

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062784.D
 Acq On : 13 Sep 2024 12:26
 Operator : JU/RC
 Sample : P3952-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0B51

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-BG090924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062784.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.363	43	47	53	rVB4	13684	19852	0.79%	0.101%
2	3.486	66	68	70	rBV3	2971	3124	0.12%	0.016%
3	3.768	112	116	128	rBV	55789	97199	3.86%	0.494%
4	4.097	170	172	174	rBV2	3268	3202	0.13%	0.016%
5	5.014	323	328	331	rBV5	1715	3020	0.12%	0.015%
6	7.064	672	677	687	rVB	56761	95764	3.80%	0.487%
7	7.229	699	705	717	rBV	136342	223575	8.87%	1.137%
8	7.434	733	740	750	rVB	162849	276004	10.95%	1.403%
9	7.899	813	819	828	rBV	154478	255236	10.13%	1.298%
10	8.604	932	939	950	rBV	175008	297728	11.81%	1.514%
11	9.050	1008	1015	1022	rBV	176069	311561	12.36%	1.584%
12	9.491	1084	1090	1093	rBV3	1483	2574	0.10%	0.013%
13	9.614	1106	1111	1114	rVB2	2793	3460	0.14%	0.018%
14	9.773	1131	1138	1152	rBV	147389	256260	10.17%	1.303%
15	10.313	1222	1230	1239	rBV	262003	457613	18.16%	2.327%
16	10.466	1253	1256	1261	rVB3	1680	2789	0.11%	0.014%
17	10.689	1286	1294	1304	rVB	218467	396637	15.74%	2.017%
18	10.824	1309	1317	1328	rBV	210344	382348	15.17%	1.944%
19	12.282	1557	1565	1568	rBV3	3711	6874	0.27%	0.035%
20	13.421	1753	1759	1766	rVB2	17147	27617	1.10%	0.140%
21	13.933	1839	1846	1856	rVB	574373	811897	32.22%	4.128%
22	14.221	1888	1895	1902	rVV2	611547	932392	37.00%	4.741%
23	14.526	1940	1947	1954	rBV	419107	648618	25.74%	3.298%
24	14.726	1974	1981	1989	rBV2	58552	105026	4.17%	0.534%
25	14.973	2014	2023	2042	rBV	727763	1376113	54.61%	6.997%
26	15.260	2070	2072	2075	rVB3	2398	2763	0.11%	0.014%
27	15.325	2080	2083	2090	rVB7	2798	5013	0.20%	0.025%
28	15.519	2107	2116	2128	rBV	858566	1298701	51.54%	6.603%
29	15.636	2130	2136	2144	rBV	262764	393722	15.62%	2.002%
30	15.760	2153	2157	2161	rVB6	4953	6410	0.25%	0.033%
31	15.883	2170	2178	2183	rBV	103154	146509	5.81%	0.745%
32	16.447	2271	2274	2279	rVB4	6556	8762	0.35%	0.045%
33	16.671	2309	2312	2317	rBV3	4972	7679	0.30%	0.039%
34	16.953	2357	2360	2363	rBV3	2118	2998	0.12%	0.015%
35	17.270	2407	2414	2422	rVB	616361	920160	36.51%	4.679%
36	17.370	2424	2431	2440	rVV	1137322	1663480	66.01%	8.458%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062784.D
 Acq On : 13 Sep 2024 12:26
 Operator : JU/RC
 Sample : P3952-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0B51

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-BG090924.MA.M
 Title : SVOA CALIBRATION

