

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
 Data File : BG059152.D
 Acq On : 21 Sep 2023 23:30
 Operator : MA/JU
 Sample : 04386-09
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AC9

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_G\Methods\SFAM-EPA-BG091923.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG059152.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.605	34	37	45	rVB4	9184	15651	1.74%	0.122%
2	4.039	106	111	118	rBV	79648	125946	13.99%	0.984%
3	4.251	141	147	154	rVB3	11752	18732	2.08%	0.146%
4	4.504	188	190	193	rVV4	1667	1838	0.20%	0.014%
5	4.750	228	232	236	rBV5	3663	5583	0.62%	0.044%
6	5.309	320	327	332	rBV3	7463	11280	1.25%	0.088%
7	6.196	473	478	480	rBV3	1478	2105	0.23%	0.016%
8	6.678	556	560	564	rBV2	2268	4168	0.46%	0.033%
9	7.065	623	626	630	rVV2	1437	1747	0.19%	0.014%
10	7.130	632	637	642	rBV	1396	1979	0.22%	0.015%
11	7.195	646	648	652	rVB2	2144	3161	0.35%	0.025%
12	7.412	678	685	696	rVB	133510	231446	25.71%	1.808%
13	7.612	713	719	728	rVB	115951	191794	21.31%	1.498%
14	7.811	747	753	766	rVB	135420	231774	25.75%	1.810%
15	8.082	794	799	803	rVV4	2307	3473	0.39%	0.027%
16	8.299	828	836	841	rBV	80307	138090	15.34%	1.079%
17	8.981	943	952	958	rBV2	144406	256609	28.51%	2.004%
18	9.051	961	964	967	rVB	1320	1786	0.20%	0.014%
19	9.139	973	979	988	rBV	22124	40608	4.51%	0.317%
20	9.286	999	1004	1010	rBV4	1882	4295	0.48%	0.034%
21	9.492	1032	1039	1048	rVB	149421	255106	28.34%	1.993%
22	10.209	1154	1161	1170	rVB	117270	205963	22.88%	1.609%
23	10.303	1170	1177	1181	rBV3	6032	11456	1.27%	0.089%
24	10.567	1217	1222	1226	rVB3	3348	4949	0.55%	0.039%
25	10.732	1241	1250	1258	rBV	168692	297979	33.11%	2.328%
26	11.137	1312	1319	1326	rBV	134764	234885	26.10%	1.835%
27	11.278	1337	1343	1352	rBV2	75116	145984	16.22%	1.140%
28	11.660	1405	1408	1411	rBV	2112	2395	0.27%	0.019%
29	11.836	1435	1438	1442	rBV	1577	2224	0.25%	0.017%
30	11.924	1451	1453	1457	rVB3	1692	2368	0.26%	0.018%
31	12.688	1578	1583	1588	rVB4	2801	4273	0.47%	0.033%
32	13.499	1717	1721	1725	rBV	1992	2721	0.30%	0.021%
33	13.757	1755	1765	1772	rBV	144365	240737	26.75%	1.880%
34	13.851	1777	1781	1784	rVB3	2104	2565	0.28%	0.020%
35	13.934	1791	1795	1798	rVB2	3657	5581	0.62%	0.044%
36	14.039	1809	1813	1814	rBV2	2257	2232	0.25%	0.017%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
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37	14.051	1814	1815	1818	rVV	2555	2436	0.27%	0.019%
38	14.092	1820	1822	1825	rVB2	2349	2729	0.30%	0.021%
39	14.316	1854	1860	1867	rVB	338774	466506	51.83%	3.644%
40	14.392	1869	1873	1876	rBV	1425	2036	0.23%	0.016%

41	14.510	1889	1893	1895	rBV2	1752	1951	0.22%	0.015%
42	14.627	1906	1913	1922	rVB	393660	614981	68.33%	4.804%
43	14.833	1944	1948	1951	rVB2	2236	3043	0.34%	0.024%
44	14.921	1957	1963	1971	rVV	222387	341337	37.92%	2.666%
45	15.097	1984	1993	1999	rBV	135728	217320	24.15%	1.698%

46	15.197	2009	2010	2016	rVB2	2717	3941	0.44%	0.031%
47	15.267	2018	2022	2026	rVB5	6152	10220	1.14%	0.080%
48	15.479	2056	2058	2061	rBV2	1459	2106	0.23%	0.016%
49	15.520	2063	2065	2068	rVB2	1694	1831	0.20%	0.014%
50	15.579	2072	2075	2077	rBV2	1540	1774	0.20%	0.014%

51	15.638	2079	2085	2090	rVB2	2002	3824	0.42%	0.030%
52	15.679	2090	2092	2095	rBV2	1645	2169	0.24%	0.017%
53	15.914	2125	2132	2138	rVB	473901	745235	82.80%	5.821%
54	16.014	2142	2149	2156	rBV	161899	229465	25.49%	1.792%
55	16.484	2228	2229	2234	rBV3	2231	2664	0.30%	0.021%

56	16.554	2237	2241	2243	rVB3	3248	4389	0.49%	0.034%
57	16.613	2248	2251	2252	rVB2	2269	1870	0.21%	0.015%
58	16.701	2262	2266	2269	rVV3	5736	6110	0.68%	0.048%
59	16.889	2294	2298	2301	rBV2	5379	8236	0.92%	0.064%
60	16.924	2302	2304	2307	rVB4	2400	2503	0.28%	0.020%

61	16.995	2312	2316	2321	rBV2	12465	19409	2.16%	0.152%
62	17.101	2332	2334	2335	rBV	2218	1965	0.22%	0.015%
63	17.112	2335	2336	2338	rVV	2364	1695	0.19%	0.013%
64	17.230	2353	2356	2358	rVV2	2100	2146	0.24%	0.017%
65	17.271	2361	2363	2366	rVB4	2306	1869	0.21%	0.015%

66	17.347	2371	2376	2381	rBV5	10182	16376	1.82%	0.128%
67	17.465	2392	2396	2397	rBV2	2231	2407	0.27%	0.019%
68	17.488	2397	2400	2406	rVV2	26219	42689	4.74%	0.333%
69	17.553	2408	2411	2418	rVB3	23557	35742	3.97%	0.279%
70	17.671	2425	2431	2437	rBV	274354	426085	47.34%	3.328%

71	17.770	2442	2448	2455	rBV2	512956	806671	89.62%	6.301%
72	18.229	2524	2526	2530	rBV5	8150	9977	1.11%	0.078%
73	18.381	2546	2552	2557	rBV4	24986	47780	5.31%	0.373%
74	18.434	2559	2561	2565	rVV4	18625	25245	2.80%	0.197%
75	18.493	2566	2571	2584	rVB	191564	284643	31.63%	2.223%

76	18.663	2598	2600	2605	rVB6	3380	4215	0.47%	0.033%
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77	18.775	2615	2619	2621	rBV4	12981	19839	2.20%	0.155%
78	18.881	2636	2637	2639	rBV2	2367	1727	0.19%	0.013%
79	19.087	2668	2672	2674	rBV4	8060	10830	1.20%	0.085%
80	19.392	2721	2724	2728	rVB2	18346	20602	2.29%	0.161%
81	19.427	2728	2730	2735	rBV6	15060	17276	1.92%	0.135%
82	19.615	2758	2762	2764	rBV5	18783	29303	3.26%	0.229%
83	19.662	2768	2770	2777	rVV5	35787	45631	5.07%	0.356%
84	19.750	2781	2785	2792	rBV3	36293	47215	5.25%	0.369%
85	19.909	2807	2812	2813	rBV4	14661	25706	2.86%	0.201%
86	20.044	2830	2835	2842	rBV	614146	900055	100.00%	7.031%
87	20.491	2905	2911	2915	rBV	49741	76655	8.52%	0.599%
88	21.537	3084	3089	3104	rVB2	149668	297017	33.00%	2.320%
89	21.930	3153	3156	3160	rBV4	22182	35818	3.98%	0.280%
90	21.995	3161	3167	3173	rVV	218317	392315	43.59%	3.064%
91	22.389	3232	3234	3238	rVB3	31839	33797	3.75%	0.264%
92	22.765	3293	3298	3301	rBV	78636	126359	14.04%	0.987%
93	22.806	3301	3305	3323	rVB3	132249	312993	34.77%	2.445%
94	23.722	3457	3461	3467	rVB4	26028	47267	5.25%	0.369%
95	23.916	3490	3494	3502	rVB4	72131	153430	17.05%	1.198%
96	24.404	3568	3577	3585	rBV2	331747	800277	88.91%	6.251%
97	24.492	3587	3592	3603	rBV5	64095	193167	21.46%	1.509%
98	25.291	3718	3728	3741	rVB2	293897	870624	96.73%	6.801%
99	25.538	3762	3770	3785	rVB2	143421	451688	50.18%	3.528%
100	26.695	3958	3967	3981	rVB3	234411	771471	85.71%	6.026%

