

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
 Data File : BP006358.D
 Acq On : 27 Jul 2021 10:51
 Operator : CG/JU
 Sample : M3063-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AT8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP006358.D

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.299	33	36	44	rVB	53890	70757	0.25%	0.053%
2	3.705	101	105	118	rBV	360653	568579	2.00%	0.428%
3	4.022	153	159	168	rVB	624992	872981	3.07%	0.657%
4	5.811	456	463	468	rBV	64798	102198	0.36%	0.077%
5	6.958	652	658	666	rBV2	124457	209732	0.74%	0.158%
6	7.128	680	687	693	rBV	885017	1380105	4.86%	1.039%
7	7.175	693	695	703	rVB	52786	77379	0.27%	0.058%
8	7.275	705	712	715	rBV2	44827	81045	0.29%	0.061%
9	7.316	715	719	729	rVB	449000	689158	2.42%	0.519%
10	7.510	746	752	763	rBV3	35803	97836	0.34%	0.074%
11	7.787	788	799	805	rVV	829467	1390324	4.89%	1.047%
12	7.852	805	810	814	rVV	188021	290208	1.02%	0.219%
13	7.899	814	818	824	rVB	109564	169960	0.60%	0.128%
14	8.152	855	861	868	rBV	155924	253278	0.89%	0.191%
15	8.322	886	890	901	rVB3	29439	56114	0.20%	0.042%
16	8.487	912	918	927	rVV	463844	744349	2.62%	0.561%
17	8.628	931	942	948	rVB2	118071	252999	0.89%	0.191%
18	8.887	981	986	989	rBV	64926	95628	0.34%	0.072%
19	8.940	989	995	1006	rVB	1035952	1757117	6.18%	1.323%
20	9.063	1011	1016	1021	rVV	64596	97494	0.34%	0.073%
21	9.228	1036	1044	1055	rBV6	30469	88864	0.31%	0.067%
22	9.657	1110	1117	1130	rVB	380021	622739	2.19%	0.469%
23	9.822	1141	1145	1153	rVB	42426	66162	0.23%	0.050%
24	9.969	1165	1170	1178	rBV2	72462	151141	0.53%	0.114%
25	10.187	1201	1207	1216	rVB	631072	1035223	3.64%	0.780%
26	10.281	1216	1223	1229	rVB3	30796	67939	0.24%	0.051%
27	10.357	1230	1236	1244	rBV	551794	884519	3.11%	0.666%
28	10.493	1255	1259	1263	rVV3	43709	73356	0.26%	0.055%
29	10.569	1265	1272	1278	rVV	1165441	1926788	6.78%	1.451%
30	10.628	1278	1282	1287	rVB5	35400	62000	0.22%	0.047%
31	10.710	1287	1296	1311	rVB2	987270	2069910	7.28%	1.559%
32	11.181	1373	1376	1380	rVV2	42531	73395	0.26%	0.055%
33	11.681	1455	1461	1467	rBV	113221	195334	0.69%	0.147%
34	11.969	1504	1510	1516	rVB4	81922	184398	0.65%	0.139%
35	12.098	1523	1532	1544	rBV2	356466	803877	2.83%	0.605%
36	12.451	1585	1592	1598	rVV3	130107	233922	0.82%	0.176%

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Smoothing : OFF

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
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37	12.540	1602	1607	1613	rVV5	28476	75339	0.27%	0.057%
38	12.598	1613	1617	1623	rVB6	33789	68808	0.24%	0.052%
39	12.840	1650	1658	1662	rVV3	194308	389044	1.37%	0.293%
40	12.904	1665	1669	1674	rVB	165077	241983	0.85%	0.182%

41	13.116	1696	1705	1709	rBV	506355	801526	2.82%	0.604%
42	13.245	1719	1727	1732	rVB2	129233	218432	0.77%	0.164%
43	13.322	1737	1740	1744	rVB2	37286	56170	0.20%	0.042%
44	13.381	1744	1750	1753	rBV	150497	230069	0.81%	0.173%
45	13.428	1753	1758	1762	rBV	384539	535239	1.88%	0.403%

46	13.545	1772	1778	1782	rBV	198460	332454	1.17%	0.250%
47	13.581	1782	1784	1789	rVB	89944	123751	0.44%	0.093%
48	13.834	1821	1827	1833	rBV	2512073	3462364	12.18%	2.607%
49	14.022	1856	1859	1864	rVB4	49138	64808	0.23%	0.049%
50	14.104	1866	1873	1884	rBV	2739380	4006834	14.10%	3.017%

51	14.187	1884	1887	1889	rBV2	63226	75490	0.27%	0.057%
52	14.304	1904	1907	1914	rVB2	37743	58371	0.21%	0.044%
53	14.410	1919	1925	1934	rVV2	1698569	2648551	9.32%	1.995%
54	14.557	1945	1950	1951	rVV4	52545	82584	0.29%	0.062%
55	14.616	1951	1960	1967	rVB8	146641	489827	1.72%	0.369%

56	14.739	1977	1981	1984	rBV3	44689	66261	0.23%	0.050%
57	14.886	2002	2006	2011	rBV	64293	95098	0.33%	0.072%
58	14.951	2011	2017	2021	rBV2	217994	365126	1.28%	0.275%
59	14.992	2021	2024	2027	rVV	156478	223471	0.79%	0.168%
60	15.034	2027	2031	2038	rVB2	719857	1068169	3.76%	0.804%

