

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
 Data File : BG061866.D
 Acq On : 3 Jul 2024 16:21
 Operator : MA/JU
 Sample : P2964-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4CT0

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-BG070224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG061866.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.482	29	33	38	rBV4	10490	15647	1.30%	0.106%
2	3.758	75	80	86	rBV	20347	37695	3.13%	0.255%
3	3.911	104	106	111	rBV2	17409	27698	2.30%	0.187%
4	4.016	120	124	129	rVB5	12823	17658	1.46%	0.119%
5	4.087	134	136	142	rVB3	3883	6312	0.52%	0.043%
6	4.245	160	163	168	rVB4	5670	8150	0.68%	0.055%
7	4.763	246	251	254	rBV5	3296	5614	0.47%	0.038%
8	4.810	255	259	263	rVV4	11229	18587	1.54%	0.126%
9	4.851	263	266	272	rVB2	12596	17374	1.44%	0.117%
10	5.062	298	302	304	rBV4	3053	3377	0.28%	0.023%
11	5.086	304	306	310	rVB3	2806	3408	0.28%	0.023%
12	5.127	310	313	315	rBV2	3052	4246	0.35%	0.029%
13	5.162	315	319	326	rVV2	38565	59552	4.94%	0.403%
14	5.250	332	334	338	rBV4	3638	5420	0.45%	0.037%
15	5.362	352	353	357	rBV	2294	3315	0.27%	0.022%
16	6.155	482	488	492	rBV4	9155	15730	1.30%	0.106%
17	7.248	666	674	683	rVV	169742	288988	23.97%	1.954%
18	7.418	697	703	711	rBV	115870	187738	15.57%	1.270%
19	7.624	731	738	744	rBV	151819	262184	21.75%	1.773%
20	7.765	758	762	763	rVB	2960	3235	0.27%	0.022%
21	8.094	812	818	825	rVV	162067	293366	24.33%	1.984%
22	8.147	825	827	833	rVB4	11085	15451	1.28%	0.104%
23	8.329	857	858	862	rBV	2706	3170	0.26%	0.021%
24	8.664	913	915	920	rVB	3088	3481	0.29%	0.024%
25	8.799	929	938	945	rBV	187912	329480	27.33%	2.228%
26	8.916	956	958	962	rVB2	4742	3890	0.32%	0.026%
27	9.263	1007	1017	1026	rVB	151197	281924	23.38%	1.906%
28	9.645	1075	1082	1088	rBV3	8459	15332	1.27%	0.104%
29	9.810	1101	1110	1115	rBV2	26600	52896	4.39%	0.358%
30	9.986	1130	1140	1147	rBV2	145599	268098	22.24%	1.813%
31	10.520	1224	1231	1241	rVB	201446	379120	31.44%	2.564%
32	10.614	1245	1247	1252	rVB3	3941	4014	0.33%	0.027%
33	10.908	1290	1297	1305	rBV	252486	458328	38.01%	3.099%
34	11.038	1313	1319	1329	rBV	117882	219785	18.23%	1.486%
35	11.137	1334	1336	1340	rVB2	3815	3388	0.28%	0.023%
36	11.349	1370	1372	1376	rVV	2242	3626	0.30%	0.025%

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Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
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37	11.431	1379	1386	1391	rBV4	13455	24339	2.02%	0.165%
38	11.643	1414	1422	1427	rBV2	31635	64889	5.38%	0.439%
39	12.307	1530	1535	1537	rBV2	6455	8192	0.68%	0.055%
40	12.471	1561	1563	1570	rVV4	4744	9001	0.75%	0.061%

41	12.659	1592	1595	1598	rBV	2631	3829	0.32%	0.026%
42	12.965	1642	1647	1649	rBV2	4633	4527	0.38%	0.031%
43	12.988	1649	1651	1654	rVV	3301	3593	0.30%	0.024%
44	13.029	1655	1658	1662	rVB	3160	4301	0.36%	0.029%
45	13.188	1684	1685	1690	rVB5	3584	4489	0.37%	0.030%

46	13.329	1705	1709	1710	rBV2	5095	5526	0.46%	0.037%
47	13.529	1739	1743	1747	rBV3	12796	16779	1.39%	0.113%
48	14.116	1838	1843	1849	rVV	434293	628583	52.14%	4.251%
49	14.357	1880	1884	1885	rBV3	6223	7278	0.60%	0.049%
50	14.416	1887	1894	1902	rVB2	411627	669876	55.56%	4.530%

51	14.533	1910	1914	1915	rBV2	2716	3323	0.28%	0.022%
52	14.598	1923	1925	1930	rVB4	5084	7668	0.64%	0.052%
53	14.716	1938	1945	1953	rBV	434976	675475	56.02%	4.568%
54	14.774	1953	1955	1957	rBV2	4374	3610	0.30%	0.024%
55	14.904	1970	1977	1986	rBV2	220806	390258	32.37%	2.639%

56	15.062	1998	2004	2008	rBV5	12367	21150	1.75%	0.143%
57	15.115	2010	2013	2018	rVV6	9021	14262	1.18%	0.096%
58	15.485	2073	2076	2081	rVB3	7117	9806	0.81%	0.066%
59	15.526	2081	2083	2085	rBV3	4302	4636	0.38%	0.031%
60	15.550	2085	2087	2091	rVV3	3910	4088	0.34%	0.028%

61	15.656	2103	2105	2107	rBV3	3963	4582	0.38%	0.031%
62	15.709	2107	2114	2119	rVV	574808	877701	72.80%	5.935%
63	15.761	2119	2123	2125	rVB3	5895	5938	0.49%	0.040%
64	15.814	2126	2132	2137	rBV2	218952	309498	25.67%	2.093%
65	15.926	2147	2151	2152	rBV3	3543	4579	0.38%	0.031%

66	16.043	2167	2171	2172	rVV3	4023	3847	0.32%	0.026%
67	16.149	2186	2189	2190	rBV2	4187	3578	0.30%	0.024%
68	16.167	2190	2192	2194	rVV2	4197	4201	0.35%	0.028%
69	16.273	2208	2210	2213	rBV4	3037	3387	0.28%	0.023%
70	16.408	2231	2233	2234	rBV2	5003	3612	0.30%	0.024%

71	16.690	2276	2281	2282	rBV4	5452	6891	0.57%	0.047%
72	16.854	2307	2309	2311	rBV3	5495	3743	0.31%	0.025%
73	16.989	2327	2332	2335	rBV5	17386	26861	2.23%	0.182%
74	17.148	2357	2359	2364	rVB5	12232	12838	1.06%	0.087%
75	17.195	2364	2367	2370	rBV5	9252	11499	0.95%	0.078%

