

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
 Data File : BG062400.D
 Acq On : 14 Aug 2024 18:20
 Operator : MA/JU
 Sample : P3589-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E0CK3

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-BG081424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062400.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.418	19	22	28	rVB3	10464	15531	0.30%	0.076%
2	3.677	63	66	73	rBV	16764	25421	0.50%	0.124%
3	3.824	88	91	98	rBV	40579	65485	1.28%	0.320%
4	6.044	463	469	476	rBV	13155	21053	0.41%	0.103%
5	6.220	495	499	505	rVB3	4719	6592	0.13%	0.032%
6	7.113	644	651	661	rVB2	31430	61627	1.20%	0.301%
7	7.284	674	680	689	rBV	72333	122394	2.39%	0.598%
8	7.489	708	715	721	rBV2	68119	116359	2.27%	0.568%
9	7.548	721	725	737	rVB	27696	50408	0.98%	0.246%
10	7.959	788	795	801	rBV	179720	305980	5.98%	1.494%
11	8.018	801	805	812	rVB2	16223	26859	0.52%	0.131%
12	8.652	907	913	922	rBV	83206	143215	2.80%	0.699%
13	9.105	984	990	998	rBV	79831	141110	2.76%	0.689%
14	9.216	1005	1009	1015	rVV4	5637	8162	0.16%	0.040%
15	9.557	1062	1067	1073	rBV4	7952	13763	0.27%	0.067%
16	9.669	1079	1086	1093	rVB2	37965	57263	1.12%	0.280%
17	9.827	1104	1113	1119	rBV	43838	78085	1.53%	0.381%
18	10.368	1195	1205	1215	rBV	92643	163720	3.20%	0.799%
19	10.526	1226	1232	1236	rBV4	9368	15761	0.31%	0.077%
20	10.750	1263	1270	1278	rBV	280708	485884	9.49%	2.372%
21	10.879	1284	1292	1299	rBV	77080	143860	2.81%	0.702%
22	11.284	1352	1361	1366	rVB3	12253	21190	0.41%	0.103%
23	11.390	1374	1379	1384	rBV2	4030	5939	0.12%	0.029%
24	11.484	1387	1395	1408	rBV	123811	235718	4.60%	1.151%
25	13.417	1718	1724	1731	rVV	350123	531526	10.38%	2.595%
26	13.481	1731	1735	1742	rVV3	8290	14954	0.29%	0.073%
27	13.857	1795	1799	1808	rBV6	4010	7587	0.15%	0.037%
28	13.981	1814	1820	1827	rVB	252961	362114	7.07%	1.768%
29	14.227	1857	1862	1864	rBV2	23534	35916	0.70%	0.175%
30	14.268	1864	1869	1879	rVV2	247051	401768	7.85%	1.962%
31	14.486	1900	1906	1910	rVB2	9860	11455	0.22%	0.056%
32	14.574	1915	1921	1928	rVB	516279	820758	16.03%	4.007%
33	14.809	1948	1961	1975	rVV	2199173	5119576	100.00%	24.996%
34	14.909	1977	1978	1988	rVB6	4348	8074	0.16%	0.039%
35	14.985	1988	1991	1998	rBV6	8555	14744	0.29%	0.072%
36	15.167	2018	2022	2026	rBV3	4226	5892	0.12%	0.029%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
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37	15.308	2043	2046	2050	rBV3	5394	7981	0.16%	0.039%
38	15.379	2053	2058	2062	rVB3	8746	13328	0.26%	0.065%
39	15.431	2064	2067	2071	rBV	11359	14865	0.29%	0.073%
40	15.567	2081	2090	2099	rBV	375508	560203	10.94%	2.735%

41	15.672	2103	2108	2115	rBV	48214	74032	1.45%	0.361%
42	15.737	2115	2119	2123	rVB	7384	10934	0.21%	0.053%
43	15.843	2133	2137	2141	rVB4	4340	7051	0.14%	0.034%
44	15.943	2149	2154	2156	rBV4	3838	5687	0.11%	0.028%
45	16.007	2159	2165	2168	rVB4	4359	6974	0.14%	0.034%

46	16.119	2179	2184	2191	rBV2	20852	36024	0.70%	0.176%
47	16.195	2193	2197	2202	rVB7	4548	9104	0.18%	0.044%
48	16.248	2202	2206	2210	rBV4	5299	7505	0.15%	0.037%
49	16.295	2210	2214	2216	rBV	19964	21744	0.42%	0.106%
50	16.324	2216	2219	2224	rVB4	9499	13330	0.26%	0.065%

51	16.459	2238	2242	2248	rVB5	8970	14368	0.28%	0.070%
52	16.718	2281	2286	2290	rBV5	12197	20145	0.39%	0.098%
53	16.812	2297	2302	2308	rBV6	4545	9303	0.18%	0.045%
54	16.877	2308	2313	2317	rBV3	11043	15926	0.31%	0.078%
55	17.035	2337	2340	2344	rVB3	5728	7208	0.14%	0.035%

56	17.088	2344	2349	2352	rBV	17849	24181	0.47%	0.118%
57	17.129	2352	2356	2364	rVB7	9011	20961	0.41%	0.102%
58	17.317	2381	2388	2393	rVV	772478	1138638	22.24%	5.559%
59	17.358	2393	2395	2399	rVV	20744	24113	0.47%	0.118%
60	17.411	2399	2404	2412	rVB	452391	683291	13.35%	3.336%

61	17.493	2414	2418	2424	rVB7	10813	17734	0.35%	0.087%
62	17.705	2450	2454	2459	rBV6	12914	20998	0.41%	0.103%
63	17.822	2471	2474	2477	rBV3	14689	18300	0.36%	0.089%
64	17.993	2499	2503	2511	rVB	52000	73082	1.43%	0.357%
65	18.081	2514	2518	2525	rVV4	14328	20414	0.40%	0.100%

66	18.216	2535	2541	2550	rBV	397531	576029	11.25%	2.812%
67	18.516	2586	2592	2596	rBV2	18944	29978	0.59%	0.146%
68	18.768	2631	2635	2643	rVB2	17283	28883	0.56%	0.141%
69	18.880	2650	2654	2659	rBV7	6844	10516	0.21%	0.051%
70	18.950	2661	2666	2670	rBV7	6449	11823	0.23%	0.058%

71	19.015	2673	2677	2683	rVB8	8116	15315	0.30%	0.075%
72	19.079	2683	2688	2692	rBV5	17161	33904	0.66%	0.166%
73	19.162	2692	2702	2707	rBV3	193099	282788	5.52%	1.381%
74	19.262	2716	2719	2721	rBV3	9092	12540	0.24%	0.061%
75	19.297	2721	2725	2728	rVB	37415	41040	0.80%	0.200%

