	0.19%	0.016%
57	16.440	2249	2253	2255	rBV5	13221	19730	0.45%	0.039%
58	16.581	2274	2277	2282	rBV7	6878	9938	0.23%	0.020%
59	16.669	2286	2292	2298	rBV	134796	185376	4.24%	0.364%
60	16.881	2323	2328	2334	rBV4	30717	53057	1.21%	0.104%
61	16.940	2334	2338	2344	rVV3	37252	55302	1.27%	0.109%
62	17.175	2374	2378	2382	rVB6	6469	10409	0.24%	0.020%
63	17.245	2384	2390	2393	rBV6	10787	22126	0.51%	0.043%
64	17.375	2408	2412	2420	rBV10	18288	32930	0.75%	0.065%
65	17.445	2420	2424	2427	rBV6	16048	20482	0.47%	0.040%
66	17.504	2427	2434	2440	rBV	1745074	2550640	58.36%	5.014%
67	17.604	2440	2451	2457	rBV	2178568	3280773	75.06%	6.449%
68	17.692	2463	2466	2472	rVB7	15527	27121	0.62%	0.053%
69	17.840	2488	2491	2495	rBV2	41438	46922	1.07%	0.092%
70	17.975	2510	2514	2516	rBV4	10441	14870	0.34%	0.029%
71	18.010	2516	2520	2525	rVB2	40901	60666	1.39%	0.119%
72	18.245	2555	2560	2567	rBV	168086	262673	6.01%	0.516%
73	18.304	2567	2570	2574	rVV6	20984	26471	0.61%	0.052%
74	18.357	2575	2579	2589	rVV	715268	934442	21.38%	1.837%
75	18.428	2589	2591	2594	rVB	23073	27773	0.64%	0.055%
76	18.645	2624	2628	2629	rBV4	9235	13833	0.32%	0.027%

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 Title : SVOA CALIBRATION

77	18.916	2670	2674	2679	rBV2	70996	95508	2.19%	0.188%
78	18.998	2684	2688	2691	rBV6	19328	24986	0.57%	0.049%
79	19.045	2694	2696	2702	rVB5	27120	38602	0.88%	0.076%
80	19.134	2708	2711	2716	rBV	37173	48583	1.11%	0.095%
81	19.186	2717	2720	2726	rVV5	28584	44030	1.01%	0.087%
82	19.239	2727	2729	2736	rVB3	28112	32570	0.75%	0.064%