76	17.330	2386	2390	2393	rBV6	13657	18915	1.57%	0.128%
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77	17.360	2393	2395	2402	rVB7	14857	22276	1.85%	0.151%
78	17.459	2405	2412	2417	rVV	576971	923438	76.59%	6.245%
79	17.501	2417	2419	2423	rVV	83125	95763	7.94%	0.648%
80	17.553	2423	2428	2437	rVV	619240	1001315	83.05%	6.771%
81	17.630	2437	2441	2447	rVB5	34544	50904	4.22%	0.344%
82	17.900	2484	2487	2490	rBV5	14208	17819	1.48%	0.120%
83	18.053	2508	2513	2520	rVB3	101427	173393	14.38%	1.173%
84	18.170	2530	2533	2540	rVB8	22095	38415	3.19%	0.260%
85	18.300	2552	2555	2558	rBV3	19626	30453	2.53%	0.206%
86	18.329	2558	2560	2566	rVV6	31376	50464	4.19%	0.341%
87	18.517	2587	2592	2597	rBV4	94488	152949	12.69%	1.034%
88	18.576	2599	2602	2606	rVV4	69365	91951	7.63%	0.622%
89	18.617	2606	2609	2615	rVB8	36273	53062	4.40%	0.359%
90	18.799	2636	2640	2645	rBV6	61977	101451	8.41%	0.686%
91	18.975	2667	2670	2676	rVB6	42008	52388	4.35%	0.354%
92	19.328	2727	2730	2733	rBV3	52212	64517	5.35%	0.436%
93	19.363	2734	2736	2741	rVB5	46598	64364	5.34%	0.435%
94	19.504	2755	2760	2764	rVB	154246	210828	17.49%	1.426%
95	19.833	2811	2816	2829	rBV2	715684	1205671	100.00%	8.153%
96	21.361	3073	3076	3082	rBV4	89289	140108	11.62%	0.947%
97	21.719	3131	3137	3142	rBV2	524519	904987	75.06%	6.120%
98	24.040	3528	3532	3539	rVB	236106	452281	37.51%	3.059%
99	24.739	3644	3651	3658	rVB2	255293	646482	53.62%	4.372%
100	24.968	3681	3690	3701	rVB3	329062	1023936	84.93%	6.924%

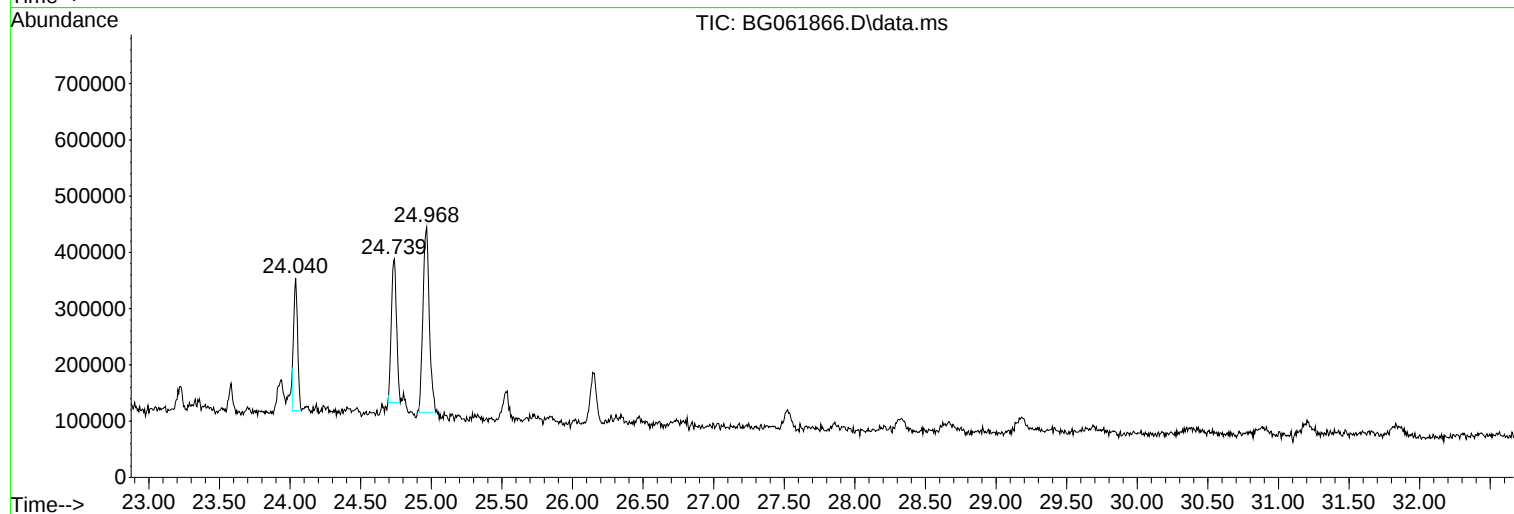
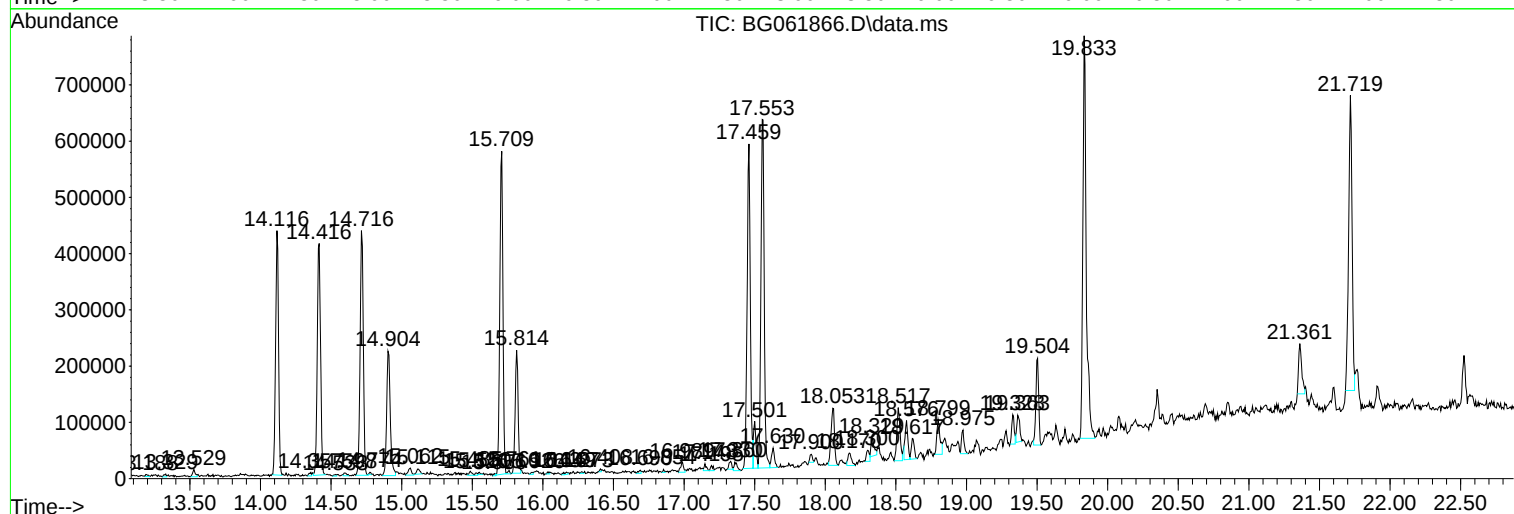
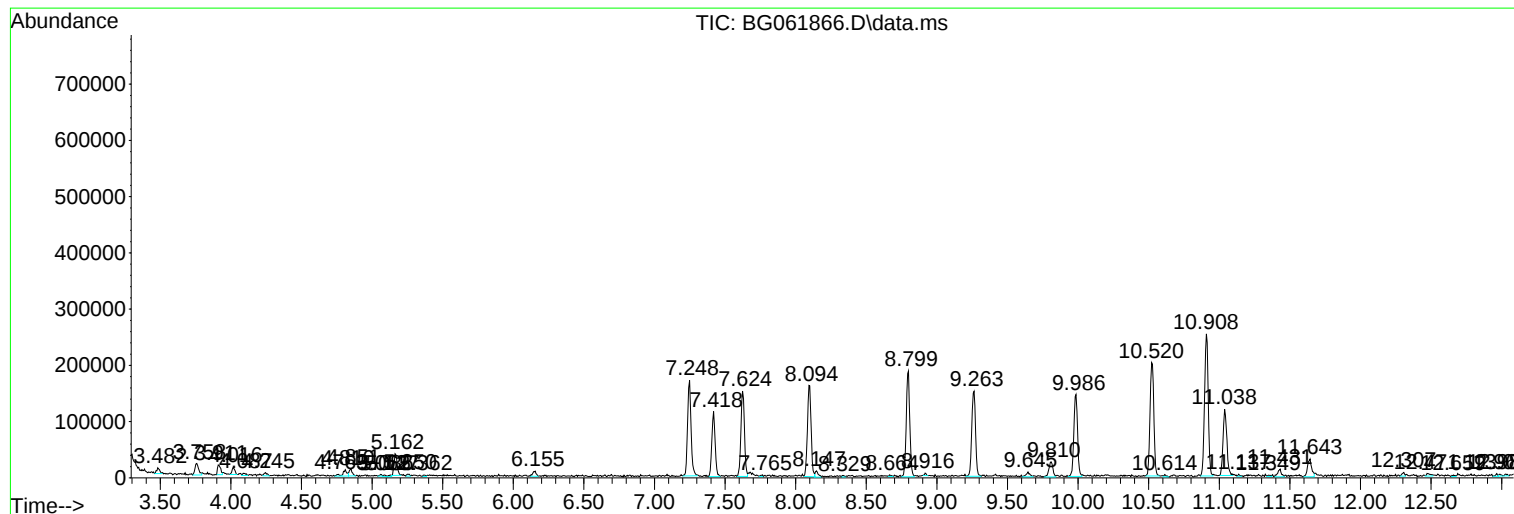
Sum of corrected areas: 14787640

