

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
 Data File : BP015200.D
 Acq On : 20 May 2023 07:00
 Operator : CG/JU
 Sample : 02689-17
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JREN9

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: BP015200.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	38	42	50	rVB	67168	89482	1.20%	0.108%
2	3.869	113	116	132	rBV	729311	1063684	14.21%	1.282%
3	3.987	132	136	140	rVB	29116	39581	0.53%	0.048%
4	4.111	153	157	164	rVB	41637	60708	0.81%	0.073%
5	4.211	170	174	178	rBV	17099	24126	0.32%	0.029%
6	4.599	235	240	244	rBV5	17046	26214	0.35%	0.032%
7	4.669	247	252	261	rVB	74548	108222	1.45%	0.130%
8	4.787	269	272	277	rVB6	8974	13664	0.18%	0.016%
9	4.969	297	303	307	rBV2	8096	14508	0.19%	0.017%
10	5.163	332	336	342	rVB	29663	41004	0.55%	0.049%
11	5.281	349	356	362	rBV10	11758	22347	0.30%	0.027%
12	5.422	376	380	388	rVB4	8797	15992	0.21%	0.019%
13	5.658	416	420	425	rVB	17386	23845	0.32%	0.029%
14	5.781	436	441	448	rVB	34531	51440	0.69%	0.062%
15	6.699	593	597	602	rBV8	5175	9017	0.12%	0.011%
16	6.781	607	611	617	rVB6	10297	17914	0.24%	0.022%
17	7.228	682	687	689	rBV	55394	76931	1.03%	0.093%
18	7.275	689	695	708	rVB	1359338	2209172	29.51%	2.663%
19	7.446	714	724	733	rBV	1099117	1723509	23.02%	2.077%
20	7.528	735	738	742	rVB3	10970	14039	0.19%	0.017%
21	7.646	749	758	768	rBV	1353104	2203794	29.44%	2.656%
22	7.846	788	792	799	rVB3	8059	15569	0.21%	0.019%
23	7.940	802	808	815	rVV	9699	18990	0.25%	0.023%
24	8.122	831	839	847	rBV	888334	1400404	18.71%	1.688%
25	8.834	952	960	977	rBV	1725627	2857922	38.18%	3.445%
26	8.963	977	982	989	rVB	14785	26745	0.36%	0.032%
27	9.299	1031	1039	1049	rBV	1306853	2207184	29.48%	2.660%
28	9.451	1061	1065	1071	rVB9	8352	13664	0.18%	0.016%
29	9.693	1099	1106	1115	rVB	170058	267771	3.58%	0.323%
30	10.028	1152	1163	1176	rBV	1188839	1974157	26.37%	2.379%
31	10.257	1195	1202	1203	rBV7	8377	12636	0.17%	0.015%
32	10.569	1247	1255	1269	rVV	1731476	2964701	39.60%	3.573%
33	10.951	1312	1320	1331	rBV	1233045	2038137	27.23%	2.457%
34	11.087	1336	1343	1358	rBV	1515026	2628768	35.12%	3.168%
35	11.481	1406	1410	1419	rVB4	12882	22971	0.31%	0.028%
36	11.728	1448	1452	1459	rBV7	7985	18501	0.25%	0.022%

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Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
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37	12.187	1525	1530	1537	rVB	127581	180229	2.41%	0.217%
38	12.428	1566	1571	1576	rBV3	12769	23143	0.31%	0.028%
39	12.528	1583	1588	1596	rVB	25213	44385	0.59%	0.053%
40	13.575	1759	1766	1771	rBV	5415	11922	0.16%	0.014%

41	13.645	1774	1778	1782	rBV2	11611	16031	0.21%	0.019%
42	14.034	1838	1844	1848	rBV8	7707	14273	0.19%	0.017%
43	14.081	1848	1852	1858	rVB5	12321	18614	0.25%	0.022%
44	14.163	1858	1866	1881	rBV	3151369	4562932	60.95%	5.500%
45	14.281	1881	1886	1891	rVB8	9832	15202	0.20%	0.018%

46	14.334	1891	1895	1901	rBV	27095	39604	0.53%	0.048%
47	14.457	1909	1916	1927	rBV	3581481	5192436	69.36%	6.258%
48	14.604	1936	1941	1953	rVB7	21951	45052	0.60%	0.054%
49	14.763	1961	1968	1975	rBV	1727723	2591521	34.62%	3.124%
50	14.951	1993	2000	2014	rBV	1018425	1619254	21.63%	1.952%

51	15.063	2017	2019	2024	rVV	22527	34131	0.46%	0.041%
52	15.110	2024	2027	2031	rVV5	17740	28919	0.39%	0.035%
53	15.169	2031	2037	2044	rVB3	70207	128354	1.71%	0.155%
54	15.704	2121	2128	2131	rBV	101269	142043	1.90%	0.171%
55	15.751	2131	2136	2143	rVV	4332182	6153501	82.20%	7.417%

56	15.798	2143	2144	2150	rVB6	12035	12535	0.17%	0.015%
57	15.869	2150	2156	2166	rBV	1075997	1435876	19.18%	1.731%
58	15.992	2173	2177	2183	rVB3	23188	35143	0.47%	0.042%
59	16.116	2191	2198	2211	rBV	1432755	1989375	26.58%	2.398%
60	16.445	2251	2254	2258	rBV5	9617	15767	0.21%	0.019%

61	16.681	2288	2294	2301	rBV	236797	325929	4.35%	0.393%
62	16.892	2325	2330	2334	rBV6	14562	22196	0.30%	0.027%
63	16.957	2337	2341	2347	rVV	85395	110780	1.48%	0.134%
64	17.186	2376	2380	2384	rBV6	11073	13717	0.18%	0.017%
65	17.392	2410	2415	2417	rBV6	10242	16657	0.22%	0.020%

66	17.451	2422	2425	2429	rVV5	10048	15135	0.20%	0.018%
67	17.510	2429	2435	2442	rVV	1943530	2783974	37.19%	3.356%
68	17.610	2446	2452	2462	rVV	4170209	6228005	83.20%	7.507%
69	17.698	2462	2467	2474	rVB2	109927	163363	2.18%	0.197%
70	17.851	2490	2493	2496	rBV5	9263	14704	0.20%	0.018%

