

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
 Data File : BG058298.D
 Acq On : 18 Jul 2023 12:54
 Operator : CG/JU
 Sample : 03444-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A46Q1

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-BG071723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG058298.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.342	40	43	48	rVB3	14006	19839	1.20%	0.099%
2	3.747	108	112	128	rBV	172240	307858	18.69%	1.538%
3	3.865	129	132	137	rVB4	11084	15222	0.92%	0.076%
4	3.976	149	151	159	rVB5	3340	5987	0.36%	0.030%
5	4.088	166	170	175	rVB2	5063	8053	0.49%	0.040%
6	4.305	203	207	214	rVB2	10809	15096	0.92%	0.075%
7	4.693	268	273	274	rVV4	2043	2770	0.17%	0.014%
8	4.717	274	277	280	rVB4	3634	5115	0.31%	0.026%
9	5.087	334	340	344	rBV3	2988	5348	0.32%	0.027%
10	7.061	669	676	688	rBV	213624	380105	23.08%	1.898%
11	7.225	697	704	714	rBV	175541	303107	18.40%	1.514%
12	7.431	732	739	749	rVV	205693	357043	21.68%	1.783%
13	7.901	811	819	826	rBV	140586	250506	15.21%	1.251%
14	8.606	930	939	953	rBV	243345	461240	28.01%	2.304%
15	9.058	1009	1016	1028	rBV	208103	379703	23.05%	1.896%
16	9.211	1037	1042	1046	rBV6	2154	3877	0.24%	0.019%
17	9.241	1046	1047	1053	rVB5	2144	3805	0.23%	0.019%
18	9.781	1132	1139	1151	rBV	169785	315276	19.14%	1.575%
19	9.875	1151	1155	1158	rVB4	2238	3580	0.22%	0.018%
20	10.322	1224	1231	1241	rBV	252096	466801	28.34%	2.331%
21	10.698	1287	1295	1303	rVV	219376	401970	24.41%	2.008%
22	10.833	1311	1318	1336	rBV	218022	437388	26.56%	2.184%
23	10.986	1341	1344	1349	rVB4	1953	2851	0.17%	0.014%
24	11.185	1372	1378	1382	rBV6	3357	6326	0.38%	0.032%
25	11.397	1411	1414	1420	rVB4	3838	5186	0.31%	0.026%
26	11.614	1448	1451	1456	rVB5	1710	3058	0.19%	0.015%
27	11.714	1463	1468	1472	rBV2	11836	18473	1.12%	0.092%
28	11.749	1472	1474	1478	rVV5	3334	4236	0.26%	0.021%
29	11.984	1510	1514	1516	rVV4	1713	2658	0.16%	0.013%
30	12.102	1529	1534	1536	rBV5	1995	2896	0.18%	0.014%
31	12.290	1563	1566	1572	rVB6	3559	6440	0.39%	0.032%
32	12.748	1639	1644	1646	rBV5	1635	2690	0.16%	0.013%
33	12.813	1650	1655	1663	rBV6	4041	8901	0.54%	0.044%
34	13.066	1693	1698	1703	rVV2	21443	32167	1.95%	0.161%
35	13.136	1703	1710	1713	rVB5	3772	6195	0.38%	0.031%
36	13.336	1737	1744	1749	rBV9	5054	11017	0.67%	0.055%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
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37	13.371	1749	1750	1755	rVV4	3358	3753	0.23%	0.019%
38	13.424	1755	1759	1764	rVV6	6575	12308	0.75%	0.061%
39	13.583	1781	1786	1790	rBV7	3458	6630	0.40%	0.033%
40	13.641	1793	1796	1799	rBV4	2563	4274	0.26%	0.021%
41	13.694	1802	1805	1809	rBV6	2236	3480	0.21%	0.017%
42	13.859	1828	1833	1838	rVB7	4060	7891	0.48%	0.039%
43	13.941	1838	1847	1854	rBV	518917	770637	46.79%	3.849%
44	14.006	1854	1858	1863	rVV	29231	42877	2.60%	0.214%
45	14.064	1863	1868	1873	rVV7	4784	8347	0.51%	0.042%
46	14.182	1883	1888	1889	rBV4	3400	5184	0.31%	0.026%
47	14.229	1889	1896	1906	rVV	586921	947891	57.55%	4.734%
48	14.364	1914	1919	1925	rVB5	9699	16531	1.00%	0.083%
49	14.429	1925	1930	1937	rBV7	4855	10391	0.63%	0.052%
50	14.534	1942	1948	1955	rVB	381813	609593	37.01%	3.044%
51	14.652	1961	1968	1975	rVB	446662	675687	41.03%	3.375%
52	14.734	1976	1982	1997	rBV	172188	334257	20.30%	1.669%
53	14.893	2005	2009	2011	rBV4	3259	4665	0.28%	0.023%
54	14.928	2011	2015	2021	rVV7	10127	17906	1.09%	0.089%
55	15.046	2031	2035	2041	rVB5	10970	17757	1.08%	0.089%
56	15.157	2050	2054	2057	rBV5	6176	10136	0.62%	0.051%
57	15.251	2065	2070	2076	rBV8	6692	16734	1.02%	0.084%
58	15.316	2078	2081	2086	rVV6	9649	15136	0.92%	0.076%
59	15.422	2096	2099	2100	rBV2	5293	4679	0.28%	0.023%
60	15.469	2103	2107	2111	rBV	22863	29330	1.78%	0.146%
61	15.527	2111	2117	2124	rBV	811628	1212872	73.64%	6.057%
62	15.645	2131	2137	2150	rVB	246262	383311	23.27%	1.914%
63	15.786	2150	2161	2169	rBV	679436	993659	60.33%	4.963%
64	15.886	2173	2178	2184	rVB	119754	179170	10.88%	0.895%
65	16.027	2201	2202	2204	rBV2	5218	3883	0.24%	0.019%
66	16.168	2224	2226	2228	rBV3	4576	4102	0.25%	0.020%
67	16.221	2232	2235	2236	rBV2	8755	8835	0.54%	0.044%
68	16.256	2236	2241	2248	rVB	72631	106394	6.46%	0.531%
69	16.391	2260	2264	2269	rVB	31883	43601	2.65%	0.218%
70	16.450	2271	2274	2278	rVV3	12020	16174	0.98%	0.081%
71	16.509	2279	2284	2292	rVB	178738	259599	15.76%	1.296%
72	16.808	2330	2335	2341	rBV	97830	151139	9.18%	0.755%
73	17.079	2376	2381	2390	rVB2	44609	74078	4.50%	0.370%
74	17.155	2390	2394	2396	rBV3	21807	28385	1.72%	0.142%
75	17.190	2396	2400	2406	rVB2	52842	78695	4.78%	0.393%
76	17.278	2409	2415	2426	rBV	503397	923409	56.07%	4.612%

