

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
 Data File : BP020195.D
 Acq On : 04 May 2024 05:33
 Operator : MA/JU
 Sample : P2234-01
 Misc :
 ALS Vial : 64 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0SA1

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020195.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.634	90	93	102	rVB	43085	74553	2.81%	0.216%
2	5.311	374	378	386	rVB4	13735	26889	1.01%	0.078%
3	5.669	432	439	445	rVB4	7215	14957	0.56%	0.043%
4	6.811	627	633	649	rBV	275522	554472	20.89%	1.608%
5	6.975	655	661	675	rVB	203686	351726	13.25%	1.020%
6	7.158	686	692	703	rBV	271664	486785	18.34%	1.412%
7	7.622	763	771	783	rBV	305688	521119	19.63%	1.512%
8	8.334	885	892	907	rBV	287634	570553	21.49%	1.655%
9	8.452	907	912	922	rVB	105185	183194	6.90%	0.531%
10	8.787	962	969	984	rBV	291421	532190	20.05%	1.544%
11	9.134	1024	1028	1035	rVB4	5572	9323	0.35%	0.027%
12	9.375	1065	1069	1081	rBV7	5436	14272	0.54%	0.041%
13	9.493	1082	1089	1103	rBV	244024	469950	17.70%	1.363%
14	9.716	1121	1127	1132	rBV5	8078	21103	0.80%	0.061%
15	9.869	1148	1153	1159	rBV3	8992	15642	0.59%	0.045%
16	10.034	1173	1181	1198	rBV	289516	683699	25.76%	1.983%
17	10.240	1210	1216	1221	rBV9	5030	10424	0.39%	0.030%
18	10.399	1234	1243	1259	rVB	428872	755577	28.46%	2.192%
19	10.563	1263	1271	1287	rBV	134862	340703	12.84%	0.988%
20	10.940	1330	1335	1340	rBV7	5917	11819	0.45%	0.034%
21	11.187	1370	1377	1386	rBV3	11390	34996	1.32%	0.102%
22	12.010	1512	1517	1525	rVV3	5948	12392	0.47%	0.036%
23	13.169	1709	1714	1723	rVB3	10145	19939	0.75%	0.058%
24	13.716	1800	1807	1820	rBV	754800	1100929	41.47%	3.194%
25	13.981	1845	1852	1870	rBV2	761983	1241634	46.78%	3.602%
26	14.293	1896	1905	1913	rVV	624650	1049062	39.52%	3.043%
27	14.369	1913	1918	1928	rVB2	13374	26915	1.01%	0.078%
28	14.569	1937	1952	1962	rBV2	154116	471948	17.78%	1.369%
29	14.734	1975	1980	1987	rBV9	10679	27189	1.02%	0.079%
30	15.181	2051	2056	2060	rVB6	6423	9597	0.36%	0.028%