83	19.398	2743	2756	2761	rBV6	98994	338046	7.73%	0.664%
84	19.445	2761	2764	2766	rBV3	70045	88666	2.03%	0.174%
85	19.469	2766	2768	2770	rBV	33904	36491	0.83%	0.072%
86	19.581	2783	2787	2795	rBV	222161	331573	7.59%	0.652%
87	19.675	2800	2803	2809	rVB4	53338	74029	1.69%	0.146%
88	19.822	2822	2828	2838	rVB	3232478	4267396	97.64%	8.388%
89	20.316	2909	2912	2914	rVB	64468	66989	1.53%	0.132%
90	20.351	2915	2918	2921	rVB	53677	55393	1.27%	0.109%
91	20.792	2991	2993	2997	rVB	84081	84985	1.94%	0.167%
92	20.963	3019	3022	3028	rVB4	81589	107669	2.46%	0.212%
93	21.251	3067	3071	3078	rBV2	347947	558763	12.78%	1.098%
94	21.528	3116	3118	3121	rVB	86235	81008	1.85%	0.159%
95	21.580	3122	3127	3136	rVB	2189276	2964856	67.83%	5.828%
96	22.180	3226	3229	3233	rBV	321767	409961	9.38%	0.806%
97	23.304	3415	3420	3426	rVB	1102265	1747935	39.99%	3.436%
98	23.969	3526	3533	3543	rVB2	2024209	4370753	100.00%	8.591%
99	24.133	3554	3561	3570	rVB	1364941	2870005	65.66%	5.641%
100	24.757	3662	3667	3678	rVB	589216	1308215	29.93%	2.571%

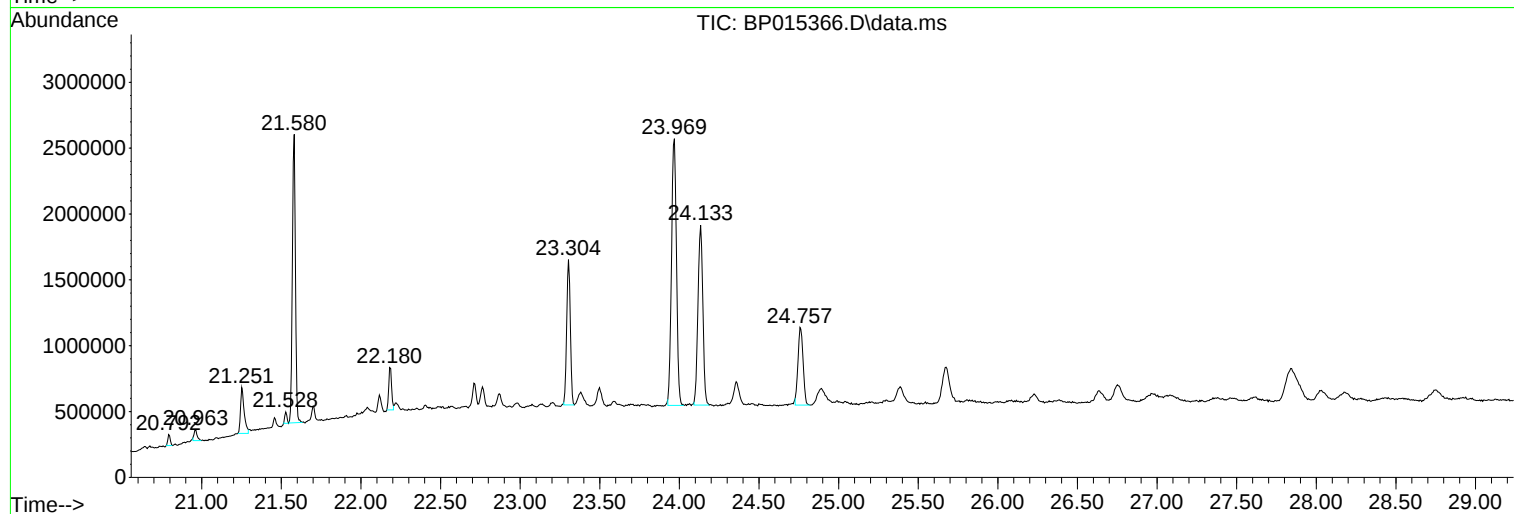
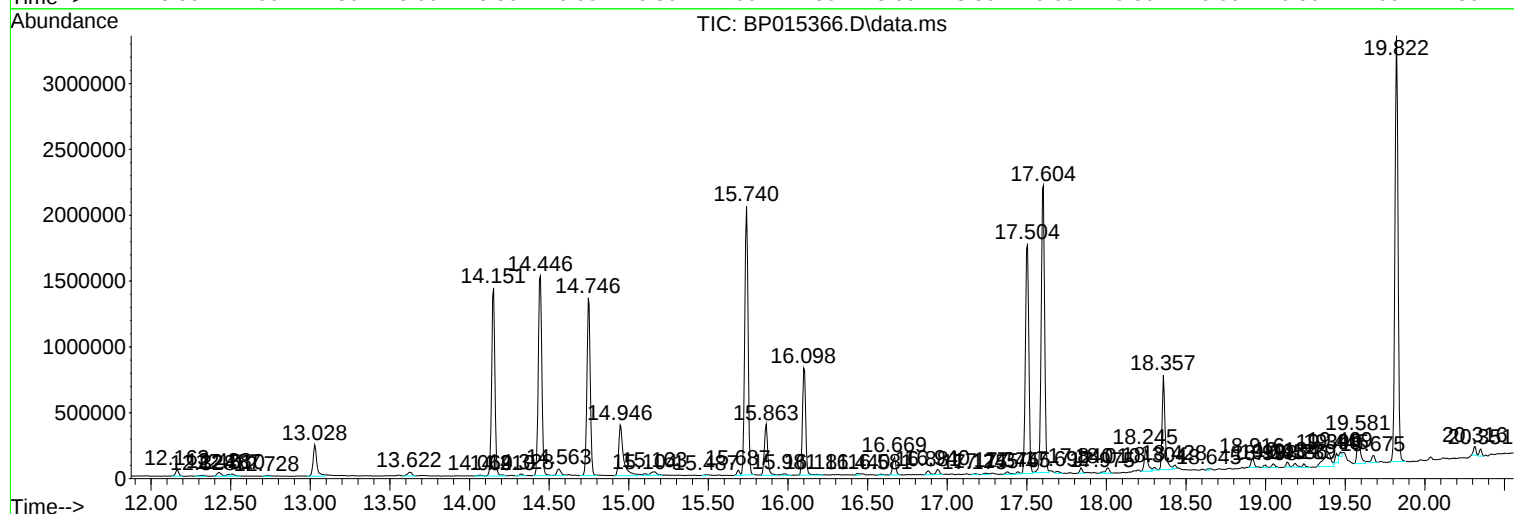
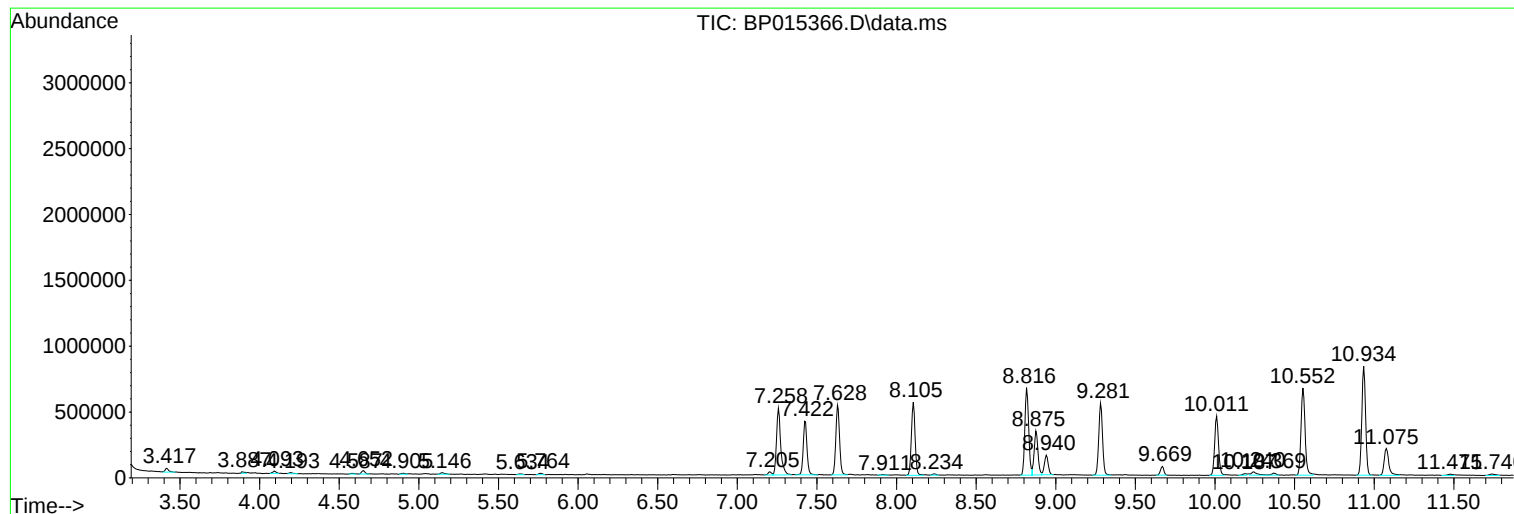
Sum of corrected areas: 50875287

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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



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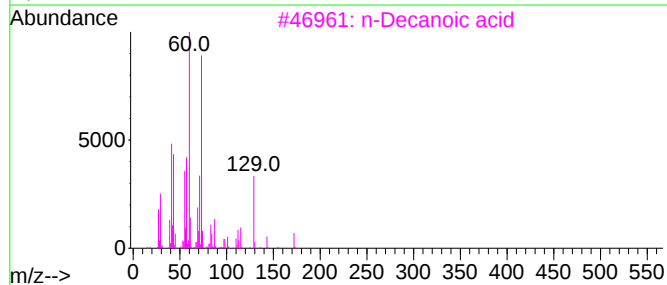
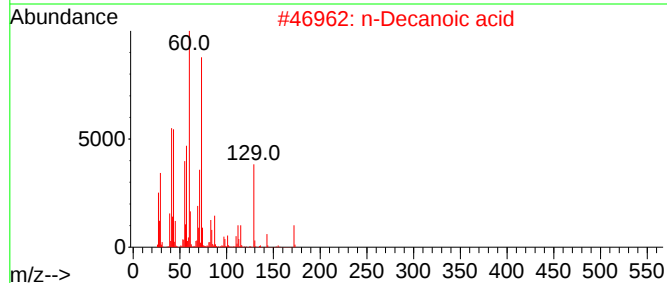
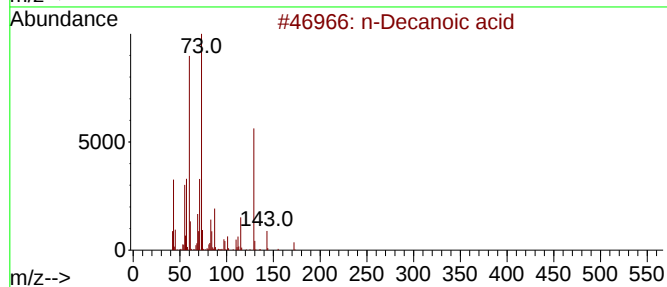
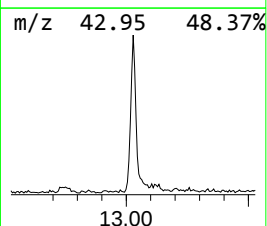