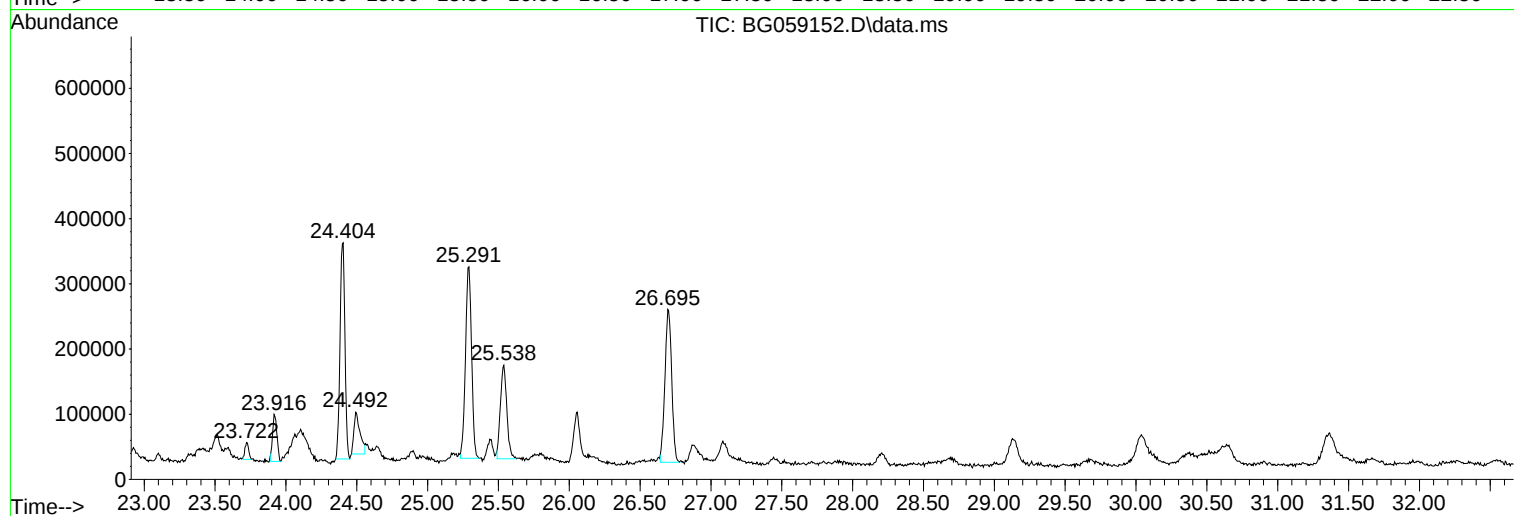
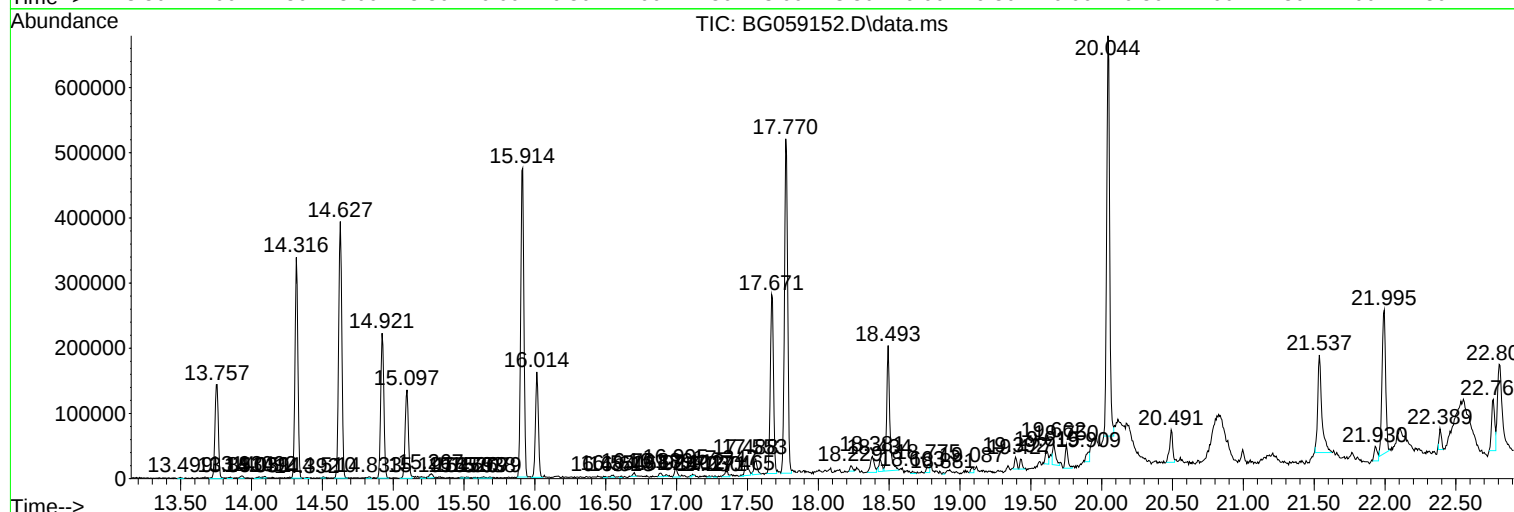
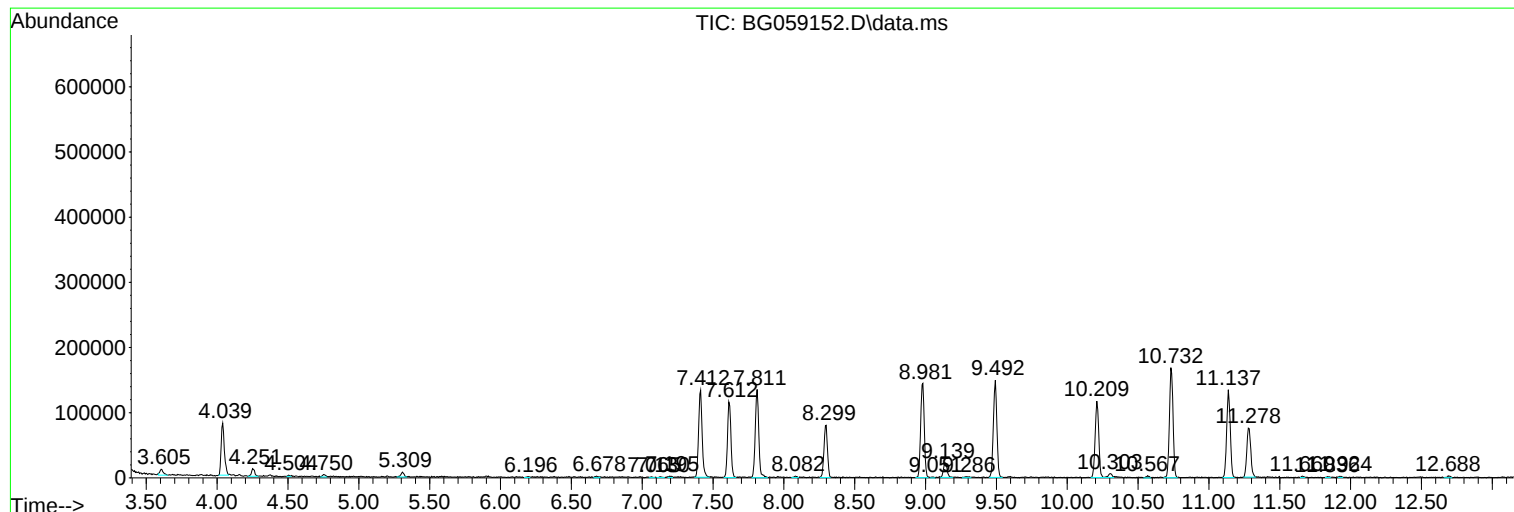
Sum of corrected areas: 12802135

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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



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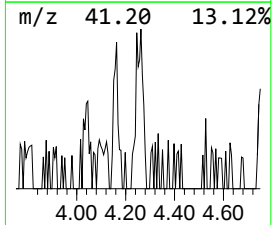
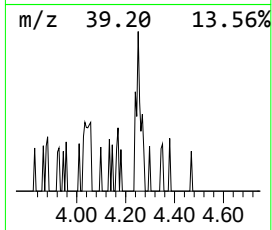
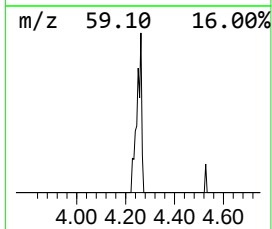
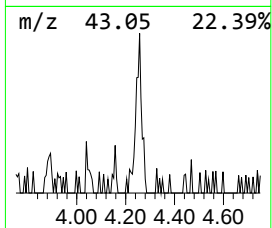
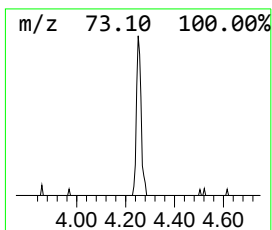
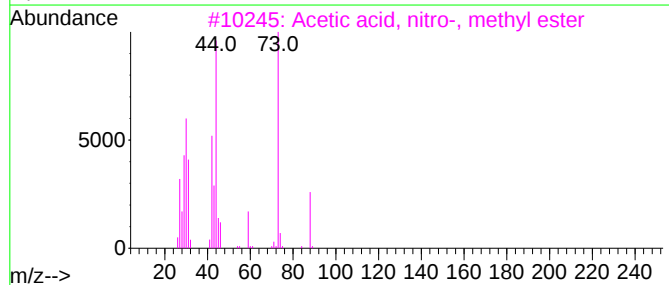
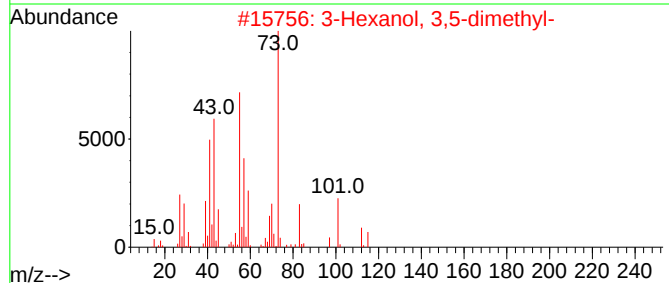
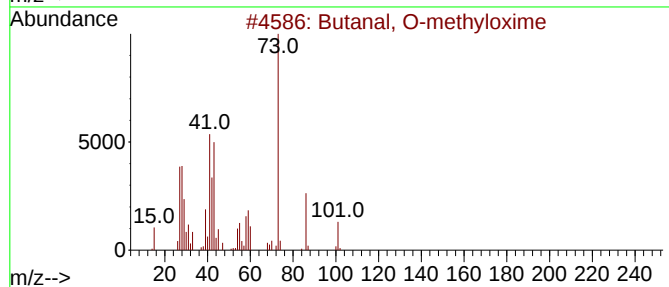
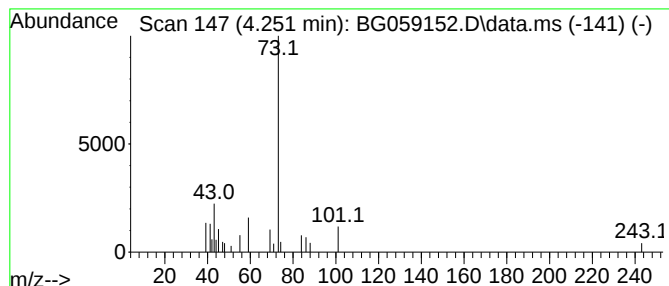
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.251	2.71 ng/ul	18732	1,4-Dichlorobenzene-d4	8.293