31	15.322	2073	2080	2087	rBV	1017770	1600723	60.30%	4.643%
32	15.381	2087	2090	2098	rVB3	13505	27495	1.04%	0.080%
33	15.469	2099	2105	2115	rBV	290249	514369	19.38%	1.492%
34	15.687	2138	2142	2149	rVB4	7055	13559	0.51%	0.039%
35	15.845	2163	2169	2174	rVB8	4928	10214	0.38%	0.030%
36	16.010	2192	2197	2203	rBV	12639	26716	1.01%	0.077%

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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-BP050124.MA.M
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37	16.092	2203	2211	2217	rVV7	12696	27028	1.02%	0.078%
38	16.416	2259	2266	2272	rBV7	5321	13714	0.52%	0.040%
39	16.540	2283	2287	2291	rBV5	9013	15011	0.57%	0.044%
40	16.792	2324	2330	2337	rVB2	21078	35406	1.33%	0.103%
41	16.857	2337	2341	2345	rBV5	6222	10403	0.39%	0.030%
42	16.945	2352	2356	2363	rVB2	18908	31445	1.18%	0.091%
43	17.039	2367	2372	2379	rBV3	27839	47129	1.78%	0.137%
44	17.128	2379	2387	2391	rBV	946036	1365419	51.44%	3.961%
45	17.169	2391	2394	2399	rVV	410912	588776	22.18%	1.708%
46	17.234	2399	2405	2418	rVB2	1131439	1790027	67.43%	5.192%
47	17.334	2418	2422	2425	rBV6	9704	18533	0.70%	0.054%
48	17.575	2456	2463	2470	rVB2	32189	63492	2.39%	0.184%
49	17.669	2475	2479	2484	rBV4	13202	27737	1.04%	0.080%
50	17.816	2501	2504	2507	rBV5	7088	10501	0.40%	0.030%
51	17.963	2523	2529	2535	rBV4	45662	78204	2.95%	0.227%
52	18.028	2535	2540	2546	rBV2	453940	714739	26.93%	2.073%
53	18.092	2546	2551	2560	rVB2	1087746	1523212	57.38%	4.418%
54	18.245	2573	2577	2581	rBV2	60541	100667	3.79%	0.292%
55	18.286	2582	2584	2592	rVB	29147	40044	1.51%	0.116%
56	18.398	2597	2603	2607	rBV6	44424	76860	2.90%	0.223%
57	18.539	2622	2627	2629	rBV5	15949	23736	0.89%	0.069%
58	18.581	2629	2634	2637	rVV	39405	59331	2.24%	0.172%
59	18.628	2637	2642	2648	rVV	118546	182051	6.86%	0.528%
60	18.686	2649	2652	2666	rVB3	29230	62090	2.34%	0.180%
61	18.786	2666	2669	2670	rBV2	13475	12776	0.48%	0.037%
62	18.886	2682	2686	2688	rBV2	12176	15422	0.58%	0.045%