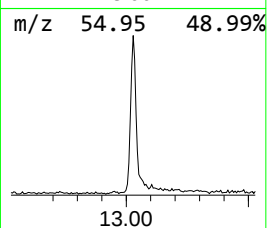
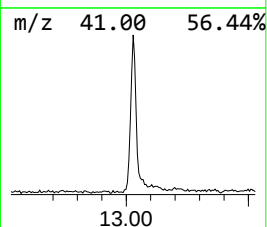
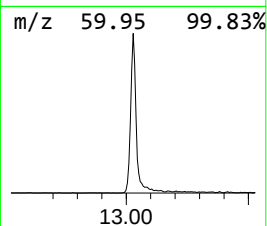
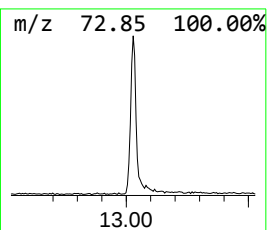
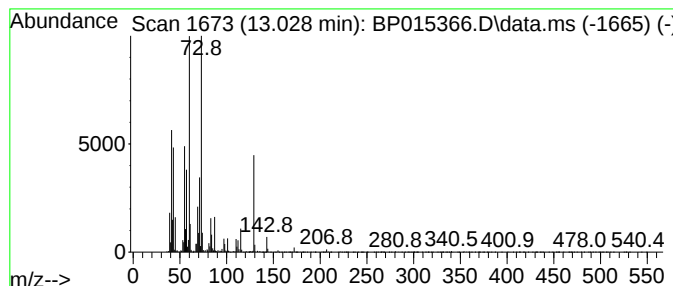
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Decanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	3.92 ng/ul	392166	Acenaphthene-d10	14.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Decanoic acid			172	C10H20O2	000334-48-5	94
2	n-Decanoic acid			172	C10H20O2	000334-48-5	94
3	n-Decanoic acid			172	C10H20O2	000334-48-5	80
4	Tetradecanoic acid			228	C14H28O2	000544-63-8	64
5	n-Decanoic acid			172	C10H20O2	000334-48-5	64



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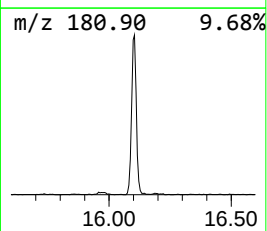
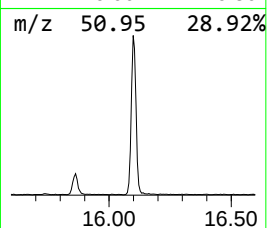
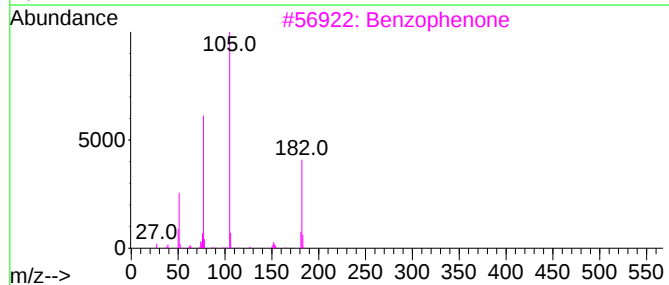
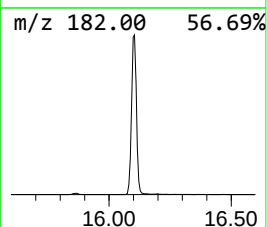
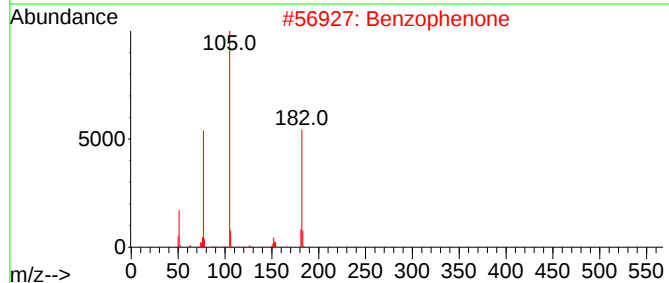
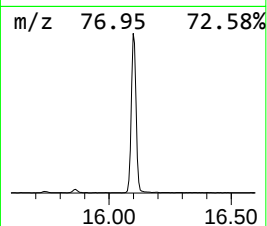
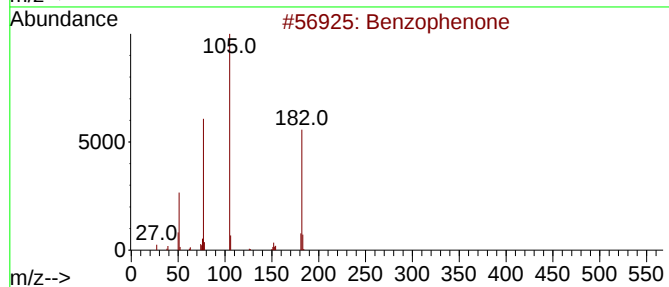
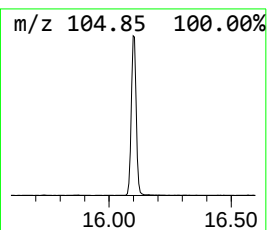
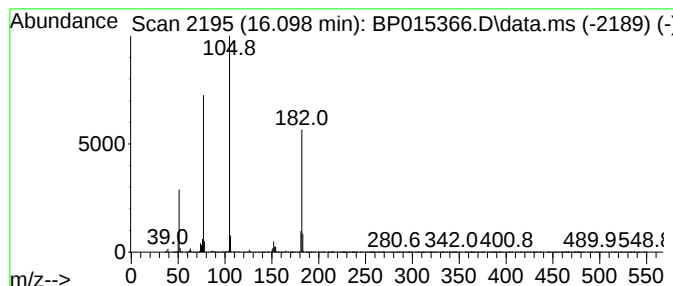
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.098	11.42 ng/ul	1143070	Acenaphthene-d10	14.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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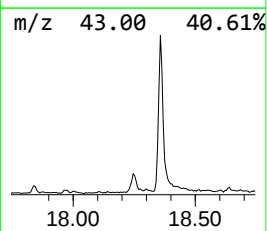
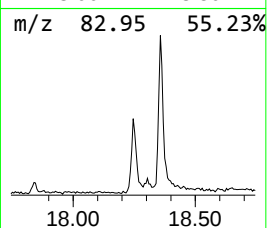
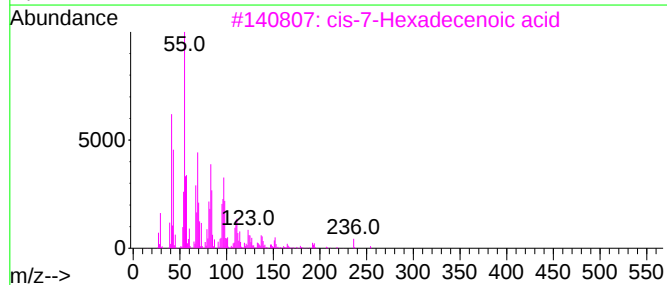
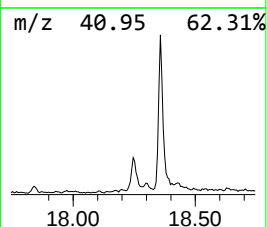
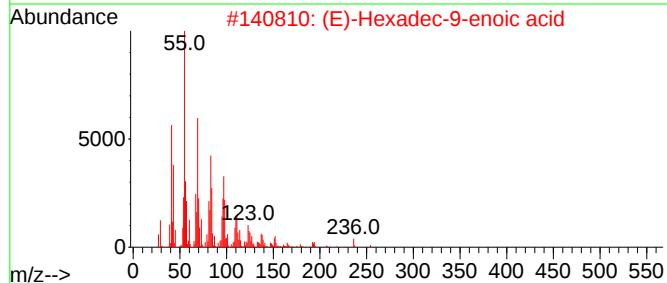
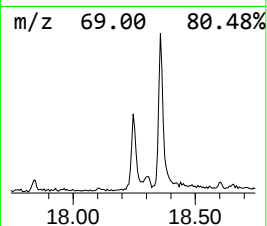
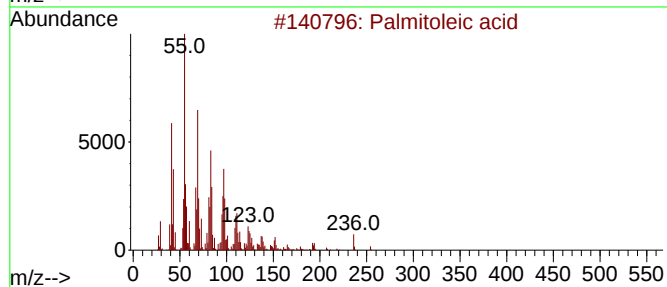
