

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
 Data File : BP015625.D
 Acq On : 20 Jun 2023 17:03
 Operator : MA/JU
 Sample : 03080-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFH6

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: BP015625.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	34	37	43	rVB	13941	17701	0.69%	0.045%
2	3.852	110	113	127	rBV	125106	230379	8.98%	0.591%
3	4.075	147	151	155	rBV2	8827	12222	0.48%	0.031%
4	4.416	205	209	214	rVB6	5455	8487	0.33%	0.022%
5	6.052	484	487	492	rVB7	3058	5556	0.22%	0.014%
6	6.305	525	530	537	rBV4	7137	15749	0.61%	0.040%
7	7.246	682	690	705	rBV	253980	470095	18.32%	1.205%
8	7.352	705	708	711	rVV5	4128	5362	0.21%	0.014%
9	7.410	711	718	730	rVB	203036	348178	13.57%	0.893%
10	7.610	746	752	761	rBV	280623	469362	18.29%	1.203%
11	7.687	761	765	768	rVB5	5619	8092	0.32%	0.021%
12	8.087	826	833	843	rBV	326928	534378	20.82%	1.370%
13	8.634	921	926	932	rVB9	3013	7299	0.28%	0.019%
14	8.722	937	941	946	rBV7	2643	4993	0.19%	0.013%
15	8.804	948	955	966	rBV	282477	520640	20.29%	1.335%
16	8.928	972	976	981	rVB2	10788	17731	0.69%	0.045%
17	9.263	1026	1033	1047	rBV	282341	508911	19.83%	1.305%
18	9.993	1149	1157	1167	rBV	253371	445871	17.37%	1.143%
19	10.075	1167	1171	1180	rVB	5979	14384	0.56%	0.037%
20	10.246	1196	1200	1205	rBV8	2911	5336	0.21%	0.014%
21	10.540	1242	1250	1268	rBV	313092	644726	25.12%	1.653%
22	10.863	1299	1305	1306	rBV6	3815	5683	0.22%	0.015%
23	10.910	1306	1313	1320	rBV	488149	817366	31.85%	2.095%
24	11.063	1331	1339	1352	rBV	197624	458151	17.85%	1.174%
25	11.604	1426	1431	1437	rVB10	2479	4981	0.19%	0.013%
26	11.728	1445	1452	1454	rBV8	3748	7109	0.28%	0.018%
27	11.916	1479	1484	1486	rBV6	4357	5244	0.20%	0.013%
28	12.304	1546	1550	1554	rBV6	4728	7889	0.31%	0.020%
29	12.498	1575	1583	1590	rVB8	6709	17715	0.69%	0.045%
30	12.569	1590	1595	1599	rBV3	7513	14908	0.58%	0.038%
31	12.787	1628	1632	1636	rBV5	6265	8380	0.33%	0.021%
32	13.245	1704	1710	1719	rVB5	3596	9729	0.38%	0.025%
33	13.534	1755	1759	1762	rBV5	3668	5458	0.21%	0.014%
34	13.610	1762	1772	1776	rVV8	14263	28901	1.13%	0.074%
35	13.687	1779	1785	1787	rVV7	7604	13773	0.54%	0.035%
36	13.710	1787	1789	1795	rVB3	10113	12623	0.49%	0.032%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	13.775	1795	1800	1804	rBV8	2817	5162	0.20%	0.013%
38	13.898	1816	1821	1822	rBV5	3936	4882	0.19%	0.013%
39	14.051	1842	1847	1851	rVB4	9500	15365	0.60%	0.039%
40	14.134	1851	1861	1867	rBV	738695	1042332	40.61%	2.672%
41	14.239	1876	1879	1883	rBV6	5982	7150	0.28%	0.018%
42	14.422	1904	1910	1927	rVV	796609	1273787	49.63%	3.265%
43	14.545	1929	1931	1936	rVV6	4451	5640	0.22%	0.014%
44	14.604	1936	1941	1946	rVB4	8357	12229	0.48%	0.031%
45	14.687	1949	1955	1956	rBV5	9618	14319	0.56%	0.037%
46	14.728	1956	1962	1970	rVV	787238	1175336	45.79%	3.013%
47	14.787	1970	1972	1980	rVV3	14193	25612	1.00%	0.066%
48	14.963	1996	2002	2023	rBV	145451	439408	17.12%	1.126%
49	15.128	2027	2030	2036	rVB4	18632	29472	1.15%	0.076%
50	15.398	2074	2076	2086	rVB4	5099	9641	0.38%	0.025%
51	15.545	2099	2101	2106	rVB3	12248	15665	0.61%	0.040%
52	15.722	2124	2131	2137	rBV	1077432	1551941	60.47%	3.978%
53	15.857	2148	2154	2166	rBV	293233	448005	17.46%	1.148%
54	16.086	2187	2193	2204	rBV	345869	493008	19.21%	1.264%
55	16.304	2228	2230	2236	rVB7	6381	9445	0.37%	0.024%
56	16.439	2245	2253	2265	rBV10	22990	76553	2.98%	0.196%
57	16.575	2272	2276	2281	rBV8	10581	17160	0.67%	0.044%
58	16.651	2284	2289	2301	rVB	87092	133594	5.21%	0.342%
59	16.833	2316	2320	2322	rBV5	4336	6855	0.27%	0.018%
60	16.857	2322	2324	2330	rVB6	13762	21059	0.82%	0.054%
61	16.904	2330	2332	2334	rBV3	4384	4703	0.18%	0.012%
62	17.016	2348	2351	2354	rBV5	8737	11183	0.44%	0.029%
63	17.151	2370	2374	2378	rVB4	15588	24216	0.94%	0.062%
64	17.210	2381	2384	2390	rBV3	21327	35839	1.40%	0.092%
65	17.304	2397	2400	2405	rVB4	10174	14982	0.58%	0.038%
66	17.357	2405	2409	2417	rBV2	41156	67851	2.64%	0.174%
67	17.428	2417	2421	2425	rBV3	31113	42079	1.64%	0.108%
68	17.486	2425	2431	2435	rVV	1109329	1544855	60.19%	3.960%
69	17.528	2435	2438	2443	rVV	204717	306545	11.94%	0.786%
70	17.586	2443	2448	2461	rVB	1244856	1773621	69.10%	4.547%
71	17.898	2498	2501	2506	rVB	29549	40093	1.56%	0.103%
72	17.951	2507	2510	2513	rBV2	14639	14731	0.57%	0.038%
73	18.128	2537	2540	2543	rVB5	16457	20826	0.81%	0.053%
74	18.233	2554	2558	2562	rVB5	30107	39688	1.55%	0.102%
75	18.286	2563	2567	2572	rBV	264404	339945	13.25%	0.871%
76	18.345	2572	2577	2582	rBV	1213700	1475099	57.47%	3.781%