63	18.922	2689	2692	2696	rVB2	21507	31038	1.17%	0.090%
64	19.004	2702	2706	2711	rBV2	25891	36451	1.37%	0.106%
65	19.057	2711	2715	2722	rBV5	42760	84151	3.17%	0.244%
66	19.163	2727	2733	2744	rBV4	73659	206905	7.79%	0.600%
67	19.281	2747	2753	2757	rBV	1339170	1883866	70.97%	5.465%
68	19.433	2774	2779	2788	rBV2	412068	580079	21.85%	1.683%
69	19.633	2806	2813	2816	rBV	1516502	2654451	100.00%	7.700%
70	19.663	2816	2818	2823	rVV	1015542	1082540	40.78%	3.140%
71	19.710	2824	2826	2829	rVB3	23254	18920	0.71%	0.055%
72	19.745	2829	2832	2836	rVB2	28875	39010	1.47%	0.113%
73	19.786	2837	2839	2843	rVB5	23816	26520	1.00%	0.077%
74	19.857	2847	2851	2856	rVB	45357	65672	2.47%	0.190%
75	19.922	2859	2862	2867	rBV4	16062	26378	0.99%	0.077%
76	20.010	2869	2877	2878	rBV3	38831	80713	3.04%	0.234%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 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 >
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77	20.180	2903	2906	2911	rVB	63084	94013	3.54%	0.273%
78	20.239	2912	2916	2921	rVB2	56174	71361	2.69%	0.207%
79	20.298	2922	2926	2930	rBV	47813	65551	2.47%	0.190%
80	20.351	2930	2935	2939	rBV2	55518	97944	3.69%	0.284%
81	20.492	2956	2959	2963	rVB	29977	38336	1.44%	0.111%
82	20.551	2965	2969	2970	rBV3	26371	32798	1.24%	0.095%
83	20.757	3000	3004	3011	rBV4	69846	131847	4.97%	0.382%
84	20.992	3041	3044	3052	rVB	75013	107327	4.04%	0.311%
85	21.175	3071	3075	3079	rBV	77489	107156	4.04%	0.311%
86	21.222	3079	3083	3086	rVV	43641	57483	2.17%	0.167%
87	21.304	3093	3097	3102	rBV4	212236	342818	12.91%	0.994%
88	21.539	3134	3137	3141	rBV	90415	139647	5.26%	0.405%
89	21.610	3142	3149	3154	rVV2	903675	1860707	70.10%	5.397%
90	21.663	3154	3158	3167	rVB	376013	661232	24.91%	1.918%
91	21.880	3191	3195	3200	rBV	71070	106197	4.00%	0.308%
92	22.533	3301	3306	3309	rBV	102019	145951	5.50%	0.423%
93	23.645	3491	3495	3502	rVB3	107548	199788	7.53%	0.580%
94	23.910	3533	3540	3548	rBV	281558	872282	32.86%	2.530%
95	24.139	3574	3579	3586	rBV2	136640	288839	10.88%	0.838%