71	17.980	2512	2515	2518	rBV5	9428	12250	0.16%	0.015%
72	18.028	2518	2523	2527	rVV	66138	82302	1.10%	0.099%
73	18.116	2535	2538	2540	rBV4	8454	9055	0.12%	0.011%
74	18.157	2542	2545	2549	rVB5	9402	14331	0.19%	0.017%
75	18.257	2556	2562	2567	rBV3	60771	120964	1.62%	0.146%

76	18.369	2576	2581	2591	rBV	748040	1017667	13.59%	1.227%
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77	18.645	2626	2628	2633	rBV6	11000	19921	0.27%	0.024%
78	18.927	2673	2676	2681	rVB	62464	77576	1.04%	0.094%
79	18.975	2681	2684	2688	rBV3	48042	58303	0.78%	0.070%
80	19.069	2697	2700	2707	rVB4	35748	54889	0.73%	0.066%
81	19.175	2715	2718	2722	rBV3	46700	60937	0.81%	0.073%
82	19.251	2728	2731	2736	rBV	40997	50693	0.68%	0.061%
83	19.304	2737	2740	2746	rVB5	49566	66311	0.89%	0.080%
84	19.410	2754	2758	2762	rBV2	67015	86833	1.16%	0.105%
85	19.474	2763	2769	2772	rBV2	73723	129033	1.72%	0.156%
86	19.551	2778	2782	2785	rBV	109595	126215	1.69%	0.152%
87	19.592	2785	2789	2800	rVV	354921	524298	7.00%	0.632%
88	19.692	2802	2806	2810	rVB	101696	134279	1.79%	0.162%
89	19.833	2824	2830	2842	rVB	5141505	7263922	97.04%	8.755%
90	19.986	2852	2856	2863	rVB	101350	124656	1.67%	0.150%
91	20.363	2916	2920	2925	rBV	89273	120732	1.61%	0.146%
92	20.669	2968	2972	2977	rBV	308365	389542	5.20%	0.470%
93	20.804	2993	2995	2999	rVB3	59695	59837	0.80%	0.072%
94	20.974	3021	3024	3031	rBV3	87366	118994	1.59%	0.143%
95	21.263	3070	3073	3080	rVB	300166	441271	5.89%	0.532%
96	21.586	3123	3128	3134	rVB	1887170	2485610	33.20%	2.996%
97	22.198	3229	3232	3235	rBV	173557	179411	2.40%	0.216%
98	23.321	3419	3423	3429	rVB	284473	457458	6.11%	0.551%
99	23.980	3527	3535	3547	rVB2	3392348	7485859	100.00%	9.023%
100	24.145	3557	3563	3575	rVB	1345123	2831049	37.82%	3.412%