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77	18.533	2605	2609	2613	rBV2	43979	70168	2.73%	0.180%
78	18.622	2621	2624	2631	rBV5	28536	53132	2.07%	0.136%
79	18.875	2664	2667	2672	rBV	52164	64075	2.50%	0.164%
80	19.222	2723	2726	2731	rBV3	60568	81961	3.19%	0.210%
81	19.327	2740	2744	2748	rBV	104054	143248	5.58%	0.367%
82	19.427	2758	2761	2763	rBV2	82304	99213	3.87%	0.254%
83	19.457	2763	2766	2767	rVV2	153432	182676	7.12%	0.468%
84	19.480	2767	2770	2777	rVV	762861	1084312	42.25%	2.780%
85	19.569	2781	2785	2792	rVV2	334889	449378	17.51%	1.152%
86	19.810	2821	2826	2829	rBV	1826968	2507604	97.70%	6.428%
87	19.839	2829	2831	2836	rVB	584104	628839	24.50%	1.612%
88	19.951	2846	2850	2855	rVB	175811	218522	8.51%	0.560%
89	20.616	2960	2963	2967	rBV4	97370	139113	5.42%	0.357%
90	21.051	3034	3037	3040	rBV	120501	146763	5.72%	0.376%
91	21.245	3066	3070	3075	rVB	556377	775410	30.21%	1.988%
92	21.569	3118	3125	3129	rBV2	1600655	2566560	100.00%	6.579%
93	21.610	3129	3132	3139	rVB	381573	479890	18.70%	1.230%
94	22.168	3223	3227	3230	rVB2	292853	361769	14.10%	0.927%
95	23.280	3411	3416	3421	rBV	964509	1507474	58.74%	3.864%
96	23.351	3422	3428	3440	rVB2	525530	1805834	70.36%	4.629%
97	23.951	3525	3530	3536	rBV2	1186663	2136114	83.23%	5.476%
98	24.121	3553	3559	3565	rVB	792706	1570846	61.20%	4.027%
99	24.733	3657	3663	3672	rVB	1112348	2370949	92.38%	6.078%
100	26.715	3994	4000	4009	rBV	439204	1261981	49.17%	3.235%

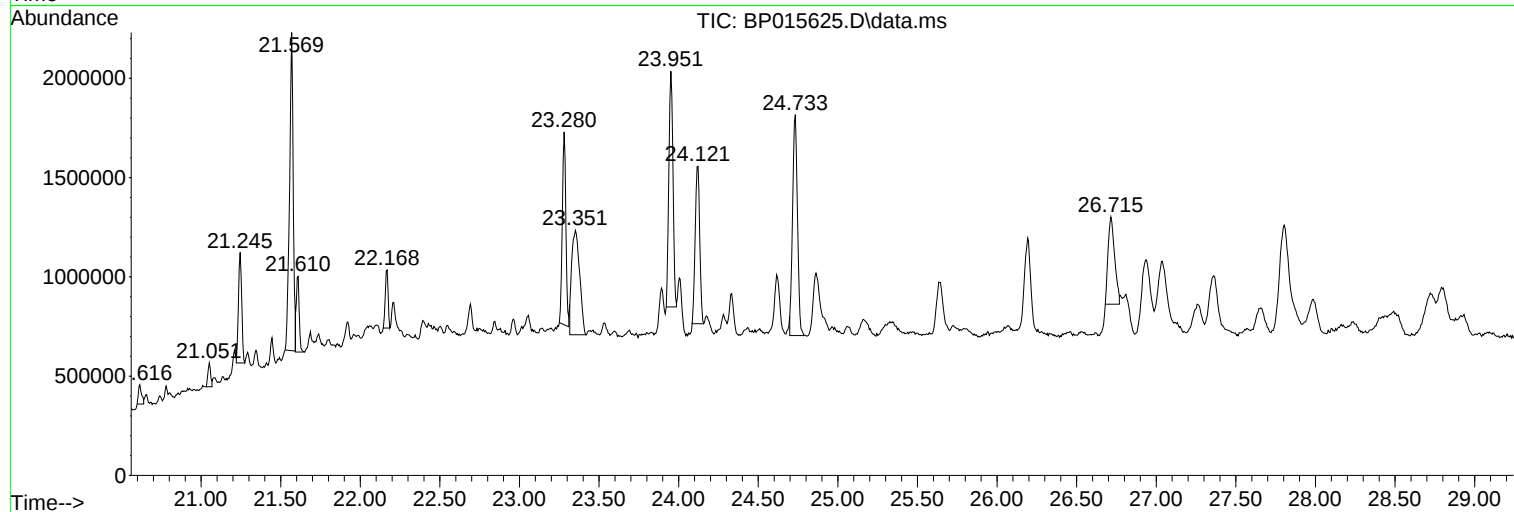
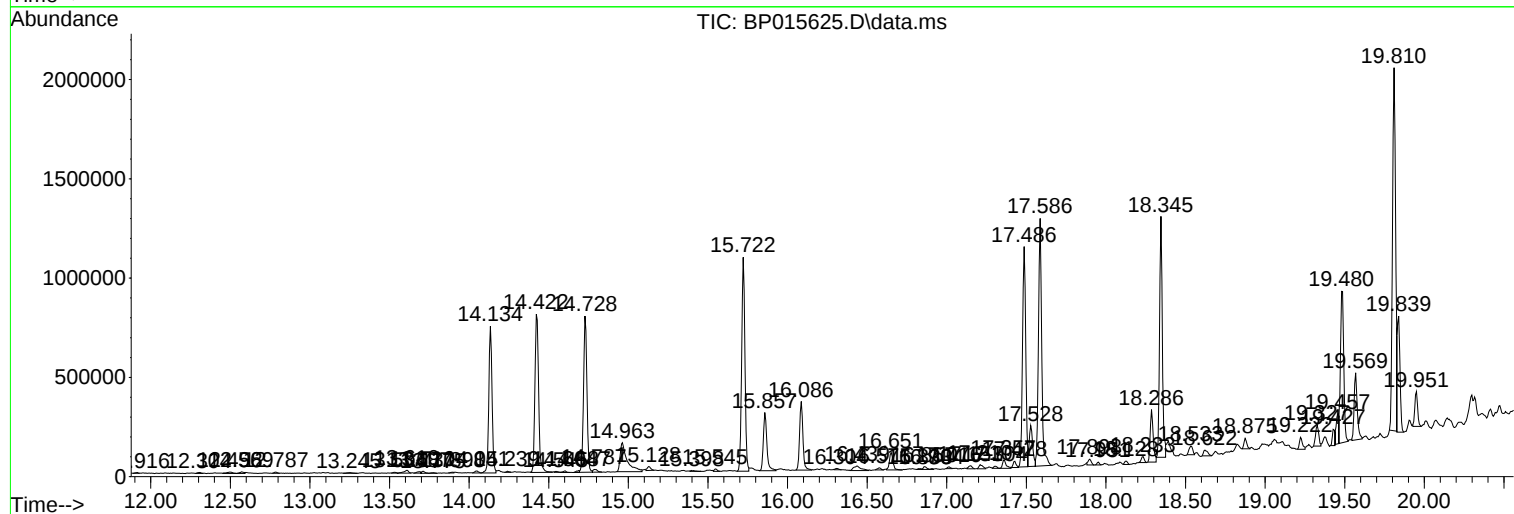
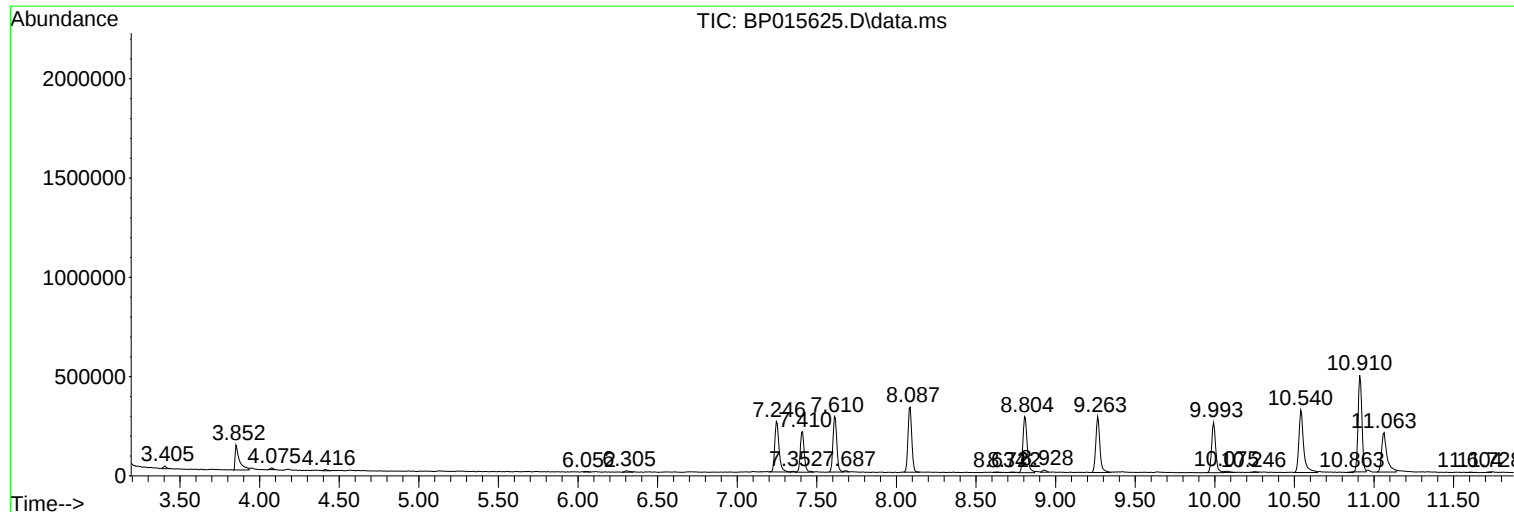
Sum of corrected areas: 39009069