96	24.233	3588	3595	3612	rVB	214018	681986	25.69%	1.978%
97	24.651	3661	3666	3672	rBV	121929	235881	8.89%	0.684%
98	24.733	3672	3680	3687	rBV2	509351	1383031	52.10%	4.012%
99	24.968	3711	3720	3728	rBV2	335460	1007124	37.94%	2.921%
100	26.363	3953	3957	3958	rBV2	132660	157484	5.93%	0.457%

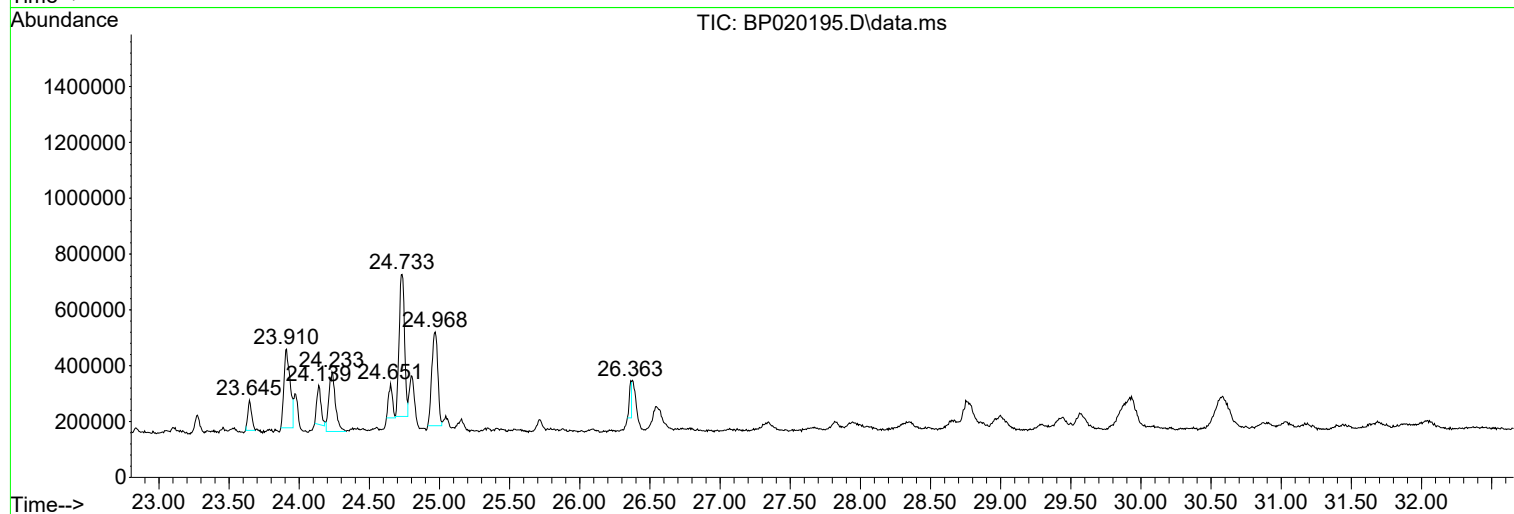
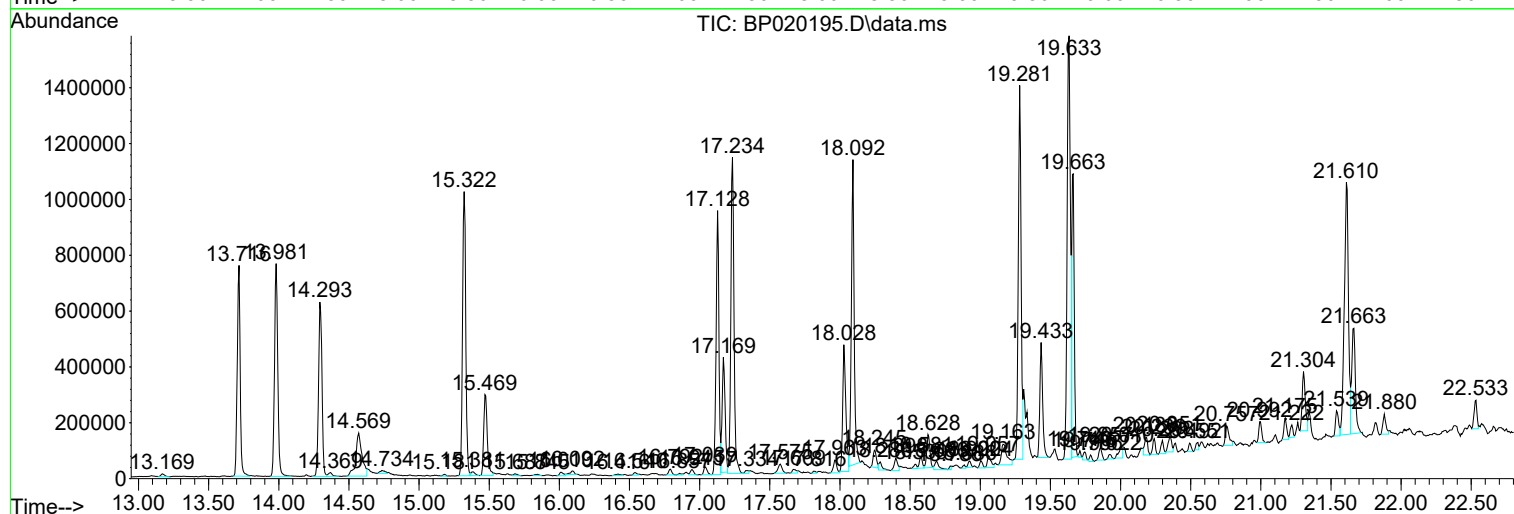
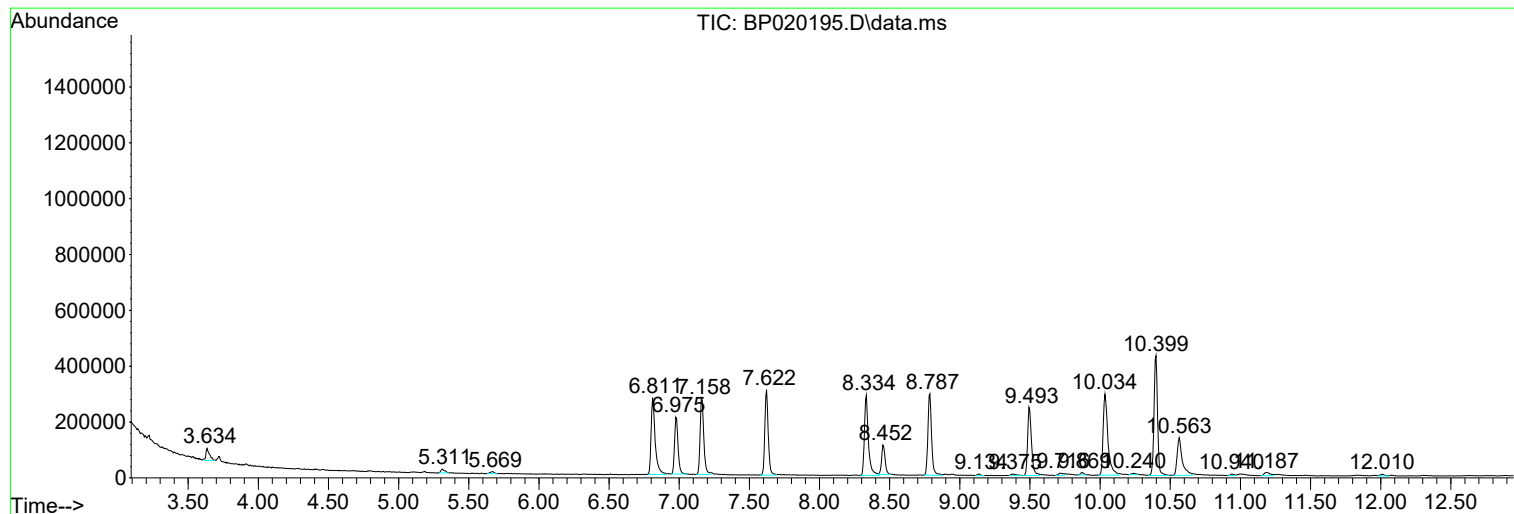
Sum of corrected areas: 34473857

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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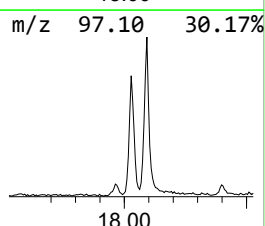
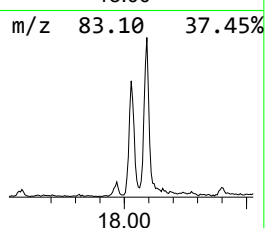
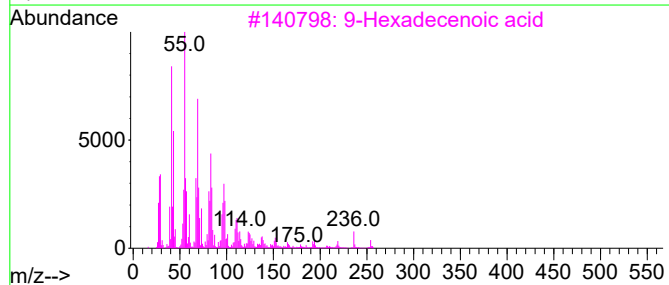
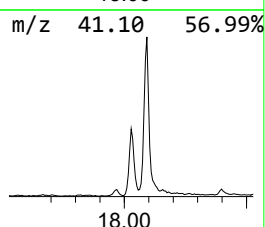
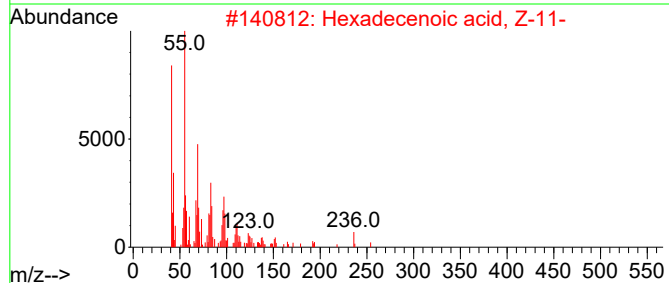
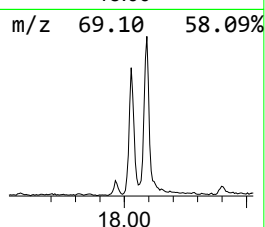
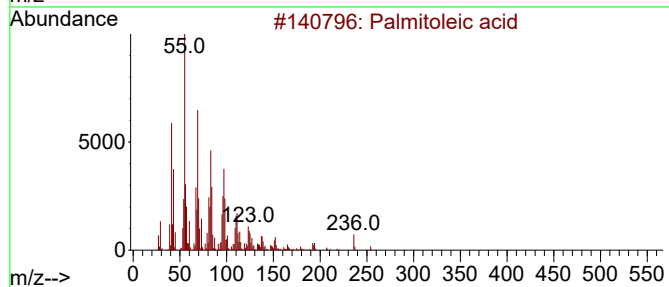
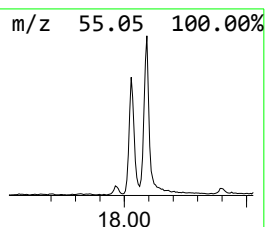
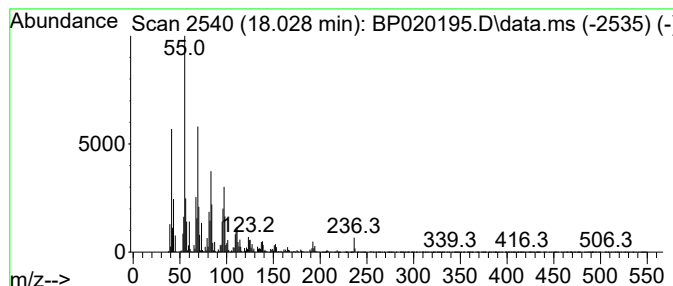
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Palmitoleic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.028	10.47 ng/ul	714739	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Palmitoleic acid	254	C16H30O2	000373-49-9	94
2			Hexadecenoic acid, Z-11-	254	C16H30O2	002416-20-8	91
3			9-Hexadecenoic acid	254	C16H30O2	002091-29-4	91
4			Palmitoleic acid	254	C16H30O2	000373-49-9	90
5			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	90



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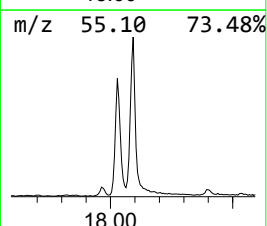
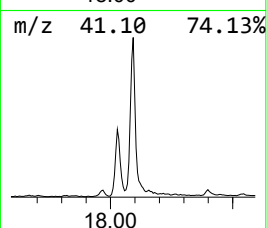
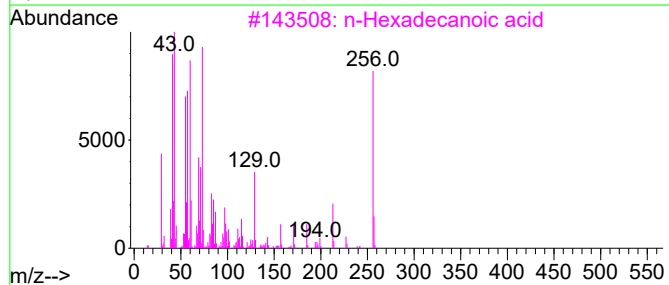
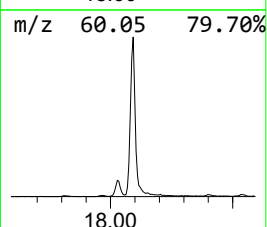
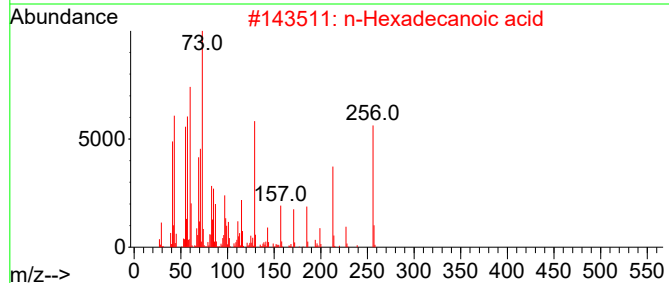
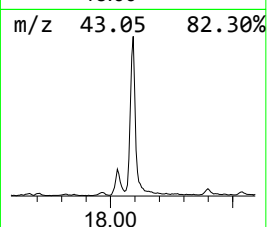
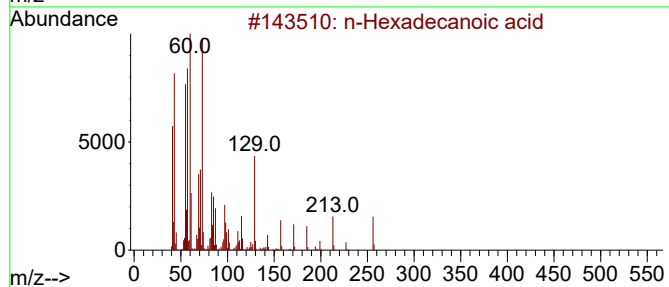
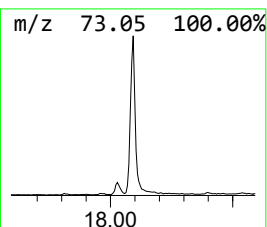
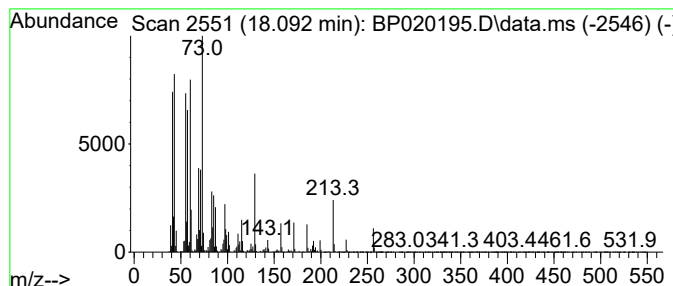
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	22.31 ng/ul	1523210	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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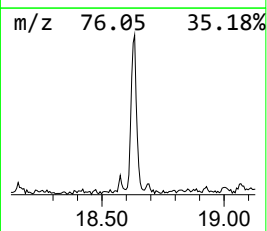