37	17.440	2440	2443	2449	rVB4	3042	4378	0.17%	0.022%
38	17.552	2458	2462	2465	rBV3	2973	4041	0.16%	0.021%
39	17.840	2508	2511	2514	rBV3	3022	4225	0.17%	0.021%
40	17.940	2525	2528	2530	rBV3	2978	3798	0.15%	0.019%
41	18.034	2541	2544	2545	rBV2	3409	3086	0.12%	0.016%
42	18.163	2562	2566	2576	rBV2	82121	129700	5.15%	0.659%
43	18.233	2576	2578	2583	rVB4	5053	5775	0.23%	0.029%
44	18.351	2595	2598	2602	rBV4	2959	3844	0.15%	0.020%
45	18.474	2615	2619	2625	rVB7	8358	11808	0.47%	0.060%
46	18.586	2635	2638	2640	rBV2	2541	2543	0.10%	0.013%
47	18.651	2647	2649	2652	rBV4	4324	3637	0.14%	0.018%
48	18.745	2662	2665	2667	rBV4	2767	3939	0.16%	0.020%
49	18.886	2686	2689	2691	rBV4	3363	3289	0.13%	0.017%
50	19.021	2707	2712	2715	rVB6	6059	9047	0.36%	0.046%
51	19.226	2741	2747	2752	rBV4	19516	31152	1.24%	0.158%
52	19.297	2754	2759	2767	rVV2	17884	34604	1.37%	0.176%
53	19.444	2780	2784	2788	rVV4	25727	34065	1.35%	0.173%
54	19.591	2806	2809	2812	rVB4	4851	6388	0.25%	0.032%
55	19.661	2815	2821	2832	rVV	1576190	2239084	88.85%	11.385%
56	19.743	2832	2835	2838	rVV4	2172	3225	0.13%	0.016%
57	19.961	2870	2872	2874	rBV2	3549	4500	0.18%	0.023%
58	20.137	2898	2902	2903	rBV4	4238	3675	0.15%	0.019%
59	20.431	2950	2952	2954	rBV3	3256	3165	0.13%	0.016%
60	20.595	2977	2980	2983	rVB5	8151	10121	0.40%	0.051%
61	21.189	3076	3081	3090	rBV4	39473	78140	3.10%	0.397%
62	21.518	3130	3137	3145	rBV	752069	1202724	47.73%	6.115%
63	21.712	3167	3170	3173	rBV5	8968	10187	0.40%	0.052%
64	22.000	3215	3219	3224	rBV3	34012	59899	2.38%	0.305%
65	22.293	3267	3269	3276	rVB7	10106	12854	0.51%	0.065%
66	22.946	3377	3380	3385	rVB6	7745	13701	0.54%	0.070%
67	23.134	3406	3412	3419	rVB	34276	65628	2.60%	0.334%
68	23.186	3419	3421	3425	rBV5	4043	4302	0.17%	0.022%
69	23.263	3432	3434	3437	rVB4	5094	4423	0.18%	0.022%
70	23.292	3437	3439	3440	rBV	5747	3383	0.13%	0.017%
71	23.715	3505	3511	3516	rBV6	13834	27028	1.07%	0.137%
72	23.892	3535	3541	3552	rBV4	31503	82586	3.28%	0.420%
73	24.367	3611	3622	3635	rBV	970566	2519999	100.00%	12.813%
74	24.579	3648	3658	3672	rVB2	527909	1409572	55.94%	7.167%
75	24.849	3702	3704	3706	rBV3	4255	2755	0.11%	0.014%
76	25.660	3838	3842	3852	rVB7	20695	50456	2.00%	0.257%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

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-BG090924.MA.M
Title : SVOA CALIBRATION

77	26.318	3952	3954	3959	rVB5	5928	6163	0.24%	0.031%
78	26.647	4008	4010	4013	rBV4	5007	6052	0.24%	0.031%
79	26.782	4026	4033	4046	rBV7	21449	73754	2.93%	0.375%
80	27.869	4216	4218	4221	rBV4	4764	4907	0.19%	0.025%
81	31.300	4800	4802	4803	rBV2	5304	3277	0.13%	0.017%
82	31.894	4901	4903	4907	rBV5	5618	9450	0.38%	0.048%