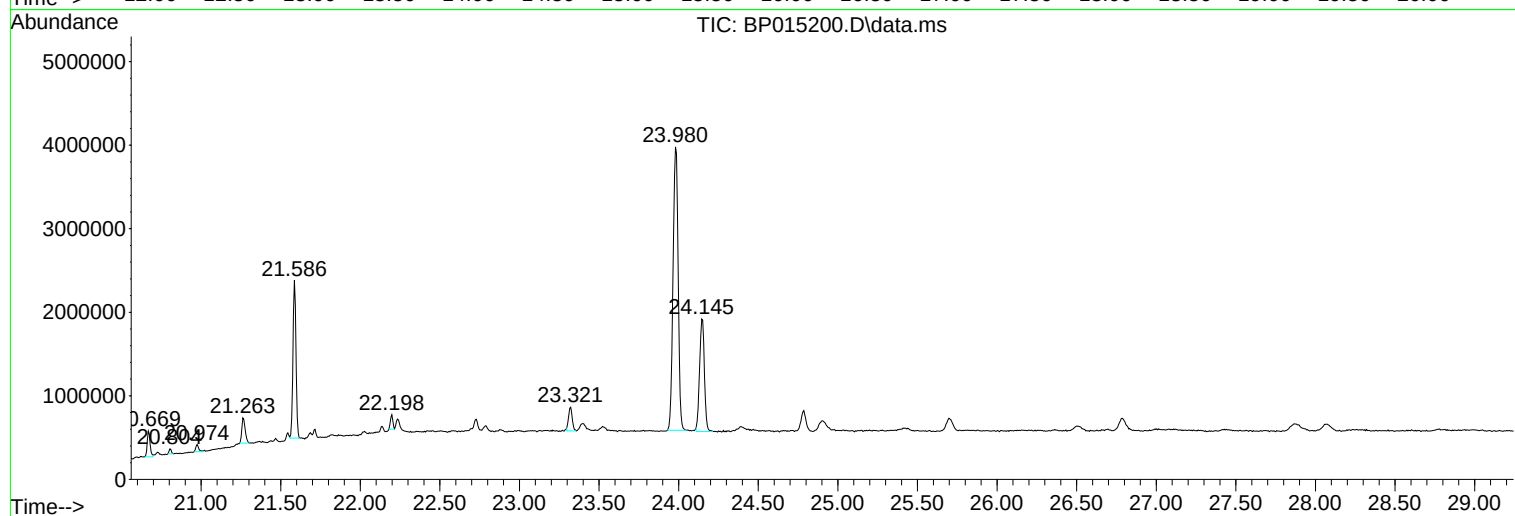
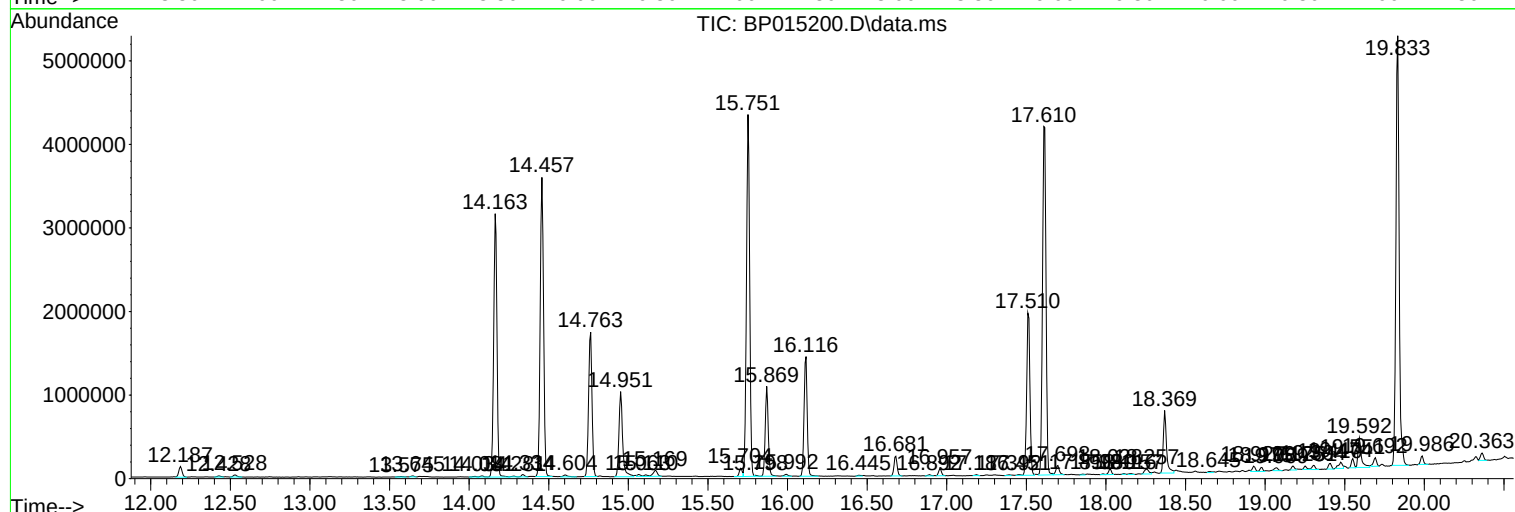
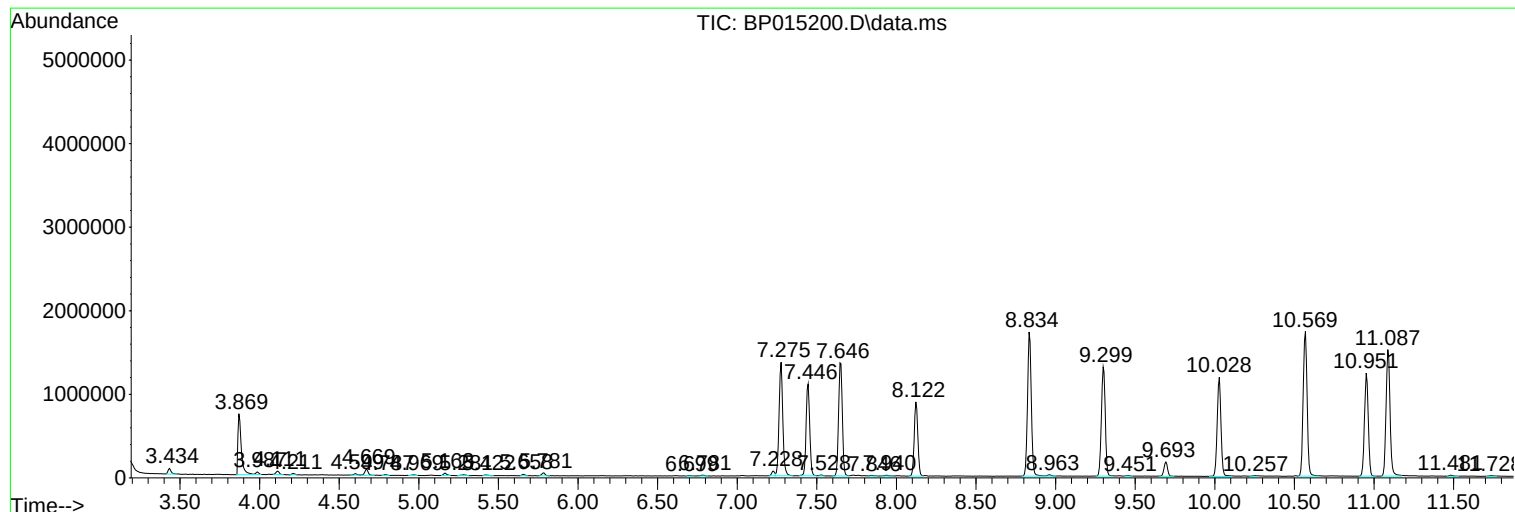
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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



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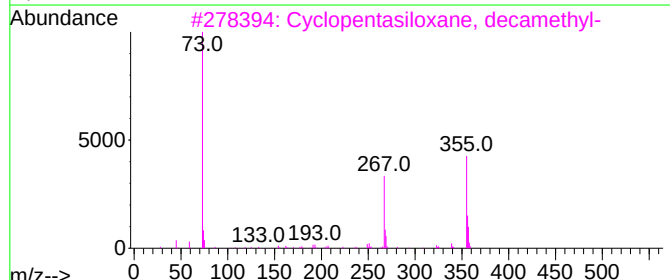
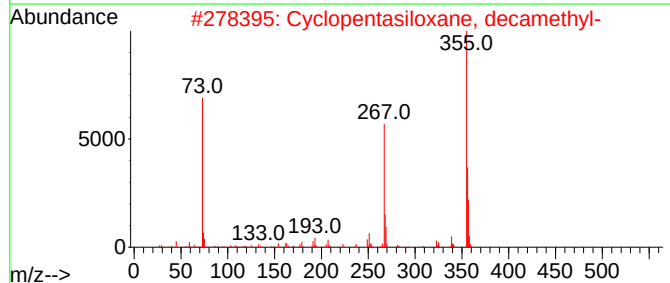
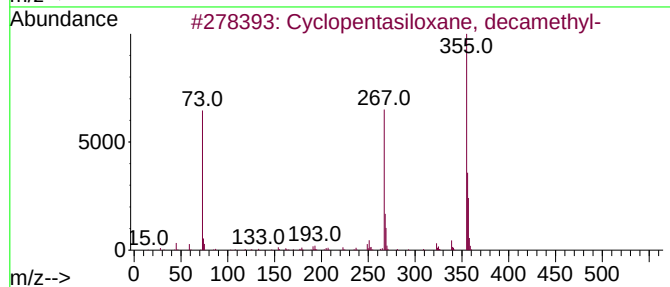
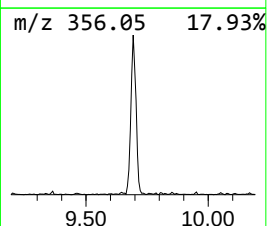
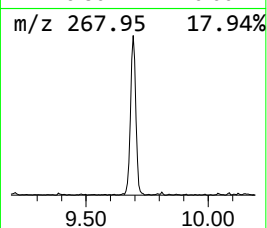
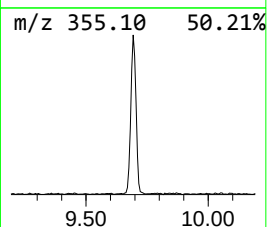
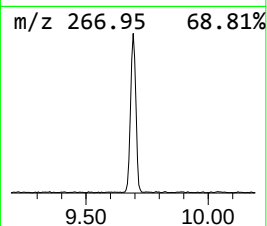
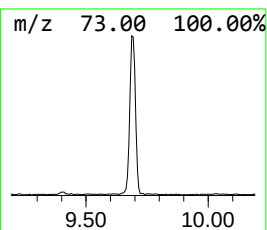
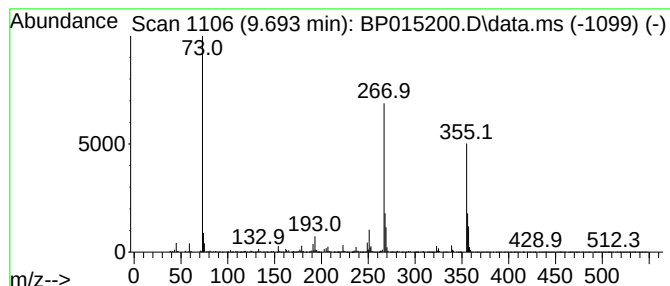
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopentasiloxane, decamet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.693	2.63 ng/ul	267771	Naphthalene-d8	10.951