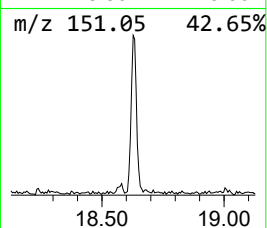
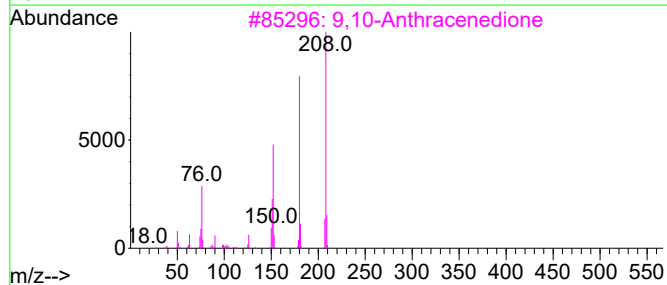
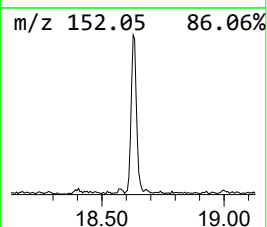
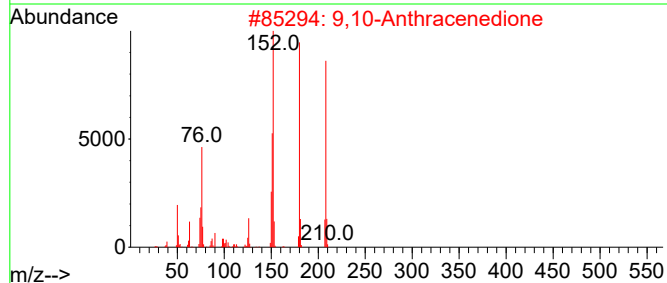
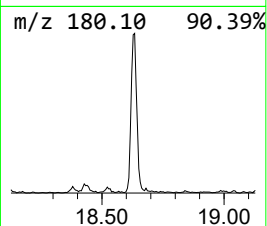
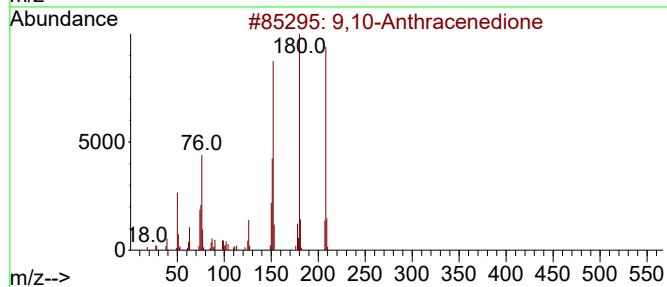
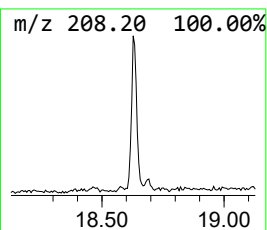
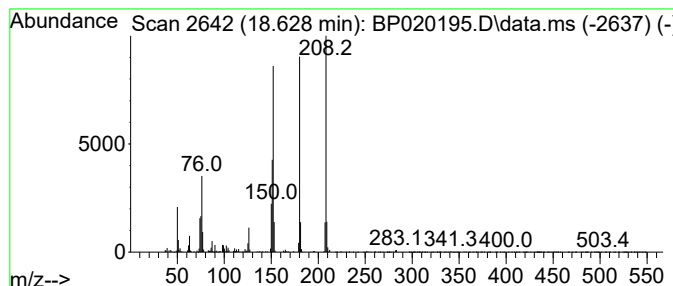
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 9,10-Anthracenedione Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.628	2.67 ng/ul	182051	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2			9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
3			9,10-Anthracenedione	208	C14H8O2	000084-65-1	95
4			9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
5			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
Operator : MA/JU
Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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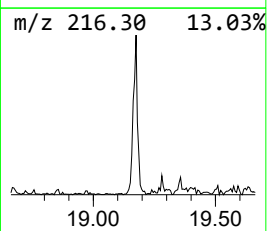
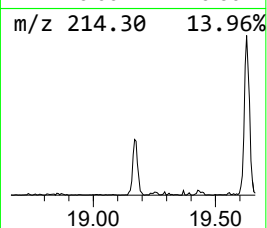
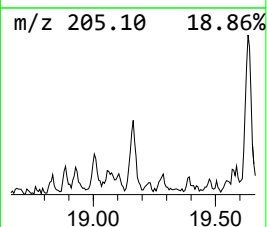
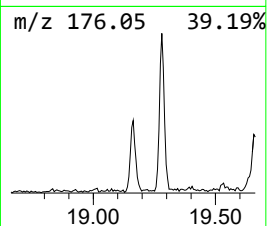
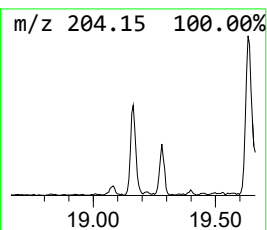
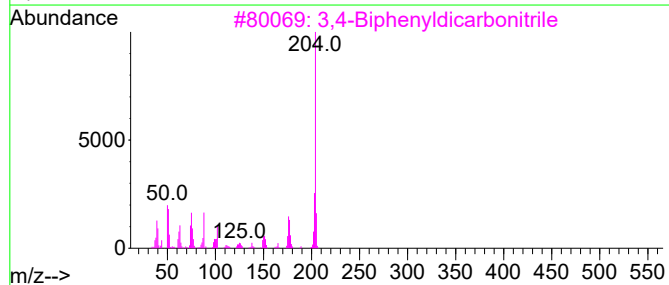
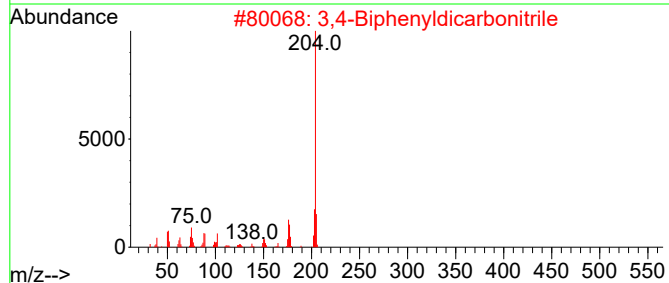
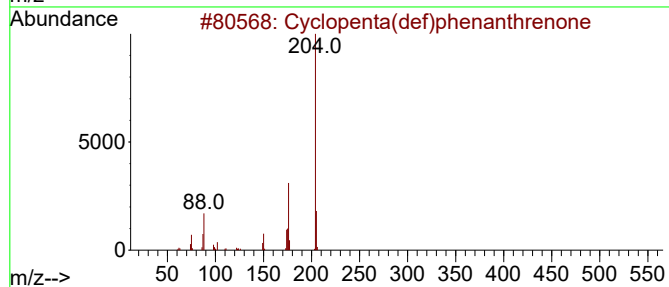
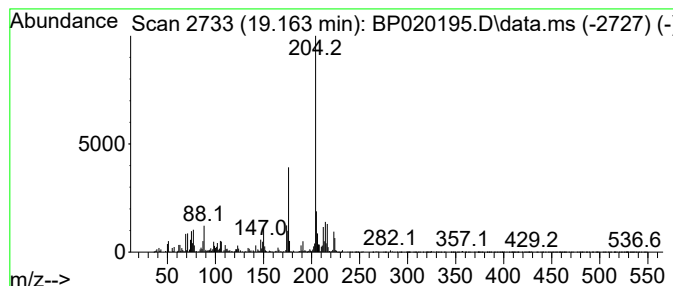
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclopenta(def)phenanthrenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.163	3.03 ng/ul	206905	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	86
2			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	47
3			3,4-Biphenyldicarbonitrile	204	C14H8N2	004128-63-6	43
4			[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	43
5			Indeno[2,1-a]indene, 5,10-dihydro-	204	C16H12	006543-29-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
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Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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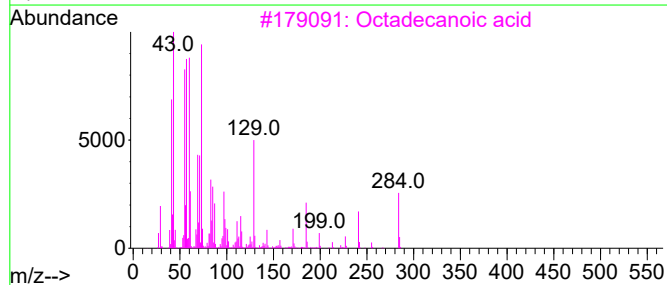