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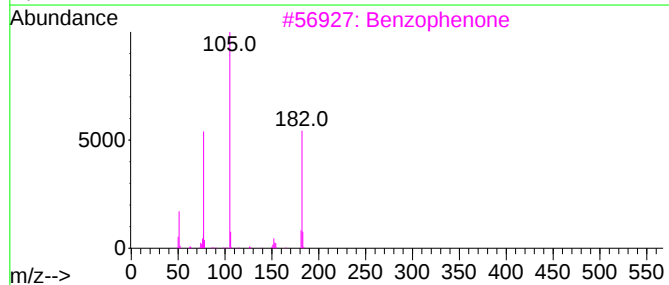
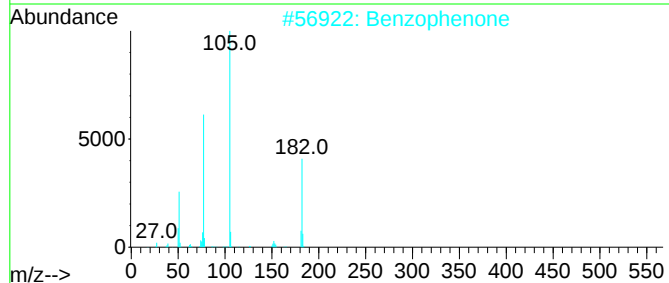
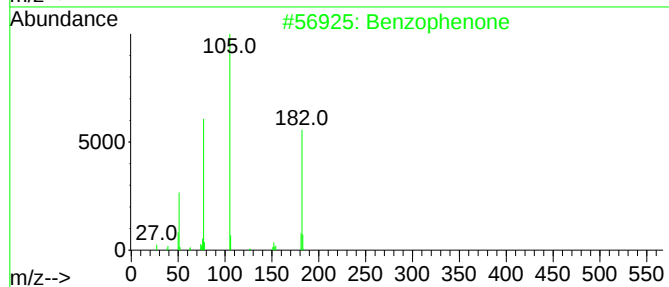
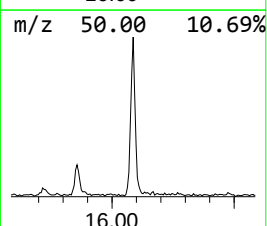
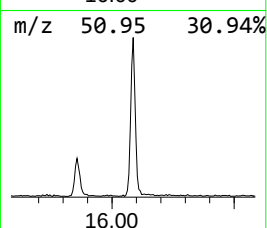
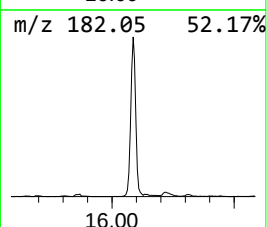
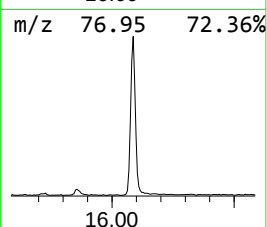
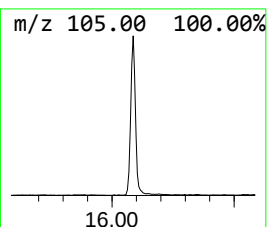
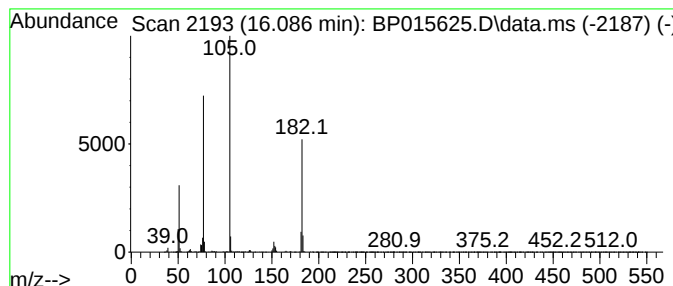
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.086	8.39 ng/ul	493008	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			1,2,4-Trioxolane, 3,3,5-triphenyl-	304	C20H16O3	023246-12-0	64