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanal, O-methyloxime	101	C5H11NO	031376-98-4	37
2			3-Hexanol, 3,5-dimethyl-	130	C8H18O	004209-91-0	23
3			Acetic acid, nitro-, methyl ester	119	C3H5NO4	002483-57-0	10
4			N-(n-Propyl)acetamide	101	C5H11NO	005331-48-6	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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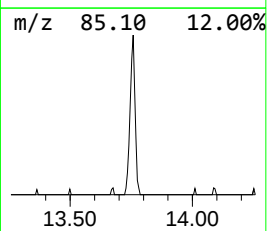
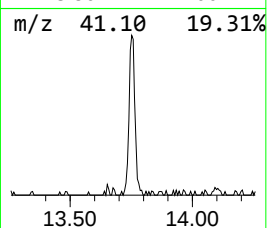
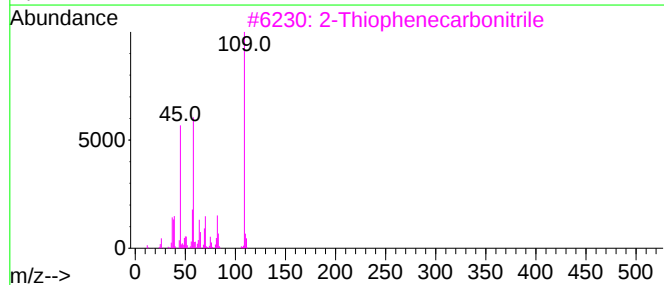
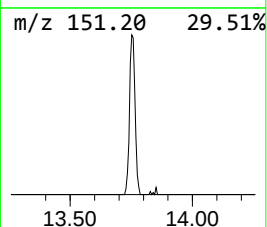
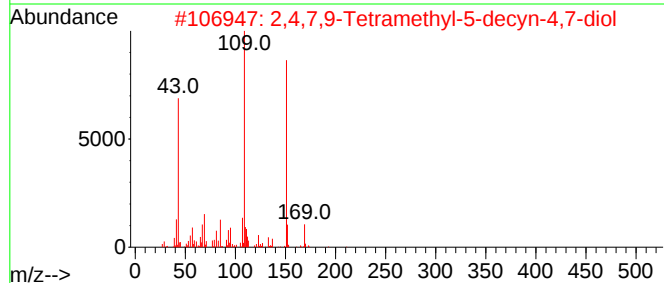
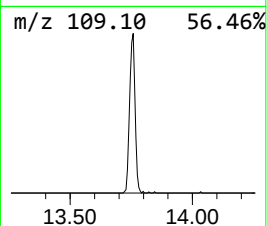
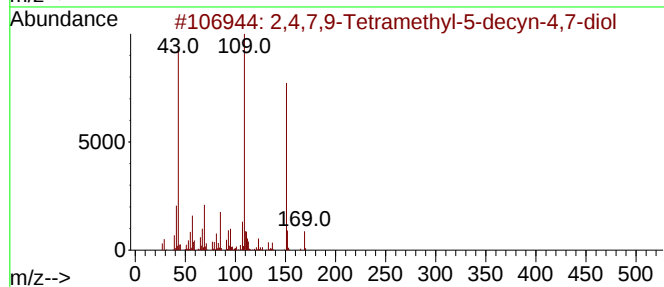
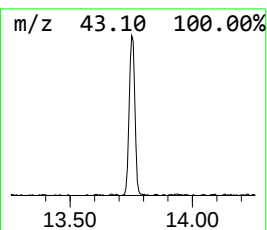
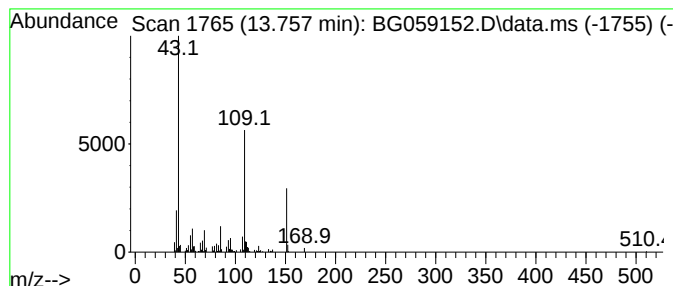
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.758	14.11 ng/ul	240737	Acenaphthene-d10	14.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	58		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	47		
3	2-Thiophenecarbonitrile	109	C5H3NS	001003-31-2	47		
4	Metacetamol	151	C8H9NO2	000621-42-1	43		
5	Acetaminophen	151	C8H9NO2	000103-90-2	43		



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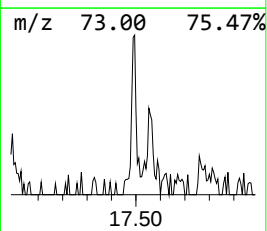
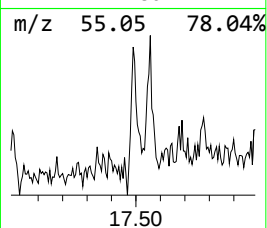
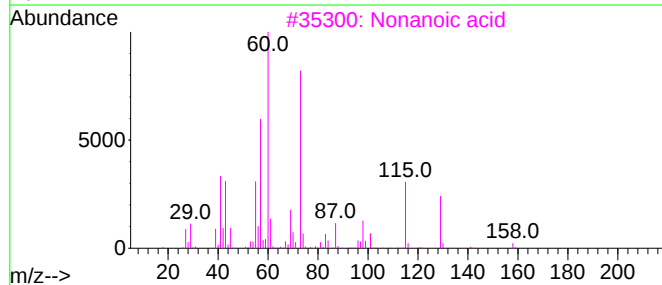
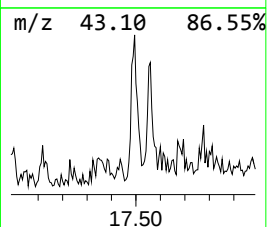
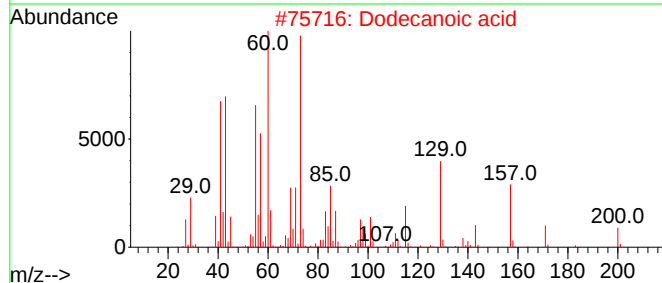
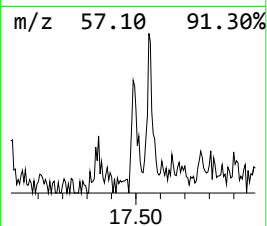
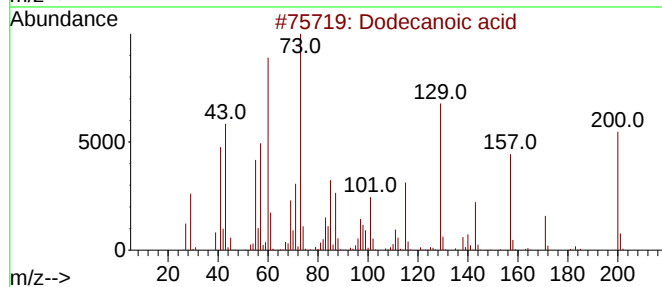
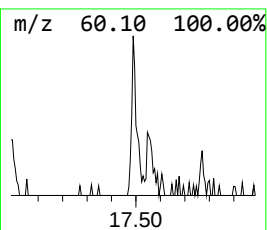
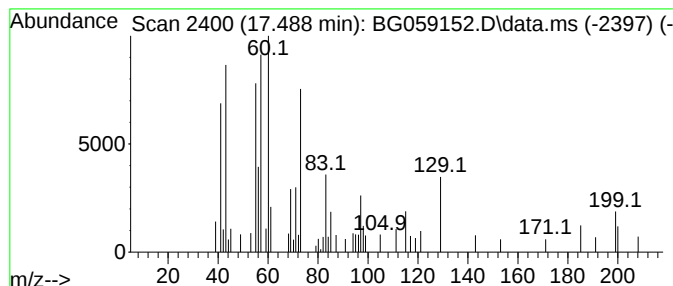
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Dodecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.488	2.00 ng/ul	42689	Phenanthrene-d10	17.671

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid	200	C12H24O2	000143-07-7	55
2			Dodecanoic acid	200	C12H24O2	000143-07-7	47
3			Nonanoic acid	158	C9H18O2	000112-05-0	47
4			n-Decanoic acid	172	C10H20O2	000334-48-5	43
5			n-Decanoic acid	172	C10H20O2	000334-48-5	43



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Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
Operator : MA/JU
Sample : 04386-09
Misc :
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Instrument :
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ClientSampleId :
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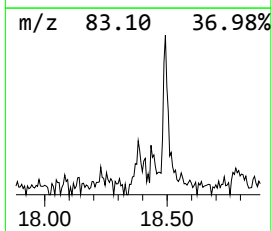
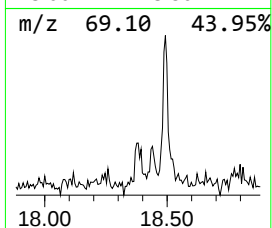
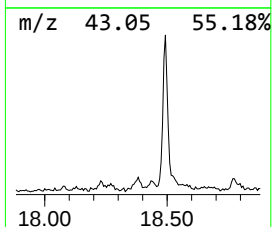
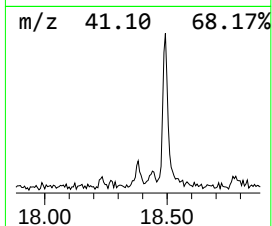
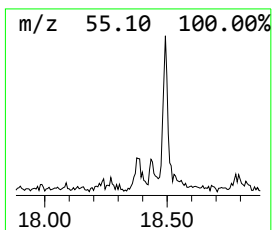
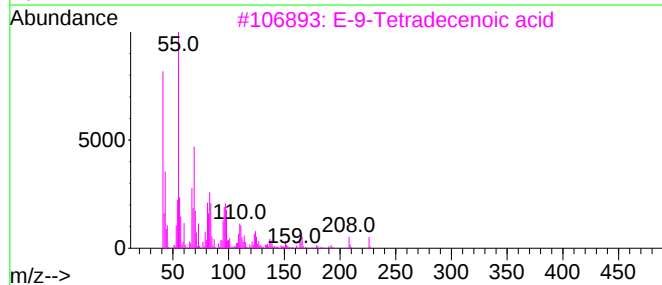
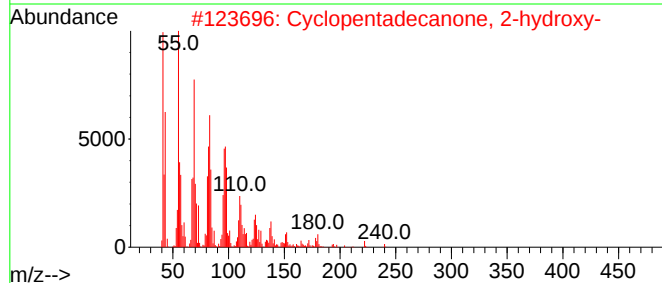
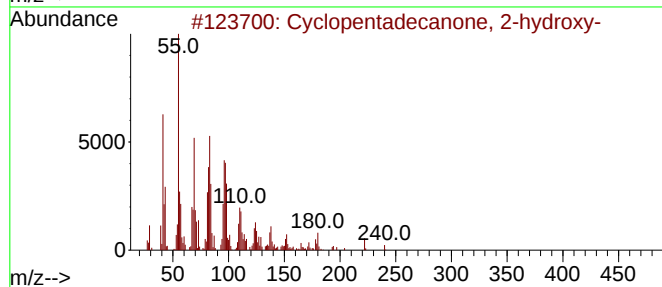
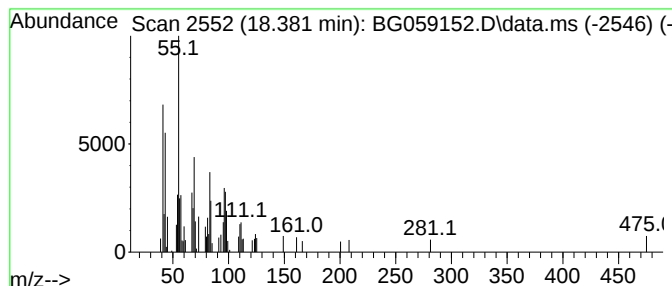
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclopentadecanone, 2-hydroxy- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.381	2.24 ng/ul	47780	Phenanthrene-d10	17.671