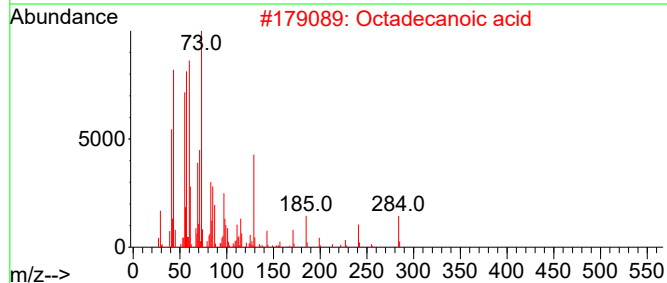
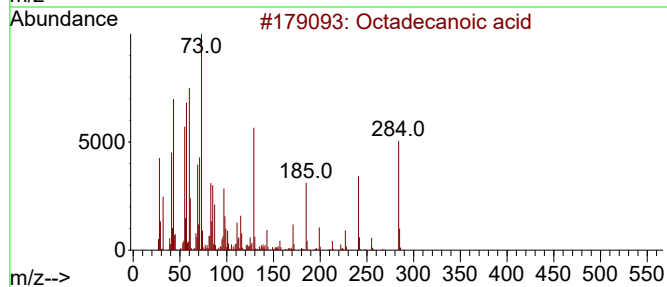
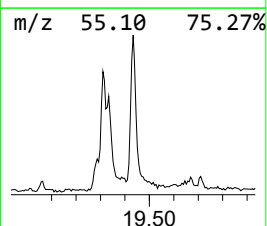
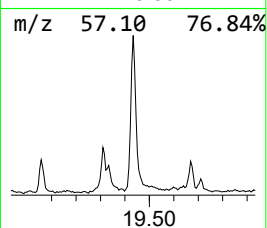
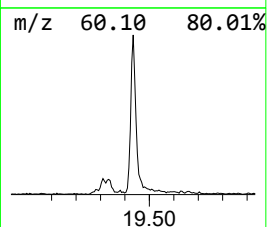
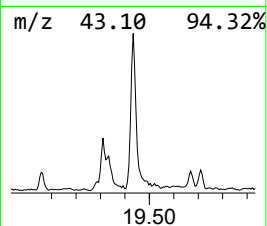
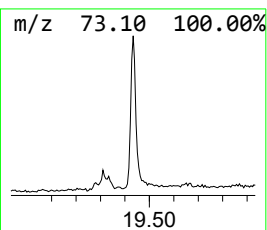
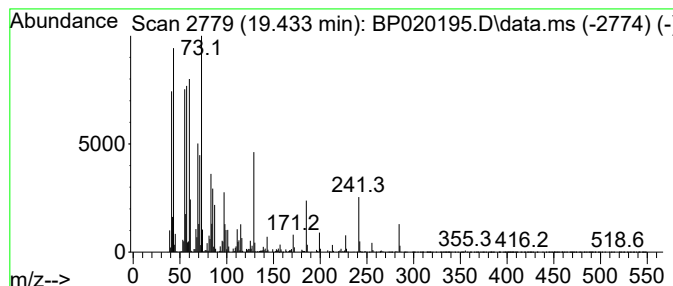
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.433	6.24 ng/ul	580079	Chrysene-d12	21.616

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	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
Operator : MA/JU
Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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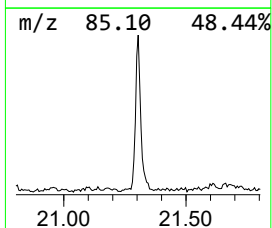
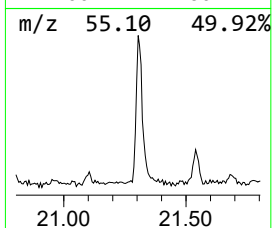
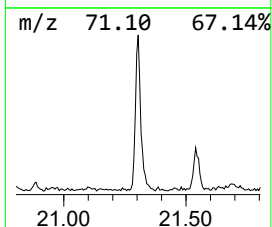
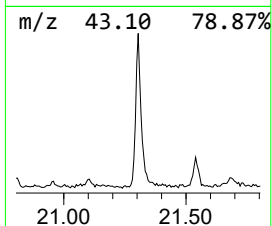
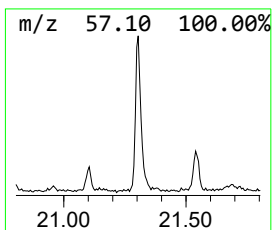
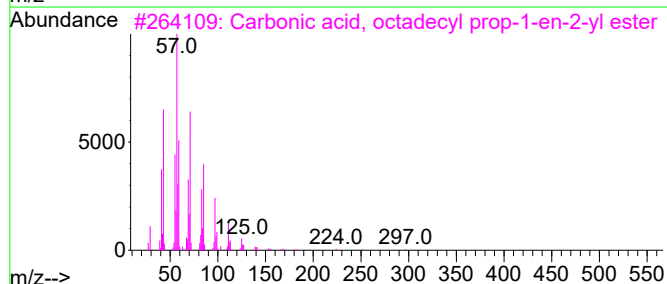
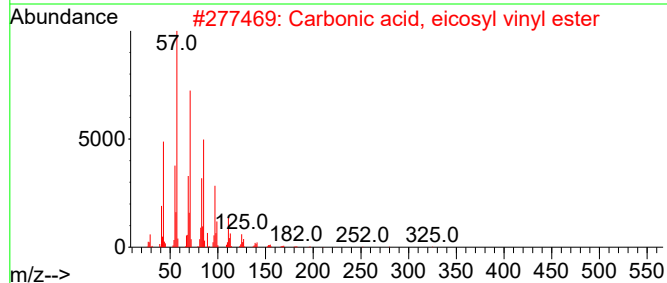
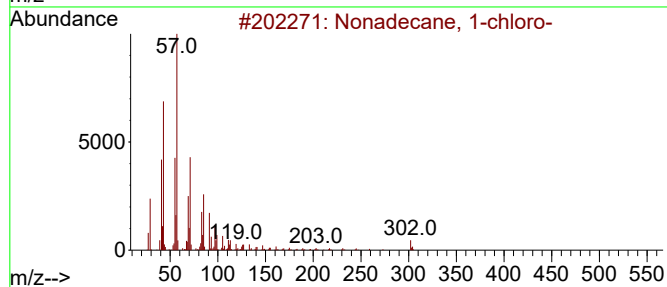
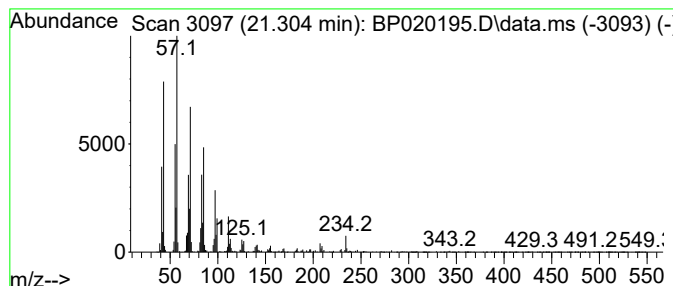
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Nonadecane, 1-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.304	3.68 ng/ul	342818	Chrysene-d12	21.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	93
2			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	91
3			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	91
4			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	91
5			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
Operator : MA/JU
Sample : P2234-01
Misc :
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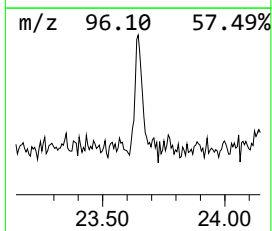
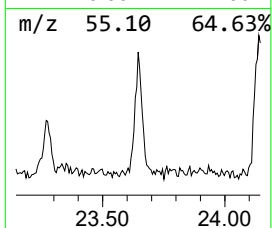
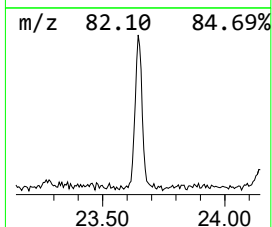
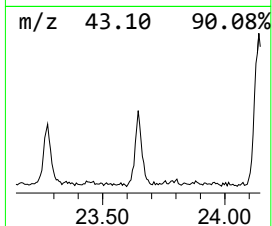
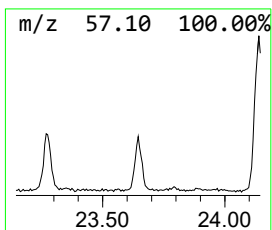
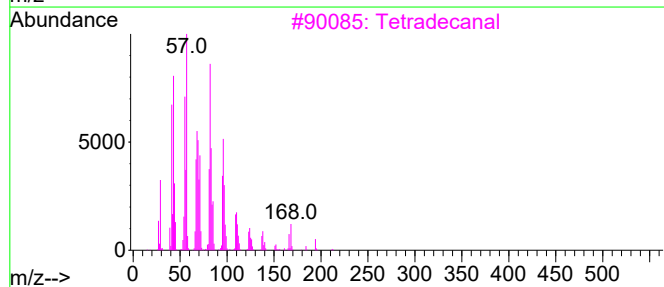
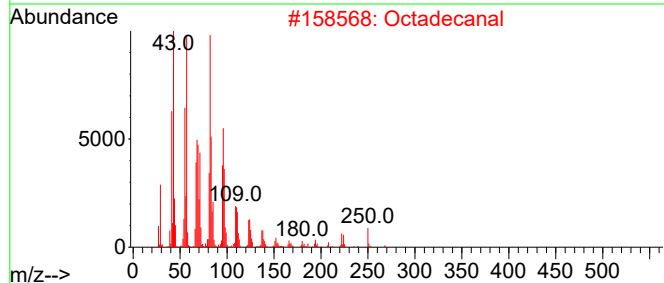
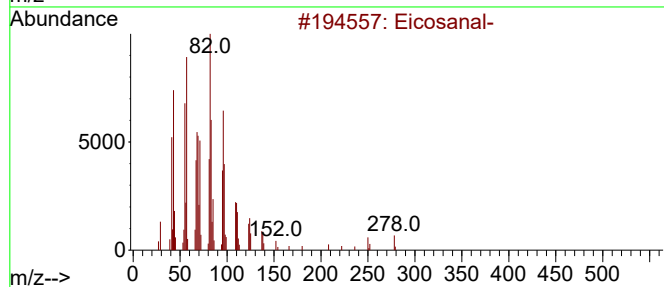