61	15.157	2048	2052	2056	rVB4	64190	88849	0.31%	0.067%
62	15.245	2061	2067	2073	rVB4	127053	230329	0.81%	0.173%
63	15.339	2081	2083	2088	rVB5	48561	56443	0.20%	0.043%
64	15.410	2088	2095	2101	rBV	3447967	5103041	17.95%	3.843%
65	15.463	2101	2104	2108	rVB2	142044	186383	0.66%	0.140%

66	15.539	2109	2117	2124	rBV	3042547	4884284	17.18%	3.678%
67	15.622	2127	2131	2136	rVV3	91316	154167	0.54%	0.116%
68	15.675	2137	2140	2149	rVB3	158078	248569	0.87%	0.187%
69	15.751	2149	2153	2161	rBV7	72693	163290	0.57%	0.123%
70	16.222	2229	2233	2238	rBV3	78614	142823	0.50%	0.108%

71	16.445	2267	2271	2276	rBV	103352	152988	0.54%	0.115%
72	16.516	2280	2283	2287	rVB2	83651	120564	0.42%	0.091%
73	16.675	2306	2310	2314	rBV3	53623	91283	0.32%	0.069%
74	16.816	2327	2334	2347	rBV	19011892	28426192	100.00%	21.407%
75	16.933	2350	2354	2360	rBV6	99788	178598	0.63%	0.134%

76	17.045	2369	2373	2380	rVB	195833	284169	1.00%	0.214%
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 Title : SVOA CALIBRATION

77	17.163	2388	2393	2399	rVB	1872239	2694443	9.48%	2.029%
78	17.263	2404	2410	2422	rBV2	4139282	5921498	20.83%	4.459%
79	17.369	2423	2428	2438	rVB	489282	770560	2.71%	0.580%
80	17.539	2451	2457	2467	rVB	2111537	2852952	10.04%	2.148%
81	17.845	2505	2509	2516	rVB2	263764	379569	1.34%	0.286%
82	18.069	2543	2547	2557	rBV2	559201	847965	2.98%	0.639%
83	18.210	2568	2571	2580	rVB6	66141	121594	0.43%	0.092%
84	19.080	2716	2719	2723	rVB3	81042	105520	0.37%	0.079%
85	19.151	2728	2731	2733	rBV	47022	69855	0.25%	0.053%
86	19.257	2744	2749	2752	rBV	11871254	13646667	48.01%	10.277%
87	19.310	2755	2758	2767	rVB	731389	884743	3.11%	0.666%
88	19.398	2768	2773	2781	rVV	3419395	4114212	14.47%	3.098%
89	19.504	2786	2791	2798	rVV	5051769	6429128	22.62%	4.842%
90	20.080	2885	2889	2896	rBV	701333	819366	2.88%	0.617%
91	20.227	2910	2914	2918	rBV	744053	808806	2.85%	0.609%
92	20.851	3016	3020	3028	rBV	1987981	2273180	8.00%	1.712%
93	20.992	3040	3044	3047	rBV	272291	336772	1.18%	0.254%
94	21.210	3078	3081	3082	rBV	152351	165250	0.58%	0.124%
95	21.263	3085	3090	3095	rVB	2571321	3310384	11.65%	2.493%
96	22.145	3236	3240	3252	rVB	792704	1158875	4.08%	0.873%
97	22.498	3296	3300	3307	rVB	473930	726863	2.56%	0.547%
98	23.457	3456	3463	3472	rVB2	3389024	6590337	23.18%	4.963%
99	23.610	3482	3489	3500	rVB2	1557939	3182055	11.19%	2.396%
100	23.780	3514	3518	3526	rVB	260153	489019	1.72%	0.368%