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	83
2			Cyclopentadecanone, 2-hydroxy-	240	C15H28O2	004727-18-8	83
3			E-9-Tetradecenoic acid	226	C14H26O2	050286-30-1	81
4			cis-Vaccenic acid	282	C18H34O2	000506-17-2	74
5			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
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ClientSampleId :
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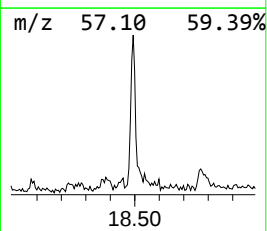
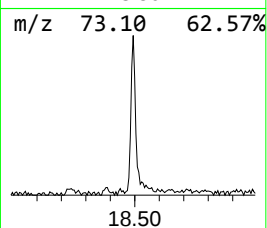
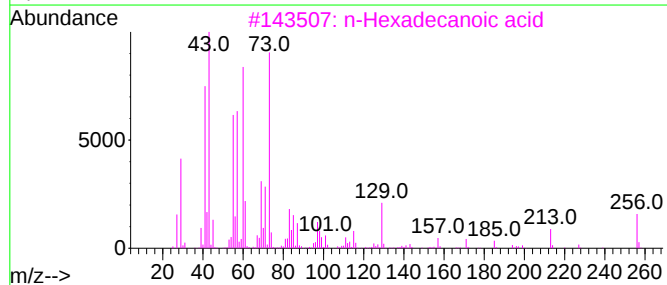
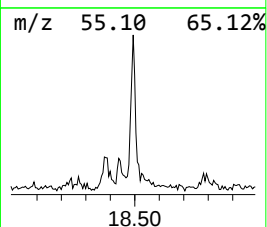
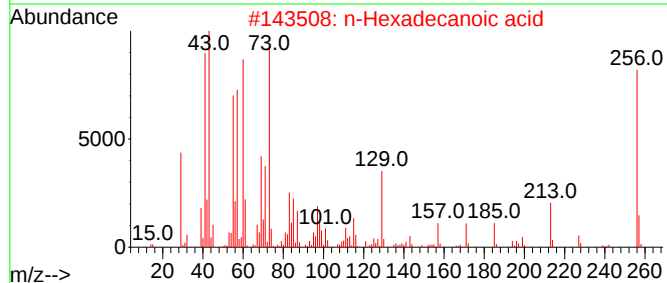
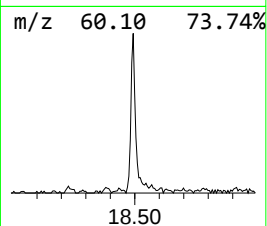
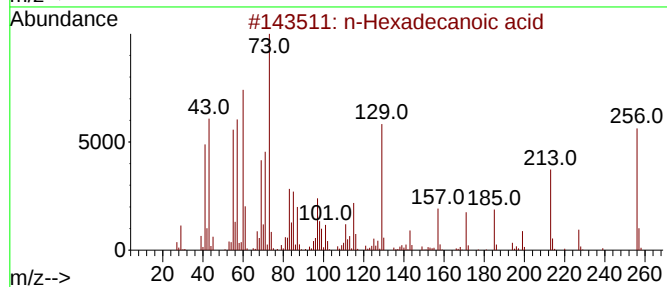
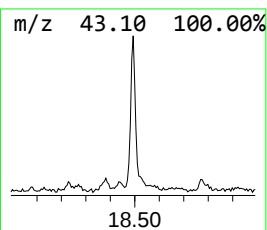
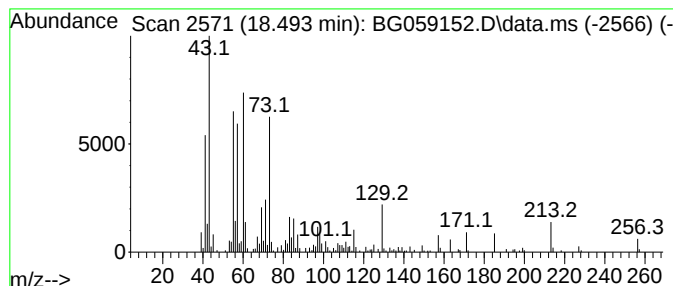
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.493	13.36 ng/ul	284643	Phenanthrene-d10	17.671

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	70
5			n-Decanoic acid	172	C10H20O2	000334-48-5	68



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Sample : 04386-09
Misc :
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Instrument :
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ClientSampleId :
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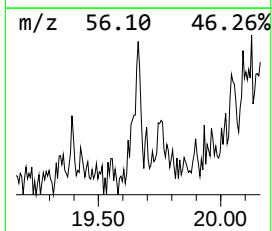
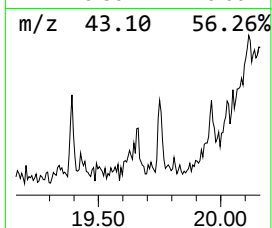
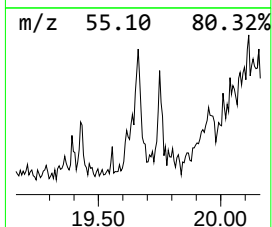
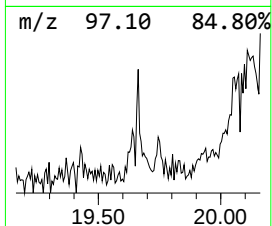
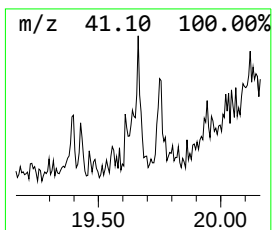
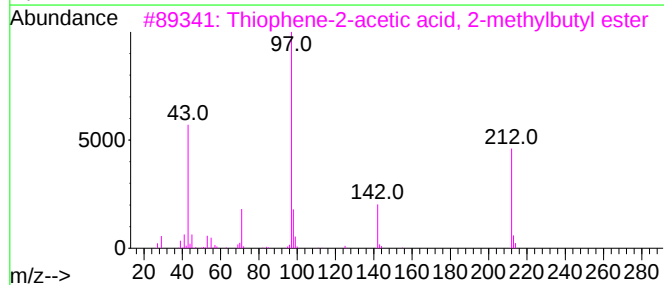
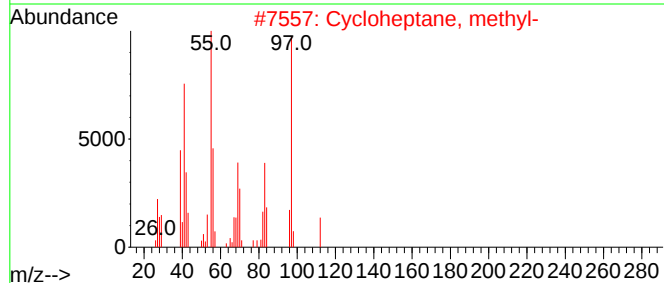
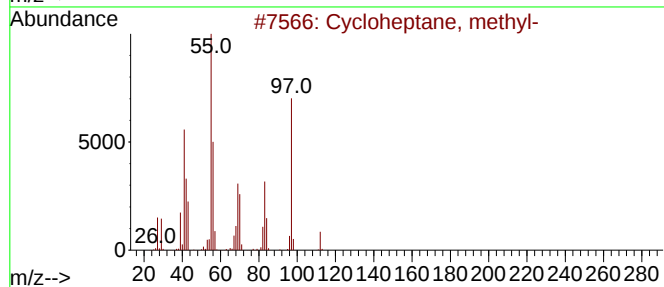
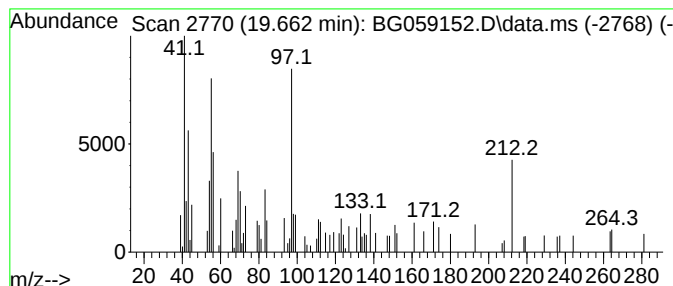
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Cyclic19.662 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.662	2.14 ng/ul	45631	Phenanthrene-d10	17.671