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77	17.378	2426	2432	2440	rVV	815213	1248610	75.81%	6.236%
78	17.455	2441	2445	2451	rVB	59123	89506	5.43%	0.447%
79	17.684	2481	2484	2488	rVB5	16748	20520	1.25%	0.102%
80	17.731	2488	2492	2495	rBV4	28659	36255	2.20%	0.181%
81	17.860	2510	2514	2521	rVB3	48922	85970	5.22%	0.429%
82	17.989	2532	2536	2543	rBV7	31674	49031	2.98%	0.245%
83	18.165	2561	2566	2580	rBV3	118626	239951	14.57%	1.198%
84	18.342	2593	2596	2602	rVV5	25526	40666	2.47%	0.203%
85	18.406	2604	2607	2610	rVV4	29671	36462	2.21%	0.182%
86	18.447	2611	2614	2618	rVB2	29173	38834	2.36%	0.194%
87	19.335	2761	2765	2770	rVB	58305	84412	5.13%	0.422%
88	19.440	2779	2783	2785	rBV4	22003	29629	1.80%	0.148%
89	19.670	2816	2822	2834	rBV	1073708	1620032	98.36%	8.091%
90	21.180	3075	3079	3092	rVB3	74755	161628	9.81%	0.807%
91	21.415	3114	3119	3125	rVB	82065	118883	7.22%	0.594%
92	21.526	3132	3138	3144	rBV2	466019	812652	49.34%	4.059%
93	21.708	3165	3169	3178	rVB	40356	71666	4.35%	0.358%
94	22.296	3265	3269	3275	rVB3	38811	61895	3.76%	0.309%
95	22.966	3380	3383	3389	rVB3	33115	54183	3.29%	0.271%
96	23.741	3510	3515	3522	rVB3	47675	107828	6.55%	0.539%
97	24.446	3626	3635	3660	rVB2	567730	1646979	100.00%	8.225%
98	24.664	3662	3672	3691	rVB	342592	996277	60.49%	4.976%
99	25.745	3852	3856	3867	rVB5	22991	60789	3.69%	0.304%
100	29.294	4458	4460	4461	rBV2	5447	4363	0.26%	0.022%

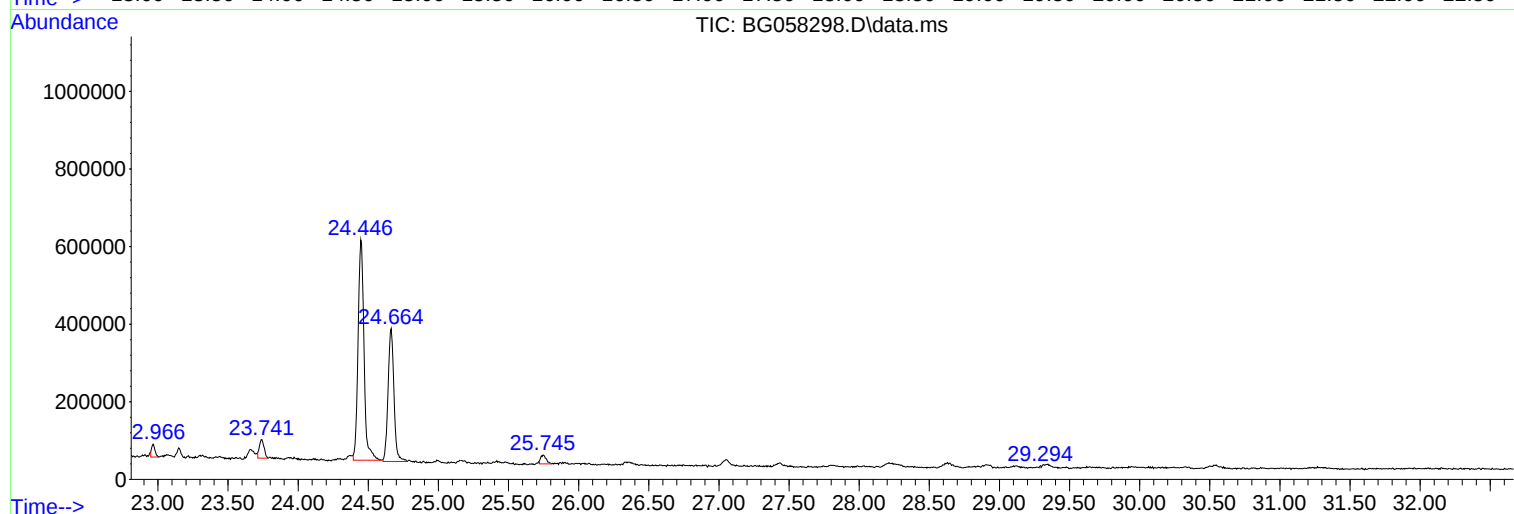
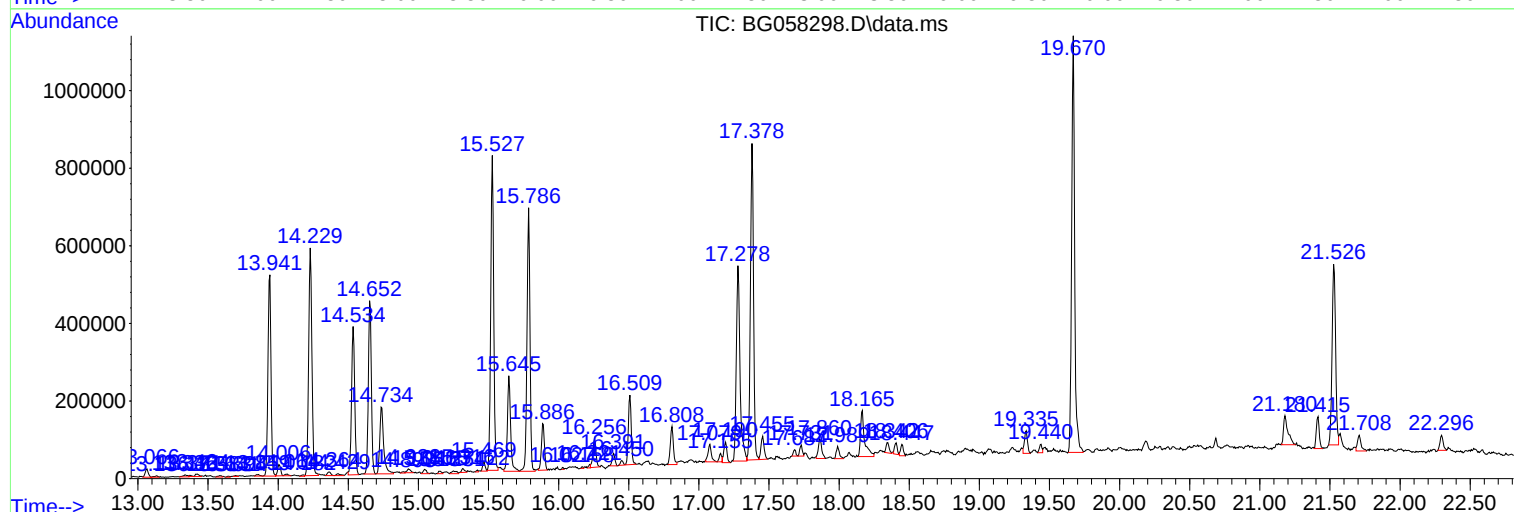
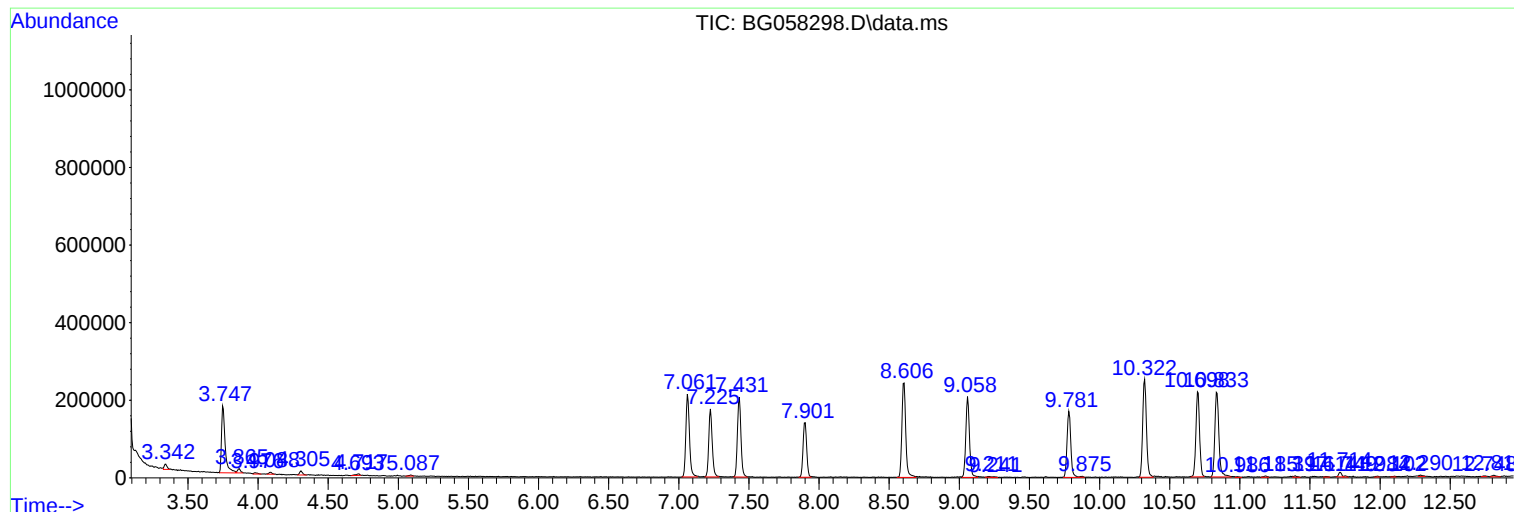
Sum of corrected areas: 20023184