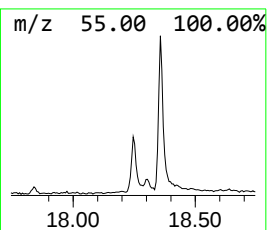
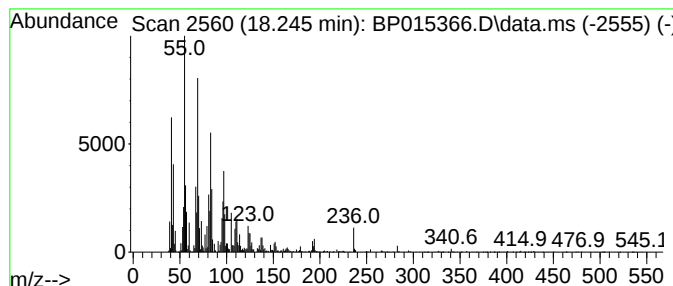
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Palmitoleic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.245	2.06 ng/ul	262673	Phenanthrene-d10	17.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	96
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	96
3			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	95
4			Palmitoleic acid	254	C16H30O2	000373-49-9	91
5			Nonadec-13-enoic acid, methyl ester	310	C20H38O2	1000191-29-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
Acq On : 31 May 2023 18:49
Operator : CG/JU
Sample : 02851-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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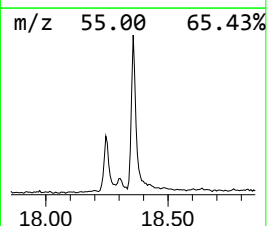
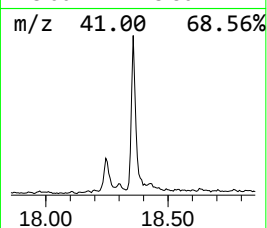
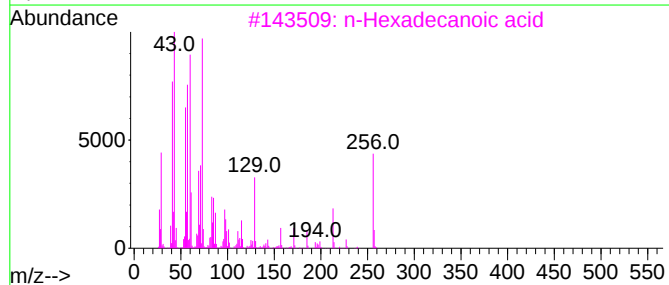
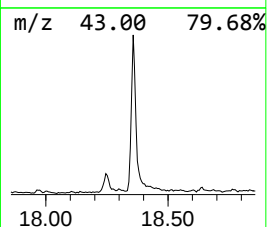
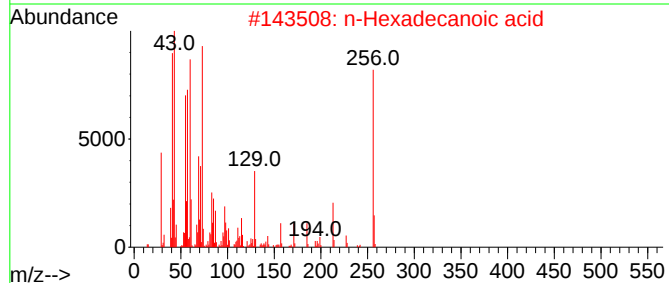
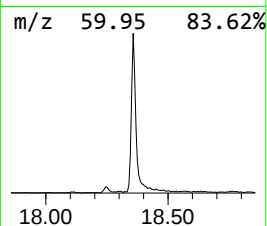
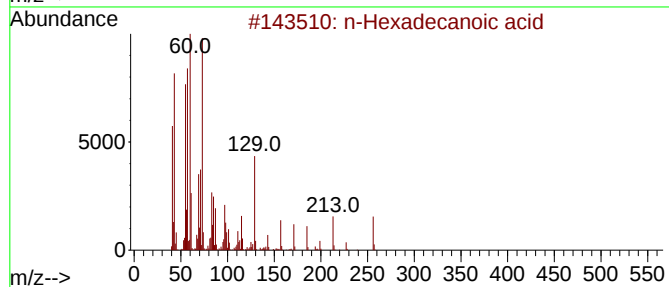
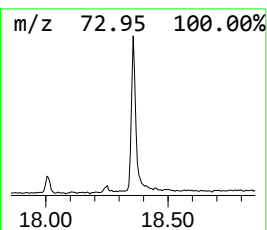
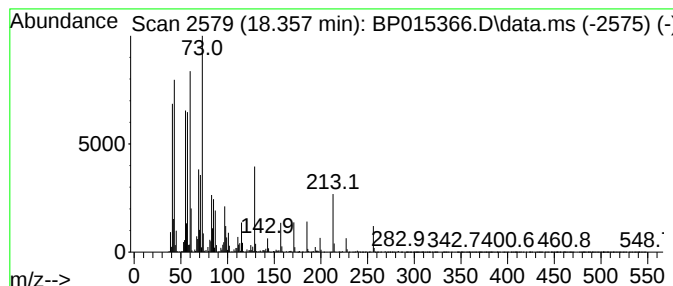
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	7.33 ng/ul	934442	Phenanthrene-d10	17.504

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	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
Acq On : 31 May 2023 18:49
Operator : CG/JU
Sample : 02851-09
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ClientSampleId :
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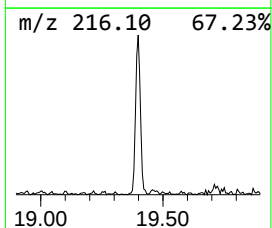