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	74
4			3-Amino-2-phenazinol ditms	355	C18H25N3OSi2	1000332-15-5	52
5			Epinephrine, (.beta.)-, 3TMS der...	399	C18H37NO3Si3	010538-85-9	49



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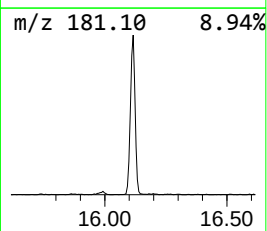
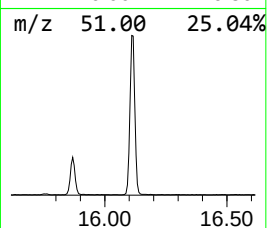
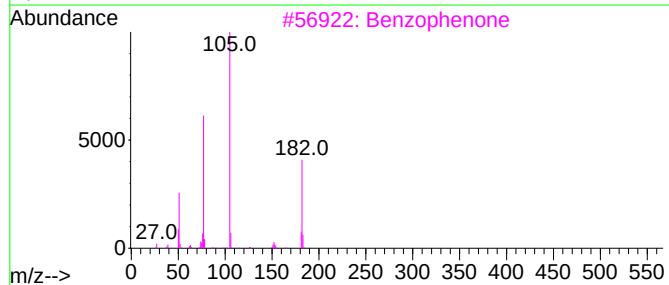
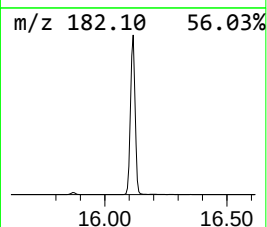
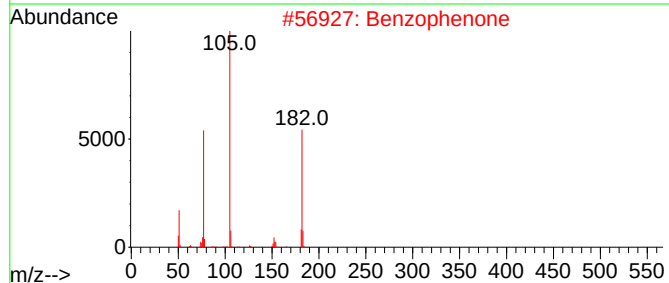
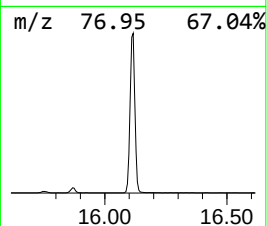
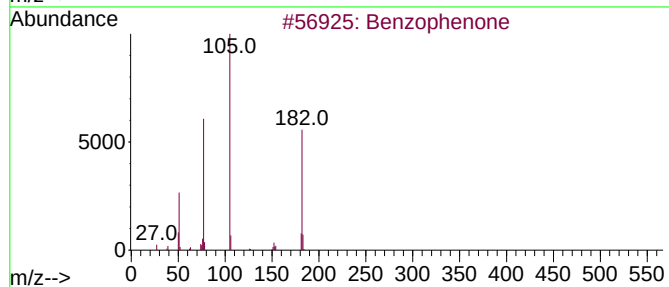
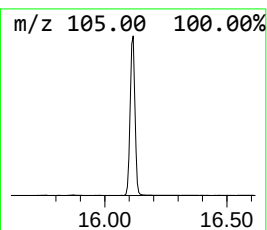
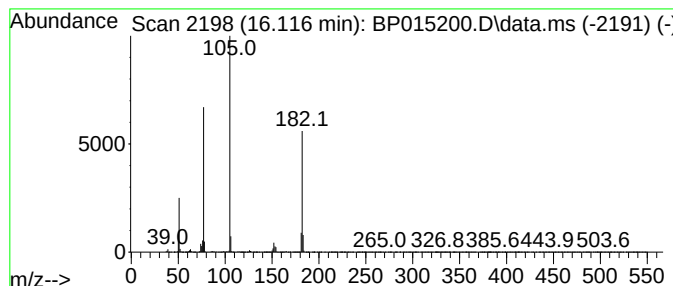
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.116	15.35 ng/ul	1989380	Acenaphthene-d10	14.763

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	90
5			Azobenzene	182	C12H10N2	000103-33-3	72