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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



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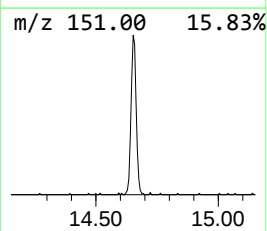
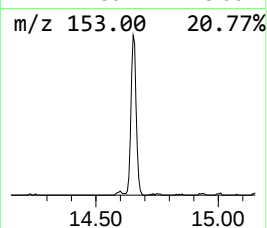
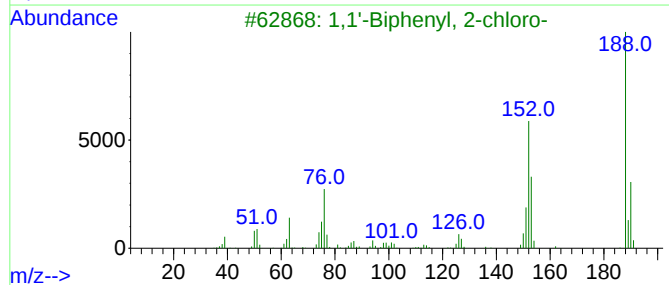
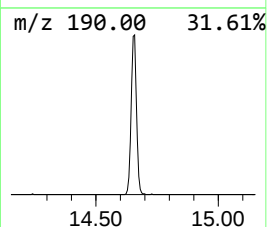
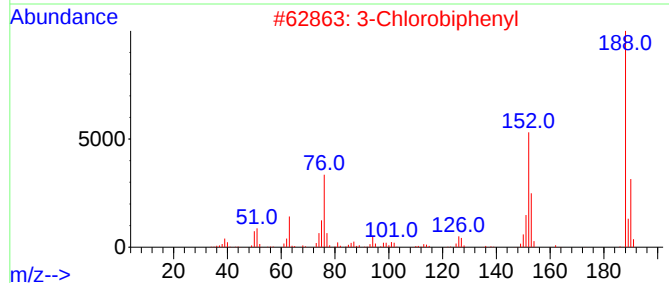
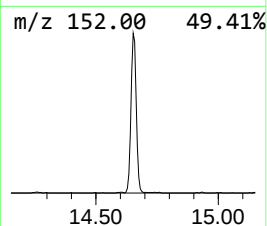
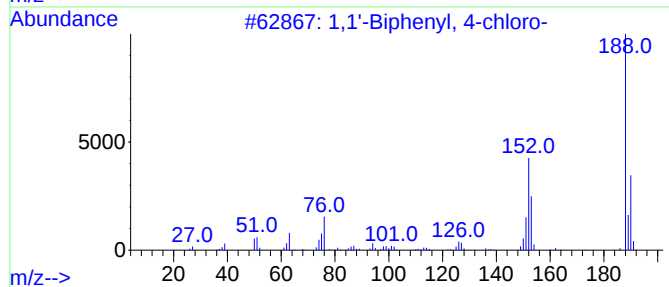
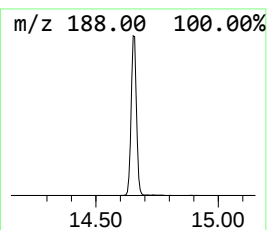
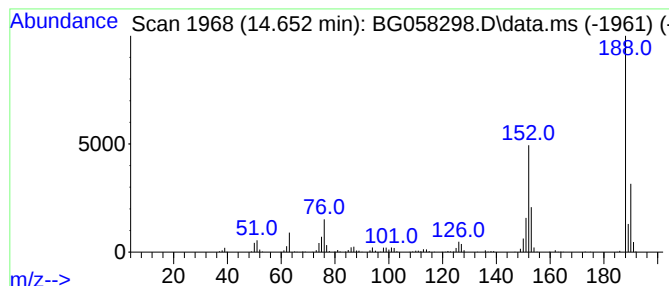
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,1'-Biphenyl, 4-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.652	22.17 ng/ul	675687	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	98
2	3-Chlorobiphenyl			188	C12H9Cl	002051-61-8	97
3	1,1'-Biphenyl, 2-chloro-			188	C12H9Cl	002051-60-7	97
4	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	95
5	1,1'-Biphenyl, 4-chloro-			188	C12H9Cl	002051-62-9	94