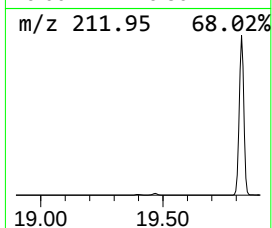
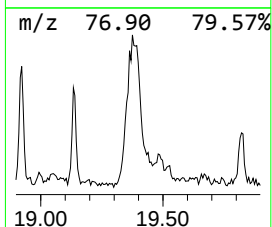
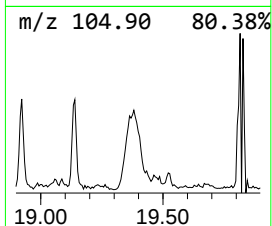
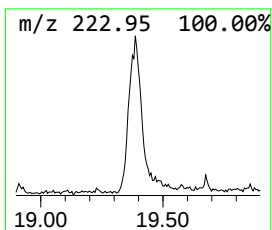
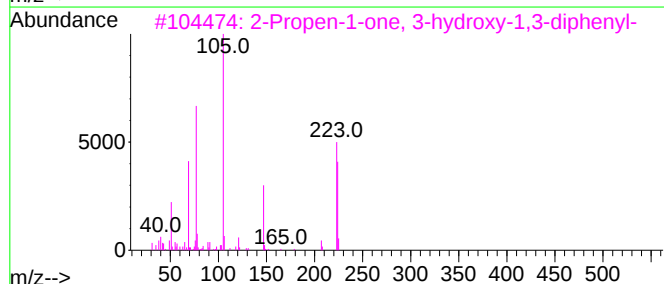
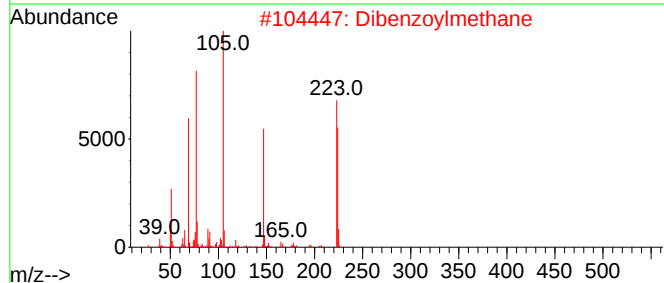
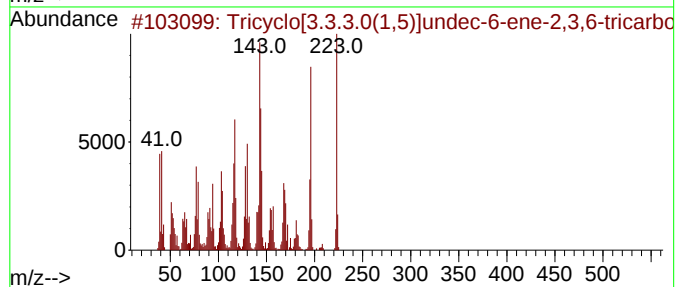
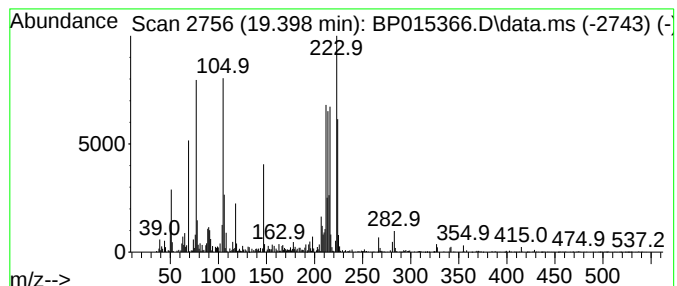
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Tricyclo[3.3.3.0(1,5)]undec... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.398	2.65 ng/ul	338046	Phenanthrene-d10	17.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[3.3.3.0(1,5)]undec-6-en...	223	C14H13N3	118894-71-6	53
2			Dibenzoylmethane	224	C15H12O2	000120-46-7	52
3			2-Propen-1-one, 3-hydroxy-1,3-di...	224	C15H12O2	001704-15-0	46
4			Dibenzoylmethane	224	C15H12O2	000120-46-7	42
5			Benzoylanthranil	223	C14H9NO2	007155-17-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
Acq On : 31 May 2023 18:49
Operator : CG/JU
Sample : 02851-09
Misc :
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ClientSampleId :
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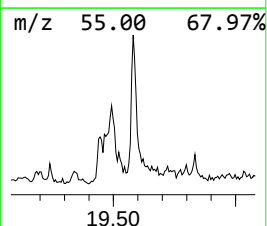
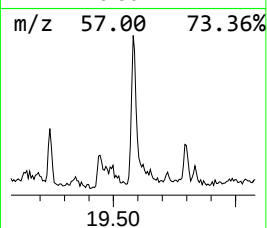
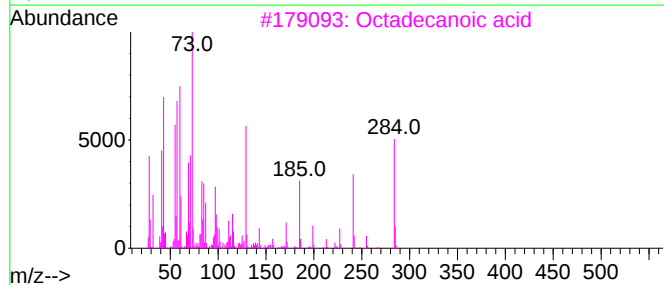
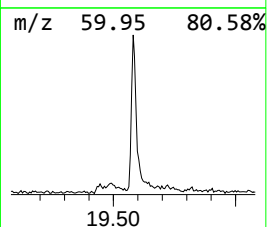
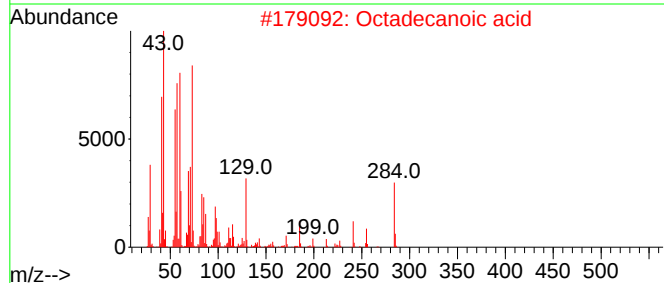
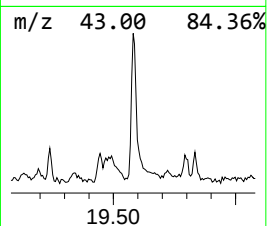
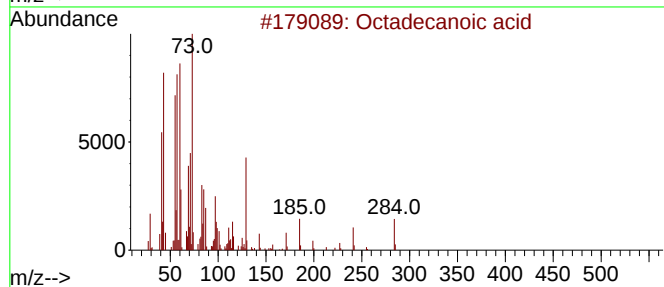
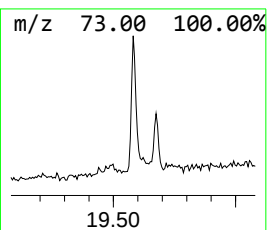
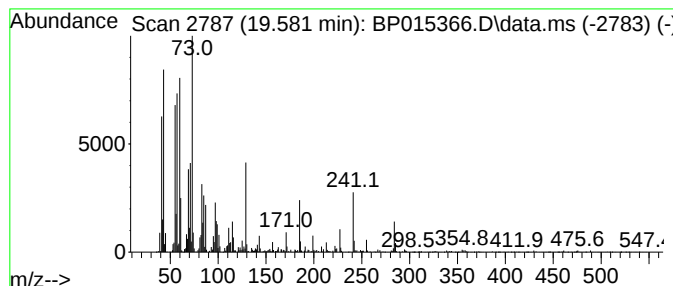