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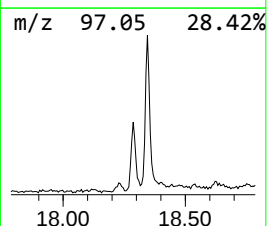
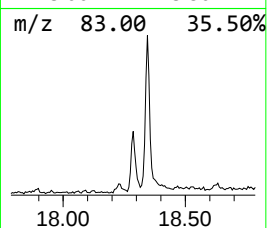
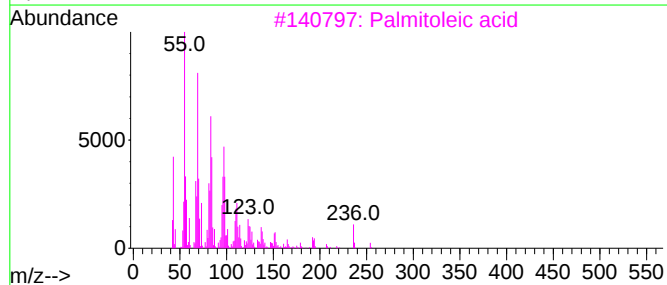
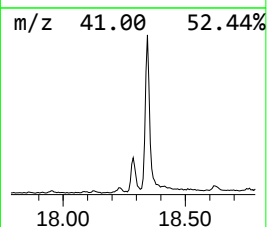
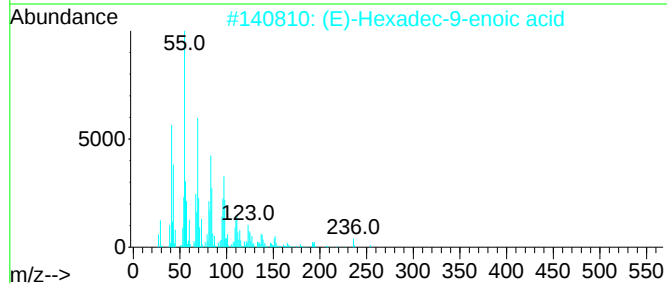
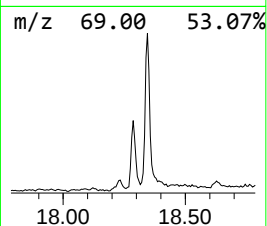
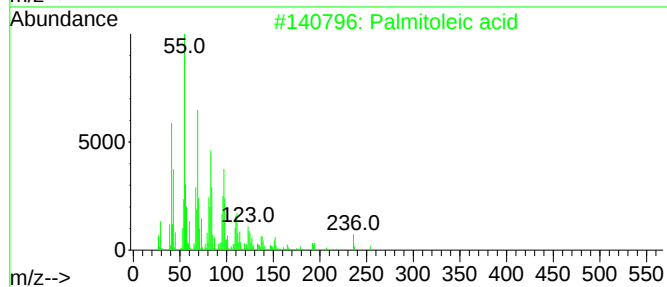
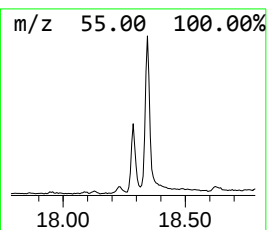
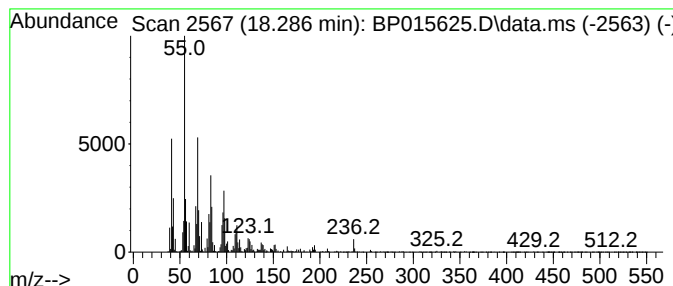
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Palmitoleic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	4.40 ng/ul	339945	Phenanthrene-d10	17.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	99
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
3			Palmitoleic acid	254	C16H30O2	000373-49-9	98
4			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	98
5			Myristoleic acid	226	C14H26O2	000544-64-9	97



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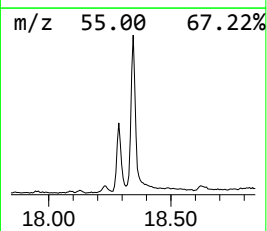
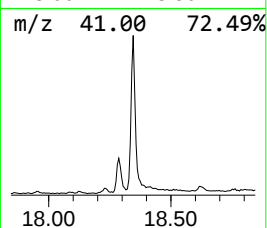
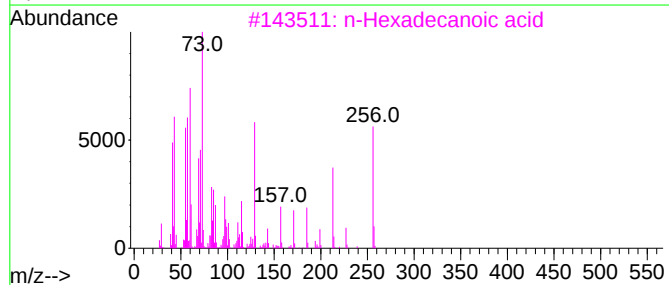
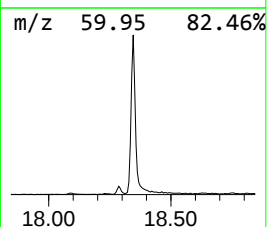
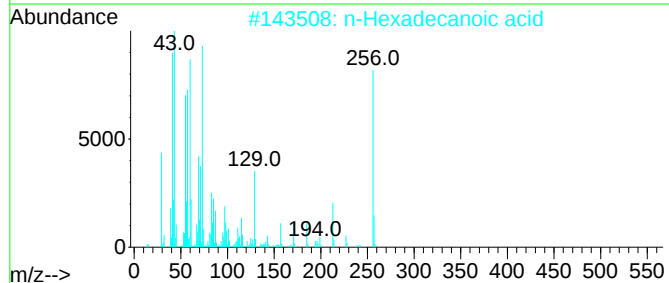
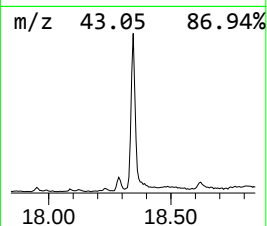
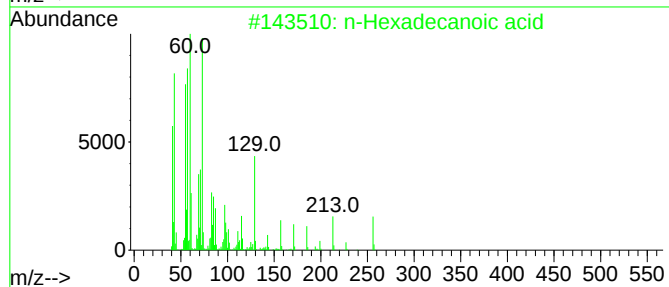
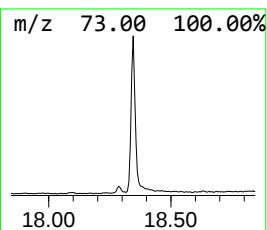
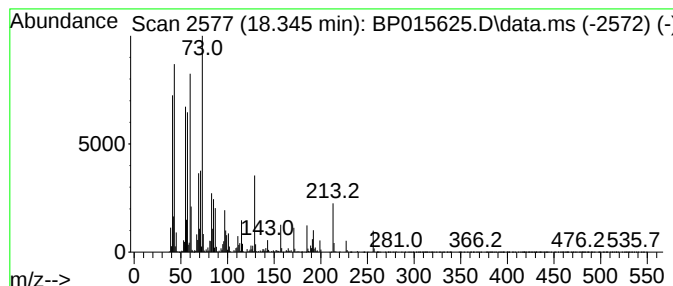
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	19.10 ng/ul	1475100	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	96		
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96		
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFH6