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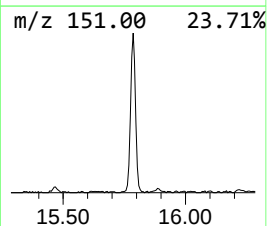
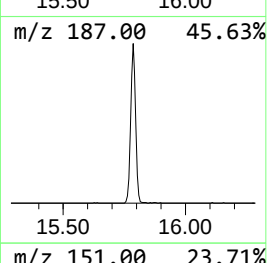
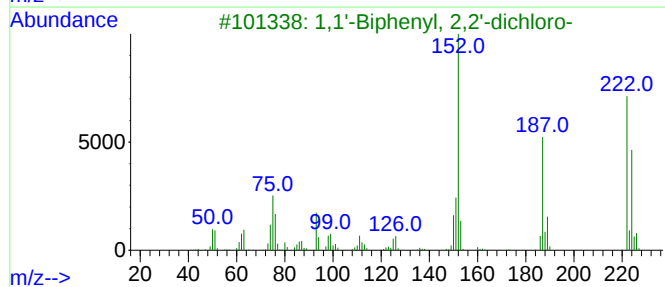
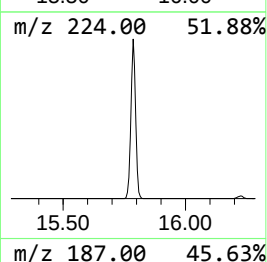
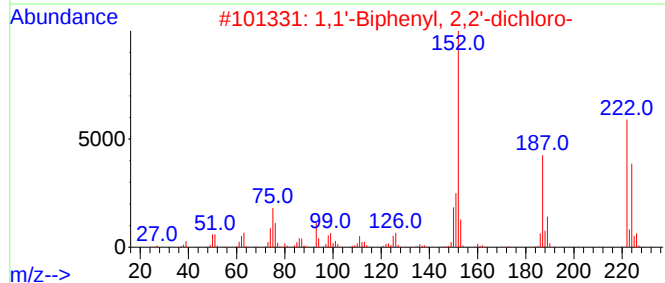
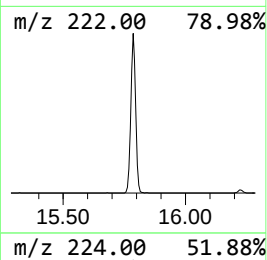
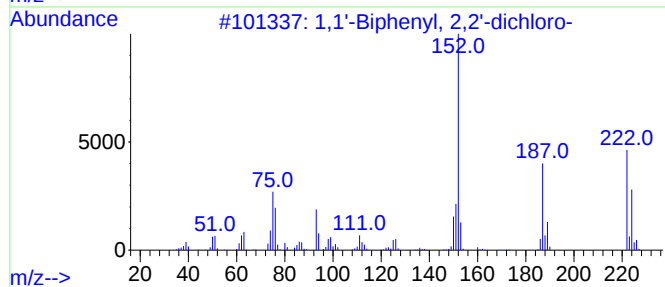
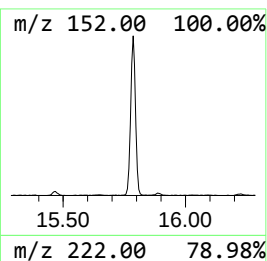
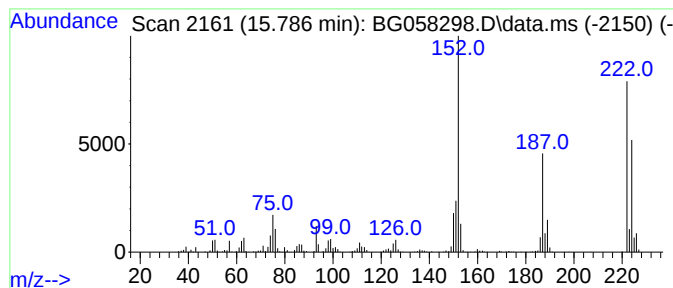
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,2'-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.786	32.60 ng/ul	993659	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99		
2	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	99		
3	1,1'-Biphenyl, 2,2'-dichloro-	222	C12H8Cl2	013029-08-8	98		
4	1,1'-Biphenyl, 3,4'-dichloro-	222	C12H8Cl2	002974-90-5	97		
5	1,1'-Biphenyl, 2,4'-dichloro-	222	C12H8Cl2	034883-43-7	97		



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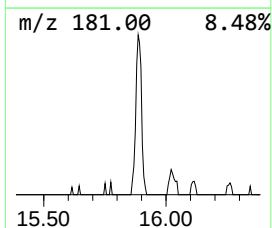
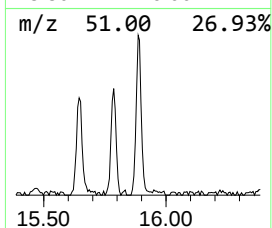
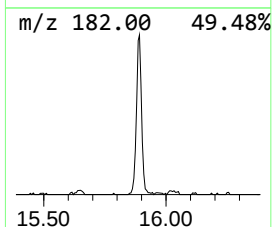
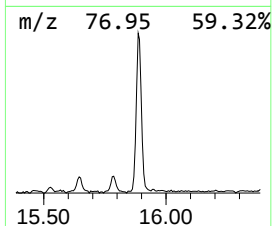
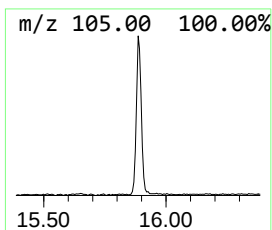
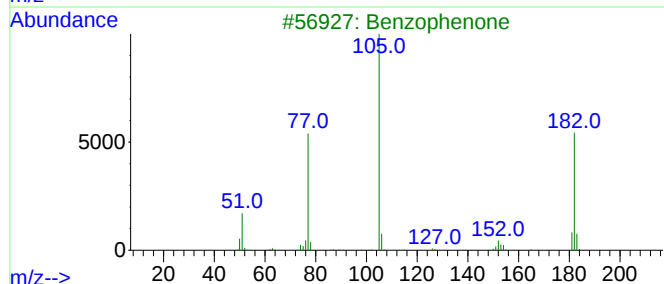
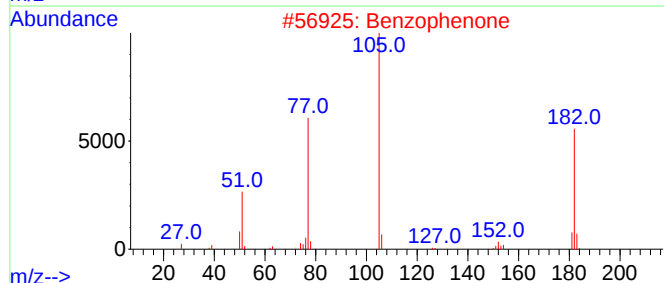
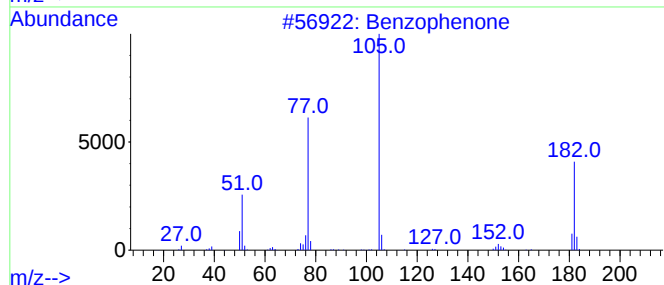
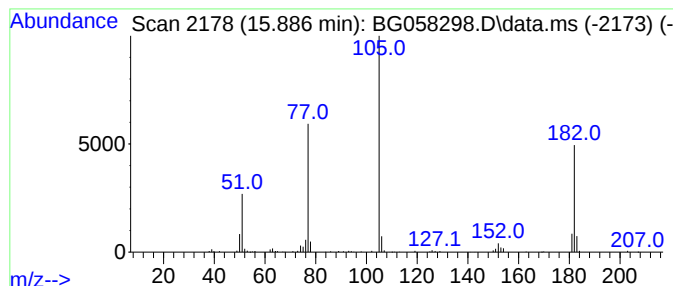
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzophenone Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.886	5.88 ng/ul	179170	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Benzophenone	182	C13H10O	000119-61-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A46Q1