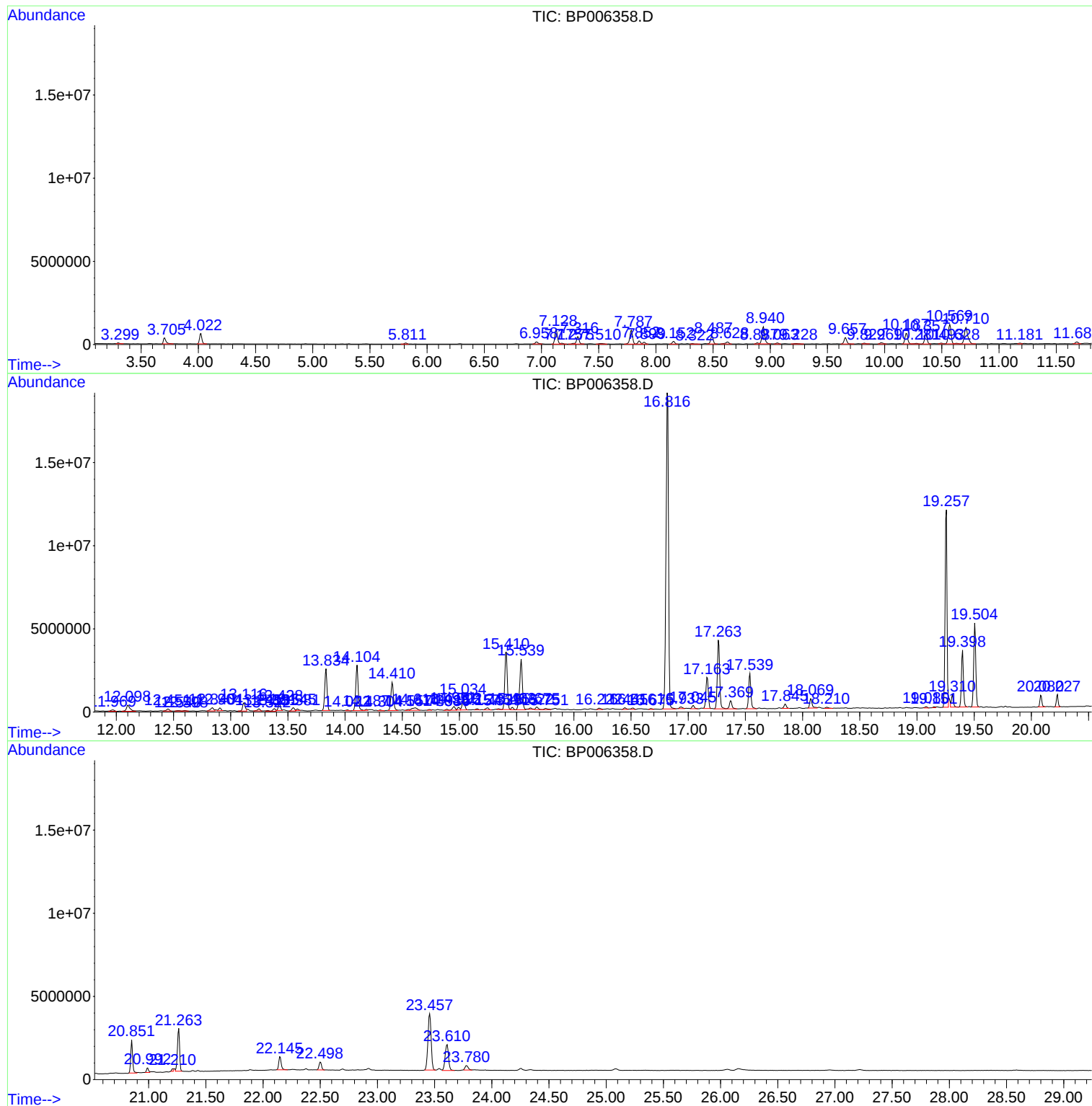
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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



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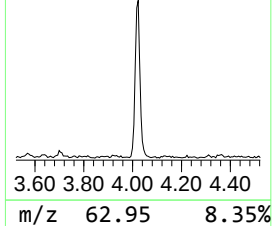
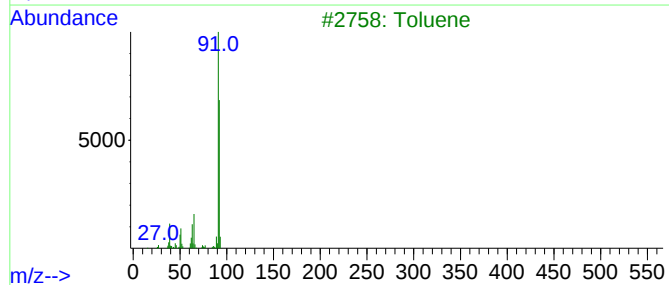
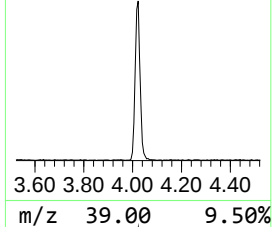
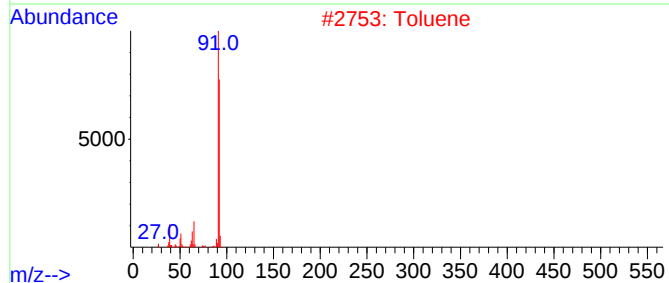
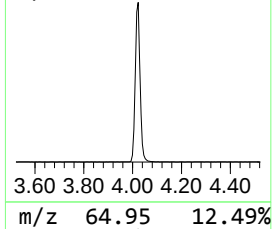
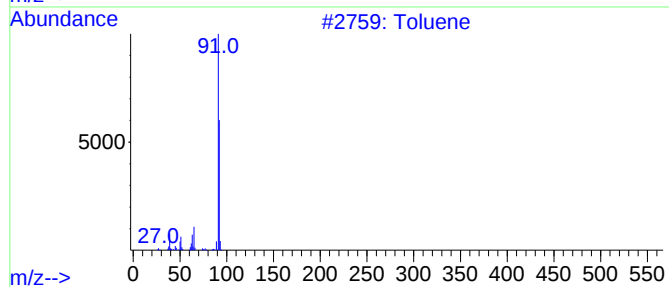
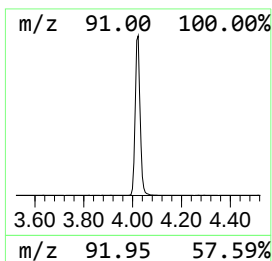
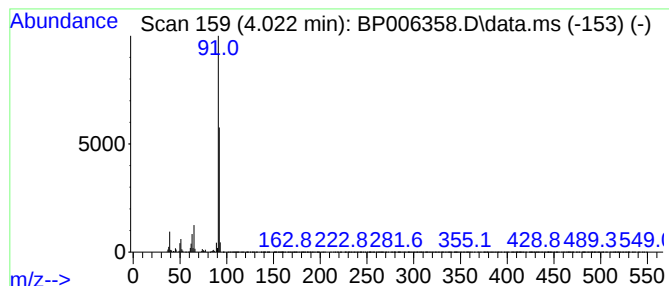
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.020	12.56 ng/ul	872981	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	95
3	Toluene			92	C7H8	000108-88-3	93
4	Toluene			92	C7H8	000108-88-3	91
5	Toluene			92	C7H8	000108-88-3	90