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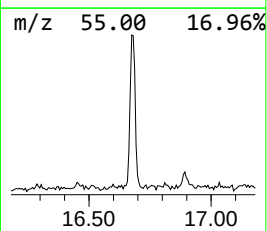
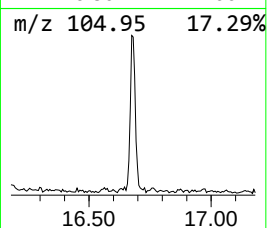
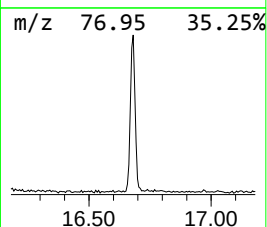
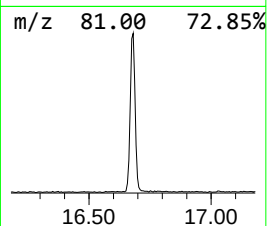
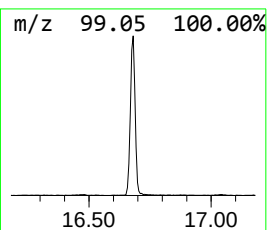
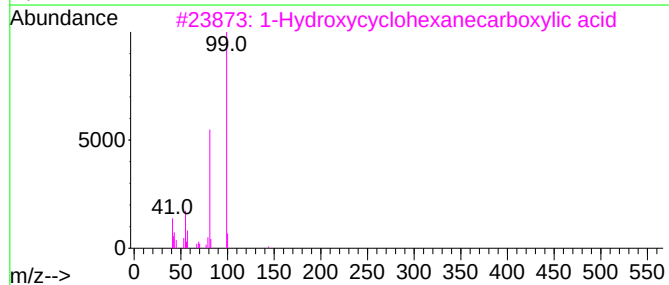
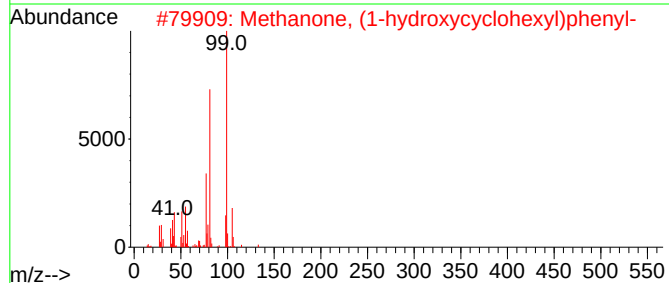
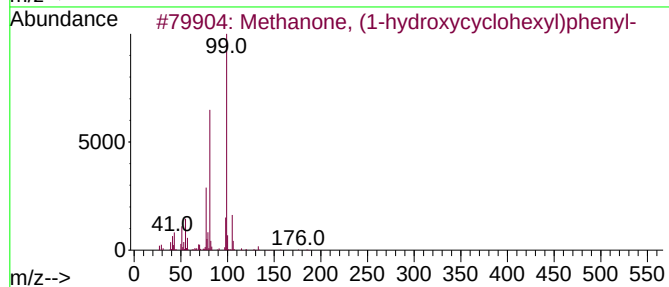
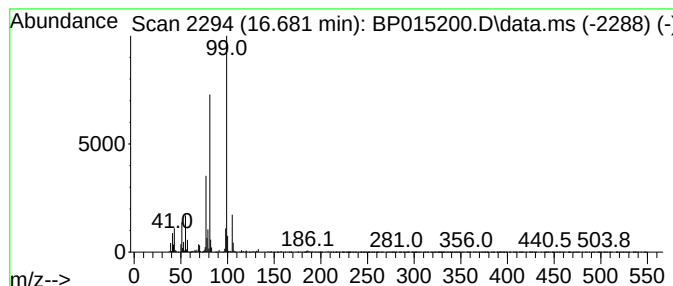
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Methanone, (1-hydroxycyclohexyl) Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.680	2.34 ng/ul	325929	Phenanthrene-d10	17.510

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	97
2			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	90
3			1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	50
4			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	50
5			Cyclohexanol, 1-(aminomethyl)-	129	C7H15NO	004000-72-0	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP015200.D
Acq On : 20 May 2023 07:00
Operator : CG/JU
Sample : 02689-17
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREN9

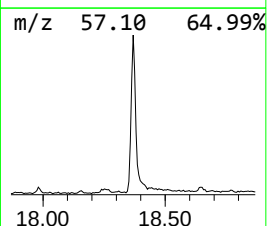
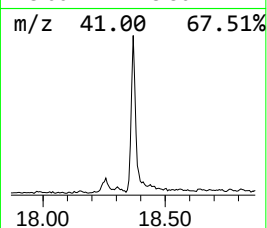
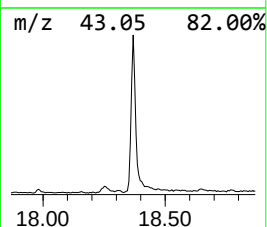
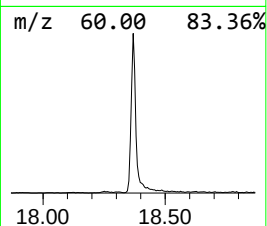
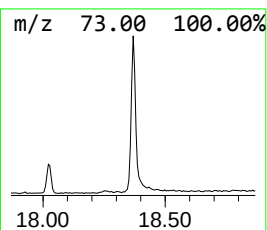
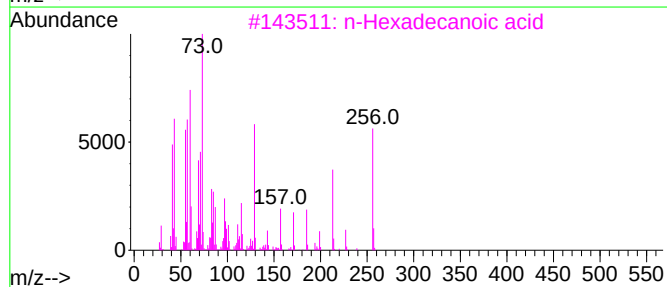
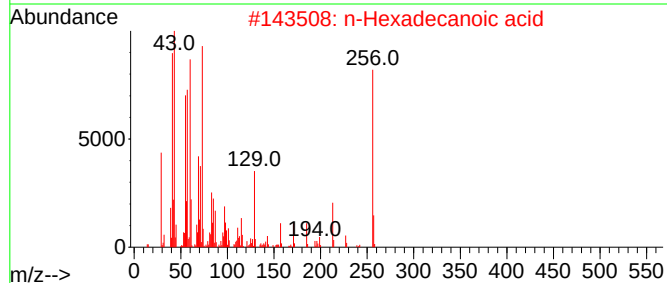
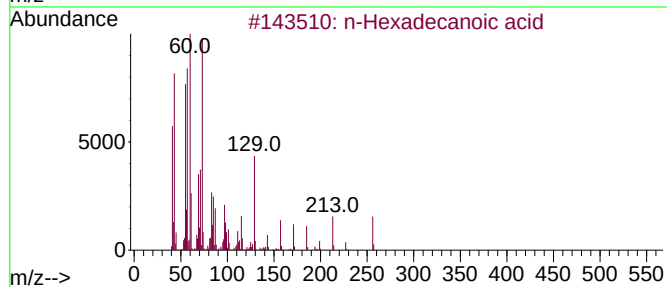
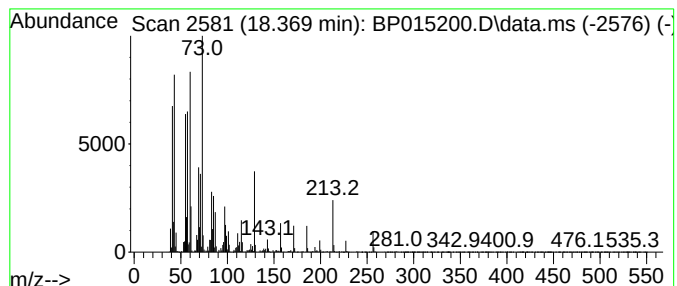
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.369	7.31 ng/ul	1017670	Phenanthrene-d10	17.510