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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



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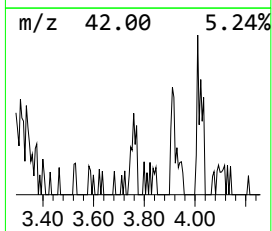
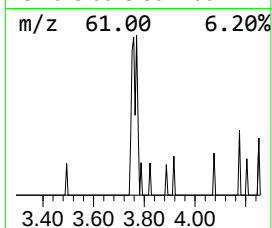
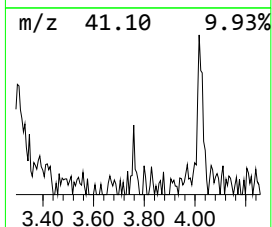
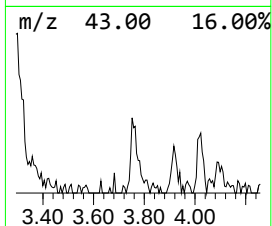
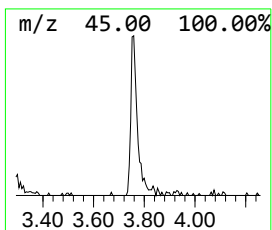
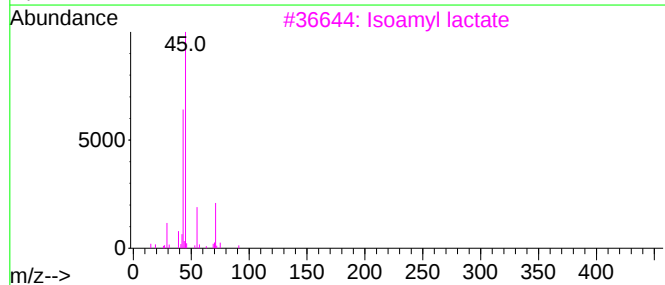
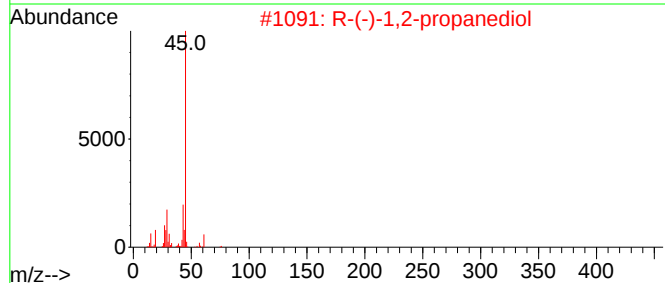
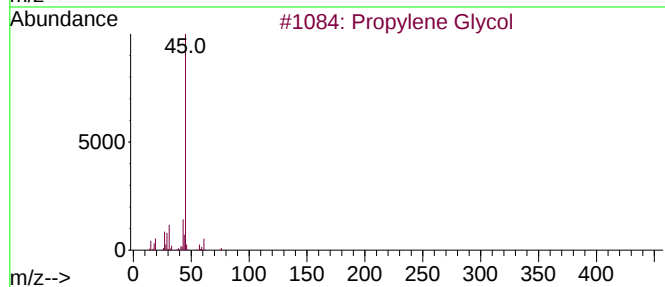
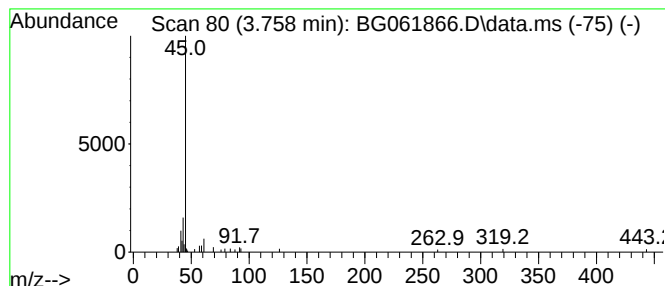
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.758	2.57 ng/ul	37695	1,4-Dichlorobenzene-d4	8.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	50
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	50
3			Isoamyl lactate	160	C8H16O3	019329-89-6	45
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	45
5			2-Butanol, 3-methyl-	88	C5H12O	000598-75-4	39