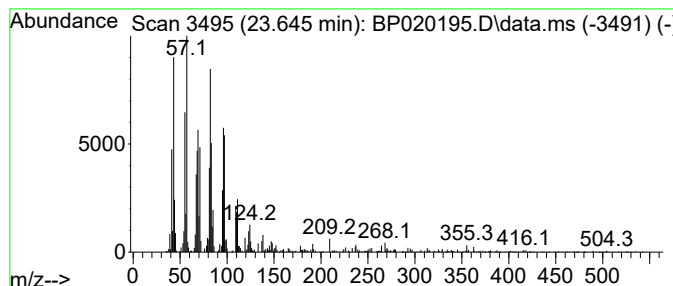
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Eicosanal- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.645	3.97 ng/ul	199788	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosanal-	296	C20H40O	002400-66-0	93
2			Octadecanal	268	C18H36O	000638-66-4	90
3			Tetradecanal	212	C14H28O	000124-25-4	90
4			1,19-Eicosadiene	278	C20H38	014811-95-1	89
5			Henicosanal	310	C21H42O	051227-32-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
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Sample : P2234-01
Misc :
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Instrument :
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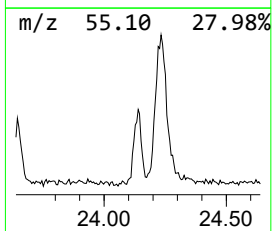
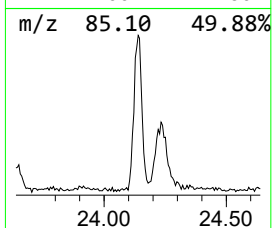
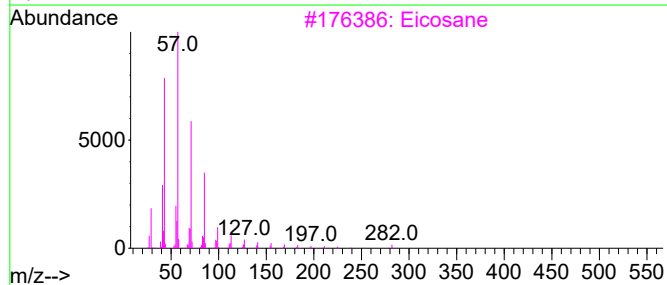
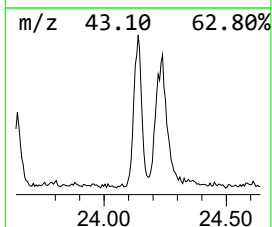
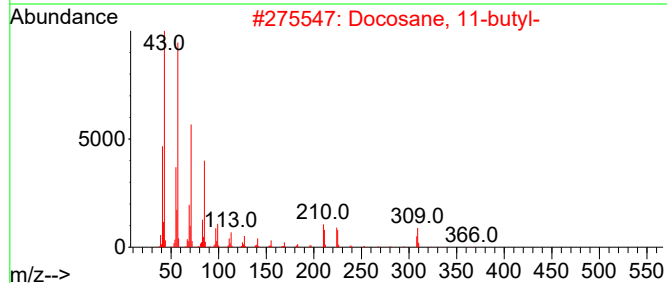
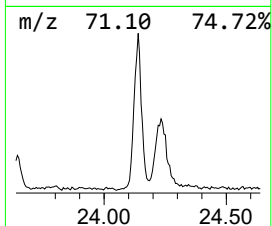
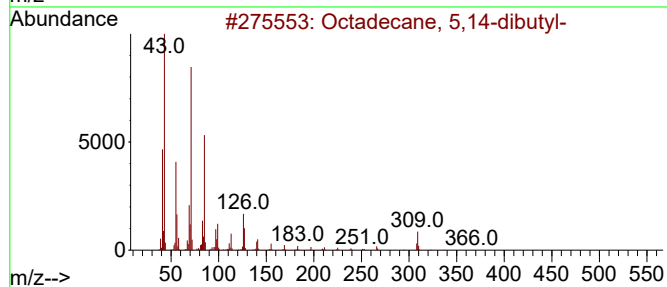
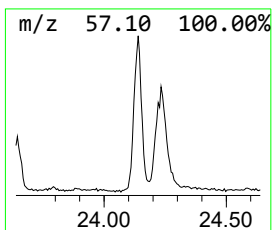
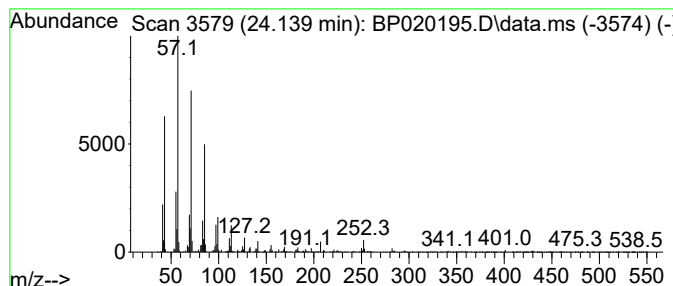
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.139	5.74 ng/ul	288839	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	93
2			Docosane, 11-butyl-	366	C26H54	013475-76-8	93
3			Eicosane	282	C20H42	000112-95-8	93
4			Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5			Pentacosane	352	C25H52	000629-99-2	93



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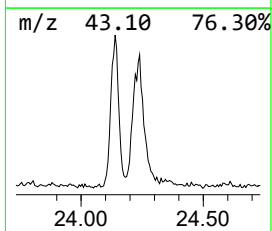
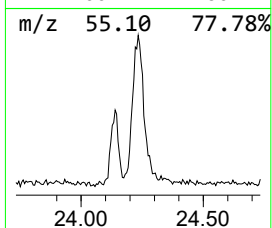
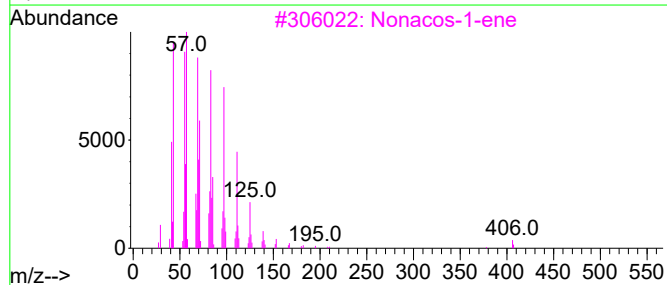
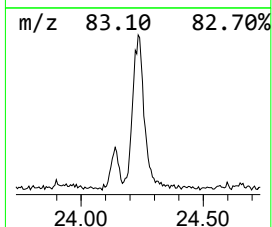
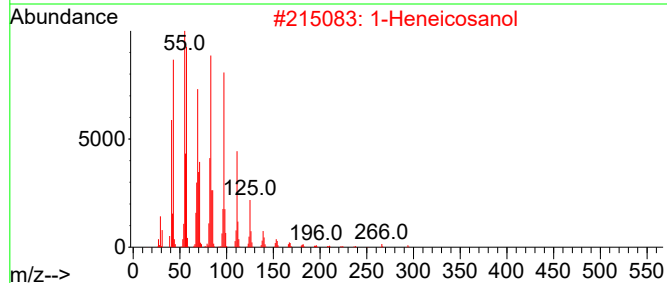
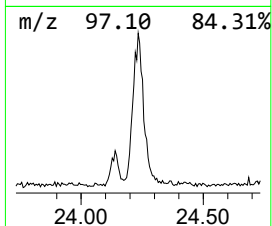
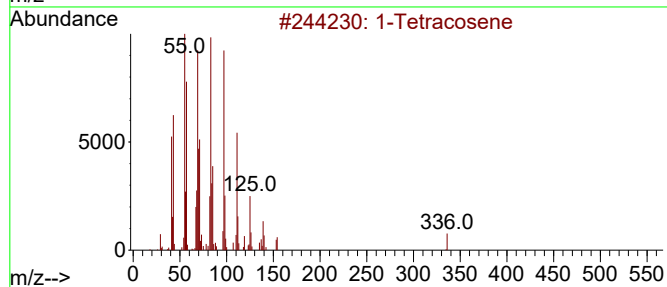
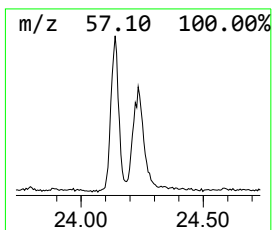
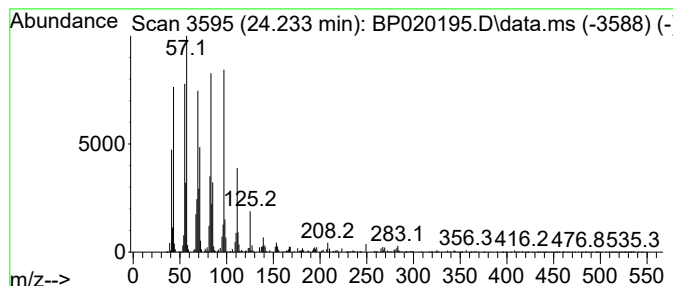
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Quant Title : SVOA CALIBRATION

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.	