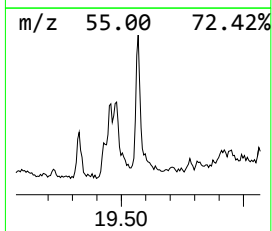
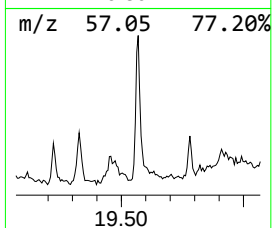
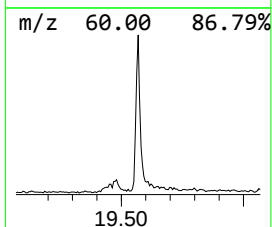
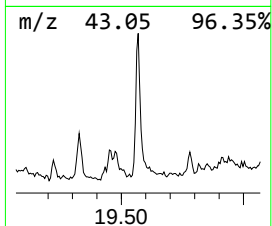
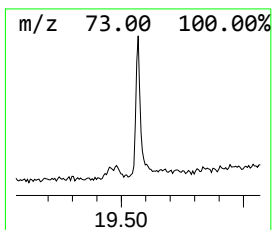
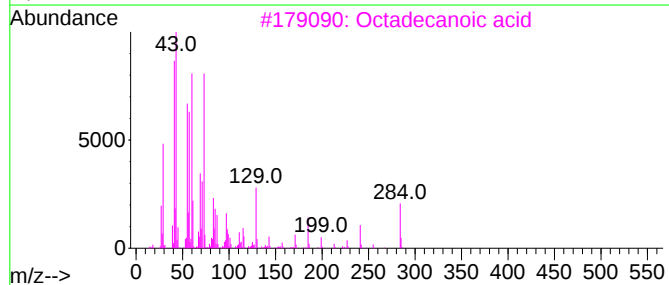
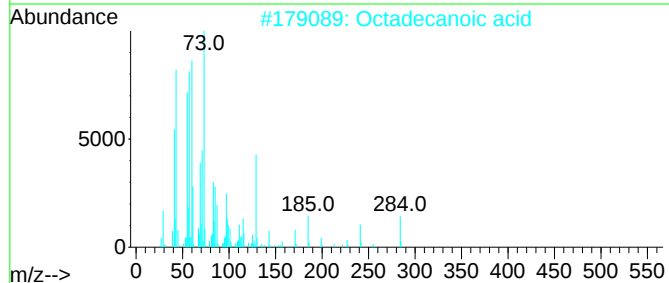
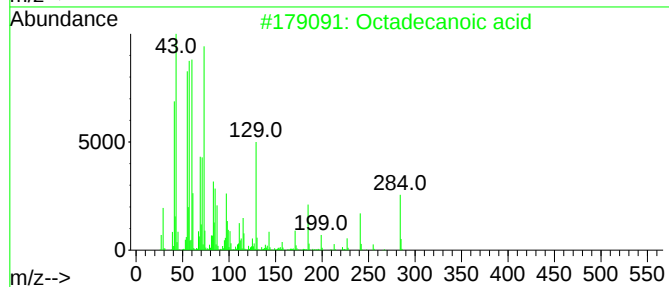
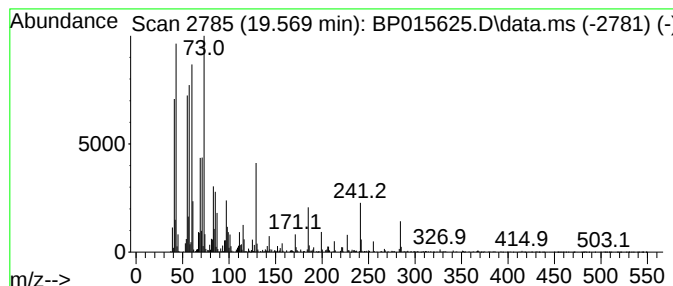
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.569	3.50 ng/ul	449378	Chrysene-d12	21.569

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	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