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cycloheptane, methyl-	112	C8H16	004126-78-7	50
2			Cycloheptane, methyl-	112	C8H16	004126-78-7	45
3			Thiophene-2-acetic acid, 2-methy...	212	C11H16O2S	1000330-96-2	38
4			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	30
5			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
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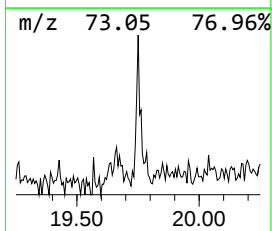
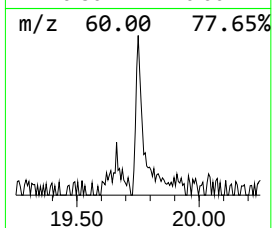
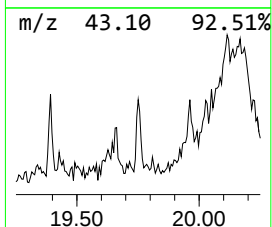
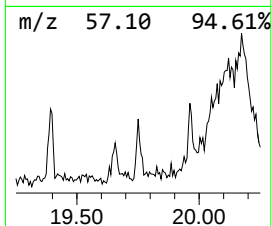
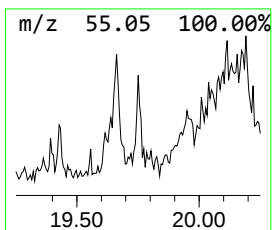
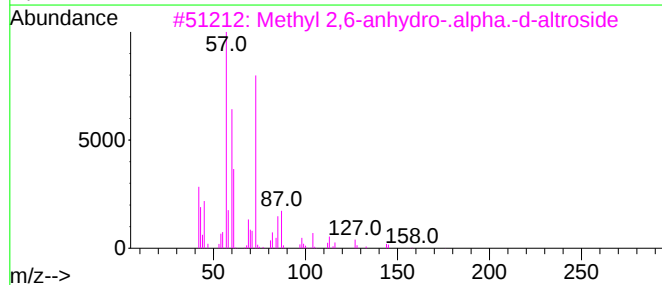
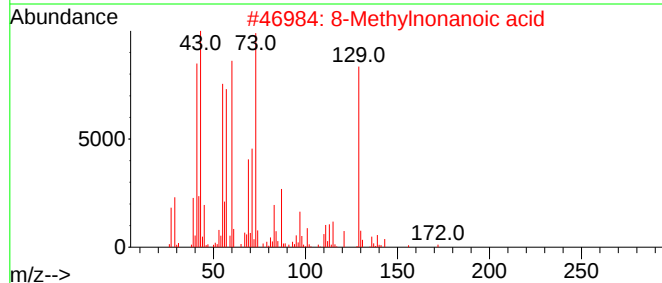
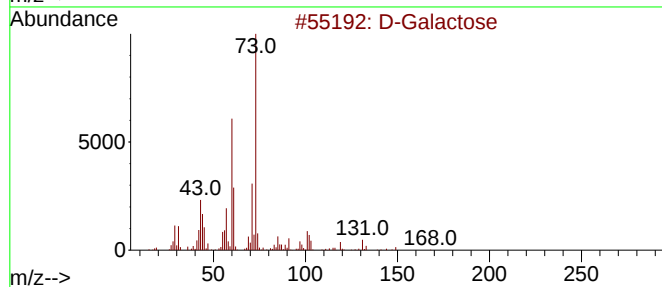
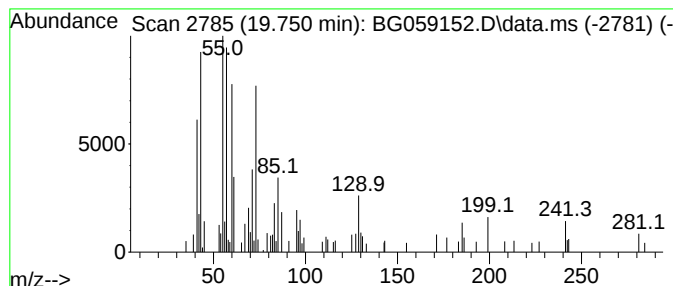
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.750	2.22 ng/ul	47215	Phenanthrene-d10	17.671

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D-Galactose	180	C6H12O6	000059-23-4	47
2			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	47
3			Methyl 2,6-anhydro-.alpha.-d-alt...	176	C7H12O5	1000130-02-0	43
4			3,5-Difluorobenzaldehyde carbamo...	199	C8H7F2N3O	163893-49-0	27
5			Decyl isopropyl ether	200	C13H28O	1000406-33-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
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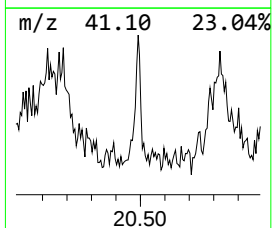
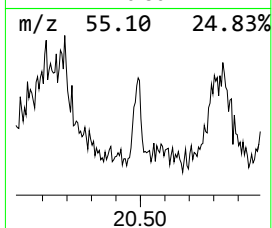
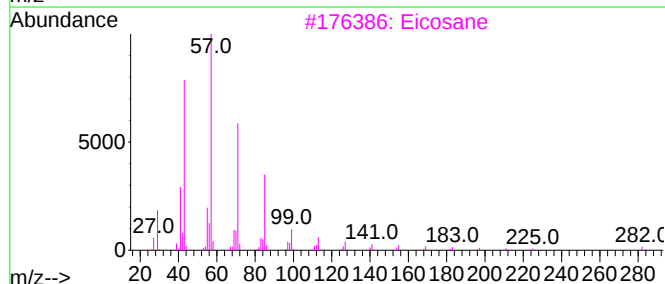
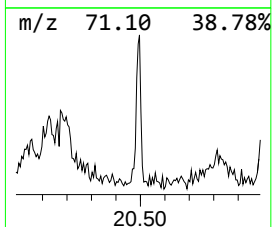
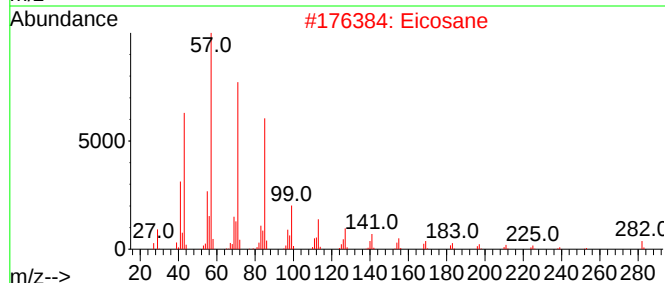
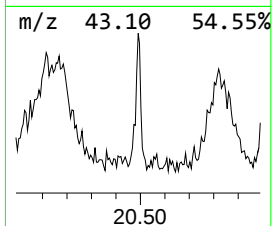
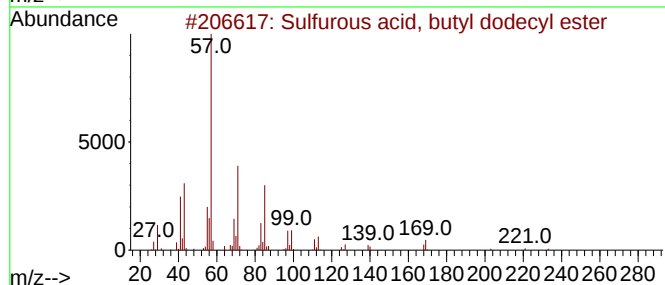
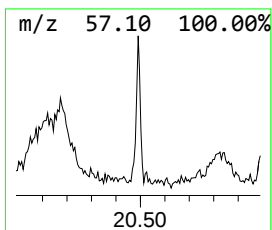
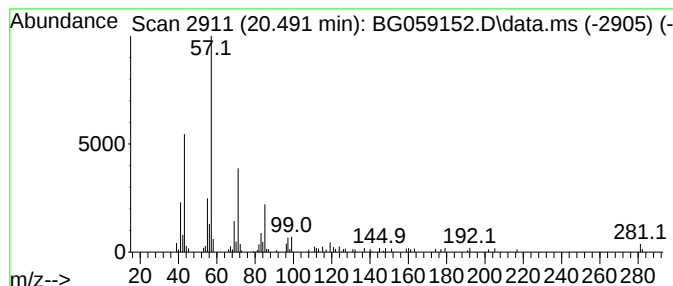
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Sulfurous acid, butyl dodec... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.491	3.91 ng/ul	76655	Chrysene-d12	21.995

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	64
2			Eicosane	282	C20H42	000112-95-8	62
3			Eicosane	282	C20H42	000112-95-8	58
4			Hexadecane	226	C16H34	000544-76-3	50
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	50



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Misc :
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ClientSampleId :
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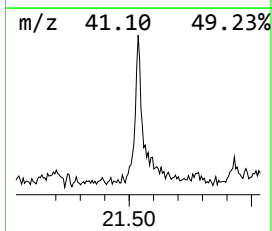
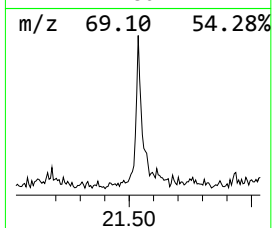
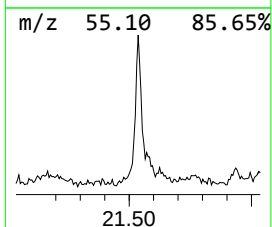
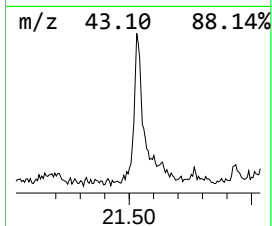
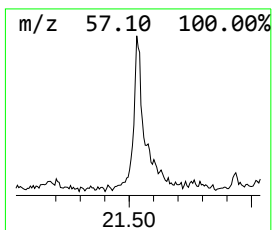
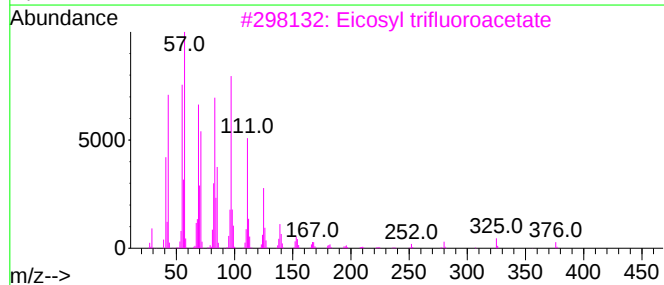
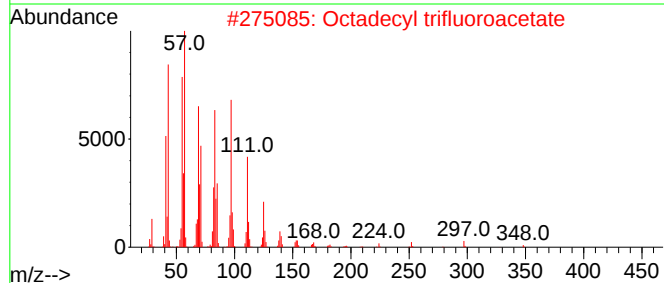
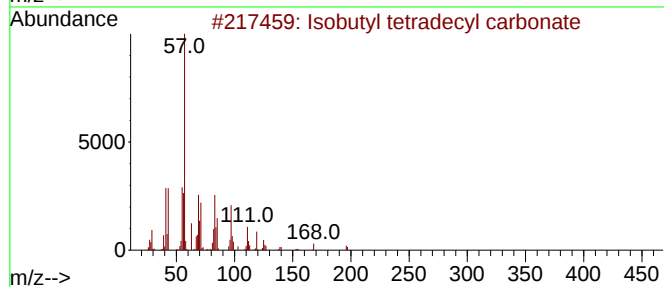
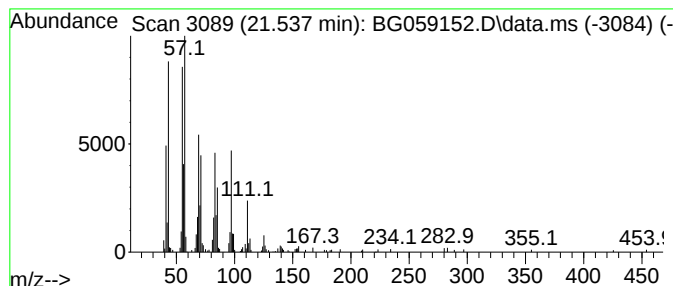
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Isobutyl tetradecyl carbonate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.537	15.14 ng/ul	297017	Chrysene-d12	21.995