76	19.338	2728	2732	2734	rBV4	30585	38471	0.75%	0.188%
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77	19.367	2734	2737	2739	rVV4	11025	13057	0.26%	0.064%
78	19.485	2753	2757	2765	rBV	129136	177324	3.46%	0.866%
79	19.638	2778	2783	2787	rVB4	20305	29739	0.58%	0.145%
80	19.696	2787	2793	2800	rVV	641790	949096	18.54%	4.634%
81	19.755	2800	2803	2808	rVB2	42073	52115	1.02%	0.254%
82	19.831	2813	2816	2819	rVB5	9302	9859	0.19%	0.048%
83	20.219	2878	2882	2884	rBV4	16877	23699	0.46%	0.116%
84	20.248	2884	2887	2891	rVB	77250	85538	1.67%	0.418%
85	20.601	2944	2947	2953	rVB5	24981	50832	0.99%	0.248%
86	20.677	2957	2960	2964	rBV6	9676	11816	0.23%	0.058%
87	20.742	2964	2971	2975	rBV2	90640	139423	2.72%	0.681%
88	20.777	2975	2977	2980	rVB2	14038	10426	0.20%	0.051%
89	20.854	2988	2990	2992	rVB2	9448	8001	0.16%	0.039%
90	21.059	3021	3025	3028	rBV6	8117	14994	0.29%	0.073%
91	21.230	3049	3054	3067	rBV2	449039	768857	15.02%	3.754%
92	21.558	3103	3110	3117	rBV	850812	1354113	26.45%	6.611%
93	21.776	3142	3147	3151	rBV	86026	128295	2.51%	0.626%
94	22.363	3243	3247	3256	rVB2	80145	143643	2.81%	0.701%
95	23.033	3356	3361	3367	rBV	59400	104919	2.05%	0.512%
96	23.215	3387	3392	3399	rVB	135645	246500	4.81%	1.203%
97	23.808	3488	3493	3500	rBV	54327	114261	2.23%	0.558%
98	24.431	3590	3599	3614	rVB2	319044	874331	17.08%	4.269%
99	24.648	3626	3636	3644	rBV	521549	1420953	27.76%	6.938%
100	25.917	3847	3852	3862	rBV5	35242	105709	2.06%	0.516%

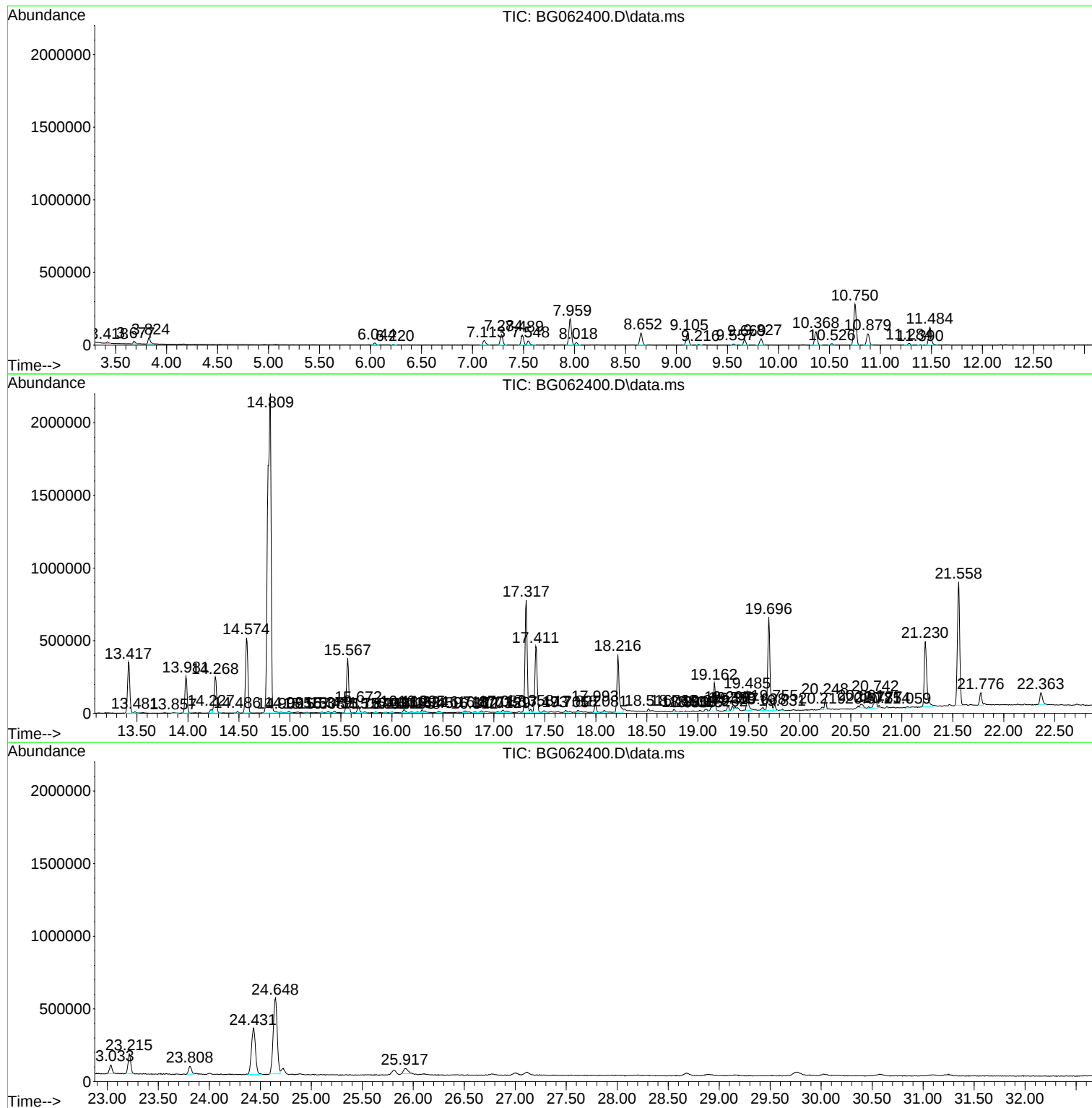
Sum of corrected areas: 20481954

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.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\BG081424\
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ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E0CK3