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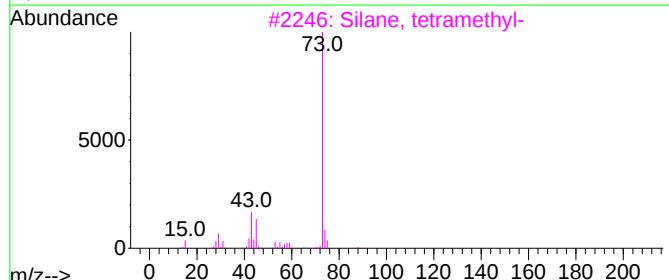
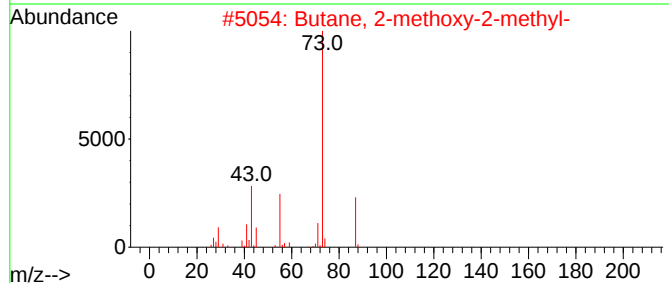
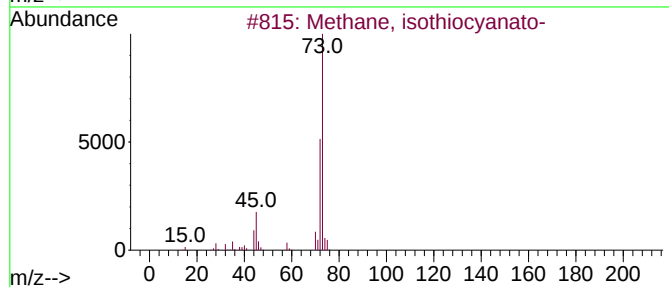
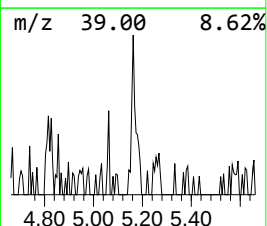
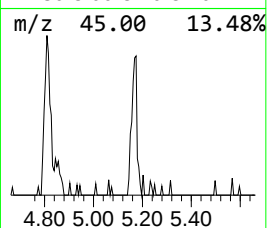
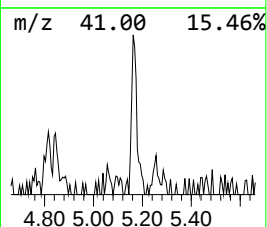
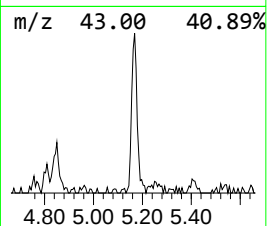
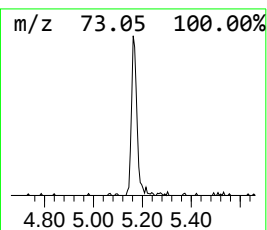
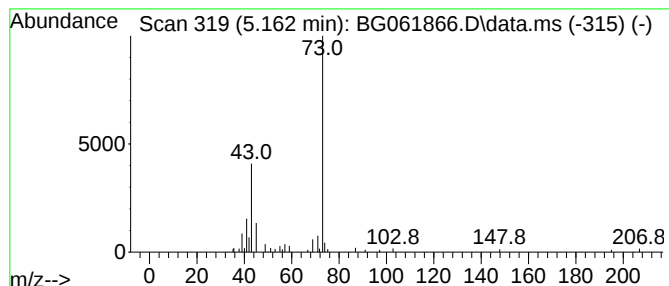
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.162	4.06 ng/ul	59552	1,4-Dichlorobenzene-d4	8.094

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methane, isothiocyanato-	73	C2H3NS	000556-61-6	37
2			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	37
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
4			7-Methoxy-3,7-dimethyloctan-2-ol...	188	C11H24O2	1000511-09-8	12
5			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	12