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	98
5			Tridecanoic acid	214	C13H26O2	000638-53-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP015200.D
Acq On : 20 May 2023 07:00
Operator : CG/JU
Sample : 02689-17
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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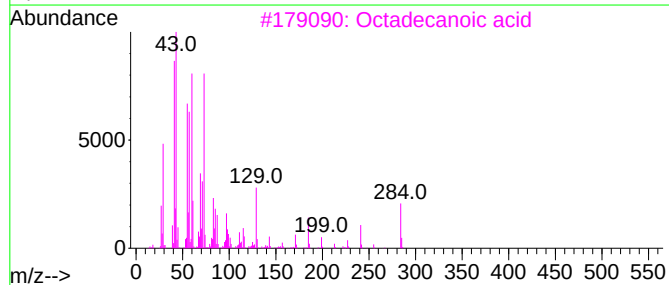
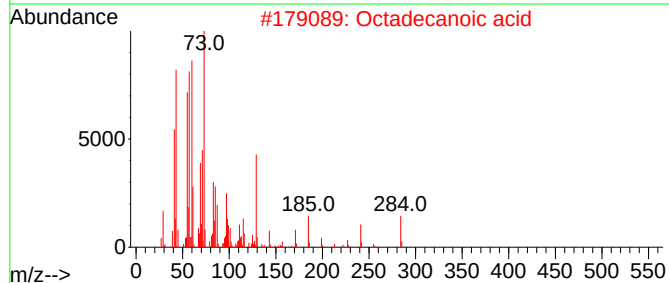
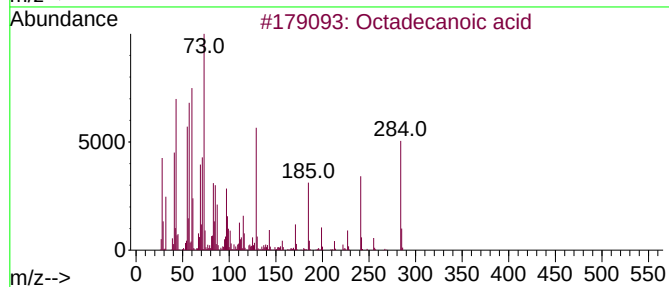
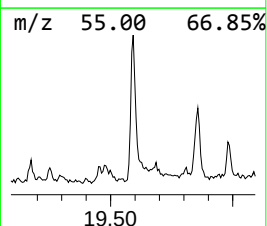
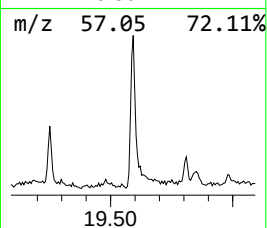
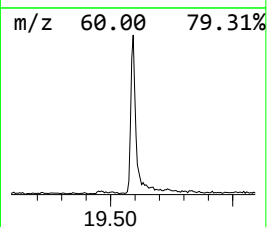
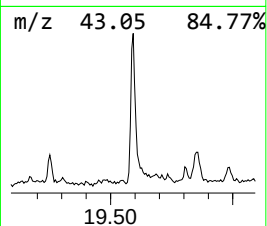
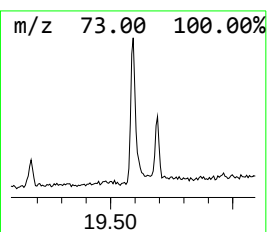
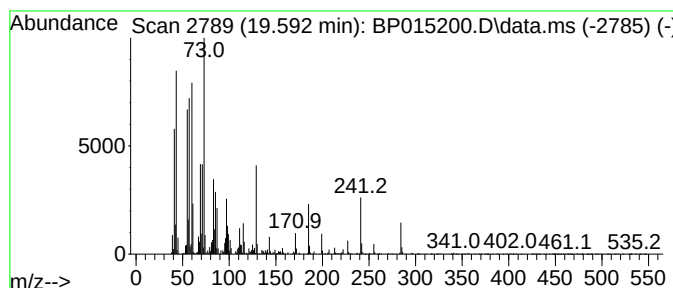
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.592	4.22 ng/ul	524298	Chrysene-d12	21.586

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			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP015200.D
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Operator : CG/JU
Sample : 02689-17
Misc :
ALS Vial : 37 Sample Multiplier: 1