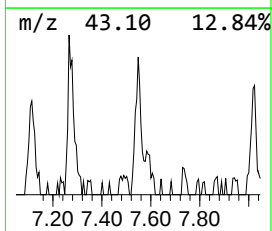
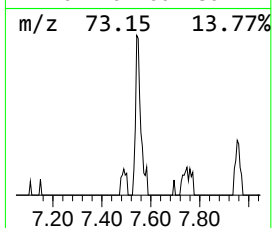
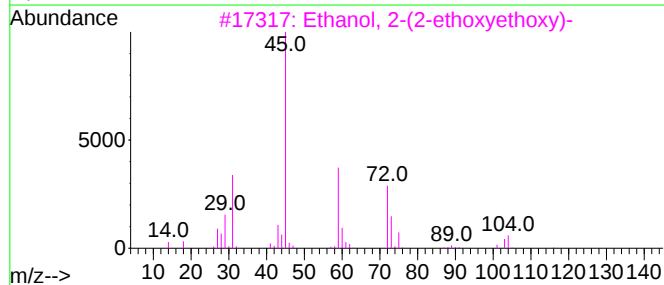
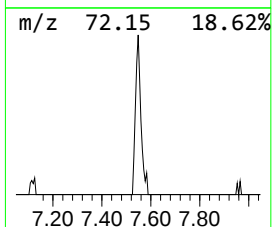
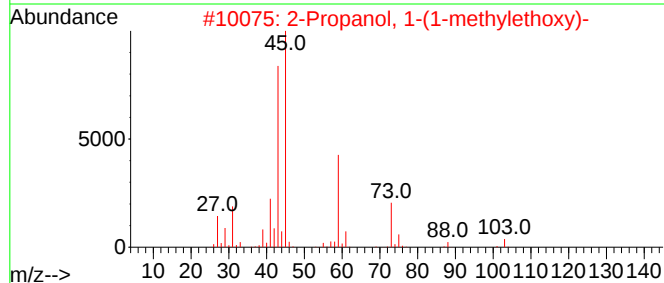
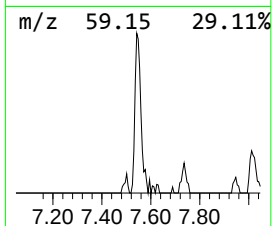
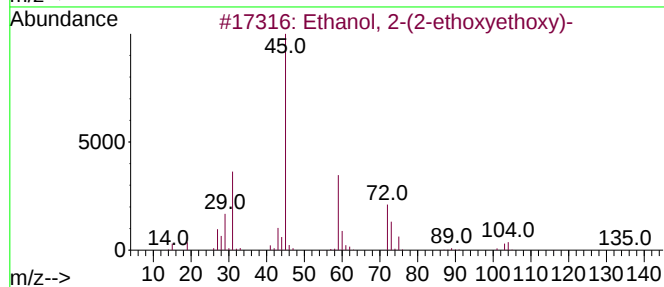
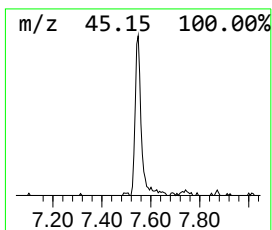
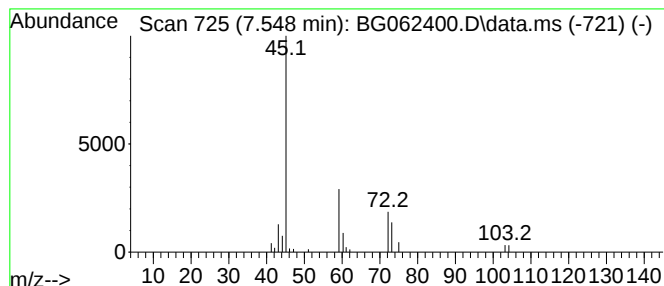
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.548	3.29 ng/ul	50408	1,4-Dichlorobenzene-d4	7.959

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	83
2			2-Propanol, 1-(1-methylethoxy)-	118	C6H14O2	003944-36-3	56
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	53
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	53
5			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	50



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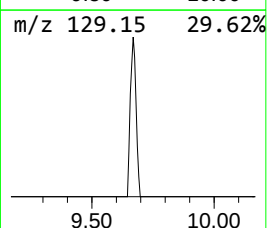
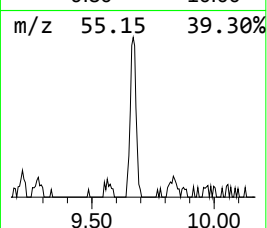
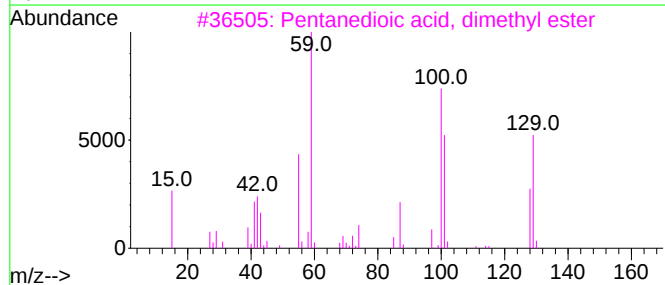
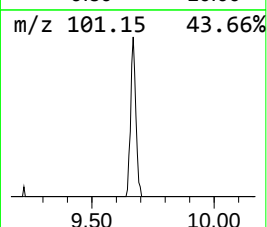
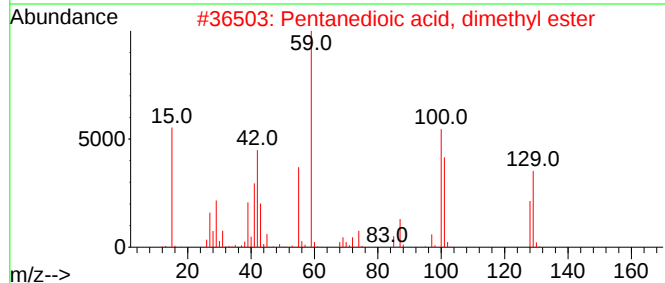
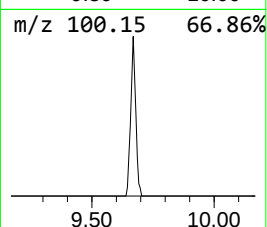
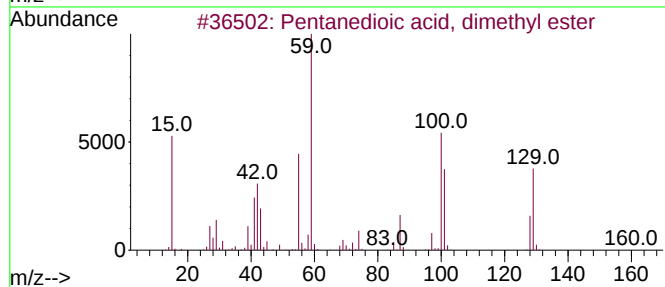
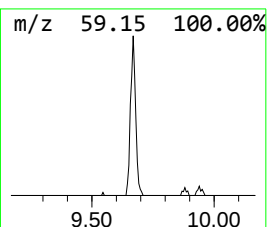
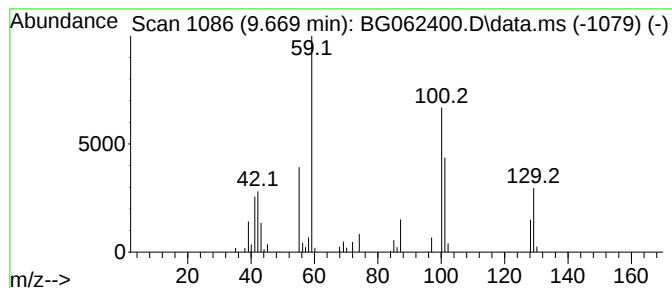
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Pentanedioic acid, dimethyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.669	2.36 ng/ul	57263	Naphthalene-d8	10.750

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	90
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
3			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	72
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	59
5			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	50