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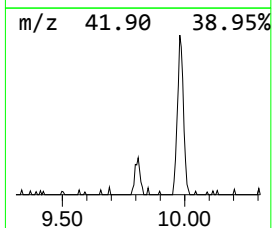
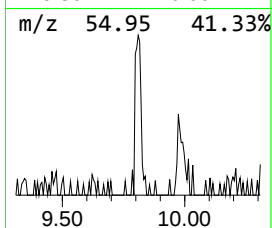
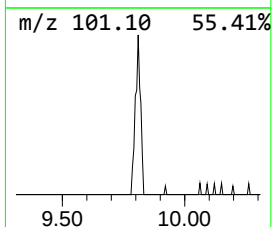
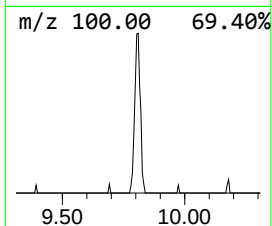
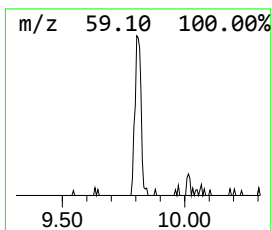
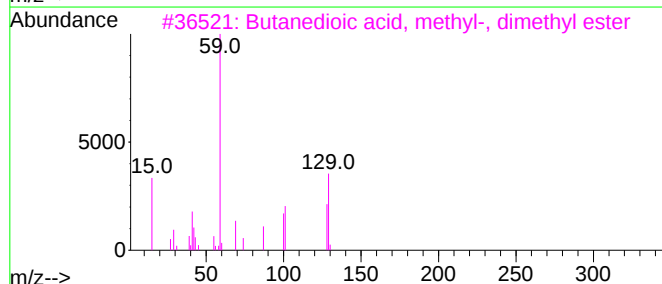
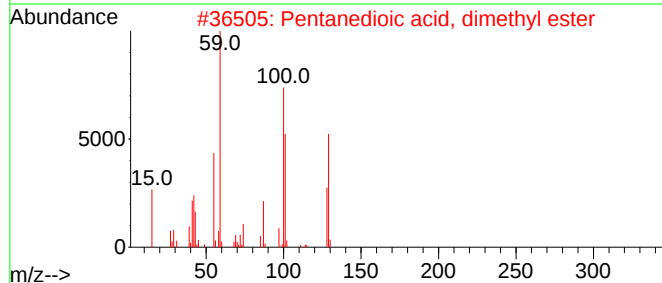
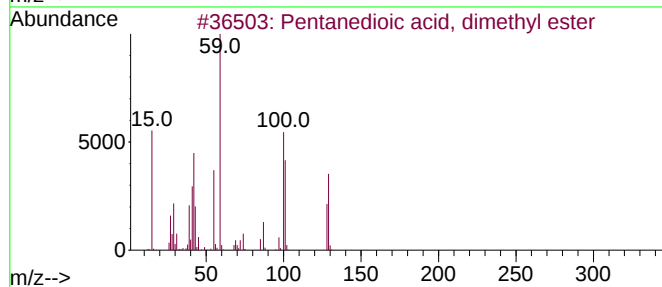
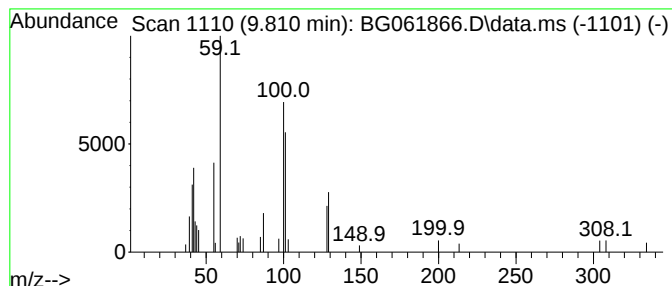
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentanedioic acid, dimethyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.810	2.31 ng/ul	52896	Naphthalene-d8	10.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	83
2			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	39
3			Butanedioic acid, methyl-, dimet...	160	C7H12O4	001604-11-1	37
4			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	34
5			Pentanedioic acid, dimethyl ester	160	C7H12O4	001119-40-0	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
Acq On : 3 Jul 2024 16:21
Operator : MA/JU
Sample : P2964-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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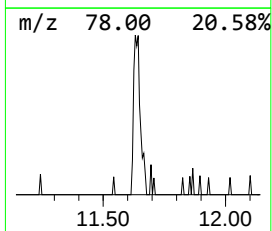
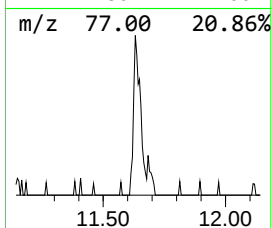
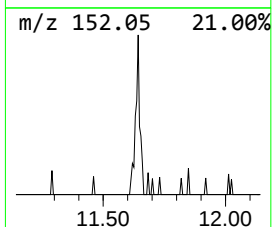
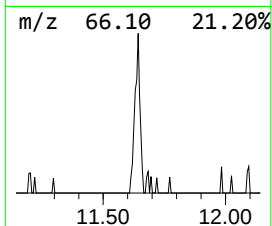
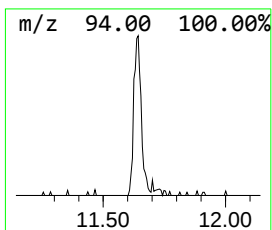
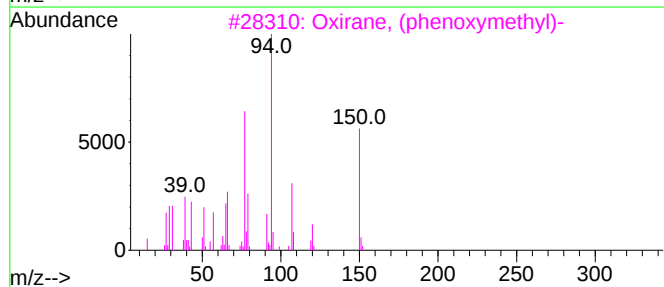
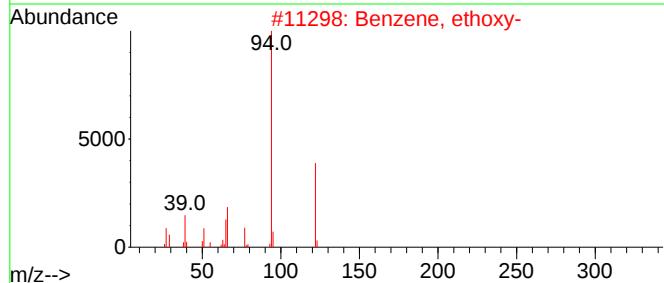
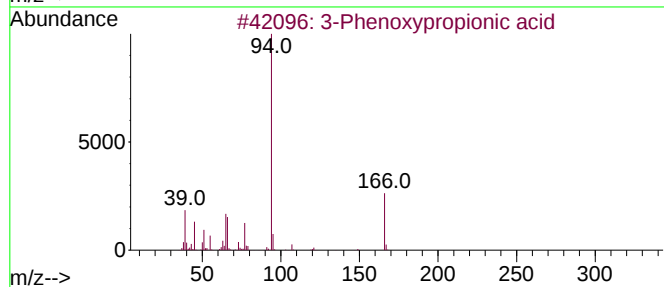
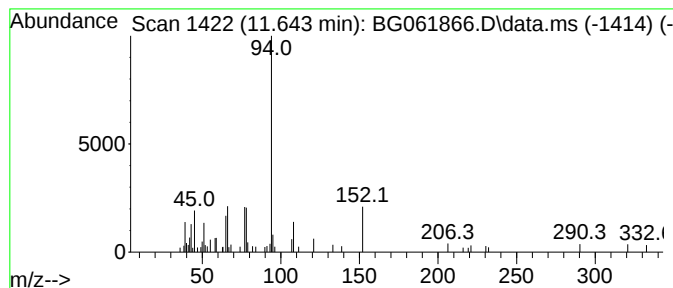
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 3-Phenoxypropionic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.643	2.83 ng/ul	64889	Naphthalene-d8	10.908