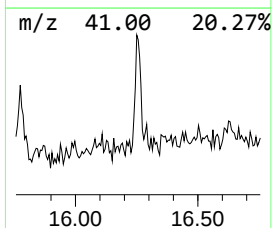
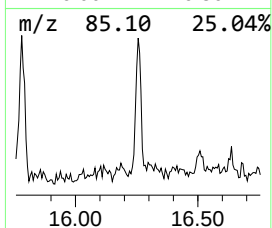
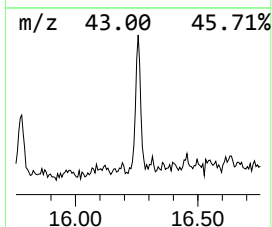
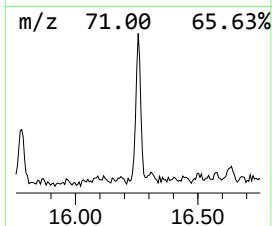
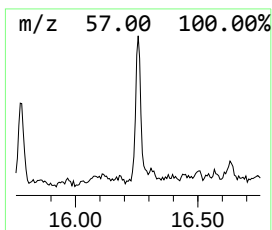
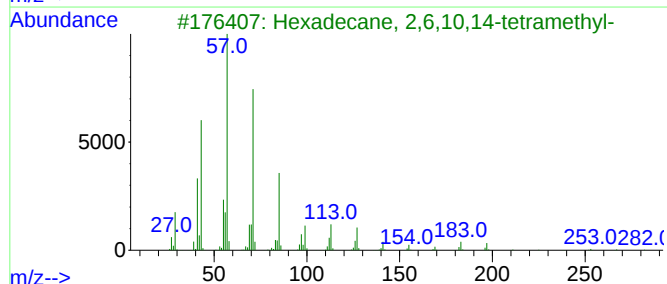
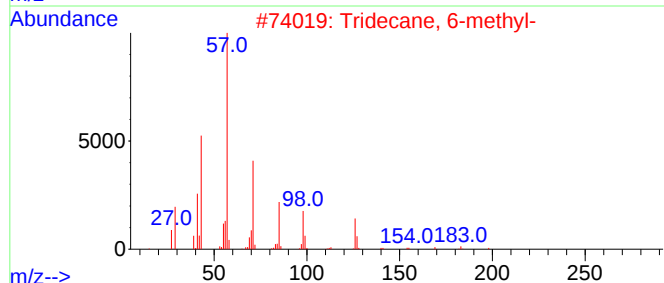
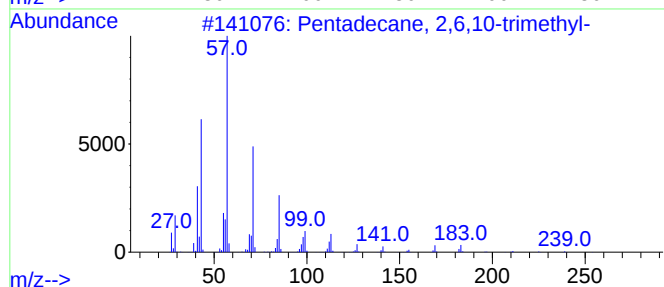
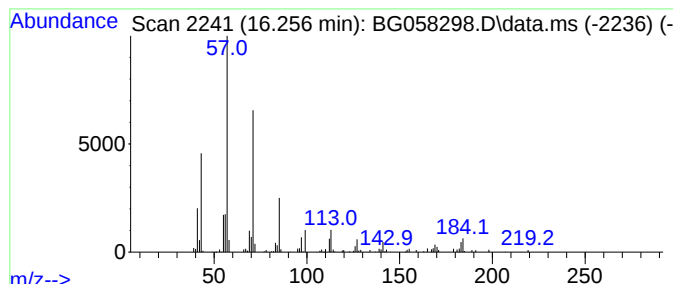
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.256	2.30 ng/ul	106394	Phenanthrene-d10	17.278