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl tetradecyl carbonate	314	C19H38O3	959275-58-2	91
2			Octadecyl trifluoroacetate	366	C20H37F3O2	079392-43-1	90
3			Eicosyl trifluoroacetate	394	C22H41F3O2	1000351-75-2	90
4			Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	90
5			Tricosyl pentafluoropropionate	486	C26H47F5O2	1000351-81-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
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Misc :
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ClientSampleId :
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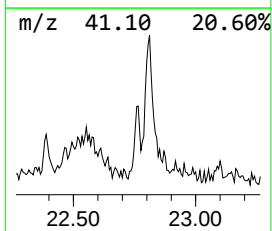
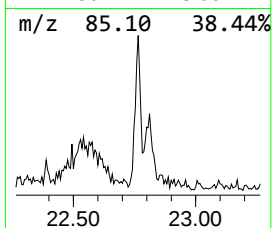
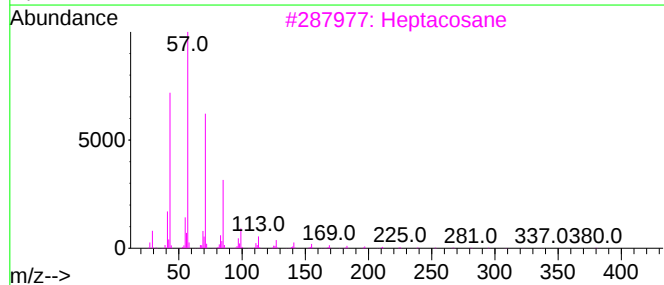
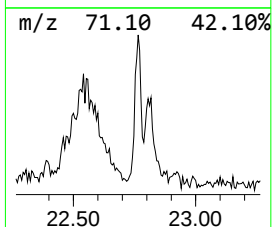
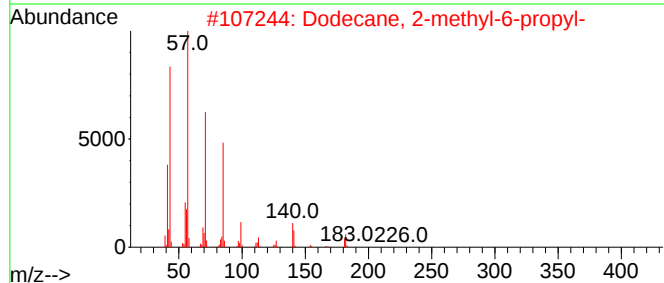
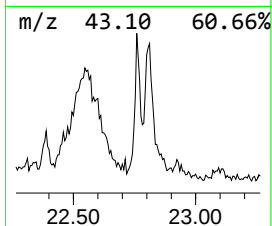
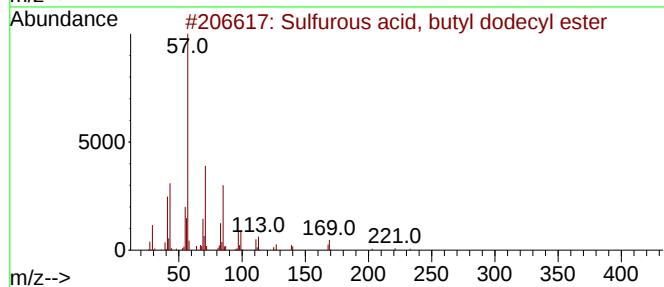
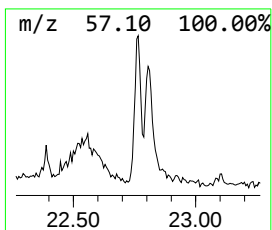
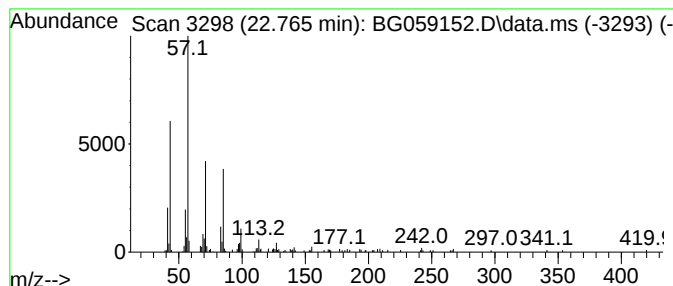
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dodecane, 2-methyl-6-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.765	6.44 ng/ul	126359	Chrysene-d12	21.995

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	83
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	78
3		Heptacosane	380	C27H56	000593-49-7	68
4		Tridecane, 3-methyl-	198	C14H30	006418-41-3	59
5		Tetracosane	338	C24H50	000646-31-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
Operator : MA/JU
Sample : 04386-09
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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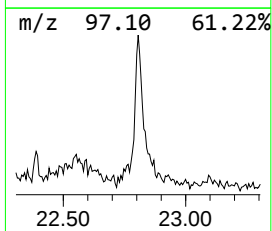
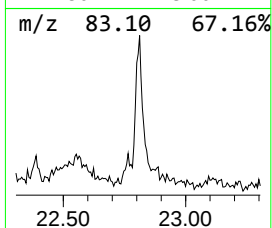
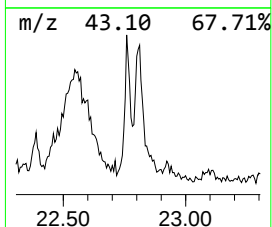
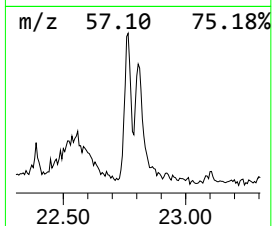
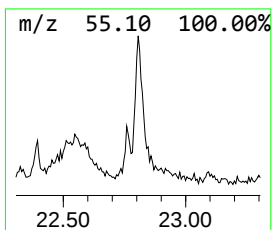
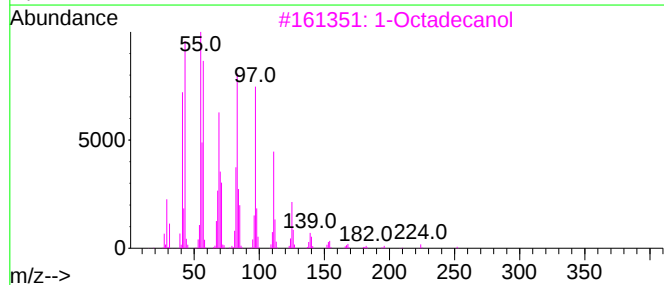
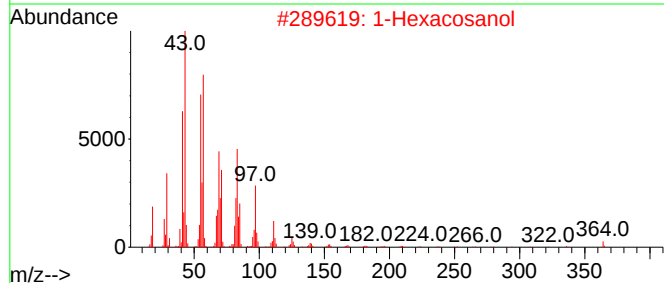
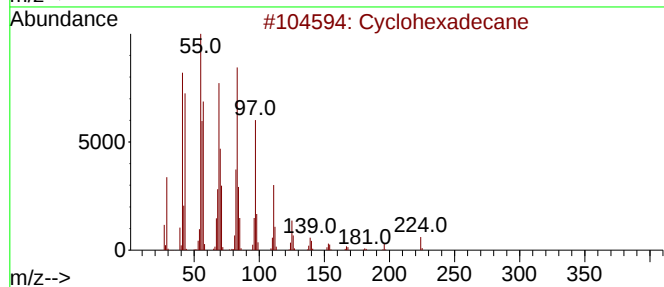
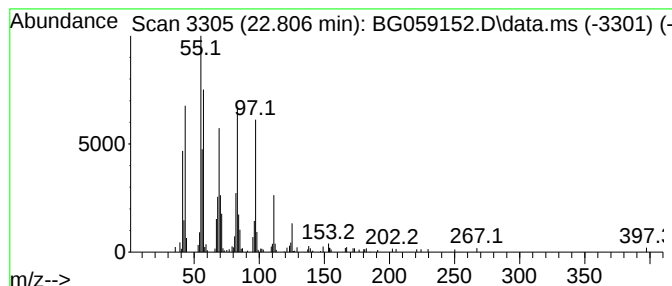
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic22.806 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.806	15.96 ng/ul	312993	Chrysene-d12	21.995