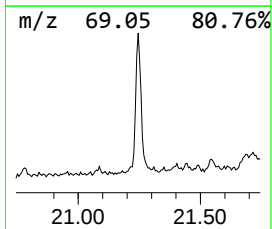
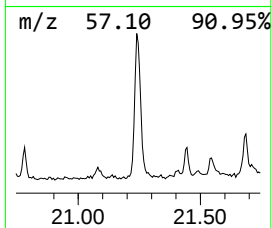
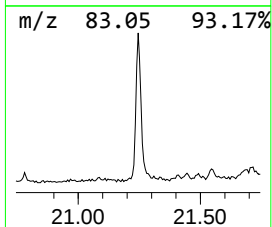
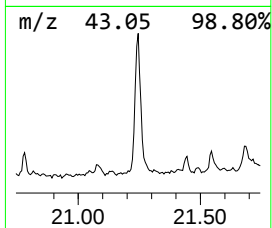
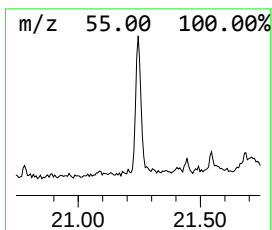
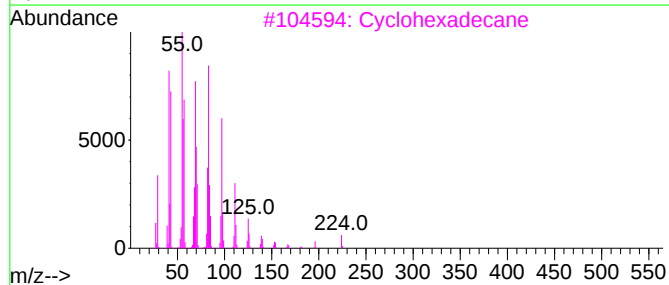
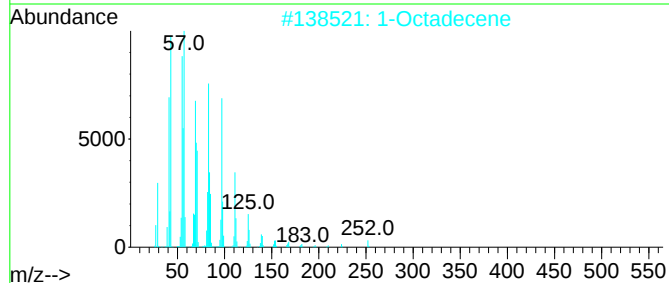
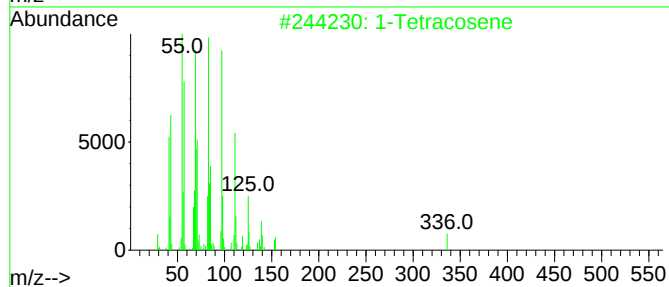
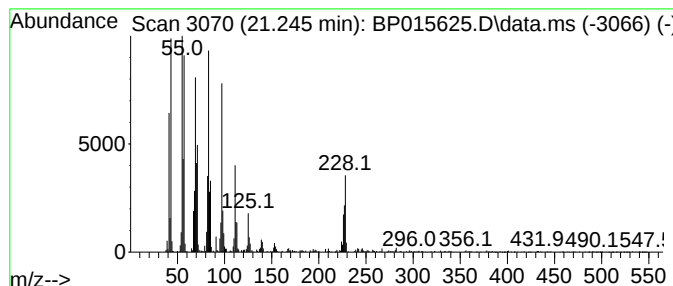
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.245	6.04 ng/ul	775410	Chrysene-d12	21.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene	336	C24H48	010192-32-2	96
2	1-Octadecene	252	C18H36	000112-88-9	95
3	Cyclohexadecane	224	C16H32	000295-65-8	94
4	1-Tricosene	322	C23H46	018835-32-0	93
5	Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

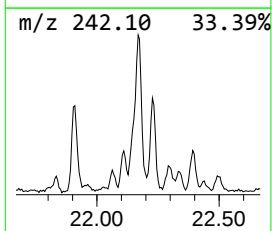
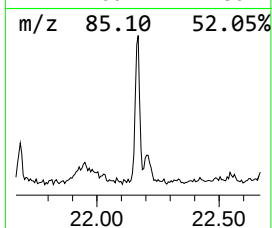
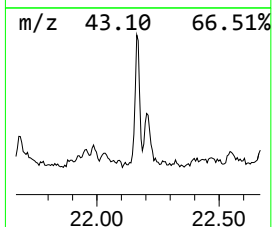
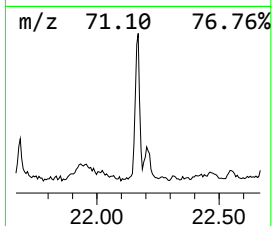
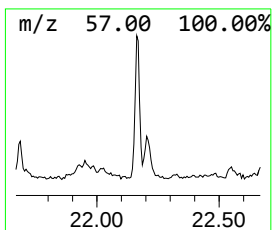
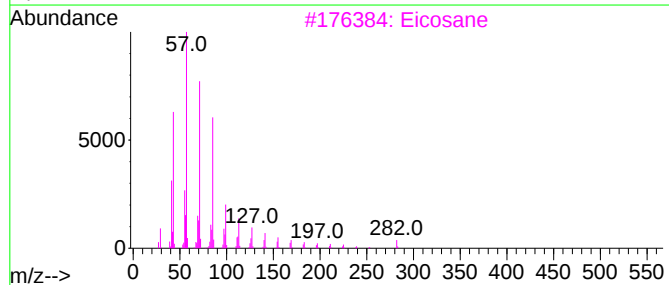
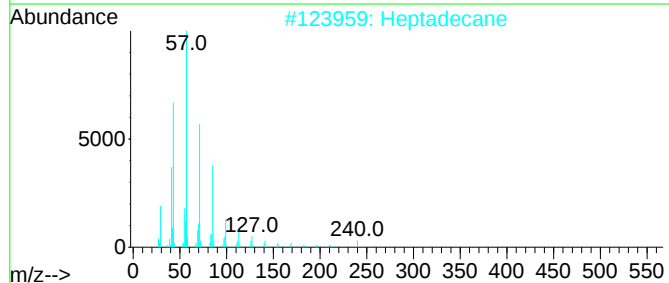
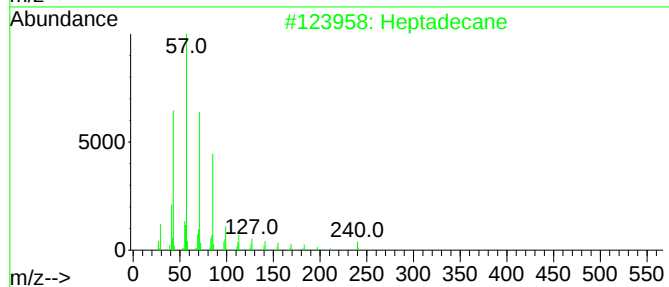
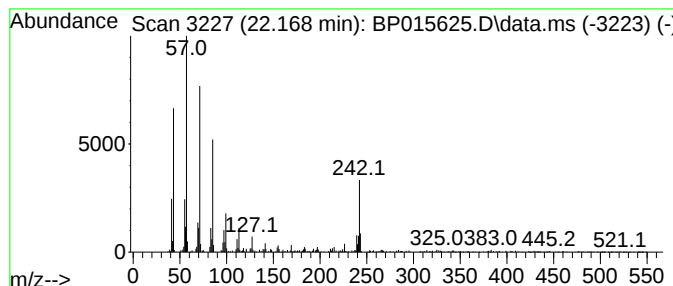
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.169	2.82 ng/ul	361769	Chrysene-d12	21.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Heptadecane	240	C17H36	000629-78-7	91
3	Eicosane	282	C20H42	000112-95-8	91
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5	Nonadecane	268	C19H40	000629-92-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