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.581	2.24 ng/ul	331573	Chrysene-d12	21.581

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	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
Acq On : 31 May 2023 18:49
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Misc :
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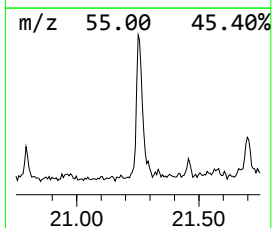
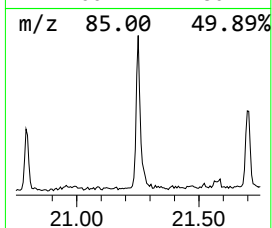
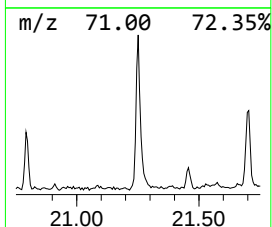
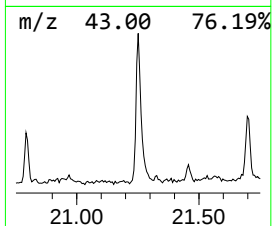
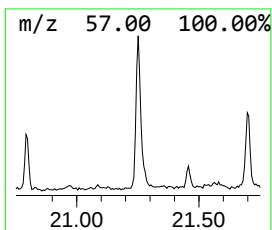
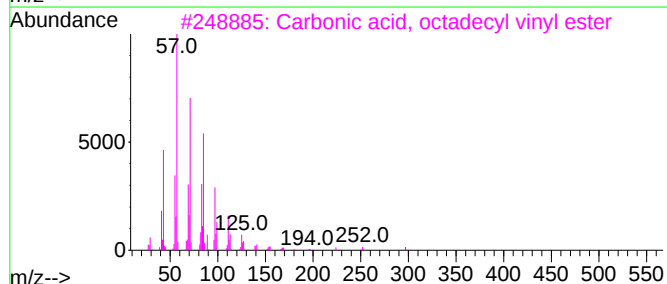
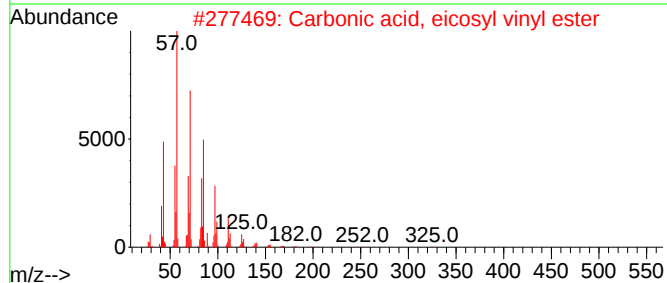
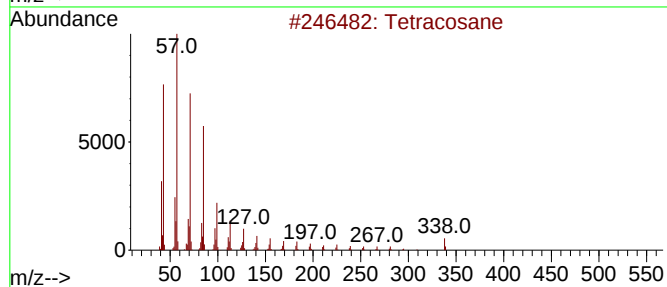
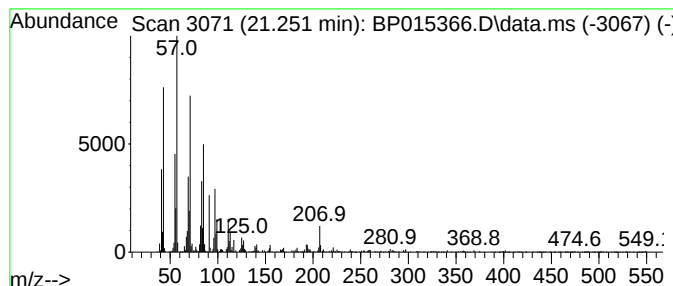
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.251	3.77 ng/ul	558763	Chrysene-d12		21.581	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	89
2	Carbonic acid, eicosyl vinyl ester		368	C23H44O3	1000382-54-3	87
3	Carbonic acid, octadecyl vinyl e...		340	C21H40O3	1000382-54-4	87
4	Carbonic acid, octadecyl prop-1-...		354	C22H42O3	1000383-11-5	87
5	Carbonic acid, hexadecyl prop-1-...		326	C20H38O3	1000382-90-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
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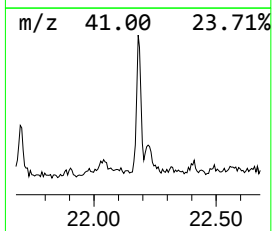
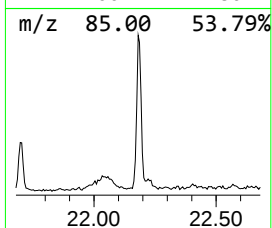
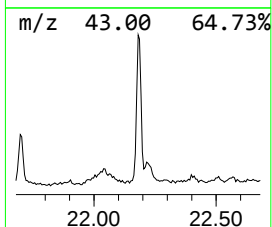
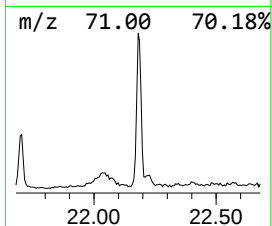
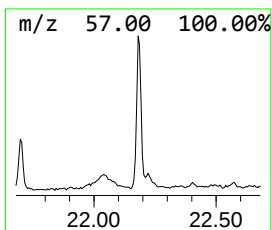
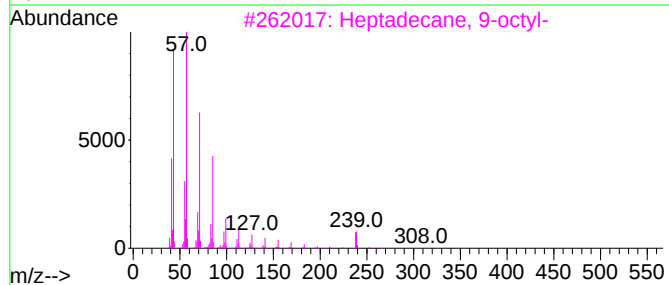
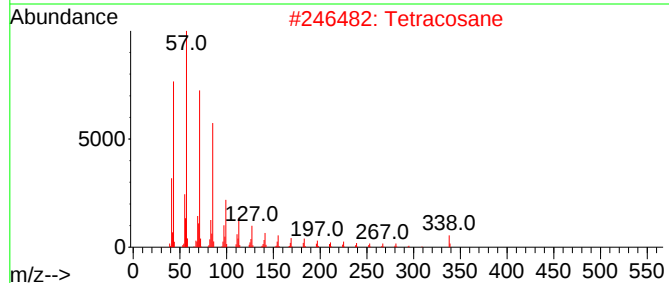
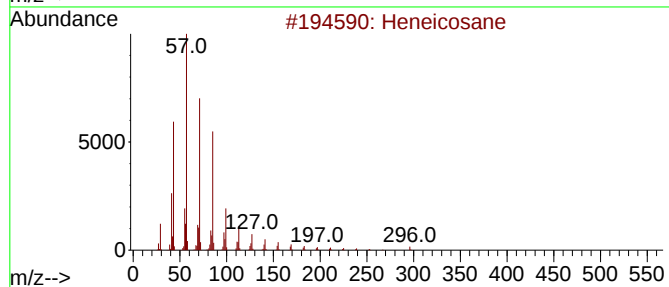