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ClientSampleId :
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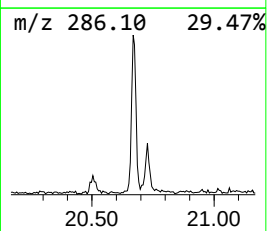
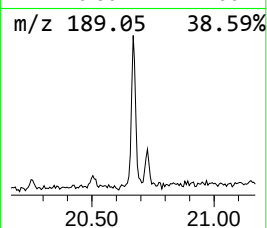
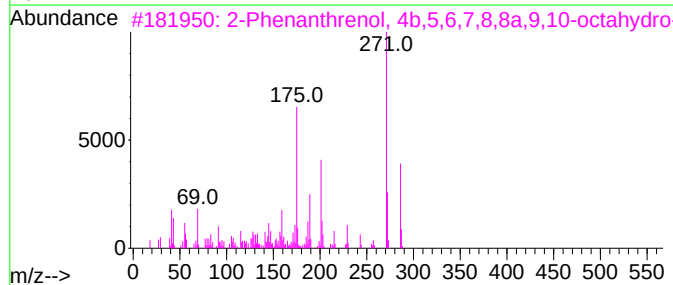
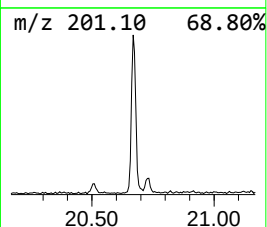
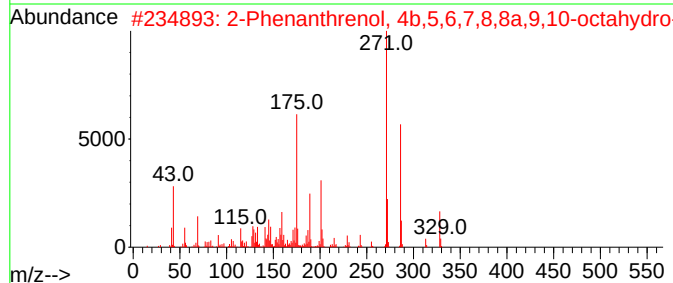
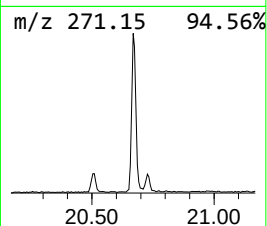
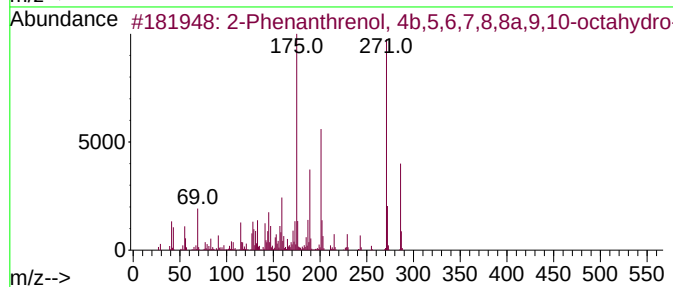
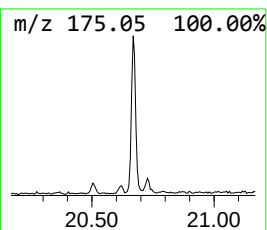
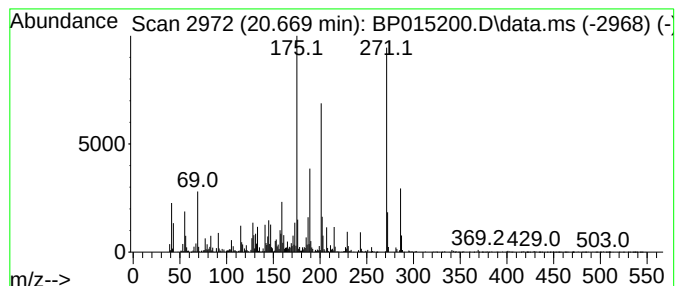
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.669	3.13 ng/ul	389542	Chrysene-d12	21.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H300	000511-15-9	99		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H3202	015340-82-6	91		
3	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H300	000511-15-9	91		
4	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H3202	015340-82-6	46		
5	(4Br,8As)-(+)-4B,5,6,7,8,8A,9,10...	346	C22H3403	1000426-72-2	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
Data File : BP015200.D
Acq On : 20 May 2023 07:00
Operator : CG/JU
Sample : 02689-17
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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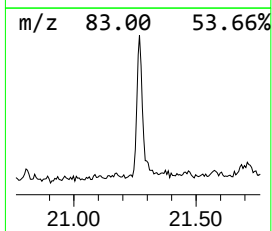
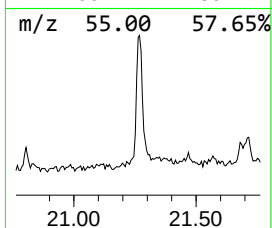
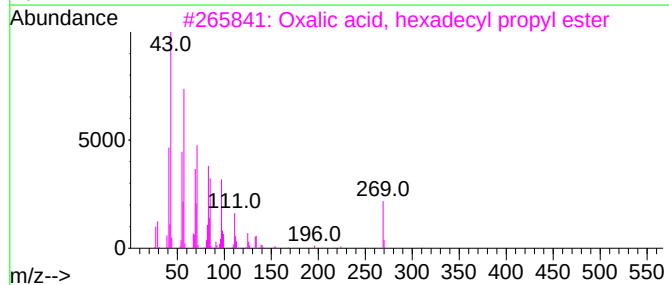
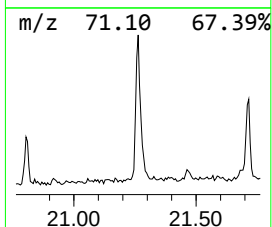
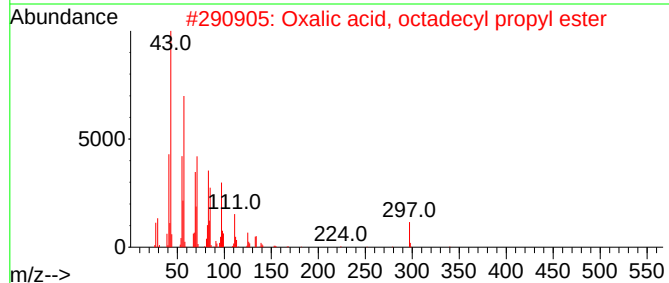
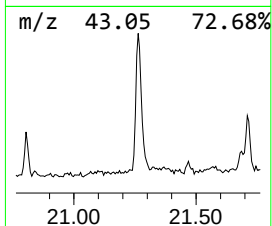
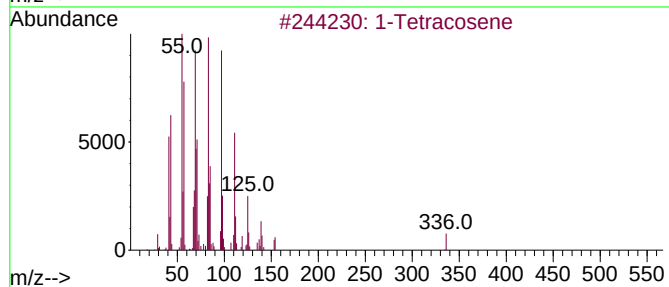
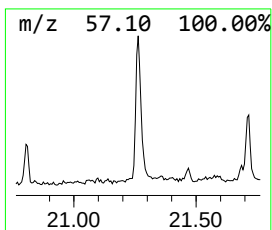
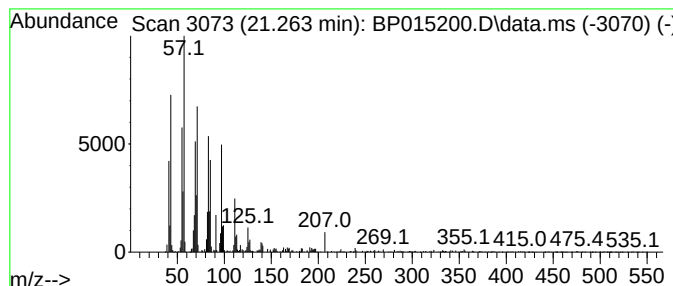
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.263	3.55 ng/ul	441271	Chrysene-d12	21.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene			336	C24H48	010192-32-2	96
2	Oxalic acid, octadecyl propyl ester			384	C23H44O4	1000309-27-0	90