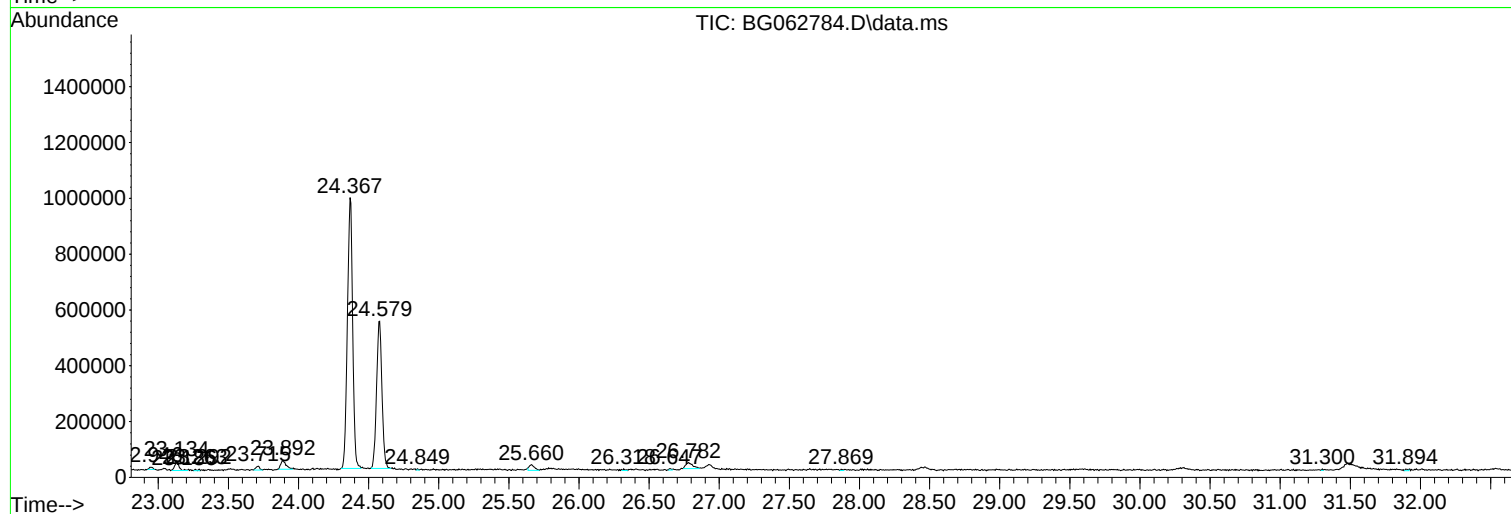
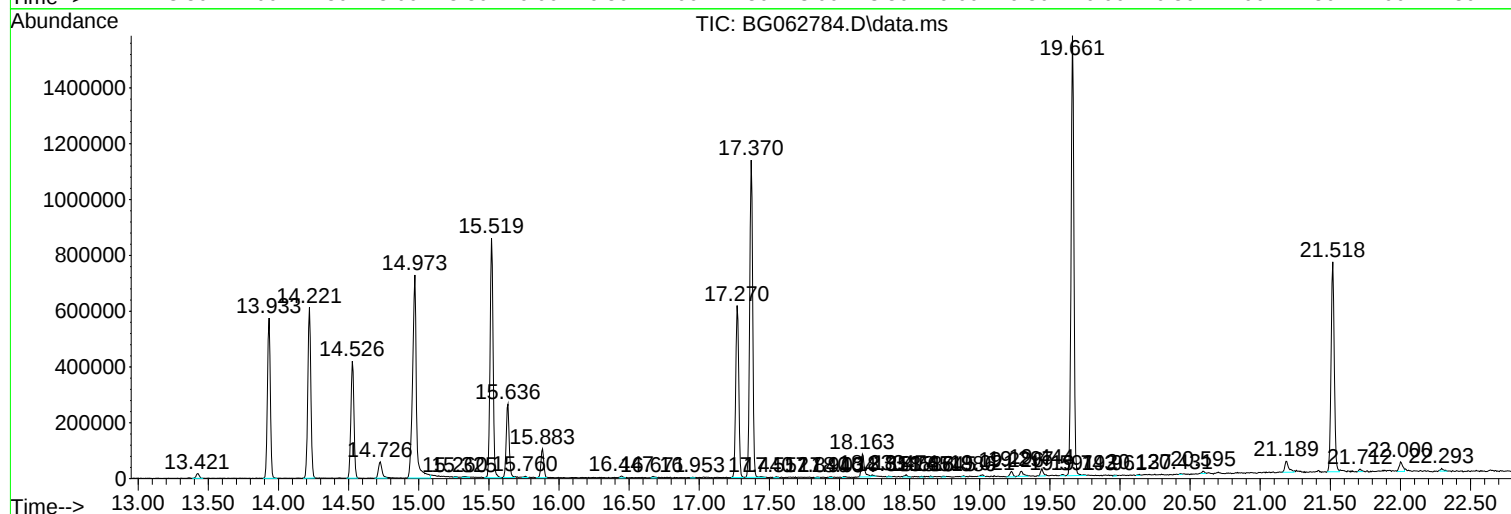