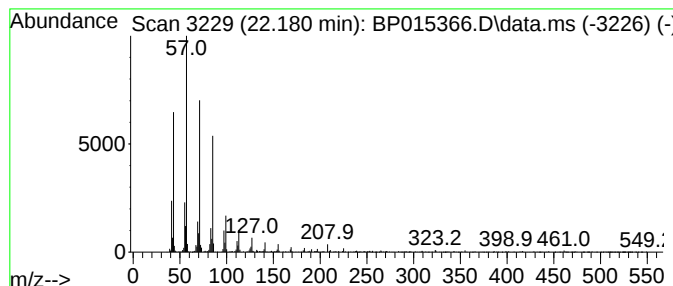
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.180	2.77 ng/ul	409961	Chrysene-d12		21.581	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	97
2	Tetracosane		338	C24H50	000646-31-1	96
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
4	Tricosane		324	C23H48	000638-67-5	93
5	2-Methylpentacosane		366	C26H54	000629-87-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
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ALS Vial : 16 Sample Multiplier: 1

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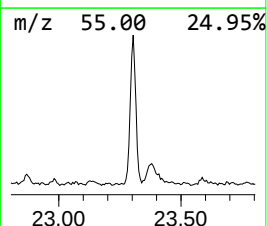
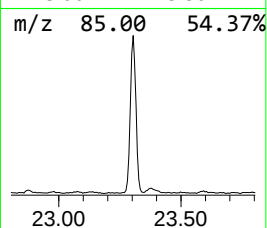
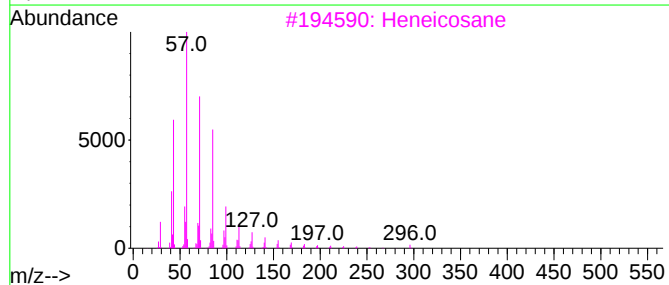
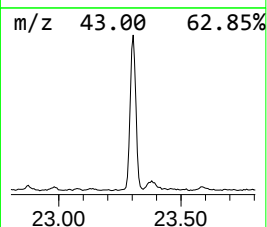
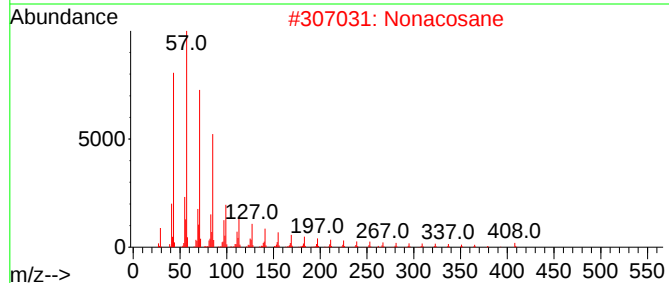
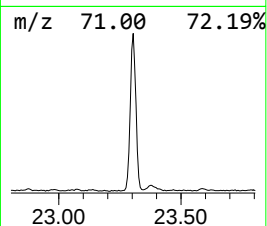
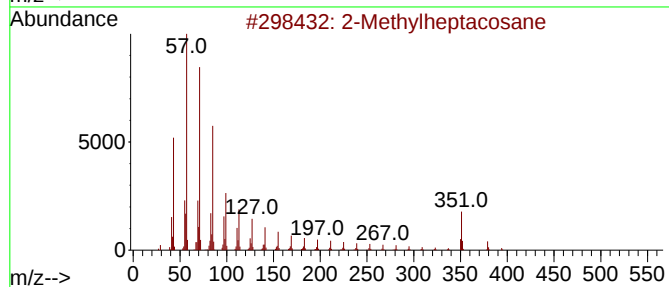
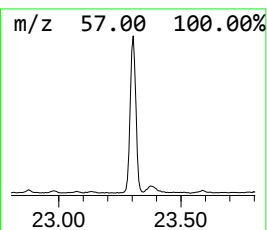
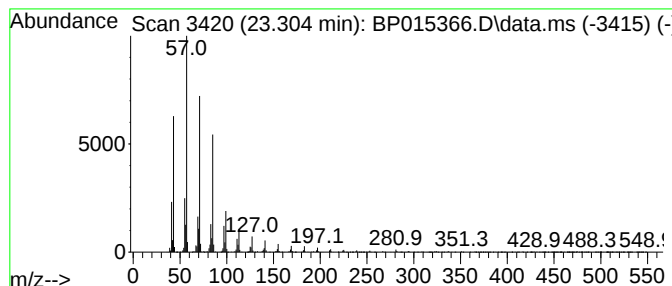
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.304	12.18 ng/ul	1747940	Perylene-d12	24.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylheptacosane			394	C28H58	001561-00-8	94
2	Nonacosane			408	C29H60	000630-03-5	93
3	Heneicosane			296	C21H44	000629-94-7	91
4	Dotriacontane, 1-iodo-			576	C32H65I	1000406-32-4	91
5	Tetracosane, 1-iodo-			464	C24H49I	1000406-32-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
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Misc :
ALS Vial : 16 Sample Multiplier: 1

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ClientSampleId :