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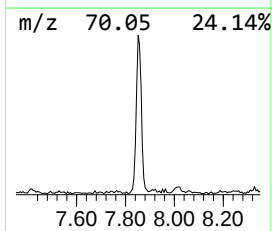
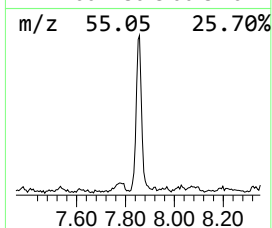
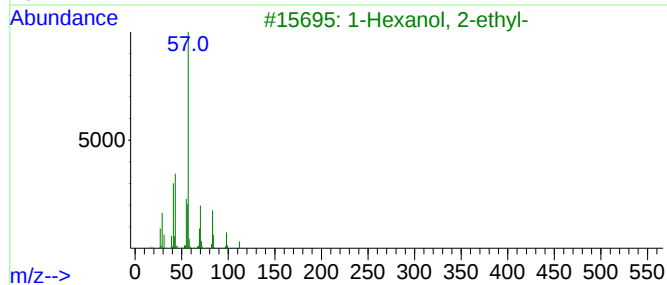
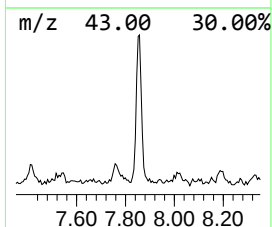
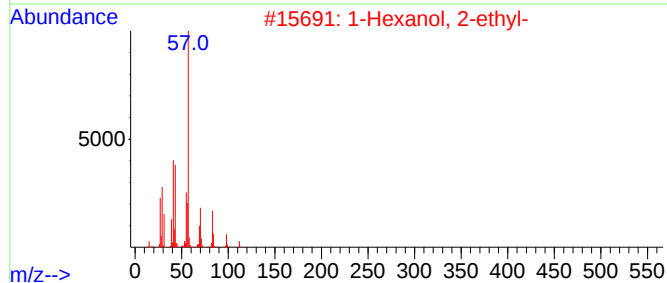
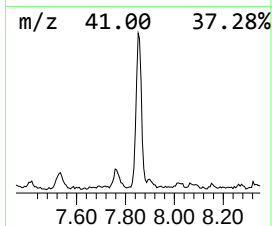
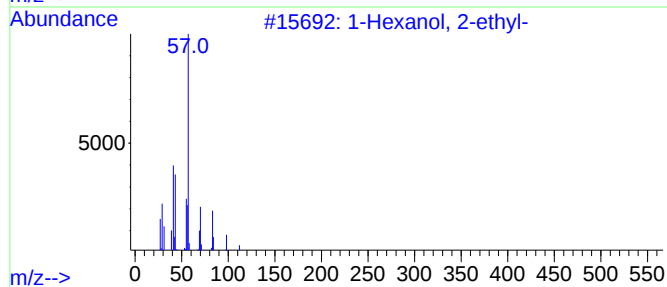
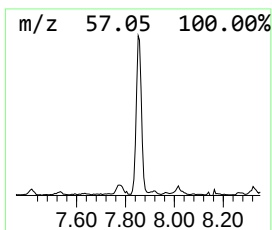
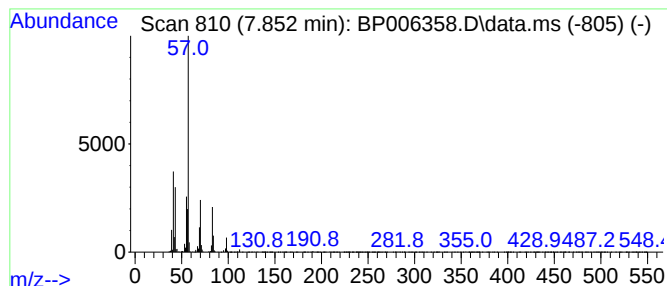
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Hexanol, 2-ethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.850	4.17 ng/ul	290208	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	83
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	64
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64