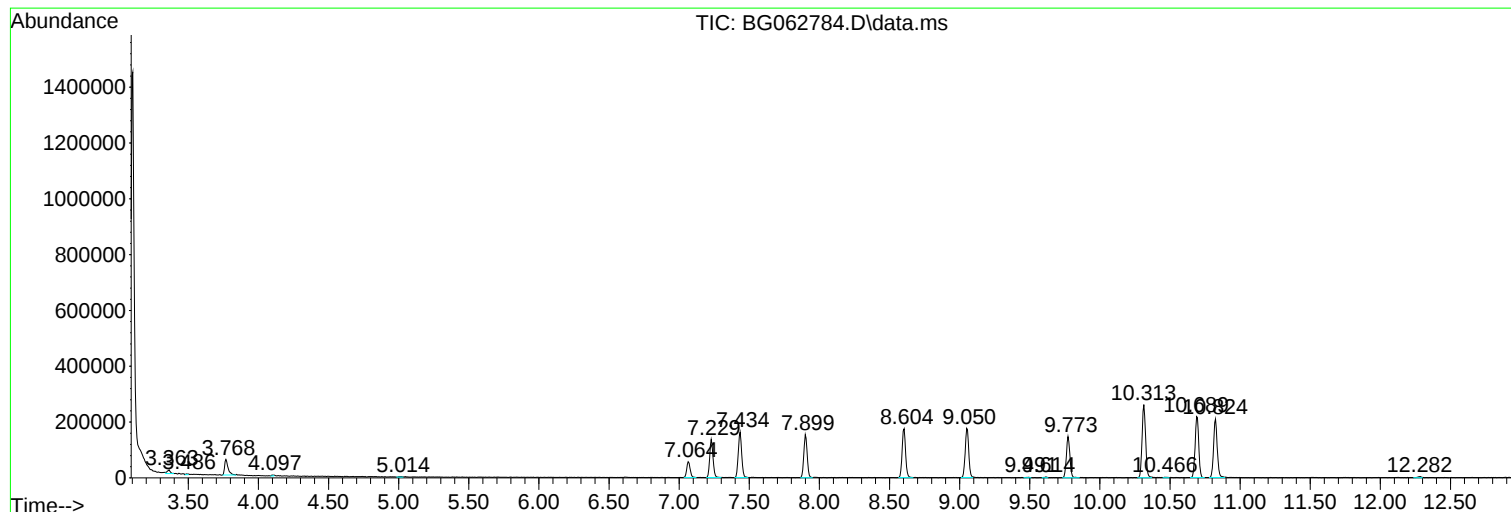
Sum of corrected areas: 19667009