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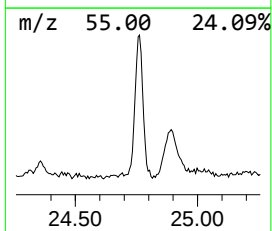
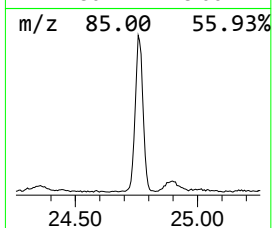
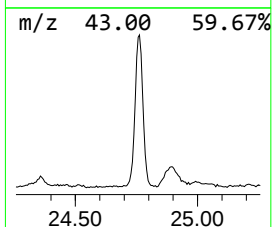
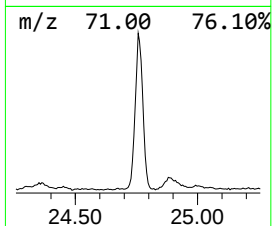
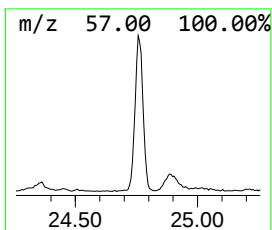
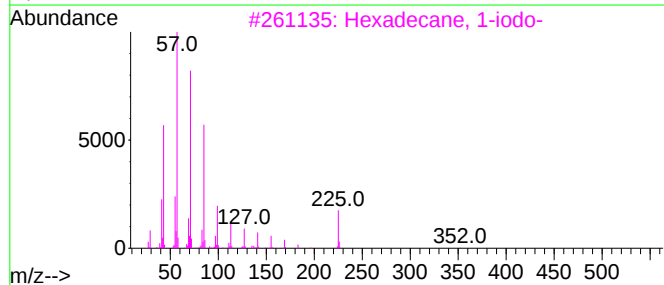
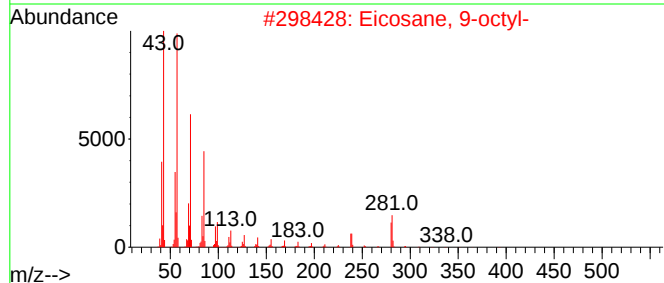
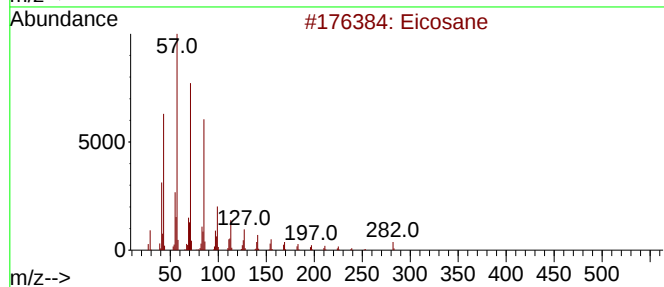
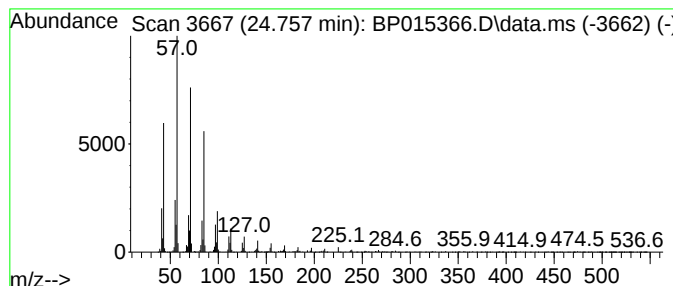
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.757	9.12 ng/ul	1308220	Perylene-d12	24.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
3			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	93
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053123\
Data File : BP015366.D
Acq On : 31 May 2023 18:49
Operator : CG/JU
Sample : 02851-09
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0FK2

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Decanoic acid	13.028	3.9	ng/ul	392166	3	14.746	2001220	20.0
Benzophenone	16.098	11.4	ng/ul	1143070	3	14.746	2001220	20.0
Palmitoleic acid	18.245	2.1	ng/ul	262673	4	17.504	2550640	20.0
n-Hexadecanoic ...	18.357	7.3	ng/ul	934442	4	17.504	2550640	20.0
Tricyclo[3.3.3....	19.398	2.6	ng/ul	338046	4	17.504	2550640	20.0
Octadecanoic acid	19.581	2.2	ng/ul	331573	5	21.581	2964860	20.0
(DEL) Alkane: S...	21.251	3.8	ng/ul	558763	5	21.581	2964860	20.0
(DEL) Alkane: S...	22.180	2.8	ng/ul	409961	5	21.581	2964860	20.0
(DEL) Alkane: S...	23.304	12.2	ng/ul	1747940	6	24.133	2870010	20.0
(DEL) Alkane: S...	24.757	9.1	ng/ul	1308220	6	24.133	2870010	20.0