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
2		Tridecane, 6-methyl-	198	C14H30	013287-21-3	74
3		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	72
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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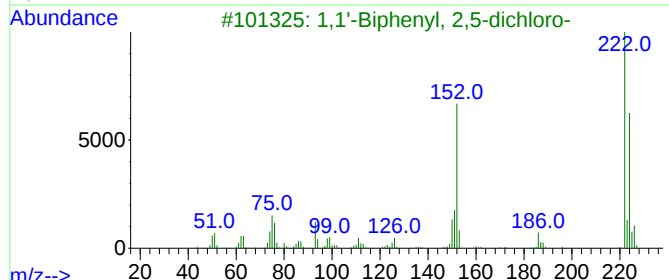
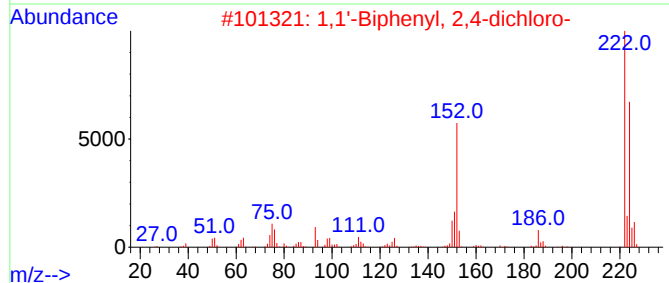
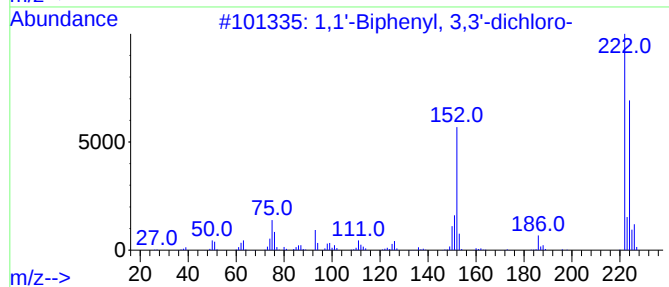
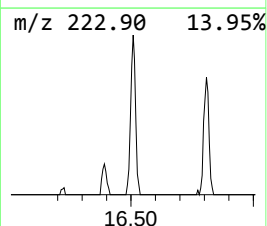
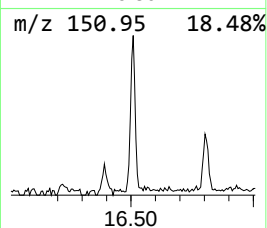
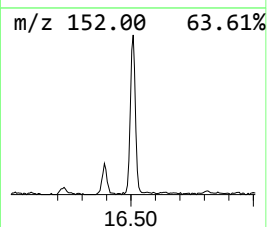
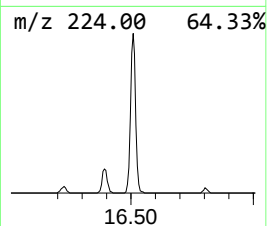
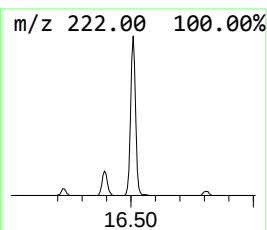
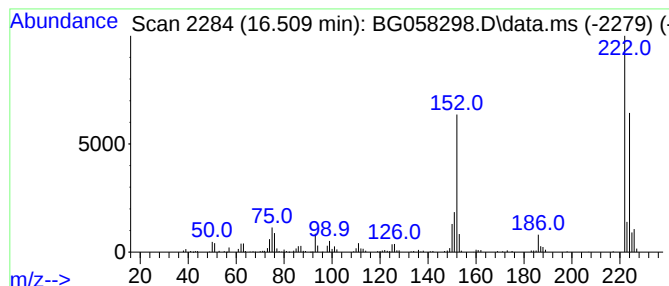
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 3,3'-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.509	5.62 ng/ul	259599	Phenanthrene-d10	17.278

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3'-dichloro-	222	C12H8Cl2	002050-67-1	99		
2	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	98		
3	1,1'-Biphenyl, 2,5-dichloro-	222	C12H8Cl2	034883-39-1	98		
4	1,1'-Biphenyl, 2,4-dichloro-	222	C12H8Cl2	033284-50-3	98		
5	1,1'-Biphenyl, 4,4'-dichloro-	222	C12H8Cl2	002050-68-2	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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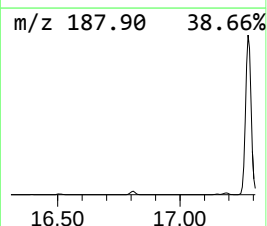
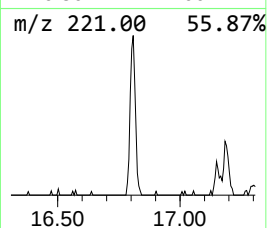
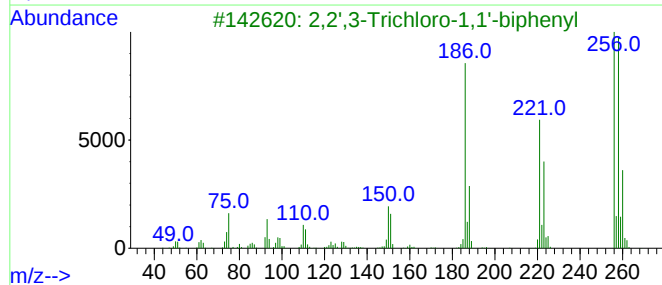
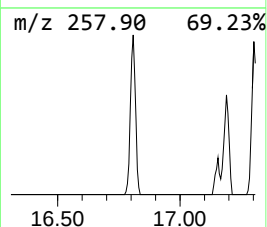
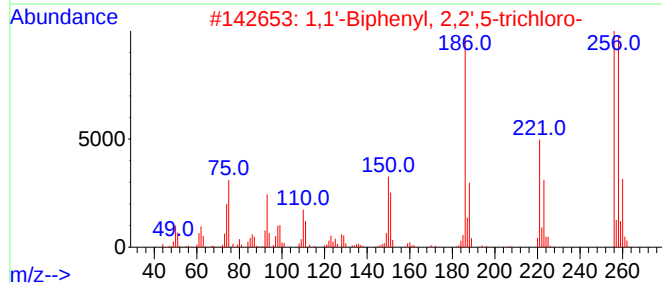
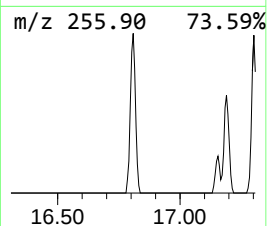
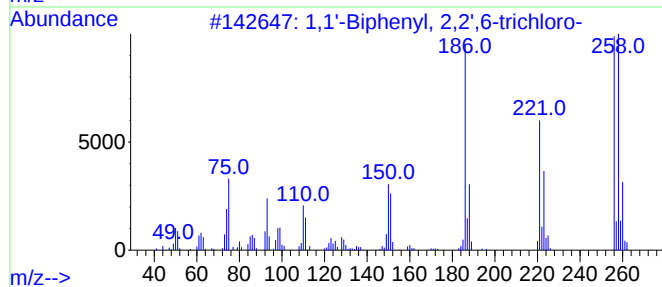
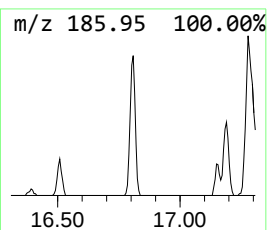
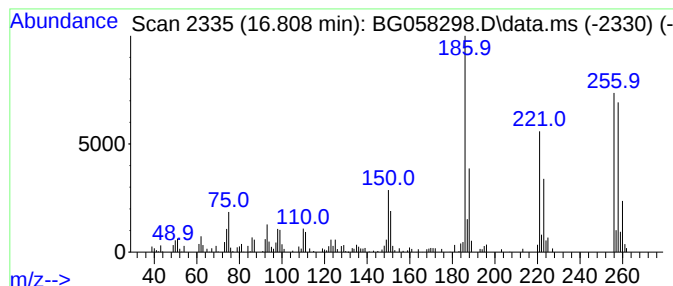
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',6-trich... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.808	3.27 ng/ul	151139	Phenanthrene-d10	17.278