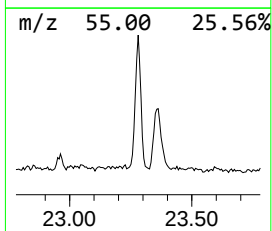
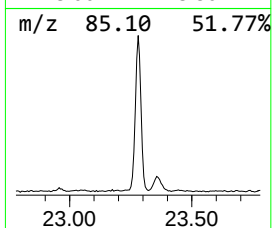
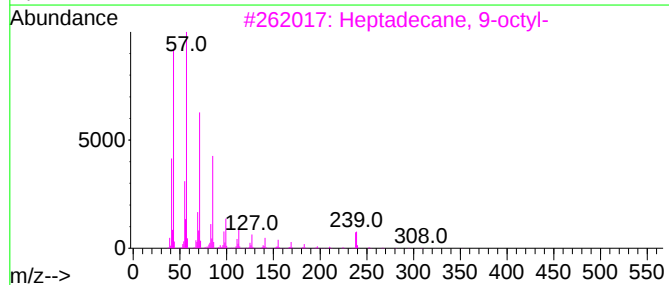
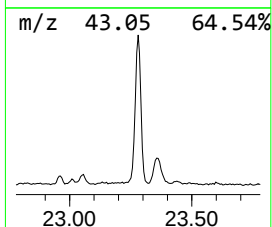
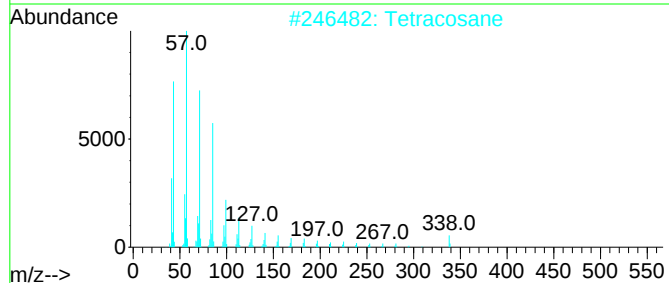
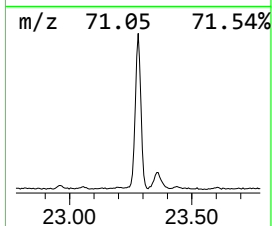
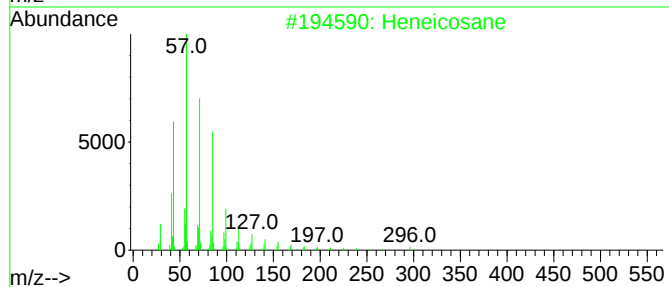
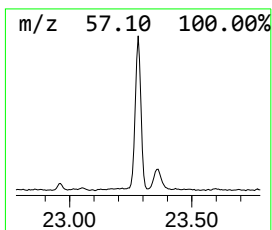
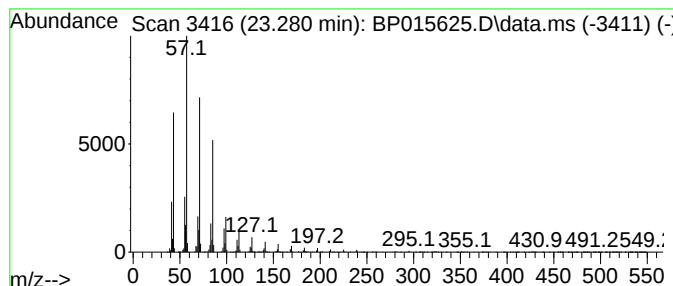
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.280	19.19 ng/ul	1507470	Perylene-d12	24.121	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	98
2	Tetracosane	338	C24H50	000646-31-1	97
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
4	Tricosane	324	C23H48	000638-67-5	94
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

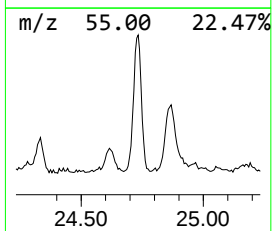
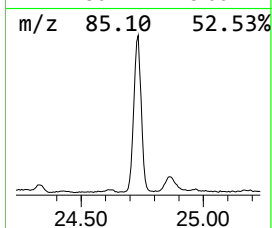
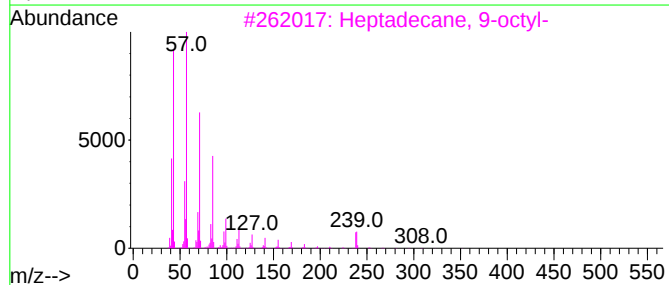
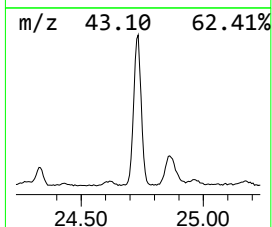
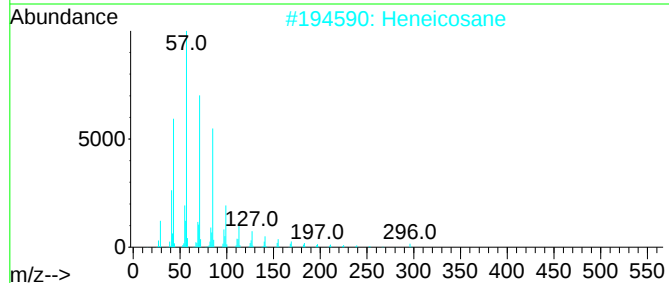
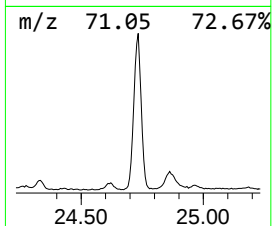
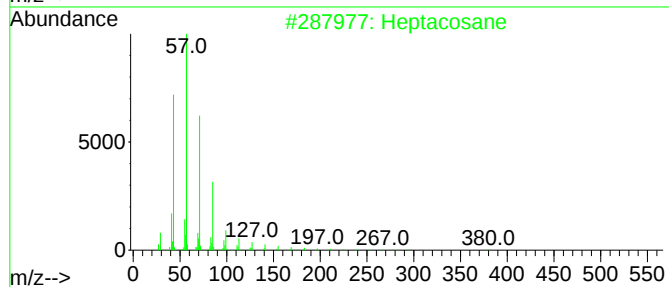
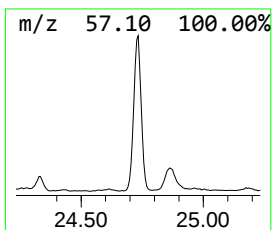
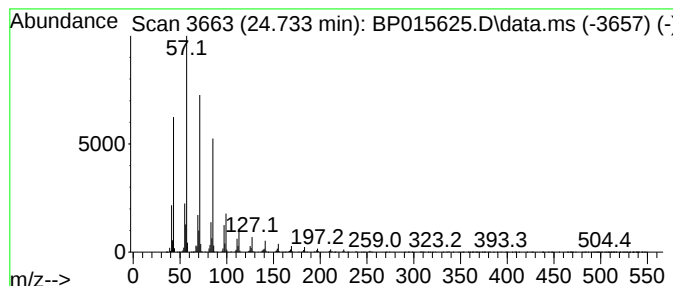
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.733	30.19 ng/ul	2370950	Perylene-d12	24.121	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptacosane	380	C27H56	000593-49-7	95
2	Heneicosane	296	C21H44	000629-94-7	91
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Tetratetracontane	619	C44H90	007098-22-8	91
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