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

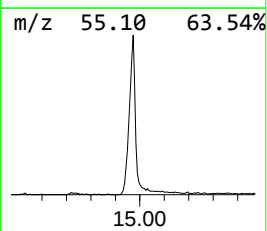
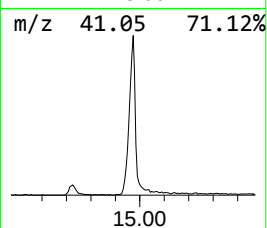
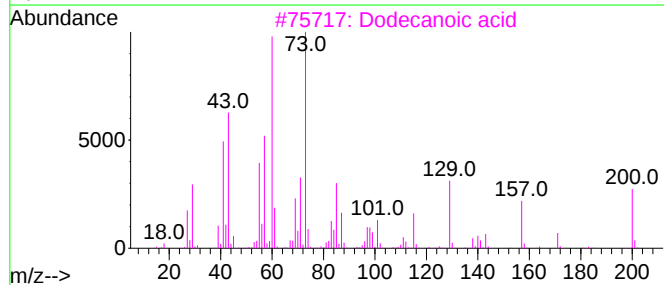
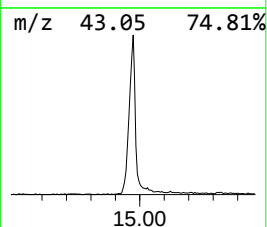
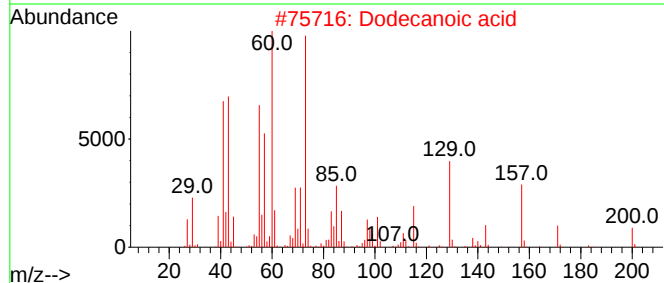
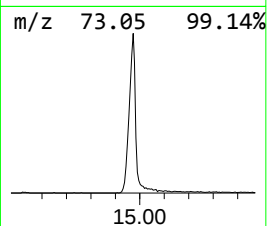
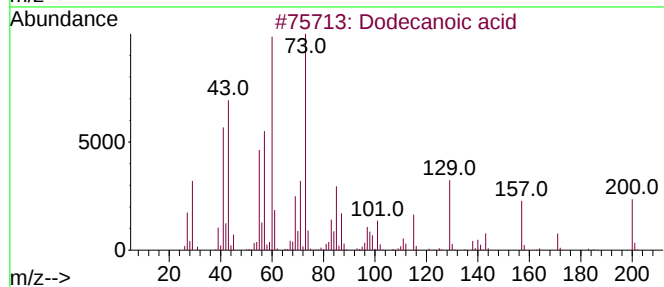
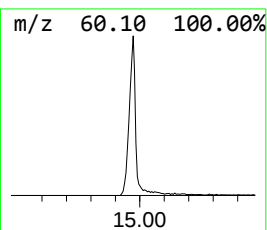
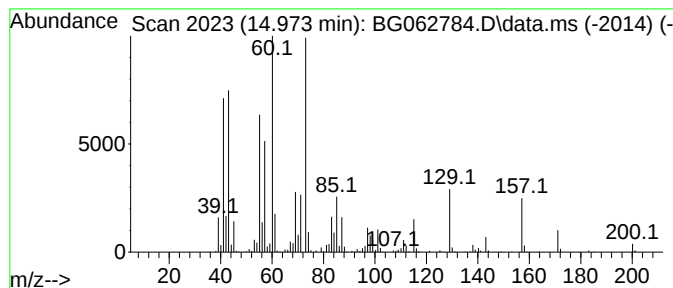
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Dodecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.973	42.43 ng/ul	1376110	Acenaphthene-d10	14.526

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecanoic acid	200	C12H24O2	000143-07-7	91
2			Dodecanoic acid	200	C12H24O2	000143-07-7	91
3			Dodecanoic acid	200	C12H24O2	000143-07-7	87
4			Dodecanoic acid	200	C12H24O2	000143-07-7	81
5			Undecanoic acid	186	C11H22O2	000112-37-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