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',6-trichloro-	256	C12H7Cl3	038444-73-4	99		
2	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	99		
3	2,2',3-Trichloro-1,1'-biphenyl	256	C12H7Cl3	038444-78-9	98		
4	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	98		
5	1,1'-Biphenyl, 2,2',5-trichloro-	256	C12H7Cl3	037680-65-2	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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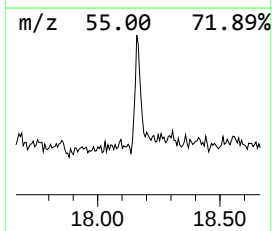
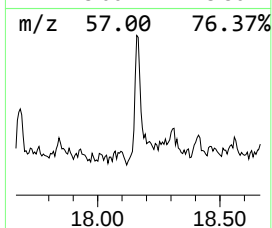
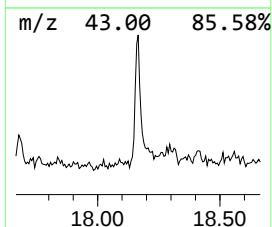
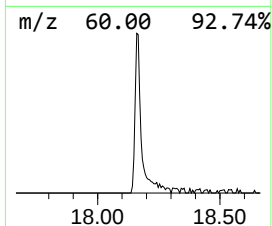
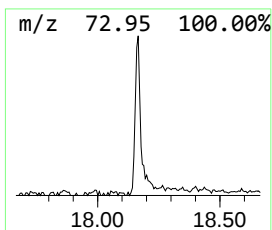
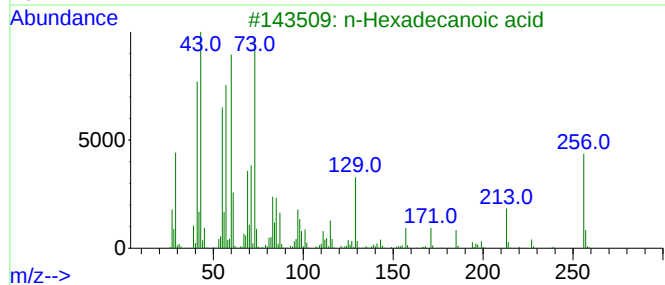
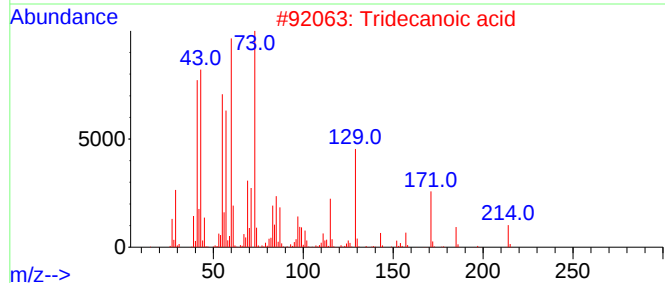
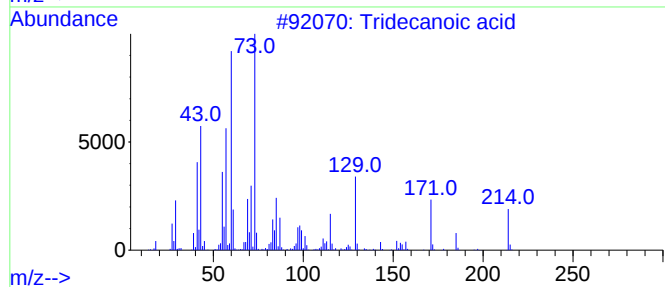
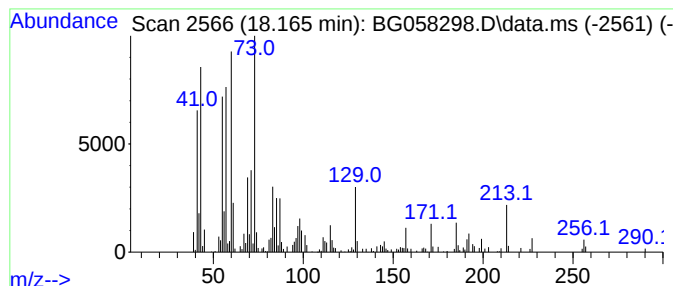
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tridecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.166	5.20 ng/ul	239951	Phenanthrene-d10	17.278

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecanoic acid	214	C13H26O2	000638-53-9	90
2			Tridecanoic acid	214	C13H26O2	000638-53-9	90
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	83
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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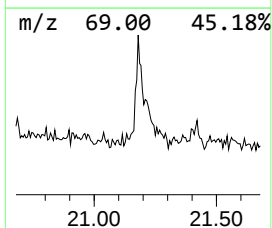
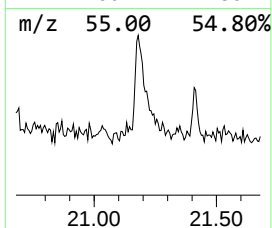
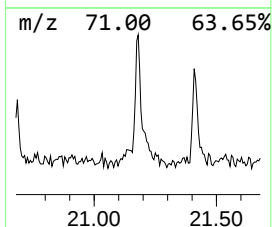
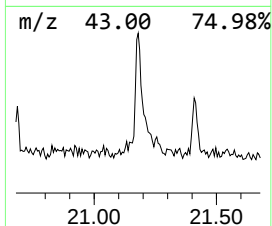
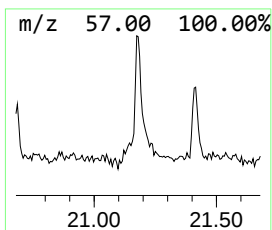
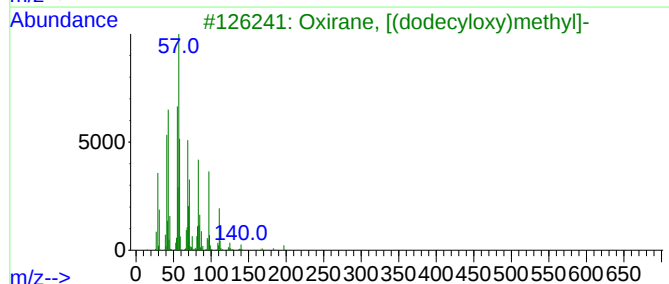
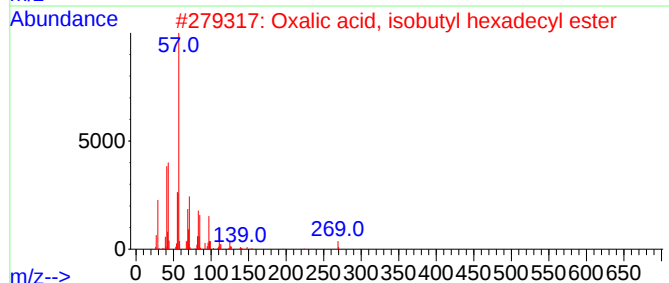
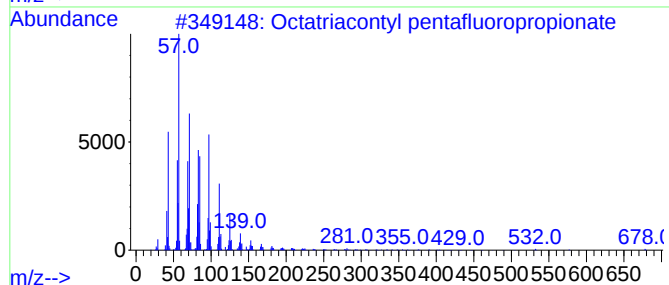
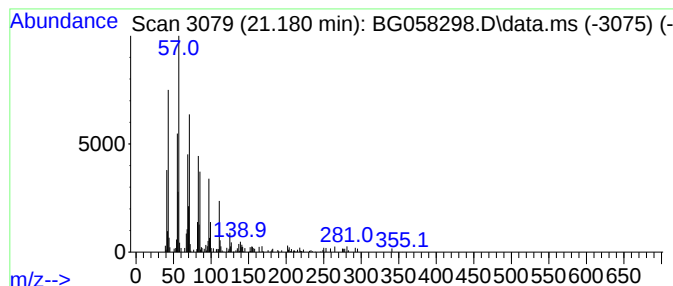
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octatriacontyl pentafluorop... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.180	3.98 ng/ul	161628	Chrysene-d12	21.526