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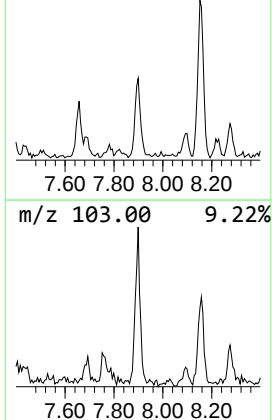
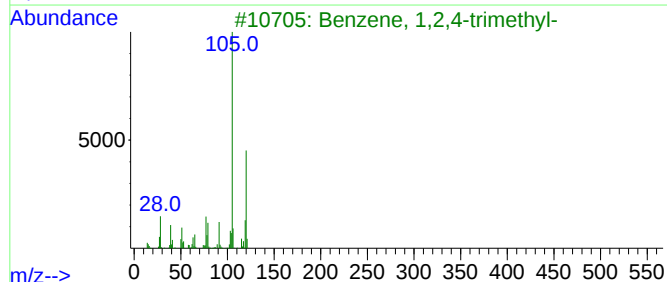
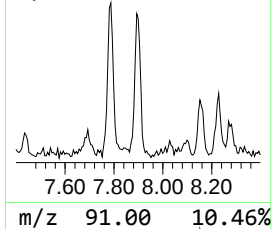
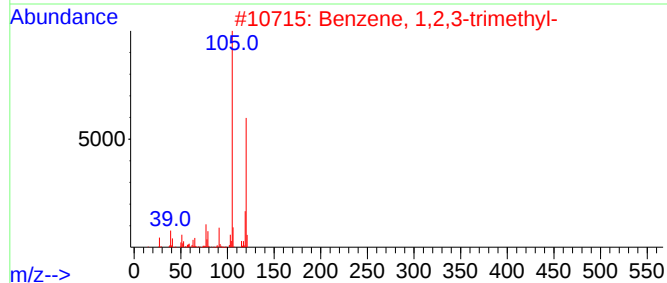
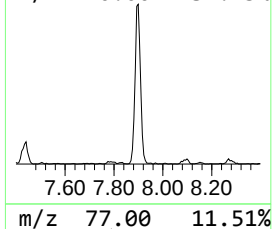
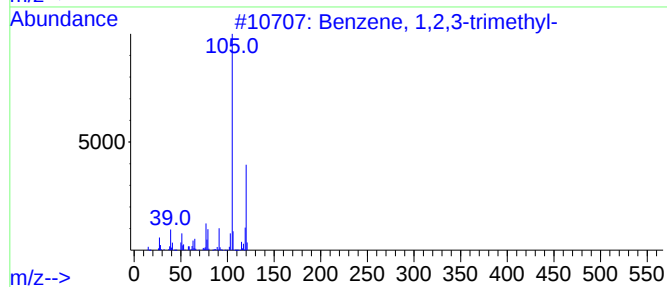
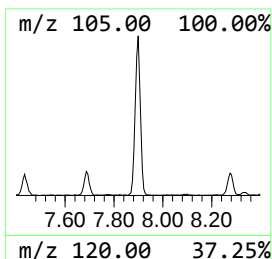
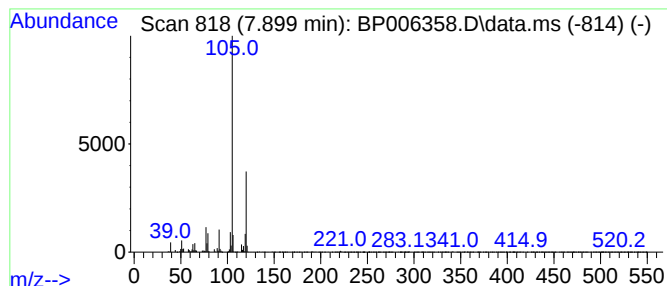
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2,3-trimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.900	2.44 ng/ul	169960	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
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Instrument :
BNA_P
ClientSampleId :
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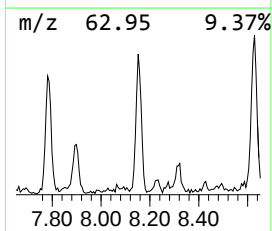
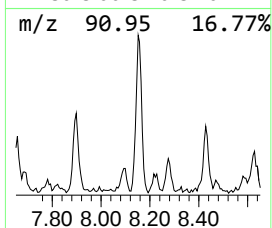
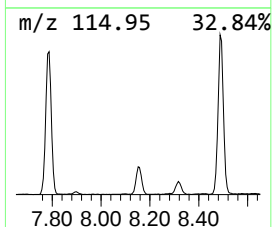
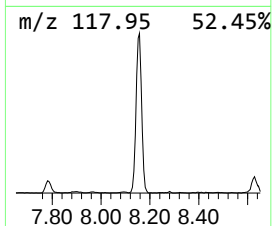
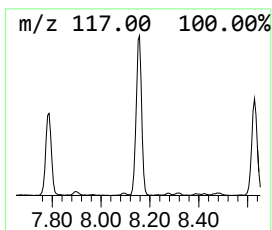
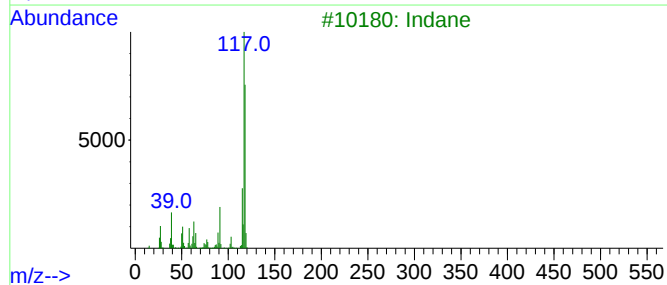
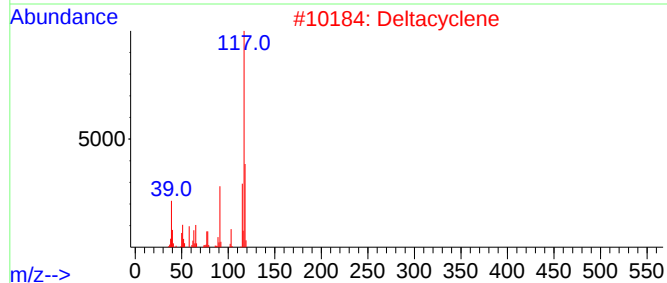
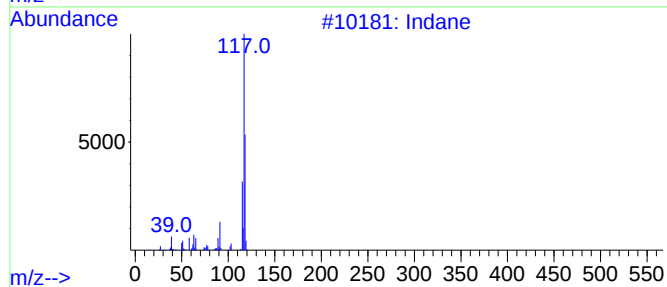
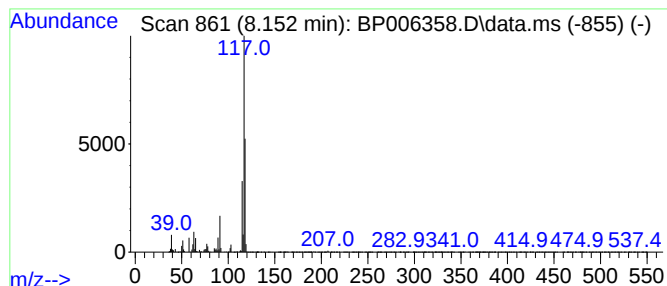
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Indane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.150	3.64 ng/ul	253278	1,4-Dichlorobenzene-d4	7.787