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	83
2		1-Hexacosanol	382	C26H54O	000506-52-5	80
3		1-Octadecanol	270	C18H38O	000112-92-5	72
4		1-Hexadecanol	242	C16H34O	036653-82-4	72
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
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Sample : 04386-09
Misc :
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Instrument :
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ClientSampleId :
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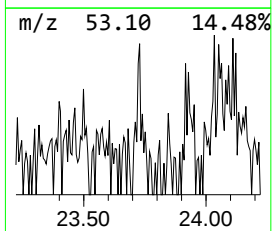
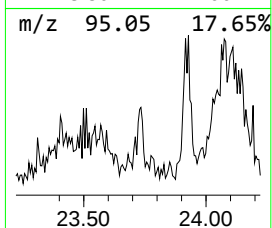
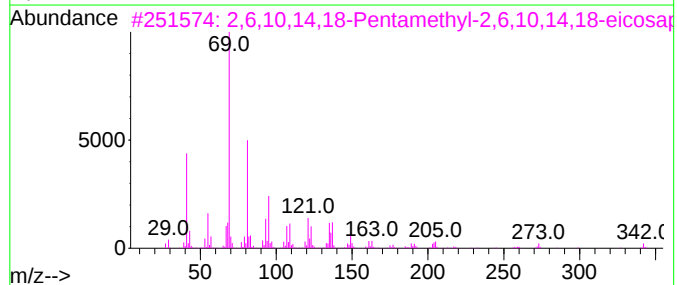
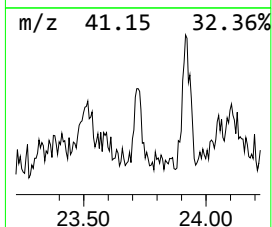
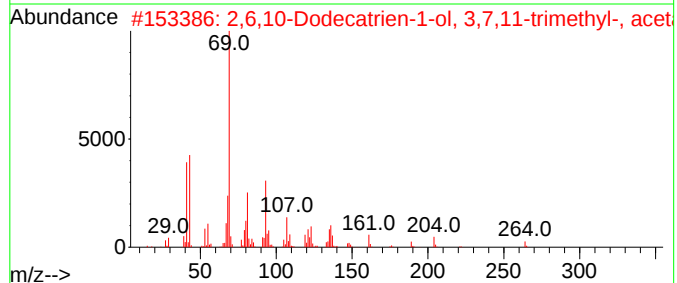
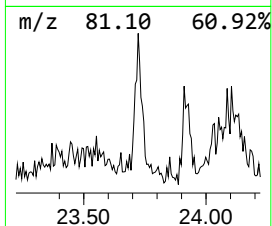
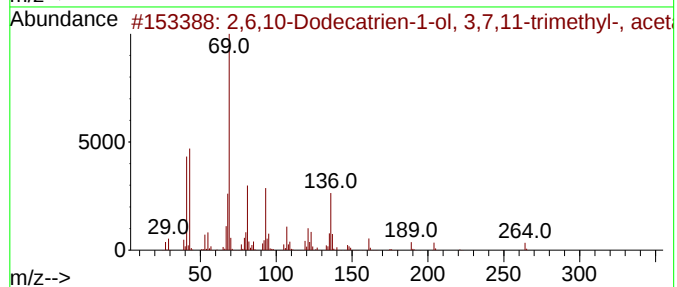
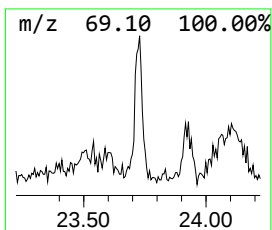
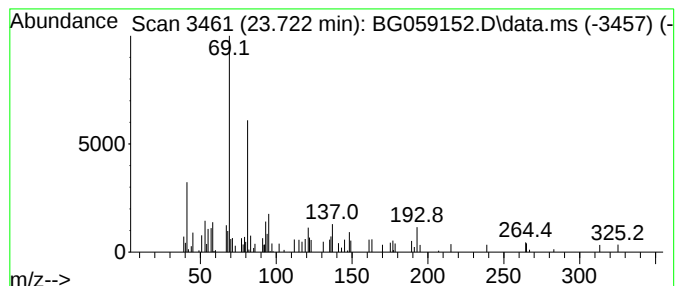
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2,6,10-Dodecatrien-1-ol, 3,... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.722	2.41 ng/ul	47267	Chrysene-d12	21.995

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	70		
2	2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	53		
3	2,6,10,14,18-Pentamethyl-2,6,10,...	342	C25H42	075581-03-2	53		
4	1-(Pyrrolidin-3-yl)pyrazole	137	C7H11N3	1000459-08-8	50		
5	Supraene	410	C30H50	007683-64-9	50		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
Operator : MA/JU
Sample : 04386-09
Misc :
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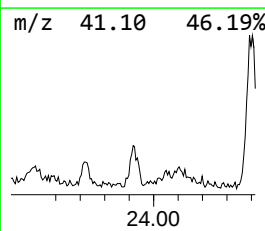
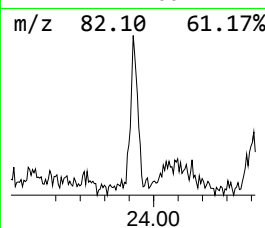
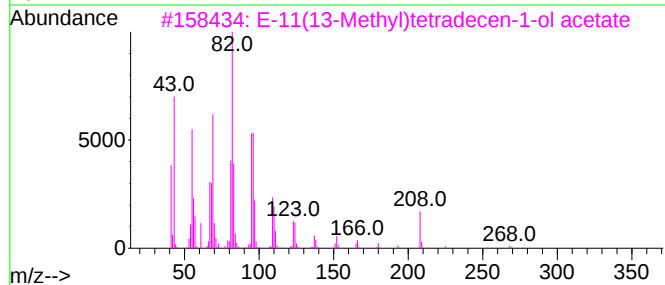
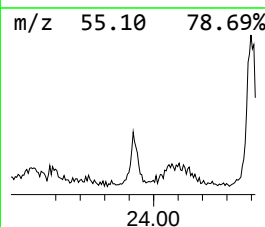
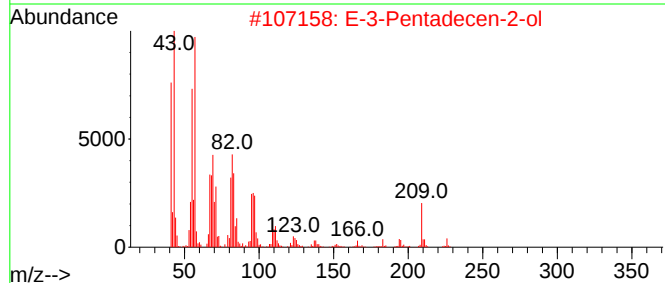
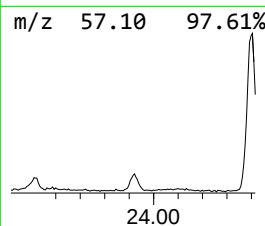
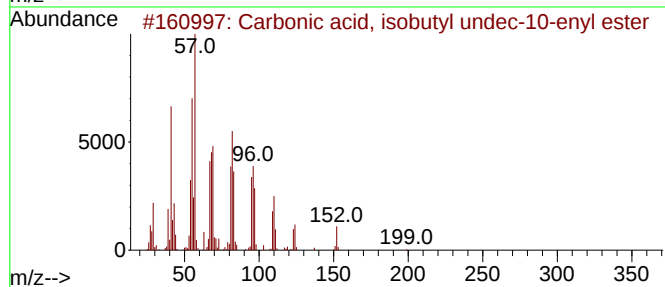
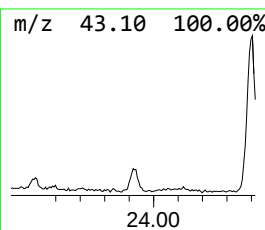
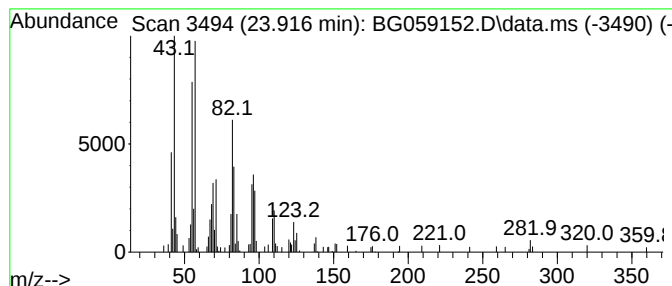
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Carbonic acid, isobutyl und... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.916	6.79 ng/ul	153430	Perylene-d12	25.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, isobutyl undec-10...	270	C16H30O3	1000314-60-8	64
2			E-3-Pentadecen-2-ol	226	C15H30O	1000130-83-8	64
3			E-11(13-Methyl)tetradecen-1-ol a...	268	C17H32O2	1000130-80-4	50
4			cis-11-Tetradecen-1-ol	212	C14H28O	034010-15-6	47
5			E-9-Tetradecen-1-ol formate	240	C15H28O2	1000130-78-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
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Sample : 04386-09
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
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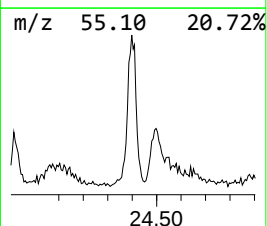
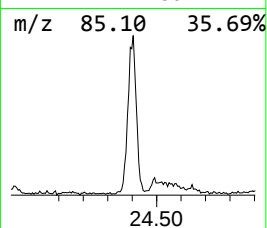
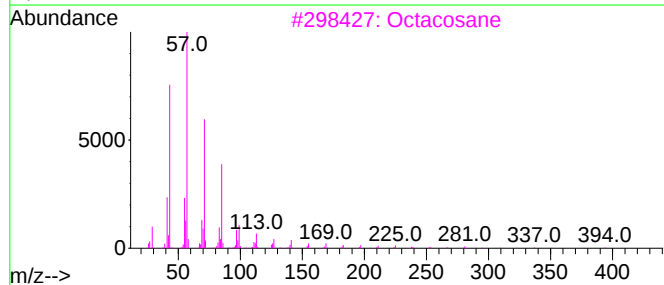
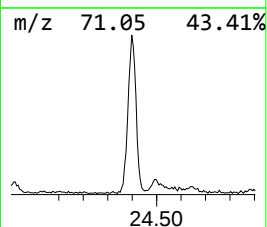
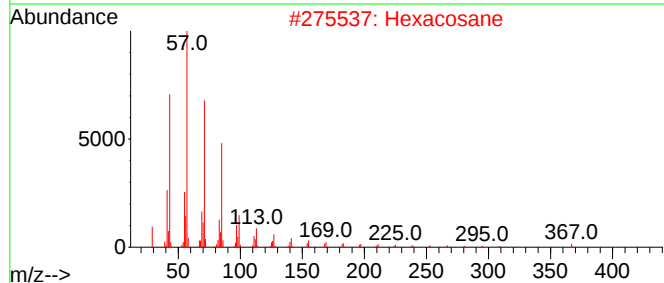
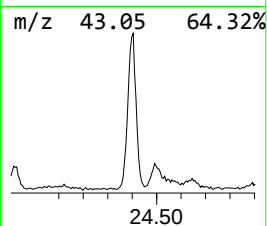
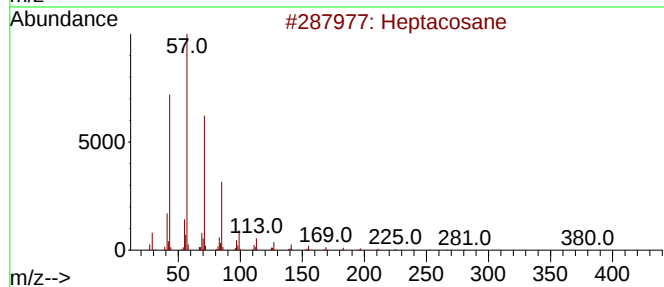
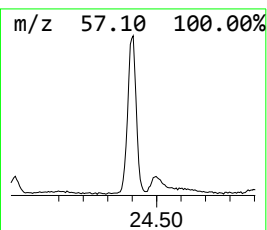
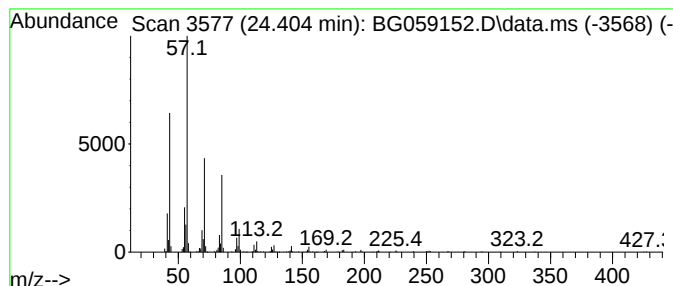
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.404	35.43 ng/ul	800277	Perylene-d12	25.538