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octatriacontyl pentafluoropropio...	697	C41H77F5O2	1000351-89-1	90
2			Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	74
3			Oxirane, [(dodecyloxy)methyl]-	242	C15H30O2	002461-18-9	74
4			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	74
5			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
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Sample : 03444-07
Misc :
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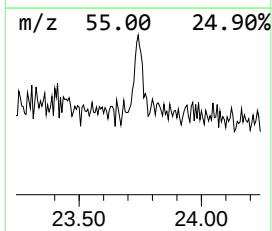
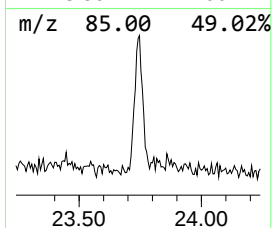
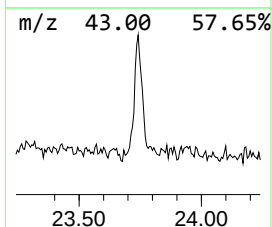
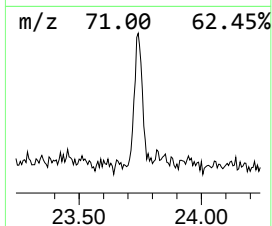
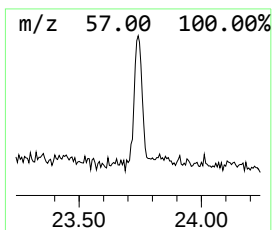
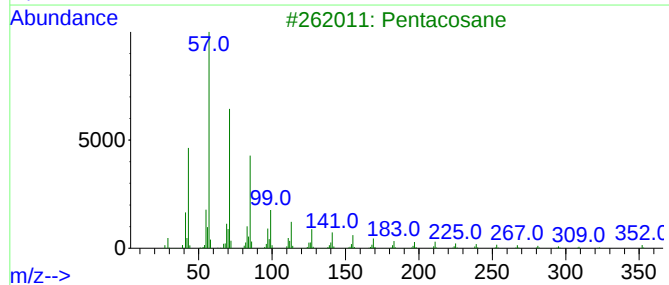
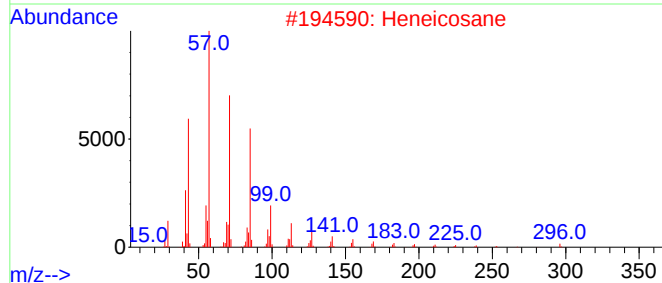
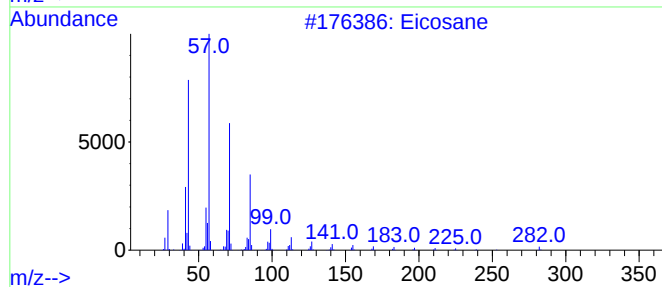
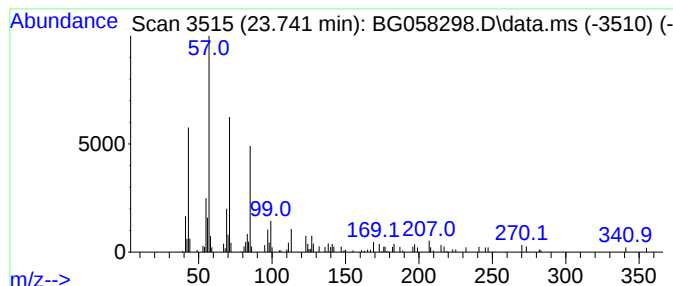
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.741	2.16 ng/ul	107828	Perylene-d12	24.664

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	90
2		Heneicosane	296	C21H44	000629-94-7	87
3		Pentacosane	352	C25H52	000629-99-2	87
4		Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	86
5		Tetratetracontane	619	C44H90	007098-22-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058298.D
Acq On : 18 Jul 2023 12:54
Operator : CG/JU
Sample : 03444-07
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.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
1,1'-Biphenyl, ...	14.652	22.2	ng/ul	675687	3	14.534	609593	20.0
1,1'-Biphenyl, ...	15.786	32.6	ng/ul	993659	3	14.534	609593	20.0
Benzophenone	15.886	5.9	ng/ul	179170	3	14.534	609593	20.0
(DEL) Alkane: S...	16.256	2.3	ng/ul	106394	4	17.278	923409	20.0
1,1'-Biphenyl, ...	16.509	5.6	ng/ul	259599	4	17.278	923409	20.0
1,1'-Biphenyl, ...	16.808	3.3	ng/ul	151139	4	17.278	923409	20.0
Tridecanoic acid	18.166	5.2	ng/ul	239951	4	17.278	923409	20.0
Octatriacontyl ...	21.180	4.0	ng/ul	161628	5	21.526	812652	20.0
(DEL) Alkane: S...	23.741	2.2	ng/ul	107828	6	24.664	996277	20.0