3	Oxalic acid, hexadecyl propyl ester			356	C21H40O4	1000309-26-9	90
4	Oxalic acid, isobutyl hexadecyl ...			370	C22H42O4	1000309-38-1	90
5	Pentadecafluorooctanoic acid, oc...			666	C26H37F15O2	1000406-04-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
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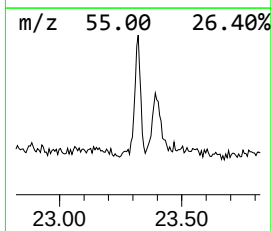
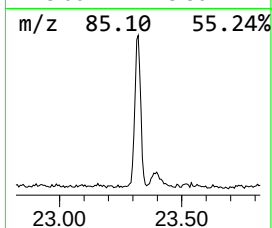
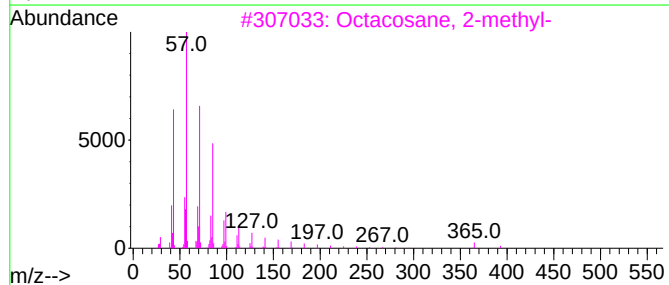
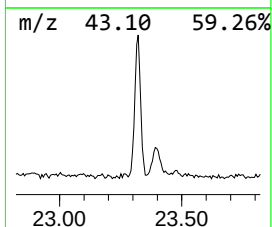
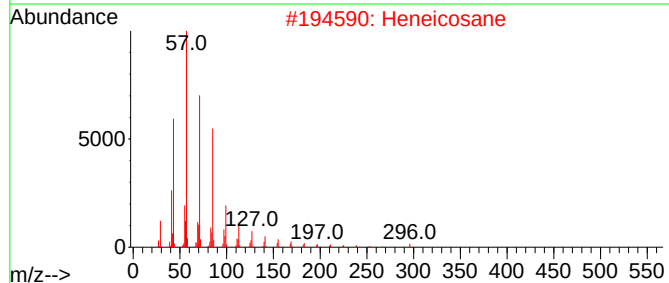
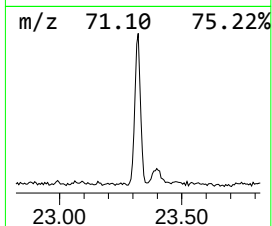
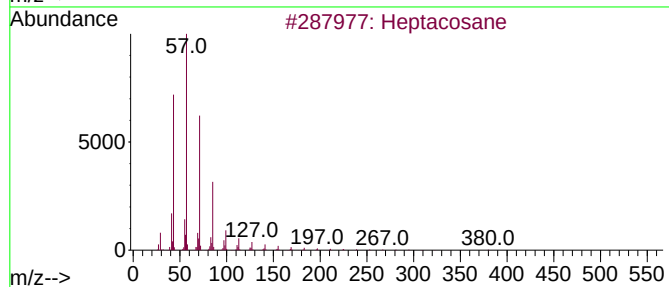
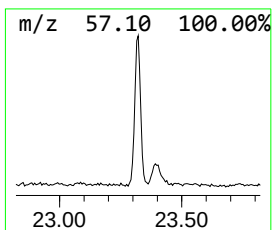
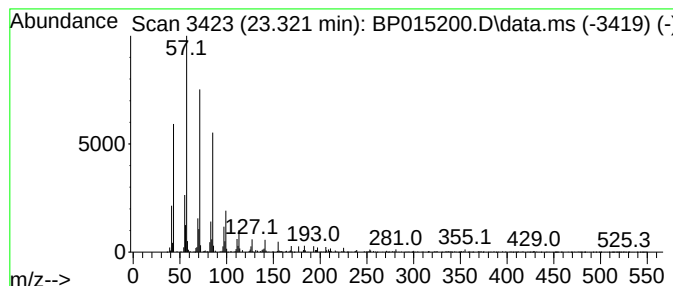
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Quant Title : SVOA CALIBRATION

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.	
23.321	3.23 ng/ul	457458	Perylene-d12		24.145	
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	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91
4	Hexacosane		366	C26H54	000630-01-3	90
5	Octacosane		394	C28H58	000630-02-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051923\
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Sample : 02689-17
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
JREN9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.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
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Benzophenone	16.116	15.4	ng/ul	1989380	3	14.763	2591520	20.0
Methanone, (1-h...	16.680	2.3	ng/ul	325929	4	17.510	2783970	20.0
n-Hexadecanoic ...	18.369	7.3	ng/ul	1017670	4	17.510	2783970	20.0
Octadecanoic acid	19.592	4.2	ng/ul	524298	5	21.586	2485610	20.0
2-Phenanthrenol...	20.669	3.1	ng/ul	389542	5	21.586	2485610	20.0
(DEL) Alkane: S...	21.263	3.5	ng/ul	441271	5	21.586	2485610	20.0
(DEL) Alkane: S...	23.321	3.2	ng/ul	457458	6	24.145	2831050	20.0