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	90
2			Deltacyclene	118	C9H10	007785-10-6	68
3			Indane	118	C9H10	000496-11-7	64
4			Indane	118	C9H10	000496-11-7	64
5			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
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Sample : M3063-01
Misc :
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Instrument :
BNA_P
ClientSampleId :
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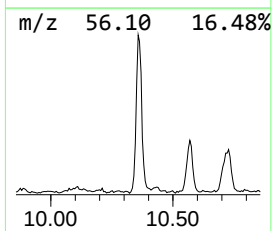
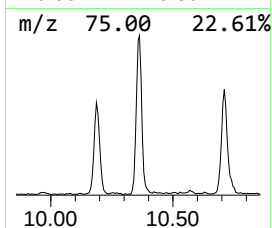
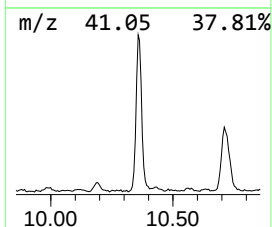
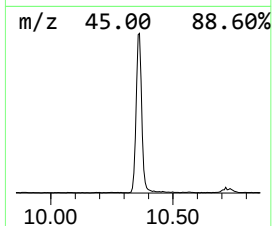
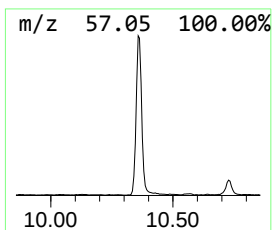
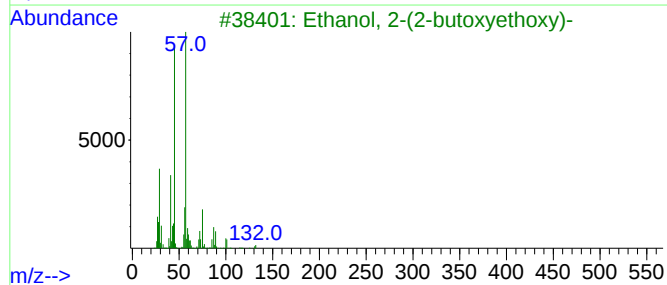
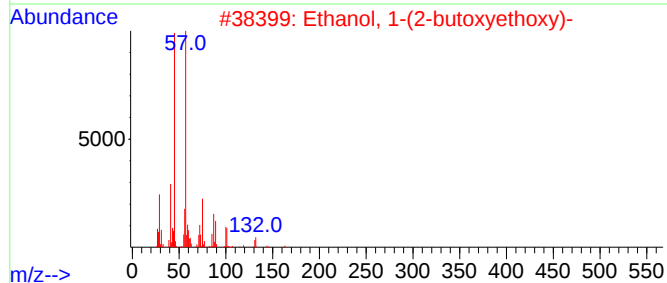
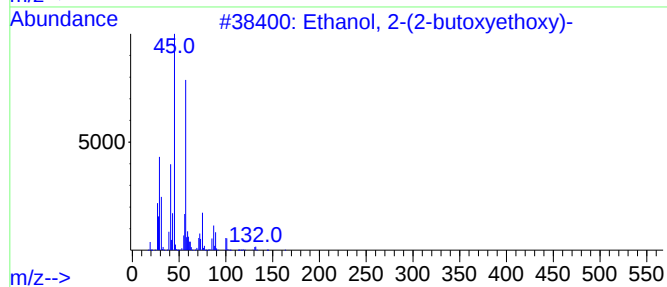
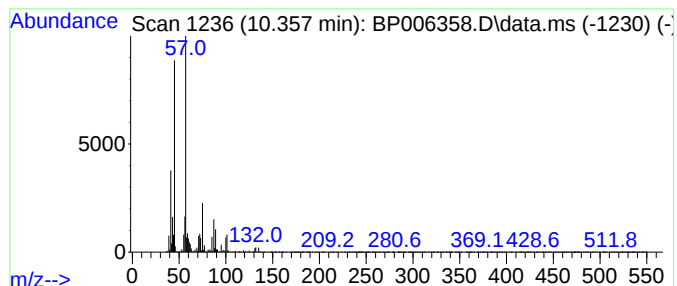
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	9.18 ng/ul	884519	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	90
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
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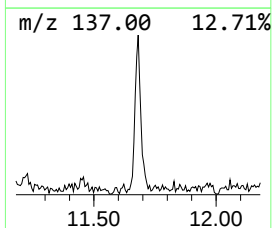