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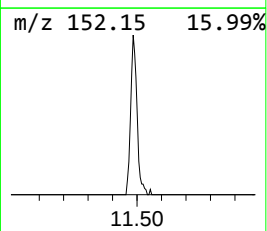
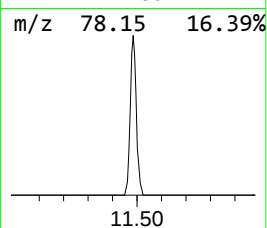
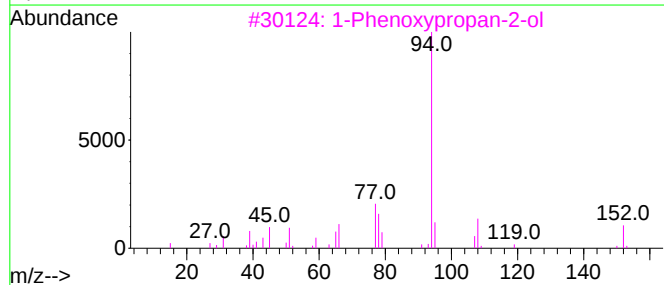
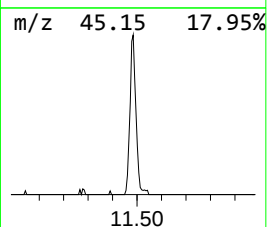
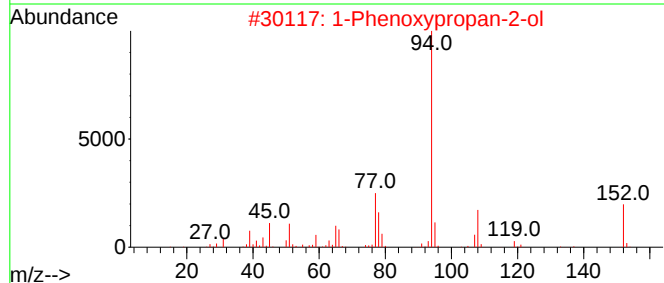
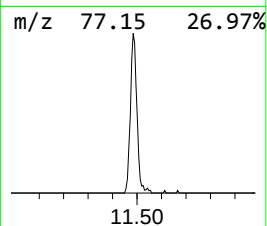
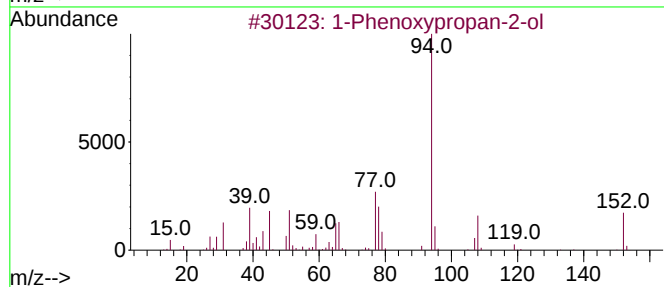
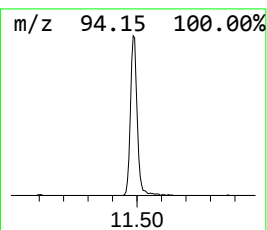
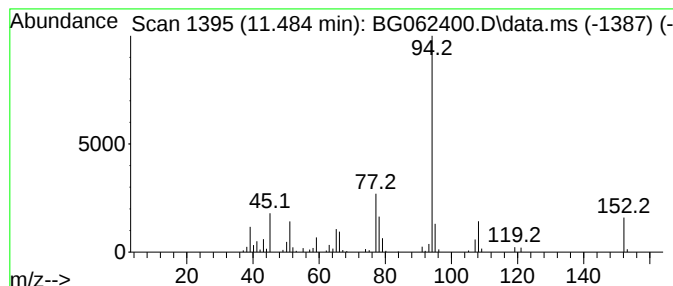
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.484	9.70 ng/ul	235718	Naphthalene-d8	10.750

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	96	
2	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	95	
3	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	94	
4	1-Phenoxypropan-2-ol		152	C9H12O2	000770-35-4	93	
5	1-Propanol, 3-phenoxy-		152	C9H12O2	006180-61-6	87	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
Operator : MA/JU
Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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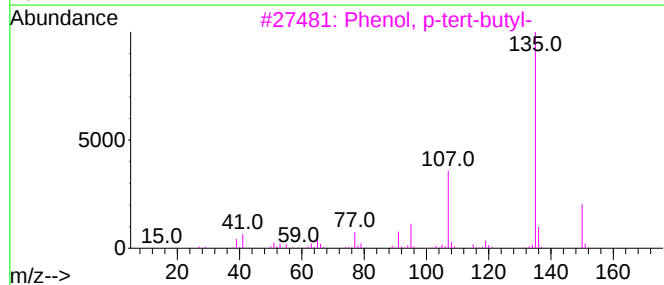
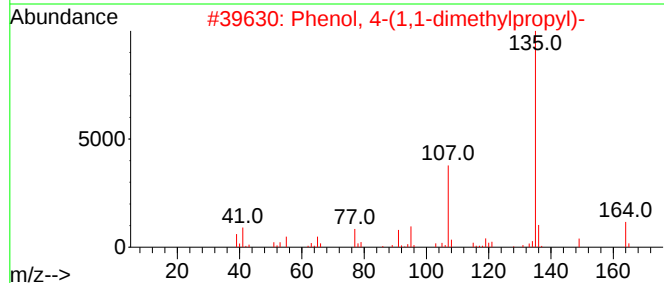
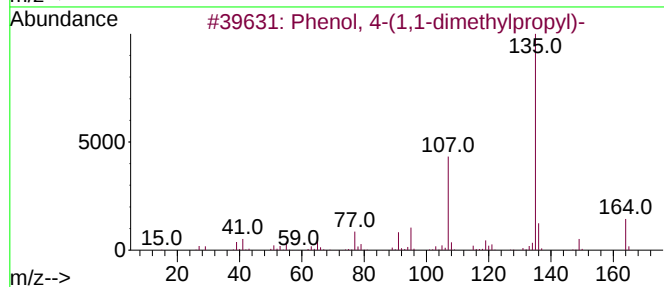
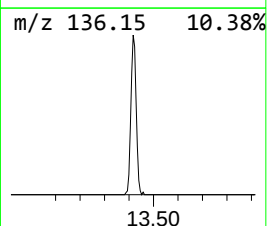
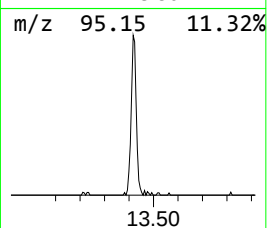
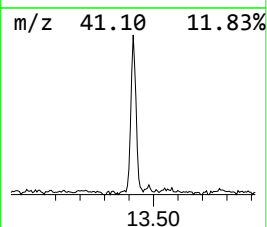
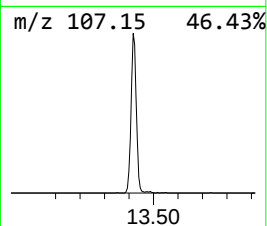
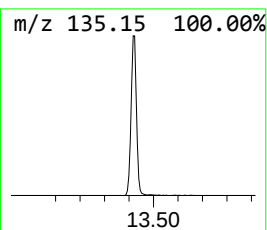
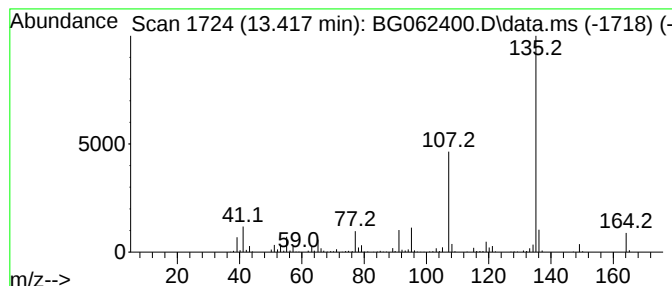
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 4-(1,1-dimethylprop... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.417	12.95 ng/ul	531526	Acenaphthene-d10	14.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	95
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	95
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	86
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	81
5			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
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Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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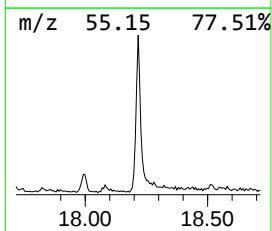
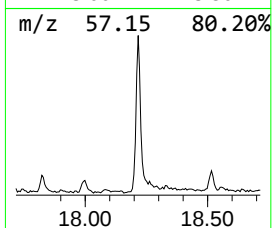
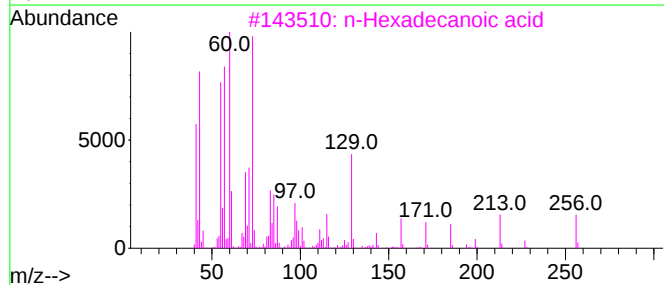
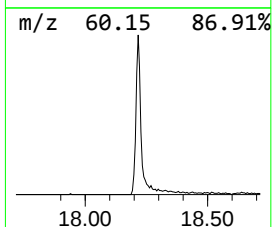
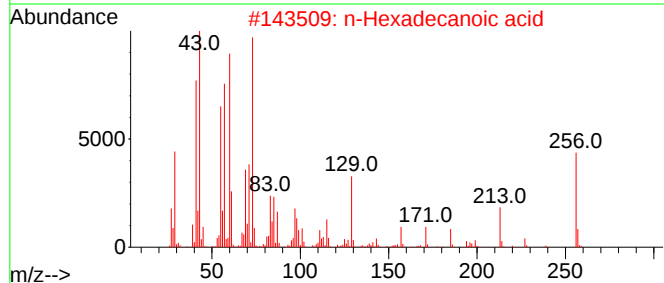
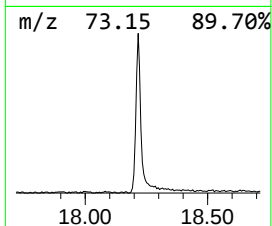
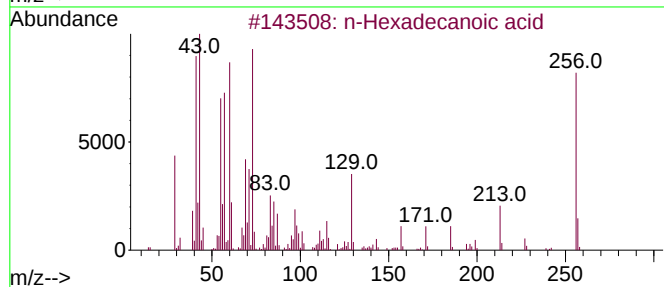
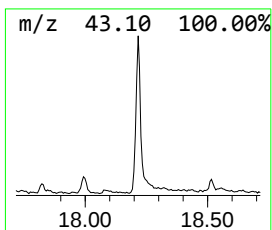
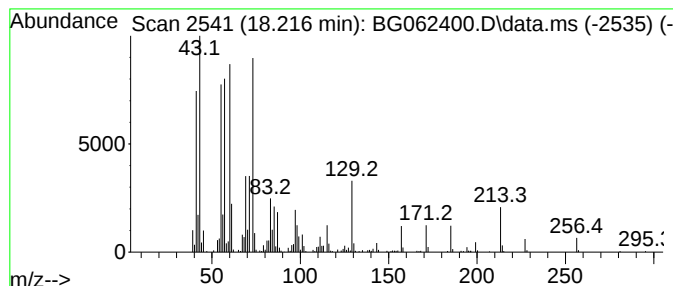
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	10.12 ng/ul	576029	Phenanthrene-d10	17.317