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenoxypropionic acid	166	C9H10O3	007170-38-9	53
2			Benzene, ethoxy-	122	C8H10O	000103-73-1	50
3			Oxirane, (phenoxymethyl)-	150	C9H10O2	000122-60-1	50
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	50
5			Benzene, ethoxy-	122	C8H10O	000103-73-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
Acq On : 3 Jul 2024 16:21
Operator : MA/JU
Sample : P2964-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A4CT0

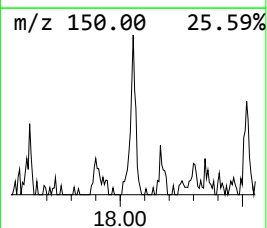
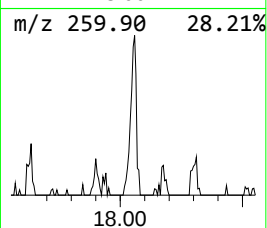
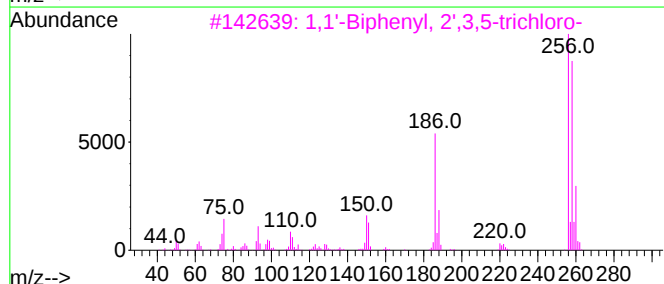
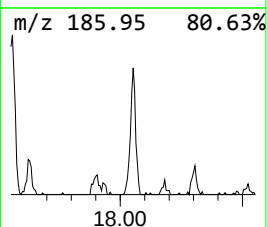
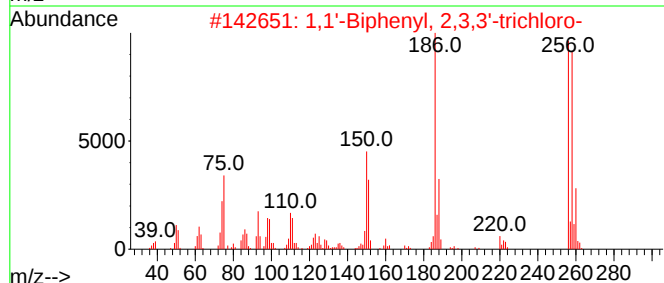
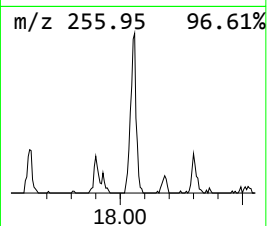
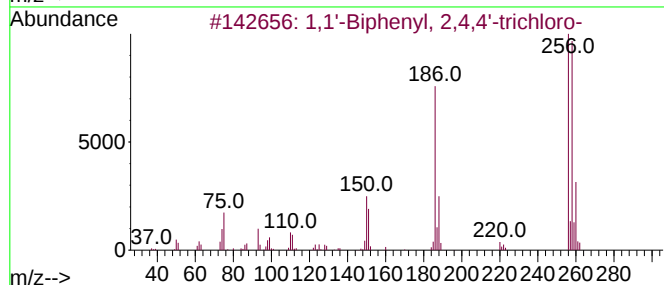
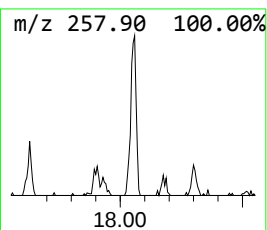
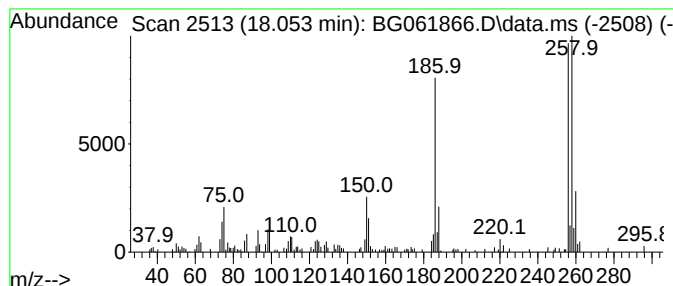
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.053	3.76 ng/ul	173393	Phenanthrene-d10	17.459