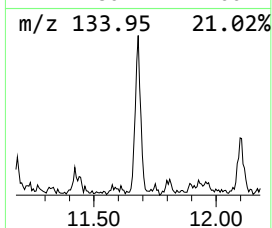
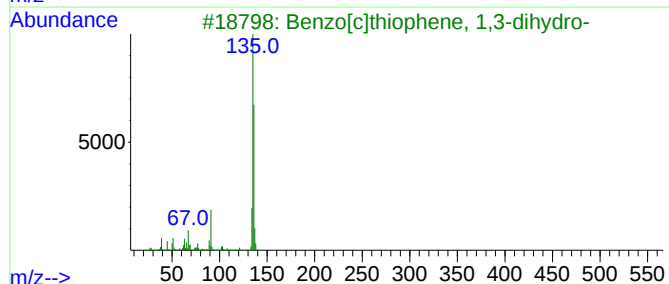
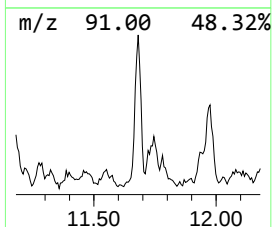
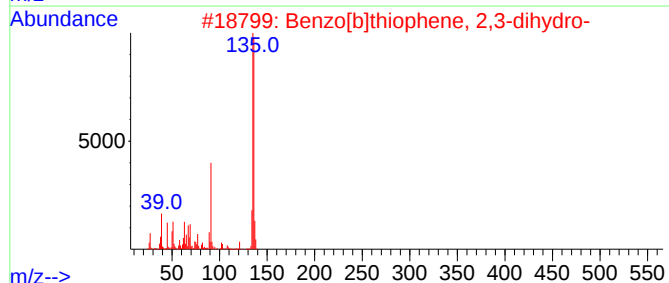
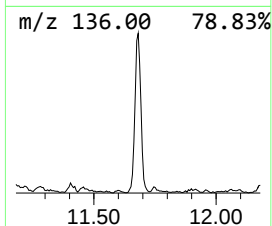
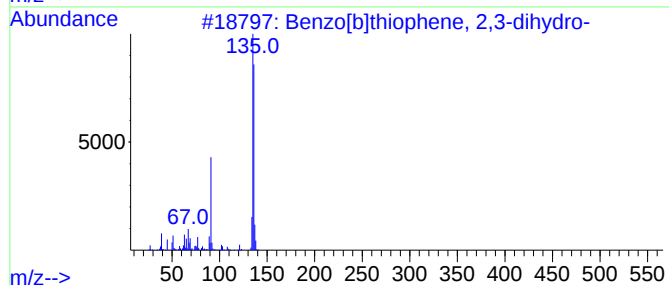
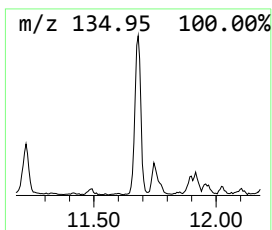
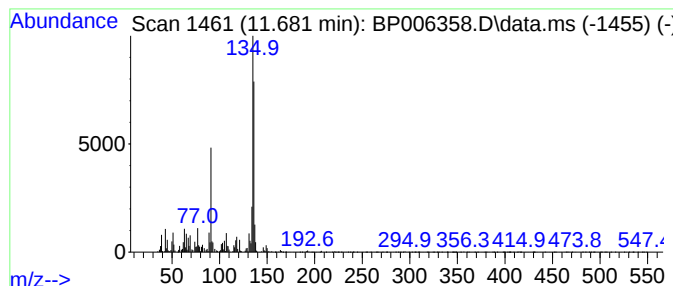
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.680	2.03 ng/ul	195334	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	95
2			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	87
3			Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	87
4			Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	83
5			Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
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Instrument :
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ClientSampleId :
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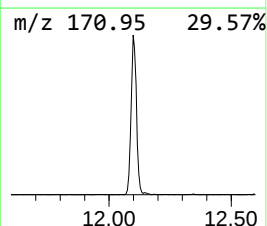
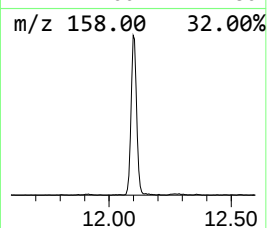
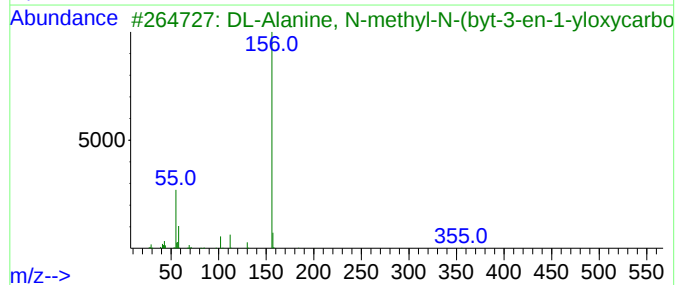
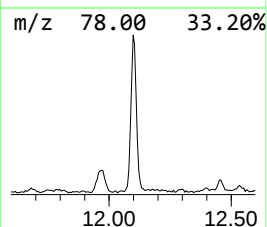
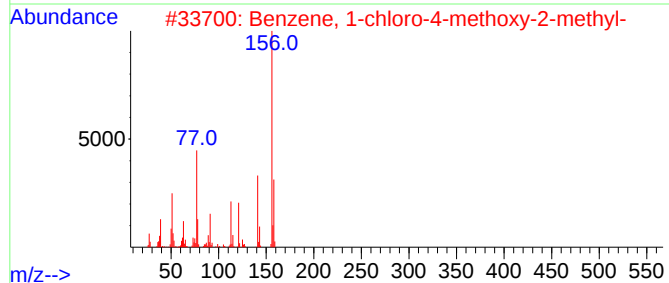