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	94
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
Operator : MA/JU
Sample : P3589-03
Misc :
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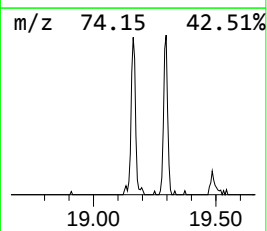
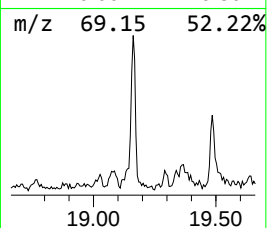
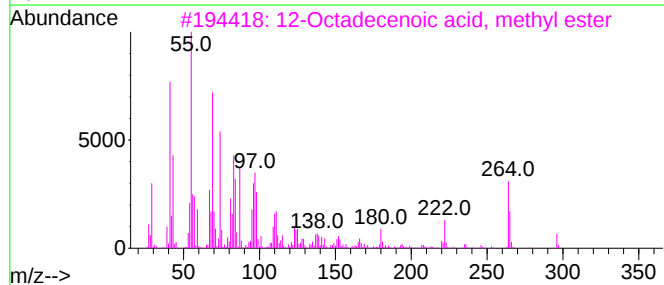
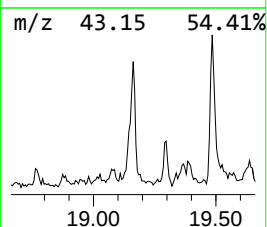
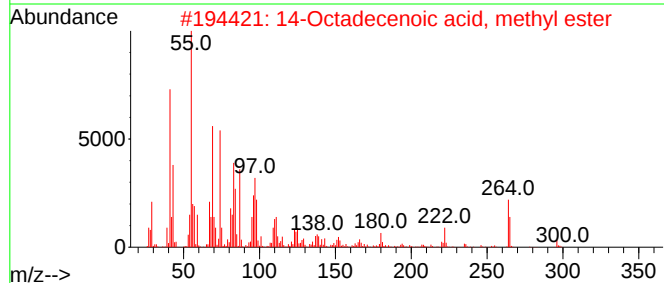
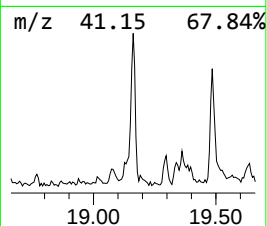
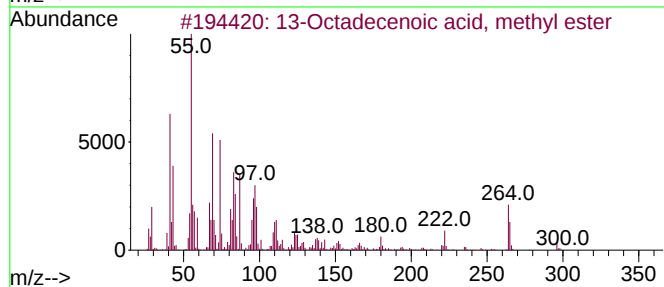
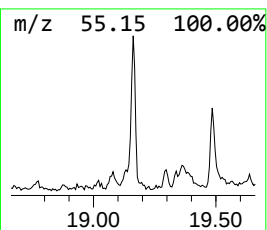
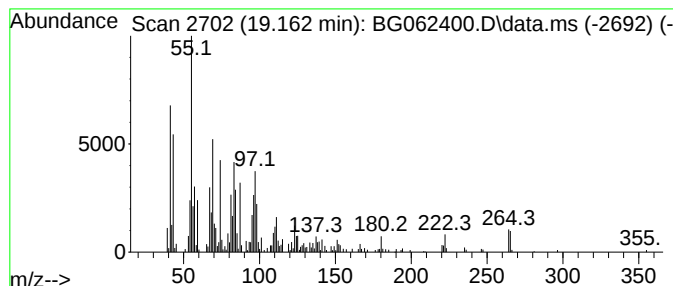
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 13-Octadecenoic acid, methyl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.162	4.97 ng/ul	282788	Phenanthrene-d10	17.317