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99	
2	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99	
3	1,1'-Biphenyl, 2',3,5-trichloro-	256	C12H7Cl3	037680-68-5	98	
4	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	98	
5	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	98	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
Acq On : 3 Jul 2024 16:21
Operator : MA/JU
Sample : P2964-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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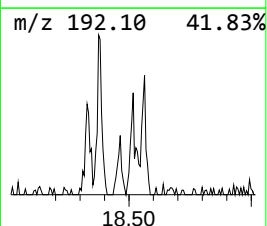
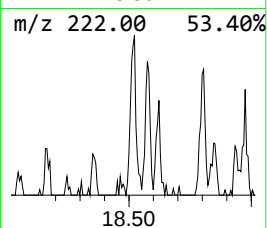
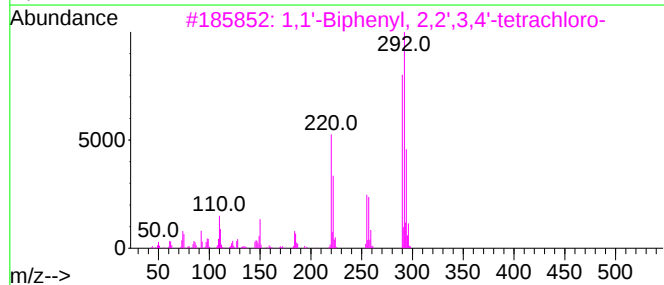
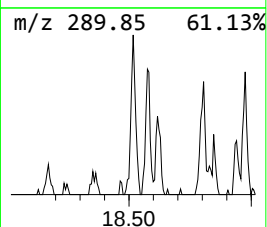
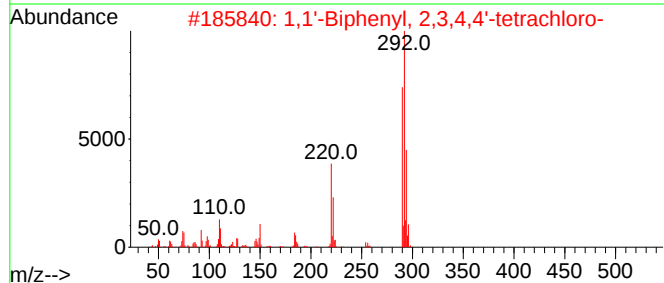
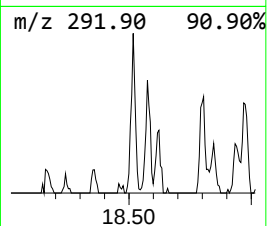
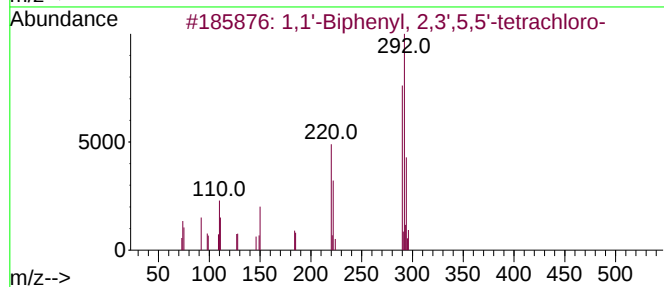
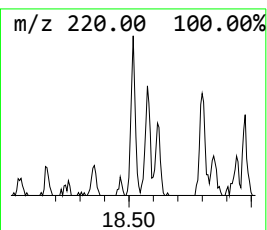
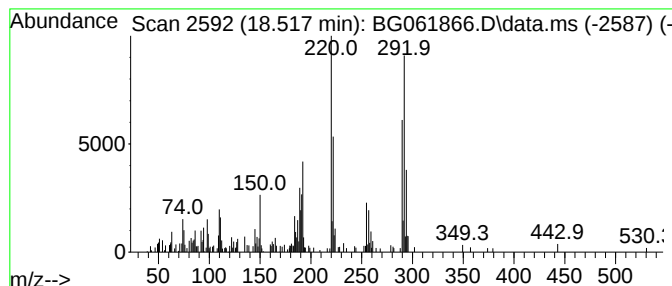
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,3',5,5'-te... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.517	3.31 ng/ul	152949	Phenanthrene-d10	17.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	95		
2	1,1'-Biphenyl, 2,3,4,4'-tetrachl...	290	C12H6Cl4	033025-41-1	93		
3	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	91		
4	1,1'-Biphenyl, 2,4,4',6-tetrachl...	290	C12H6Cl4	032598-12-2	90		
5	1,1'-Biphenyl, 2,3',4',6-tetrach...	290	C12H6Cl4	041464-46-4	90		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
Acq On : 3 Jul 2024 16:21
Operator : MA/JU
Sample : P2964-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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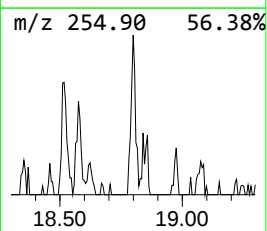
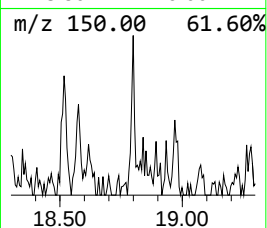
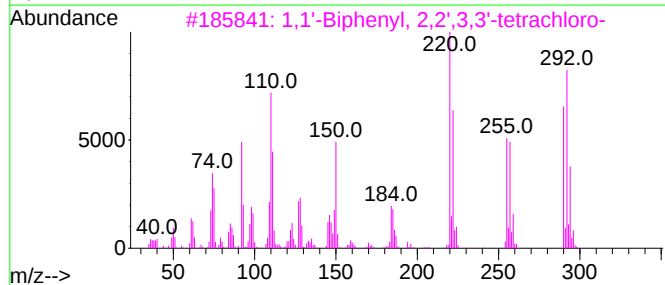
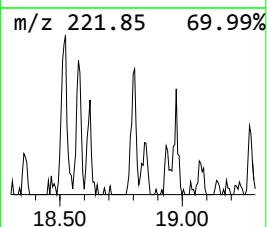
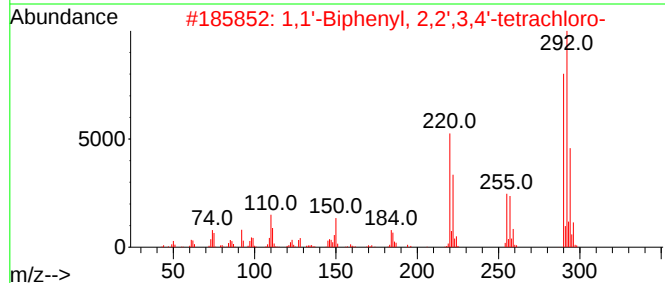
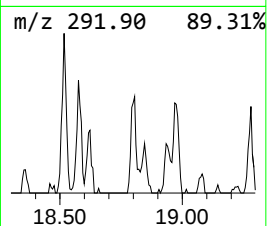
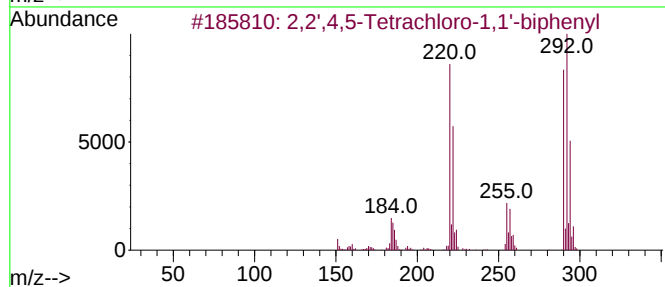
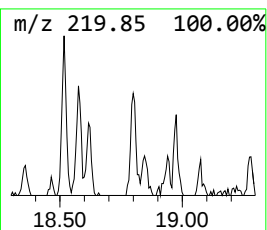
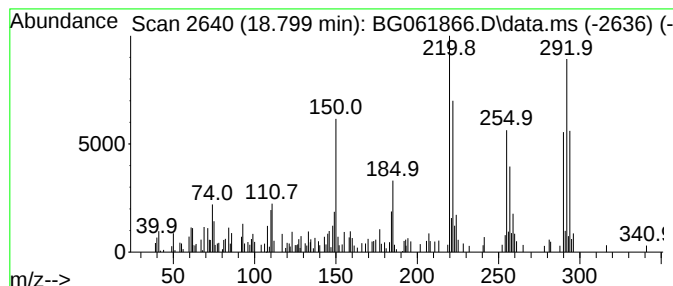
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.799	2.20 ng/ul	101451	Phenanthrene-d10	17.459