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	90
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetracosane	338	C24H50	000646-31-1	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
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Misc :
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Instrument :
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ClientSampleId :
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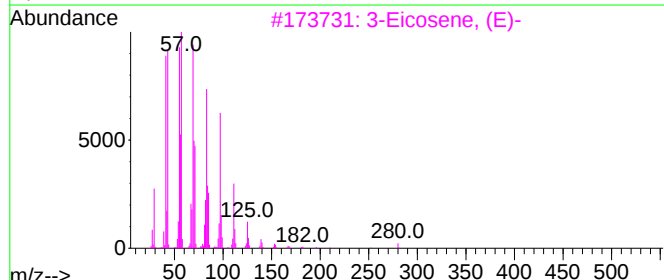
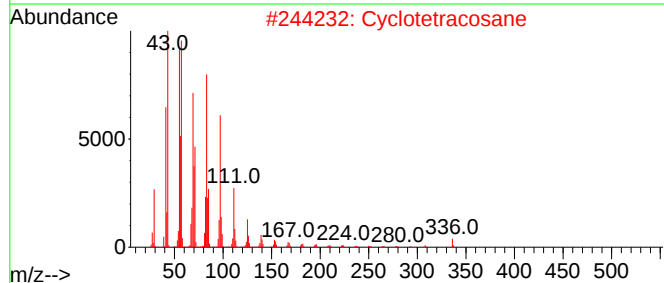
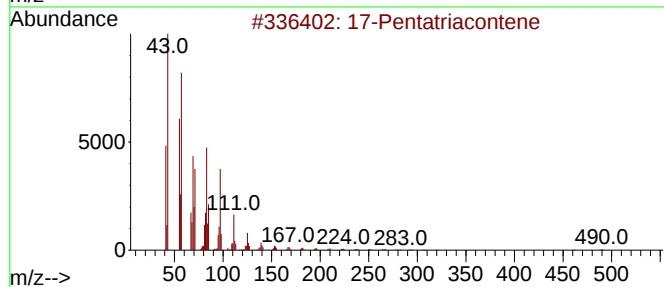
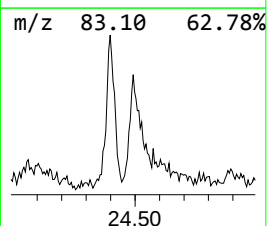
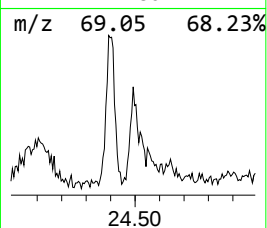
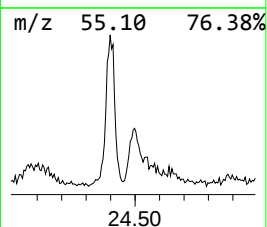
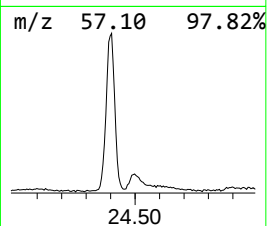
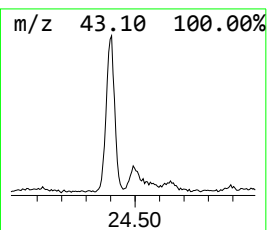
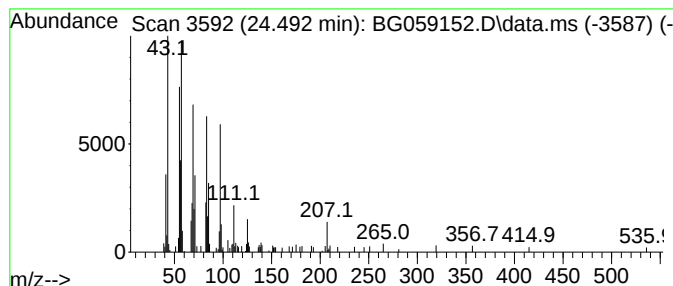
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.492	8.55 ng/ul	193167	Perylene-d12	25.538

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Pentatriacontene	491	C35H70	006971-40-0	86
2		Cyclotetracosane	336	C24H48	000297-03-0	86
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	62
4		Cyclohexanecarboxaldehyde, 3,3-d...	154	C9H14O2	065080-66-2	58
5		Cyclopentane, 2-isopropyl-1,3-di...	140	C10H20	032281-85-9	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
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Sample : 04386-09
Misc :
ALS Vial : 53 Sample Multiplier: 1

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ClientSampleId :
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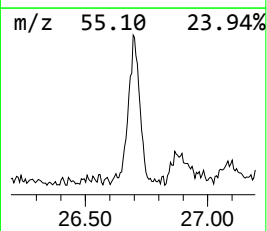
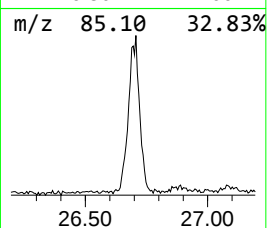
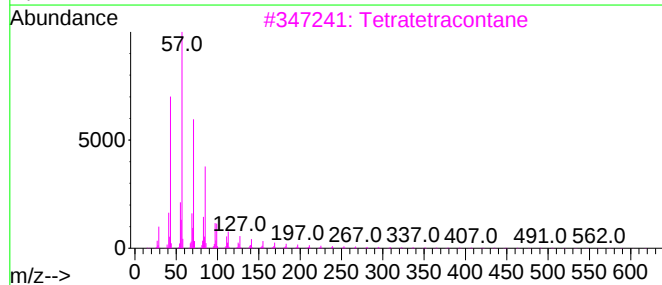
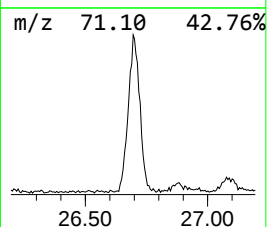
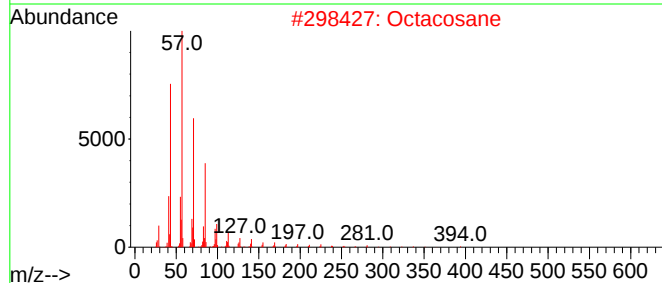
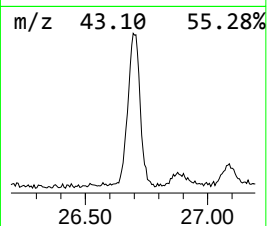
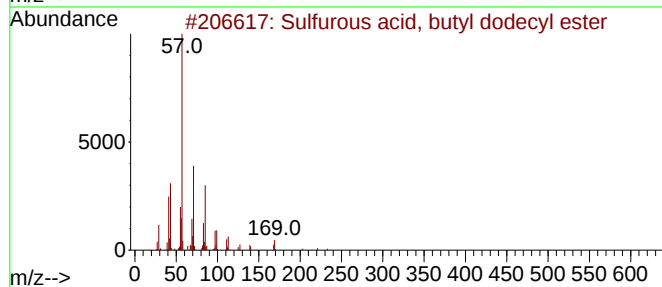
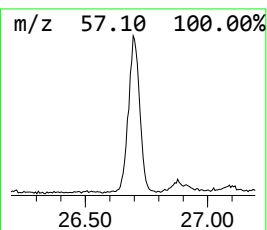
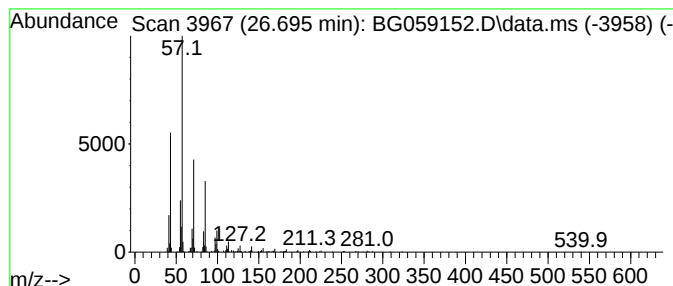
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Octacosane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.695	34.16 ng/ul	771471	Perylene-d12	25.538

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	91
2			Octacosane	394	C28H58	000630-02-4	90
3			Tetratetracontane	619	C44H90	007098-22-8	90
4			Tetracosane	338	C24H50	000646-31-1	81
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092023\
Data File : BG059152.D
Acq On : 21 Sep 2023 23:30
Operator : MA/JU
Sample : 04386-09
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AC9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091923.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
unknown-01	4.251	2.7	ng/ul	18732	1	8.293	138090	20.0
2,4,7,9-Tetrame...	13.758	14.1	ng/ul	240737	3	14.921	341337	20.0
Dodecanoic acid	17.488	2.0	ng/ul	42689	4	17.671	426085	20.0
Cyclopentadecan...	18.381	2.2	ng/ul	47780	4	17.671	426085	20.0
n-Hexadecanoic ...	18.493	13.4	ng/ul	284643	4	17.671	426085	20.0
(DEL) Alkane: C...	19.662	2.1	ng/ul	45631	4	17.671	426085	20.0
unknown-02	19.750	2.2	ng/ul	47215	4	17.671	426085	20.0
Sulfurous acid,...	20.491	3.9	ng/ul	76655	5	21.995	392315	20.0
Isobutyl tetrad...	21.537	15.1	ng/ul	297017	5	21.995	392315	20.0
Dodecane, 2-met...	22.765	6.4	ng/ul	126359	5	21.995	392315	20.0
(DEL) Alkane: C...	22.806	16.0	ng/ul	312993	5	21.995	392315	20.0
2,6,10-Dodecatr...	23.722	2.4	ng/ul	47267	5	21.995	392315	20.0
Carbonic acid, ...	23.916	6.8	ng/ul	153430	6	25.538	451688	20.0
(DEL) Alkane: S...	24.404	35.4	ng/ul	800277	6	25.538	451688	20.0
(DEL) Alkane: S...	24.492	8.6	ng/ul	193167	6	25.538	451688	20.0
Octacosane	26.695	34.2	ng/ul	771471	6	25.538	451688	20.0