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	98
2			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	96
3			12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	94
4			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	94
5			9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
Operator : MA/JU
Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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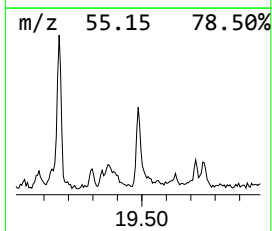
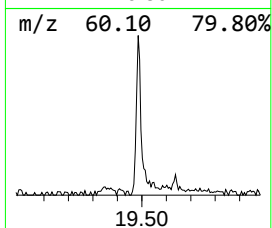
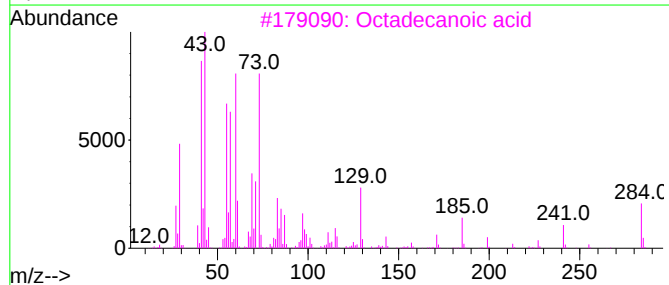
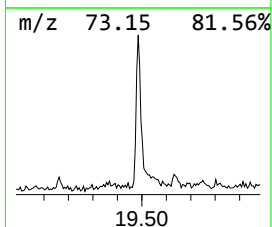
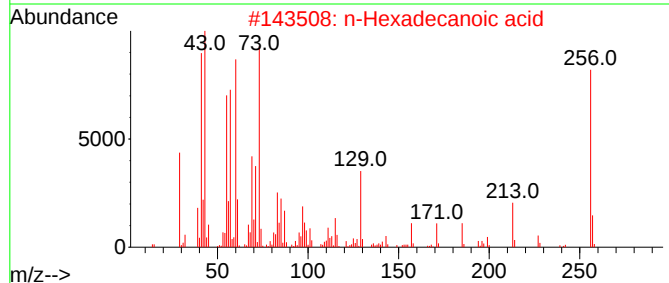
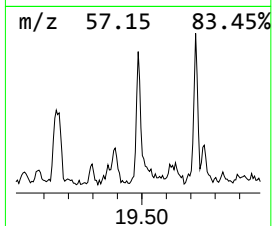
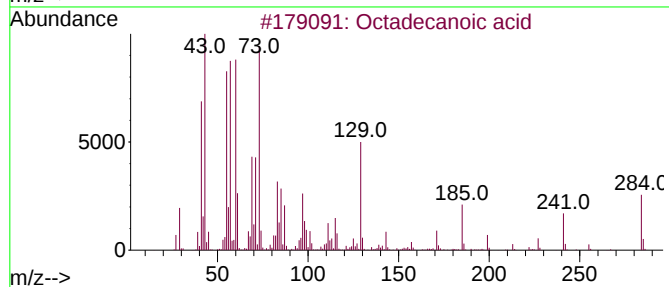
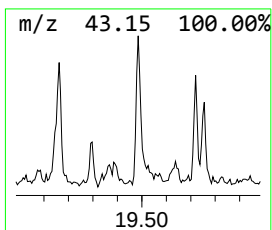
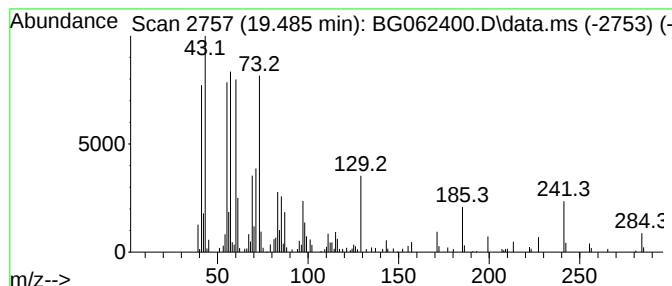
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.485	2.62 ng/ul	177324	Chrysene-d12	21.558

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	94
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
3			Octadecanoic acid	284	C18H36O2	000057-11-4	91
4			Octadecanoic acid	284	C18H36O2	000057-11-4	90
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
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Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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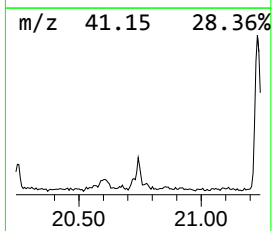
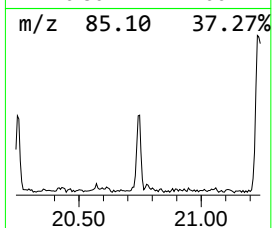
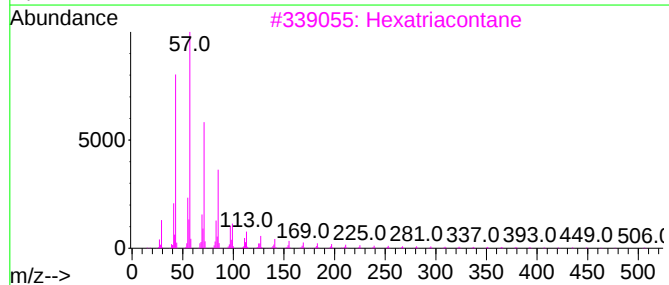
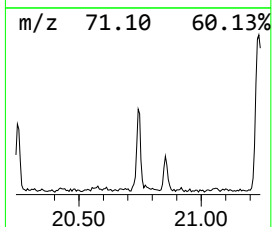
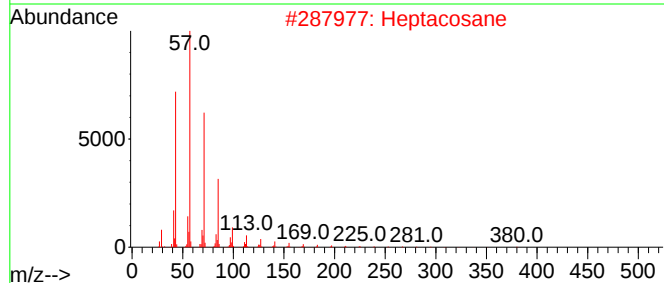
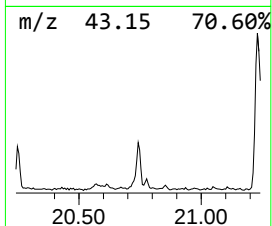
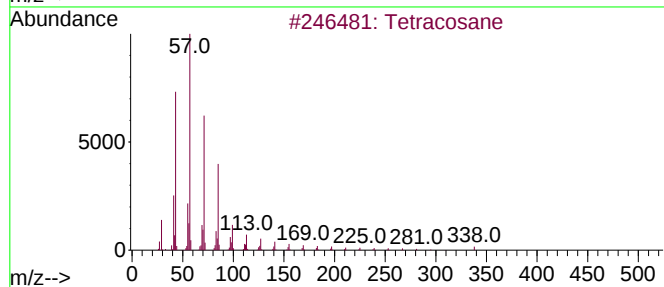
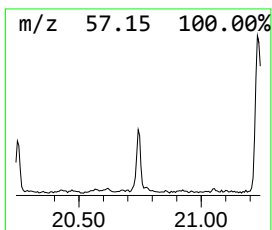
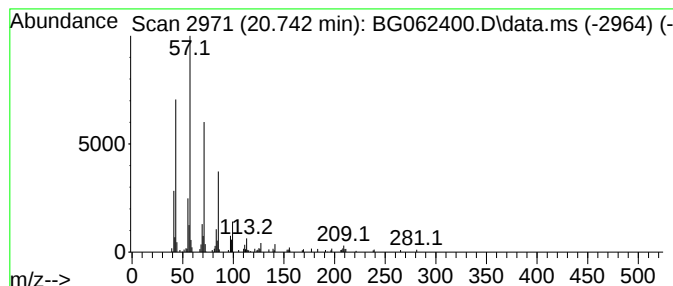
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.742	2.06 ng/ul	139423	Chrysene-d12	21.558