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	94		
2	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	94		
3	1,1'-Biphenyl, 2,2',3,3'-tetrach...	290	C12H6Cl4	038444-93-8	94		
4	2,2',3,6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-45-7	93		
5	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	92		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
Acq On : 3 Jul 2024 16:21
Operator : MA/JU
Sample : P2964-13
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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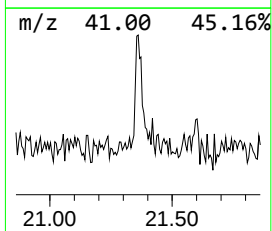
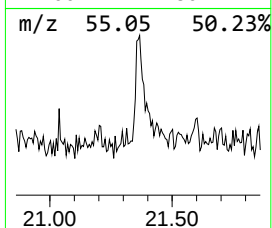
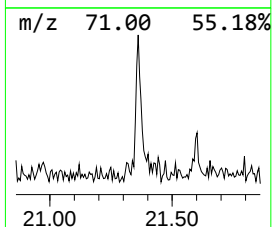
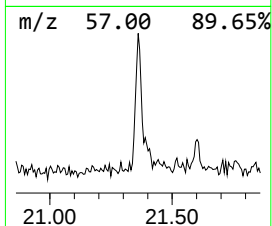
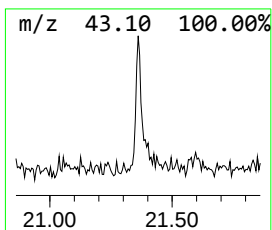
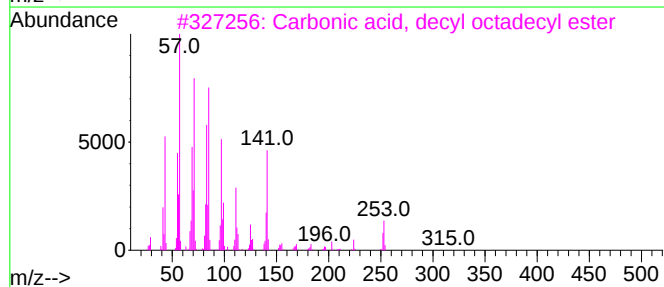
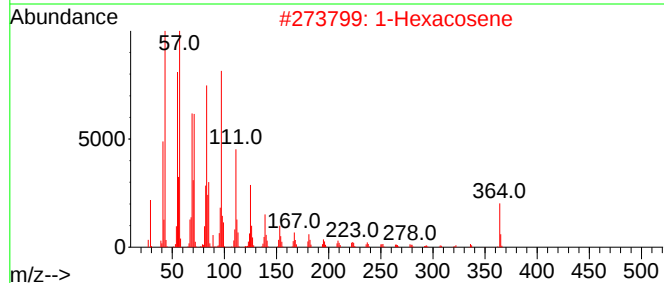
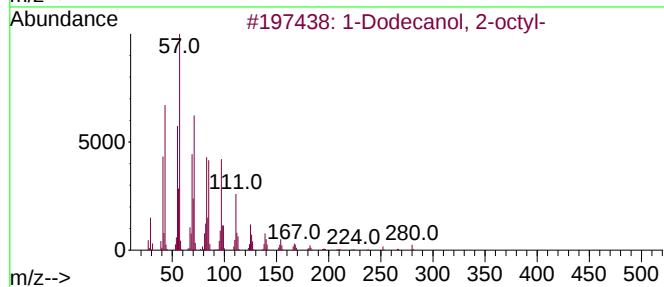
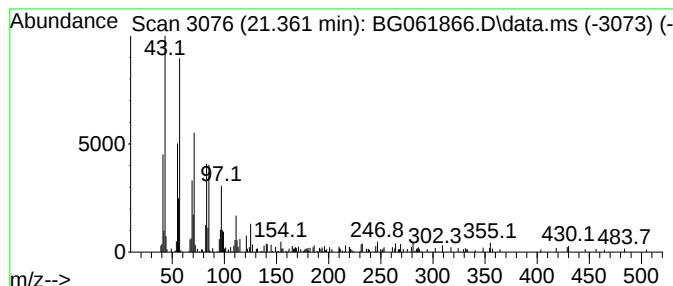
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Dodecanol, 2-octyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.361	3.10 ng/ul	140108	Chrysene-d12	21.719

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Dodecanol, 2-octyl-			298	C20H42O	005333-42-6	90
2	1-Hexacosene			364	C26H52	018835-33-1	90
3	Carbonic acid, decyl octadecyl e...			454	C29H58O3	1000383-16-7	87
4	1-Decanol, 2-octyl-			270	C18H38O	045235-48-1	87
5	Oxalic acid, cyclobutyl pentadec...			354	C21H38O4	1010309-70-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070324\
Data File : BG061866.D
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Sample : P2964-13
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-BG070224.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
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unknown-01	5.162	4.1	ng/ul	59552	1	8.094	293366	20.0
Pentanedioic ac...	9.810	2.3	ng/ul	52896	2	10.908	458328	20.0
3-Phenoxypropio...	11.643	2.8	ng/ul	64889	2	10.908	458328	20.0
1,1'-Biphenyl, ...	18.053	3.8	ng/ul	173393	4	17.459	923438	20.0
1,1'-Biphenyl, ...	18.517	3.3	ng/ul	152949	4	17.459	923438	20.0
2,2',4,5-Tetrac...	18.799	2.2	ng/ul	101451	4	17.459	923438	20.0
1-Dodecanol, 2-...	21.361	3.1	ng/ul	140108	5	21.719	904987	20.0