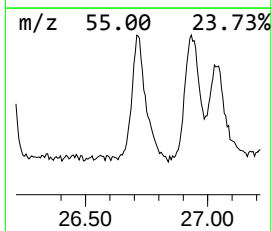
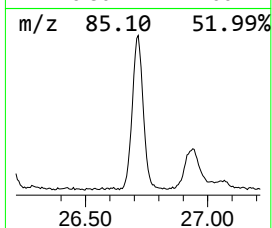
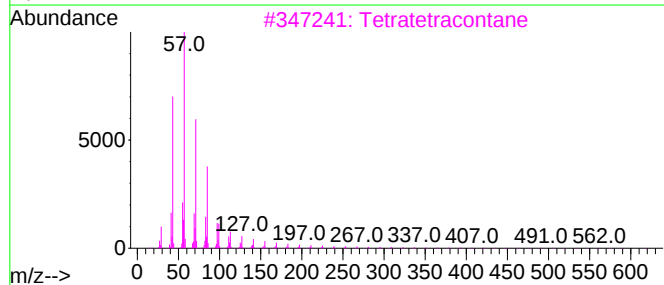
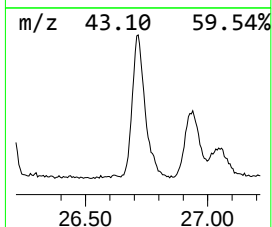
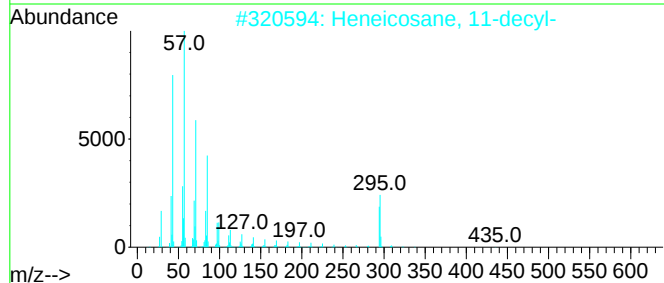
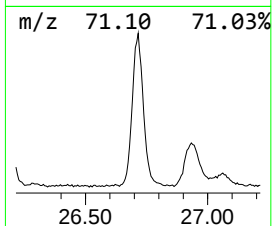
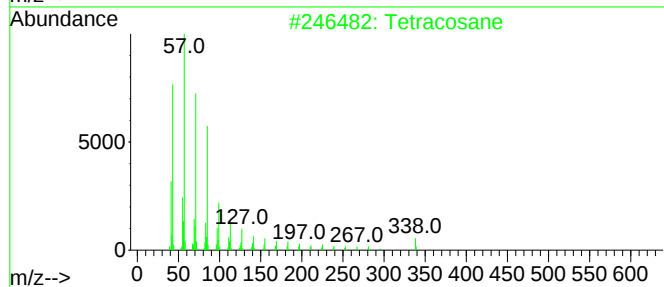
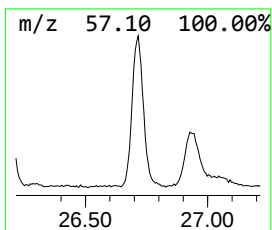
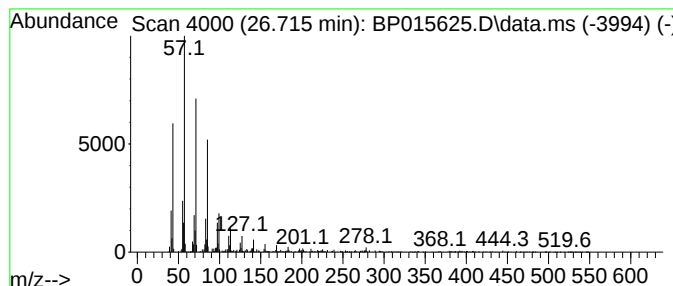
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.715	16.07 ng/ul	1261980	Perylene-d12		24.121	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetracosane		338	C24H50	000646-31-1	97
2	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	93
3	Tetratetracontane		619	C44H90	007098-22-8	93
4	Tetratriacontane		479	C34H70	014167-59-0	93
5	Tetracosane		338	C24H50	000646-31-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015625.D
Acq On : 20 Jun 2023 17:03
Operator : MA/JU
Sample : 03080-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFH6

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.086	8.4	ng/ul	493008	3	14.728	1175340	20.0
Palmitoleic acid	18.286	4.4	ng/ul	339945	4	17.486	1544860	20.0
n-Hexadecanoic ...	18.345	19.1	ng/ul	1475100	4	17.486	1544860	20.0
Octadecanoic acid	19.569	3.5	ng/ul	449378	5	21.569	2566560	20.0
(DEL) Alkane: S...	21.245	6.0	ng/ul	775410	5	21.569	2566560	20.0
(DEL) Alkane: S...	22.169	2.8	ng/ul	361769	5	21.569	2566560	20.0
(DEL) Alkane: S...	23.280	19.2	ng/ul	1507470	6	24.121	1570850	20.0
(DEL) Alkane: S...	24.733	30.2	ng/ul	2370950	6	24.121	1570850	20.0
(DEL) Alkane: S...	26.715	16.1	ng/ul	1261980	6	24.121	1570850	20.0