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	90
2		Heptacosane	380	C27H56	000593-49-7	90
3		Hexatriacontane	507	C36H74	000630-06-8	90
4		Octacosane	394	C28H58	000630-02-4	87
5		Eicosane	282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
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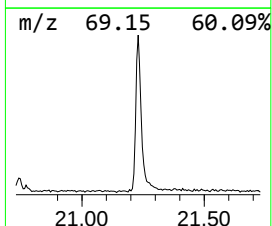
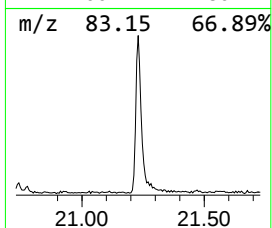
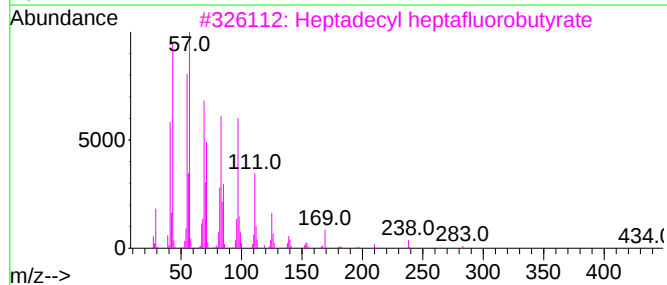
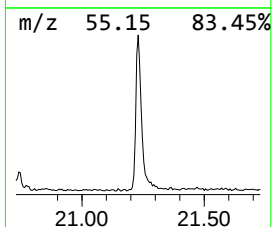
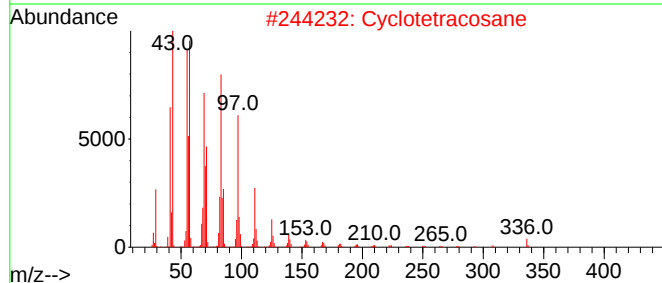
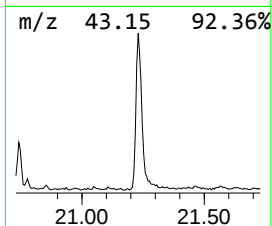
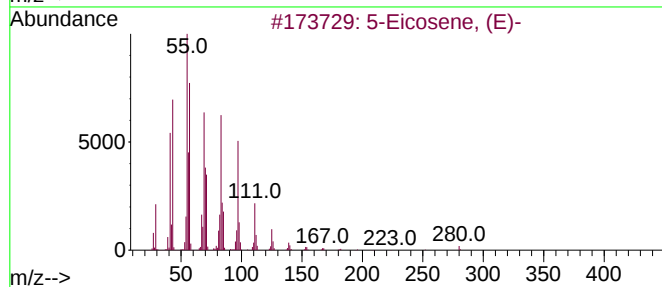
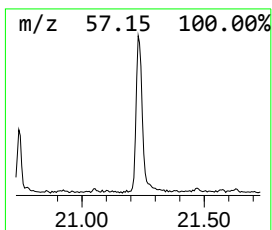
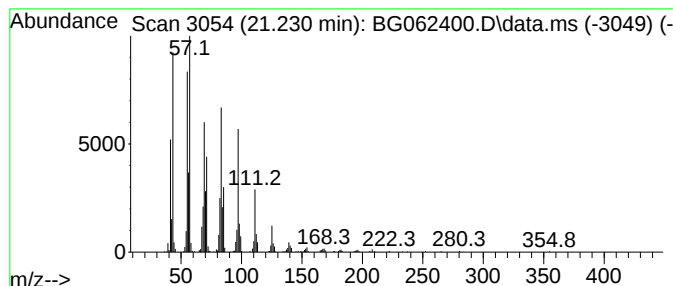
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.230	11.36 ng/ul	768857	Chrysene-d12		21.558	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	95
2	Cyclotetracosane		336	C24H48	000297-03-0	95
3	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94
4	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	91
5	n-Tetracosanol-1		354	C24H50O	000506-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
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Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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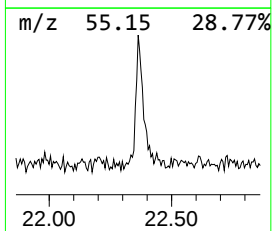
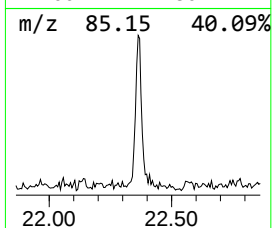
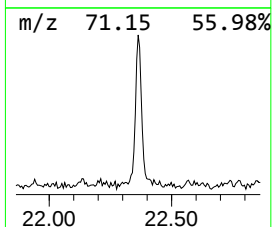
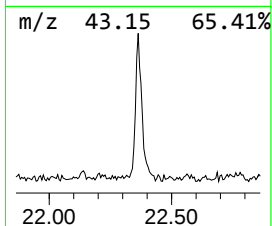
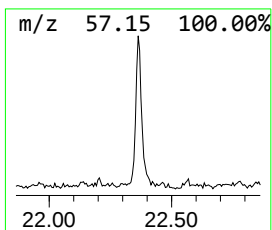
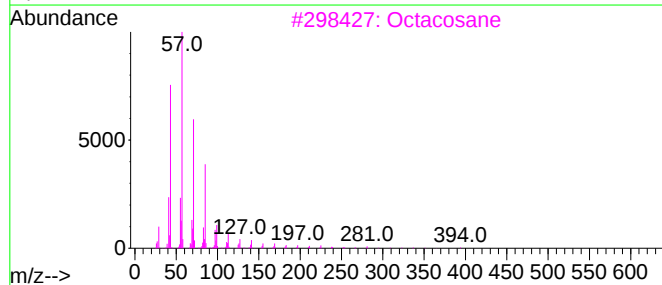
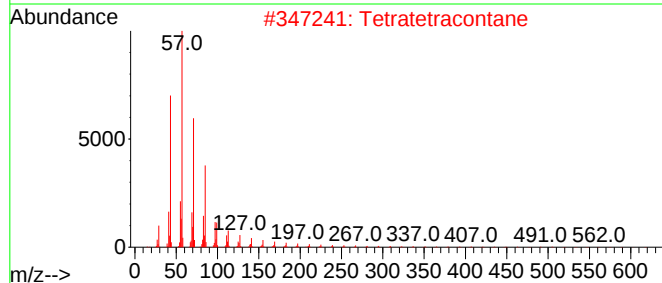
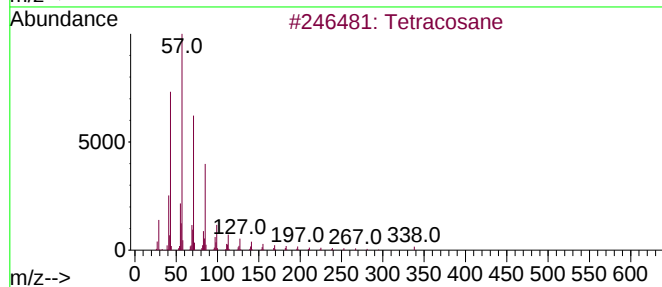
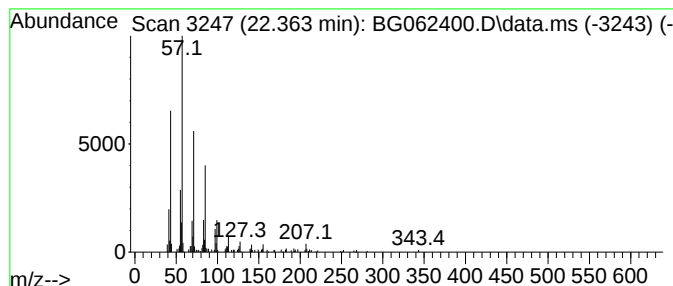
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.363	2.12 ng/ul	143643	Chrysene-d12	21.558