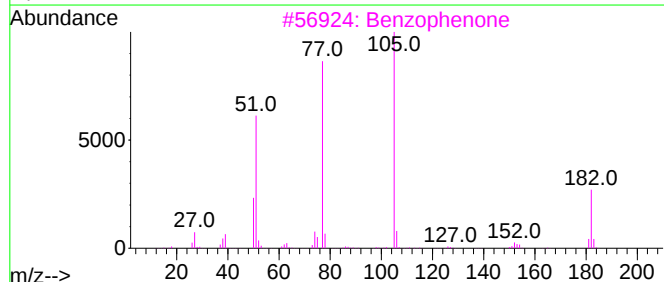
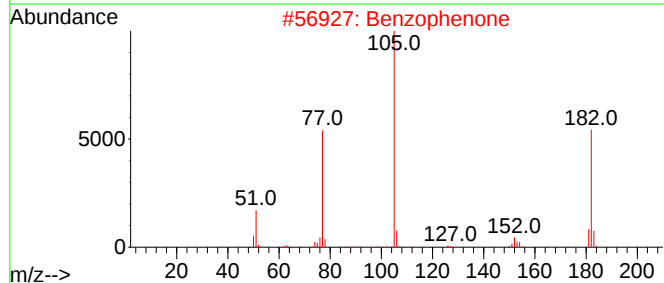
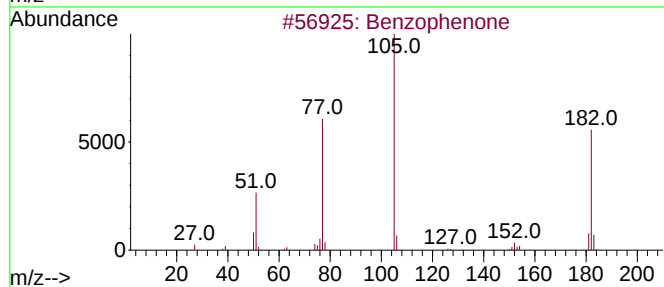
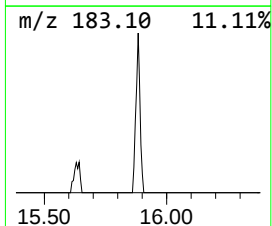
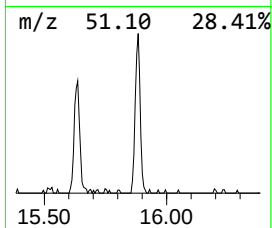
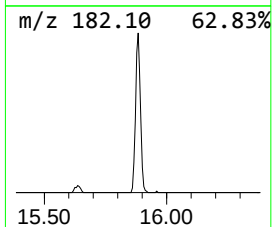
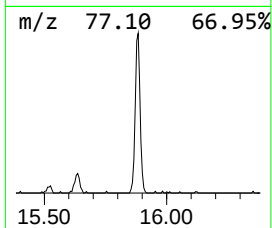
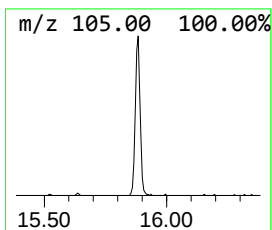
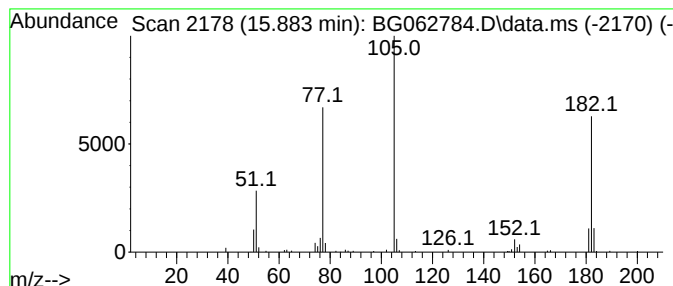
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.883	4.52 ng/ul	146509	Acenaphthene-d10	14.526

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	94
5			Benzophenone	182	C13H10O	000119-61-9	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