24.233	13.54 ng/ul	681986	Perylene-d12		24.968	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Tetracosene		336	C24H48	010192-32-2	98
2	1-Heneicosanol		312	C21H44O	015594-90-8	93
3	Nonacos-1-ene		406	C29H58	018835-35-3	91
4	1-Triacontanol		438	C30H62O	000593-50-0	89
5	n-Nonadecanol-1		284	C19H40O	001454-84-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
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Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

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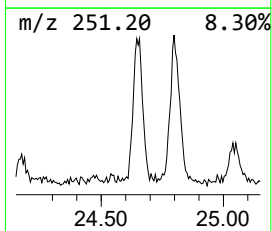
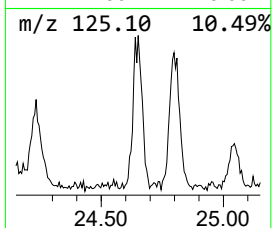
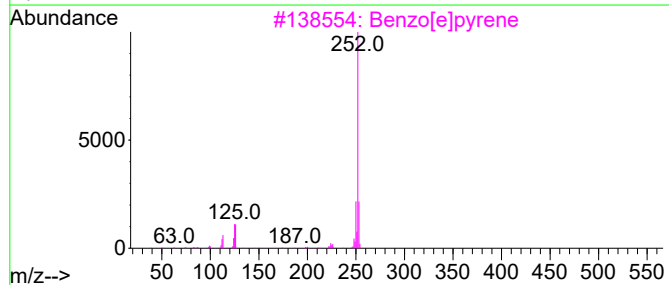
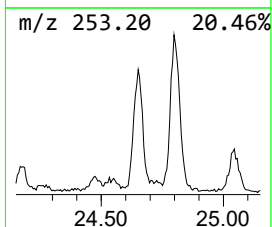
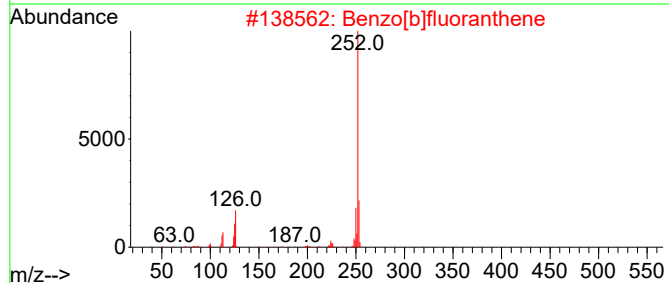
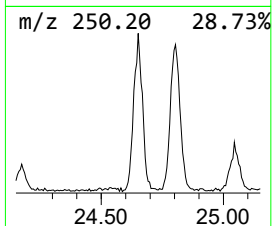
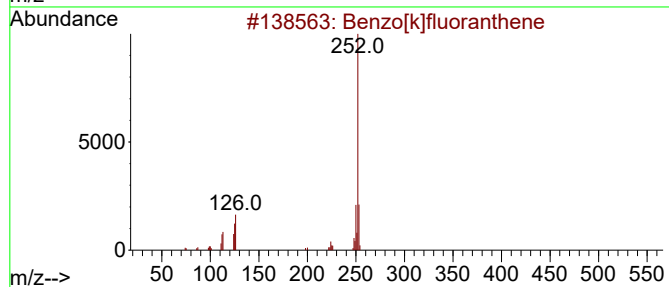
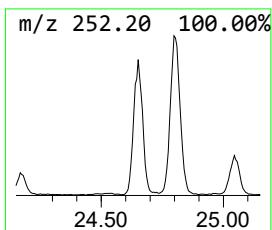
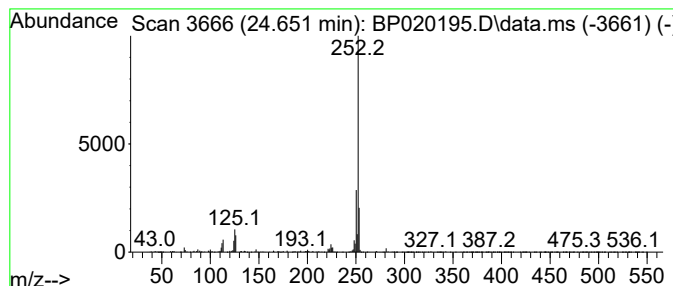
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.651	4.68 ng/ul	235881	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
2			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
3			Benzo[e]pyrene	252	C20H12	000192-97-2	90
4			Benzo[a]pyrene	252	C20H12	000050-32-8	90
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
Operator : MA/JU
Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0SA1

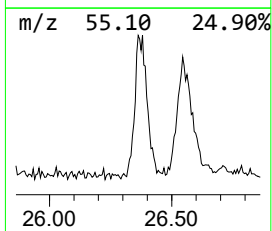
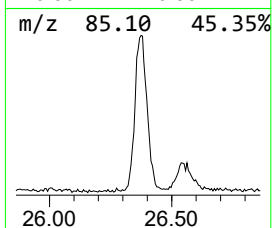
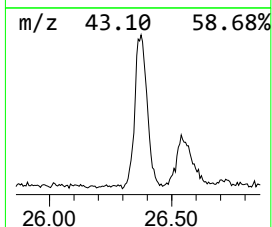
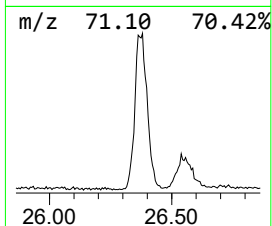
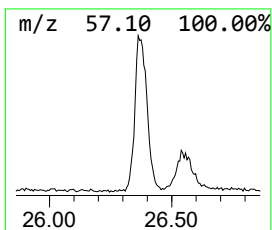
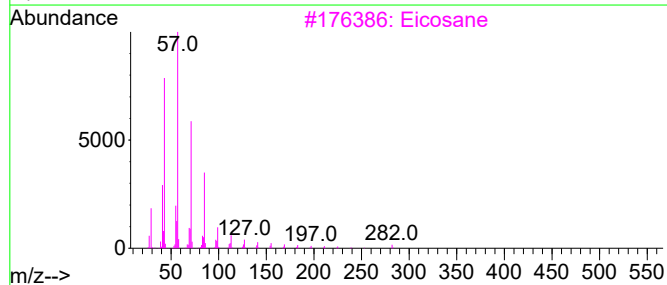
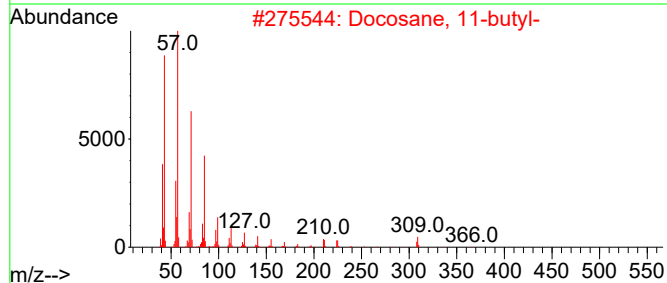
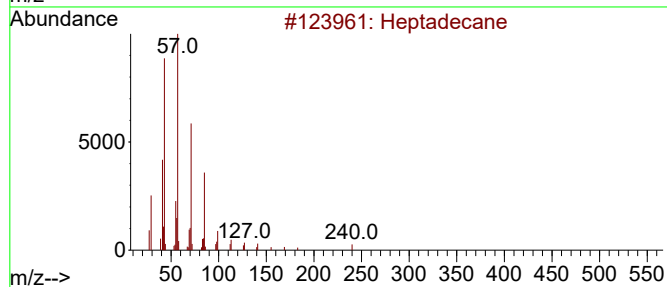
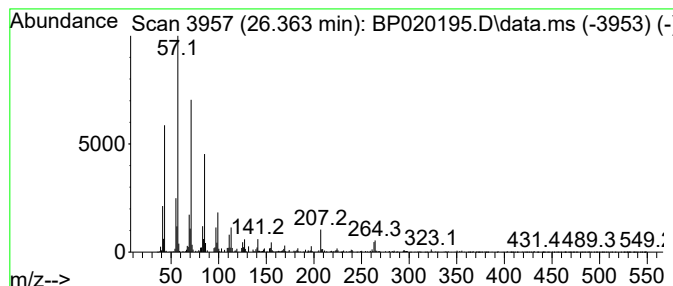
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.363	3.13 ng/ul	157484	Perylene-d12	24.968

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	93
2		Docosane, 11-butyl-	366	C26H54	013475-76-8	91
3		Eicosane	282	C20H42	000112-95-8	90
4		Heptacosane	380	C27H56	000593-49-7	87
5		Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050224\
Data File : BP020195.D
Acq On : 04 May 2024 05:33
Operator : MA/JU
Sample : P2234-01
Misc :
ALS Vial : 64 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0SA1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.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
Palmitoleic acid	18.028	10.5	ng/ul	714739	4	17.128	1365420	20.0
n-Hexadecanoic ...	18.092	22.3	ng/ul	1523210	4	17.128	1365420	20.0
9,10-Anthracene...	18.628	2.7	ng/ul	182051	4	17.128	1365420	20.0
Cyclopenta(def)...	19.163	3.0	ng/ul	206905	4	17.128	1365420	20.0
Octadecanoic acid	19.433	6.2	ng/ul	580079	5	21.616	1860710	20.0
Nonadecane, 1-C...	21.304	3.7	ng/ul	342818	5	21.616	1860710	20.0
Eicosanal-	23.645	4.0	ng/ul	199788	6	24.968	1007120	20.0
(DEL) Alkane: S...	24.139	5.7	ng/ul	288839	6	24.968	1007120	20.0
(DEL) Alkane: S...	24.233	13.5	ng/ul	681986	6	24.968	1007120	20.0
Benzo[e]pyrene	24.651	4.7	ng/ul	235881	6	24.968	1007120	20.0
(DEL) Alkane: S...	26.363	3.1	ng/ul	157484	6	24.968	1007120	20.0