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	87
2		Tetratetracontane	619	C44H90	007098-22-8	87
3		Octacosane	394	C28H58	000630-02-4	87
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	83
5		Heptacosane	380	C27H56	000593-49-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
Operator : MA/JU
Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E0CK3

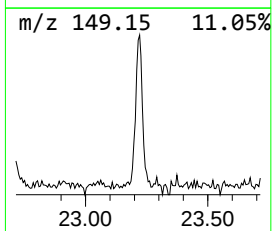
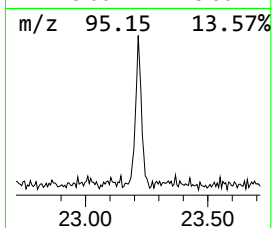
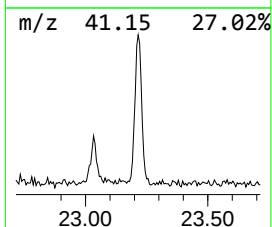
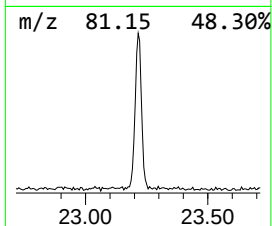
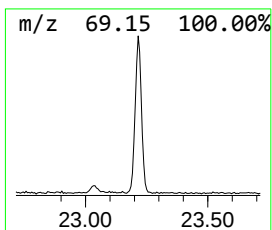
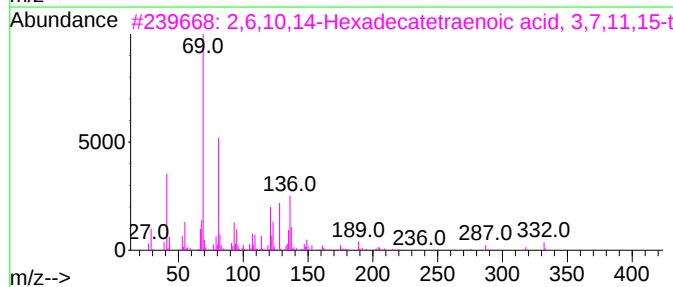
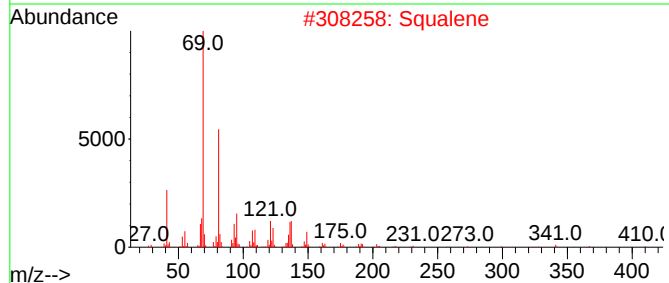
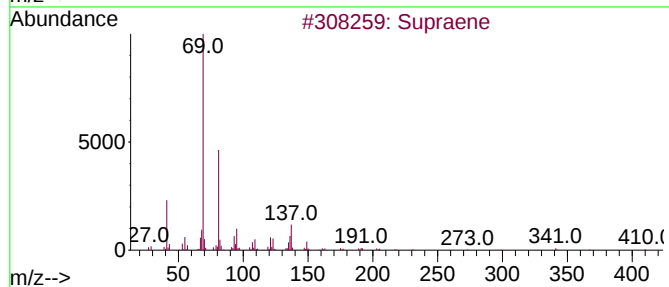
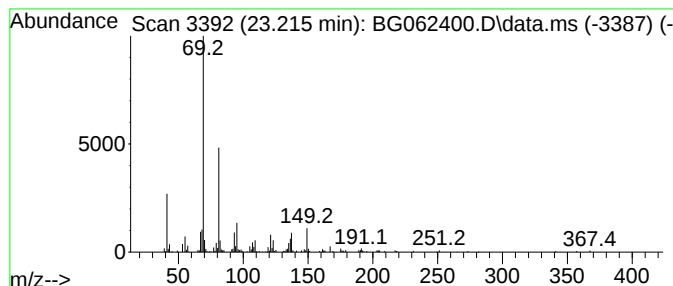
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.215	3.47 ng/ul	246500	Perylene-d12	24.648

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	90
3			2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	72
4			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	64
5			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
Data File : BG062400.D
Acq On : 14 Aug 2024 18:20
Operator : MA/JU
Sample : P3589-03
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
E0CK3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.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
Ethanol, 2-(2-e...	7.548	3.3	ng/ul	50408	1	7.959	305980	20.0
Pentanedioic ac...	9.669	2.4	ng/ul	57263	2	10.750	485884	20.0
1-Phenoxypropan...	11.484	9.7	ng/ul	235718	2	10.750	485884	20.0
Phenol, 4-(1,1-...	13.417	12.9	ng/ul	531526	3	14.574	820758	20.0
n-Hexadecanoic ...	18.216	10.1	ng/ul	576029	4	17.317	1138640	20.0
13-Octadecenoic...	19.162	5.0	ng/ul	282788	4	17.317	1138640	20.0
Octadecanoic acid	19.485	2.6	ng/ul	177324	5	21.558	1354110	20.0
(DEL) Alkane: S...	20.742	2.1	ng/ul	139423	5	21.558	1354110	20.0
(DEL) Alkane: S...	21.230	11.4	ng/ul	768857	5	21.558	1354110	20.0
(DEL) Alkane: S...	22.363	2.1	ng/ul	143643	5	21.558	1354110	20.0
Supraene	23.215	3.5	ng/ul	246500	6	24.648	1420950	20.0