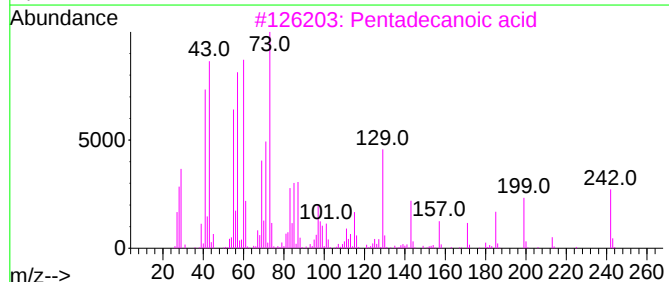
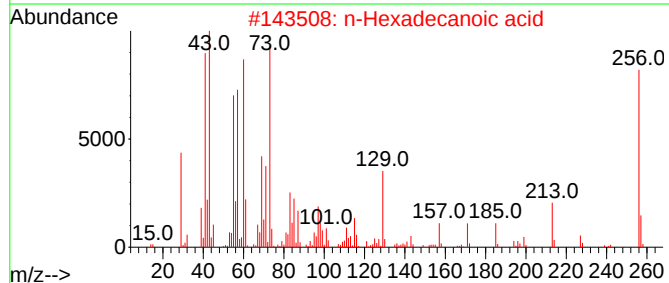
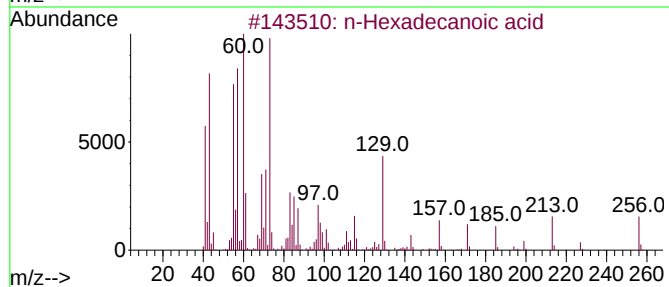
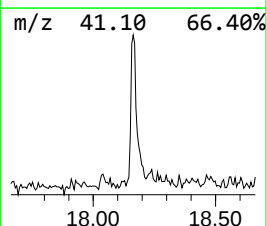
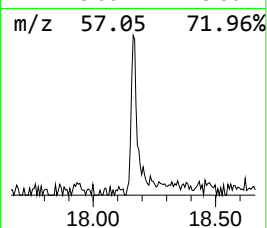
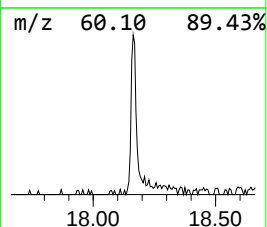
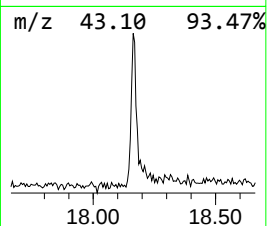
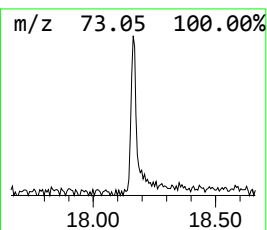
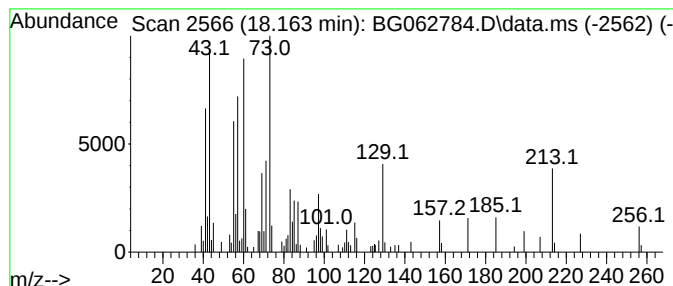
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.163	2.82 ng/ul	129700	Phenanthrene-d10	17.270

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
Data File : BG062784.D
Acq On : 13 Sep 2024 12:26
Operator : JU/RC
Sample : P3952-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0B51

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dodecanoic acid	14.973	42.4	ng/ul	1376110	3	14.526	648618	20.0
Benzophenone	15.883	4.5	ng/ul	146509	3	14.526	648618	20.0
n-Hexadecanoic ...	18.163	2.8	ng/ul	129700	4	17.270	920160	20.0