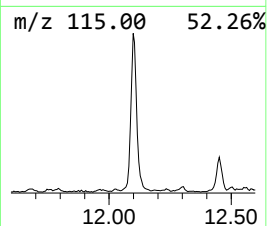
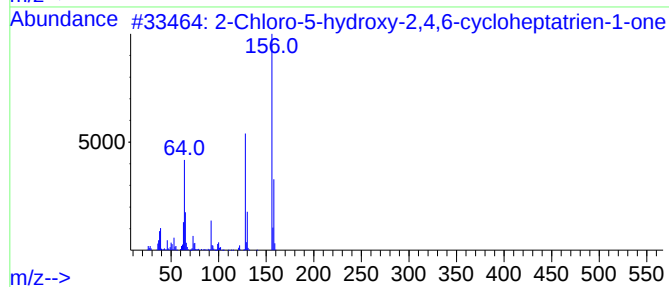
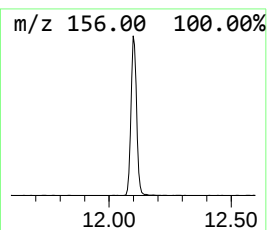
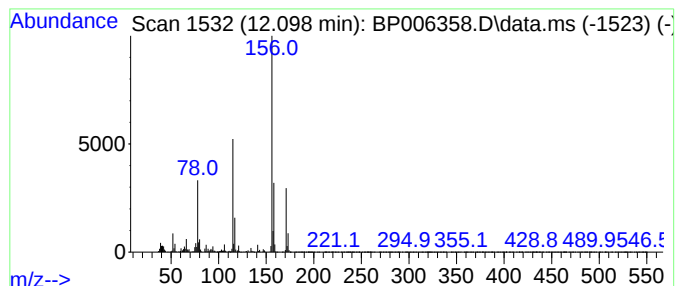
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.100	8.34 ng/ul	803877	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	38
2			Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	32
3			DL-Alanine, N-methyl-N-(byt-3-en...	355	C20H37NO4	1000392-73-6	27
4			Tebuthiuron	228	C9H16N4OS	034014-18-1	25
5			4-Chloro-3,5-dimethylphenol, O-(...	242	C12H15ClO3	1000487-35-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
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Sample : M3063-01
Misc :
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Instrument :
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ClientSampleId :
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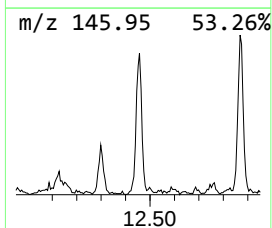
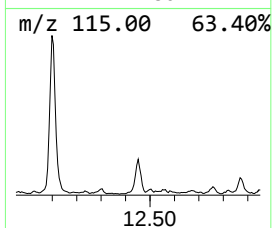
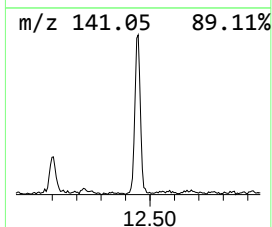
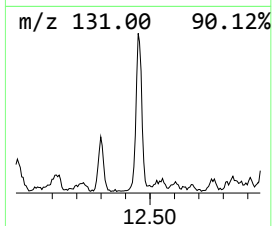
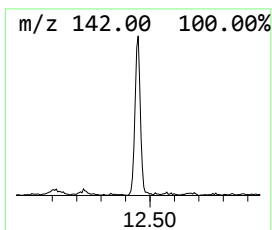
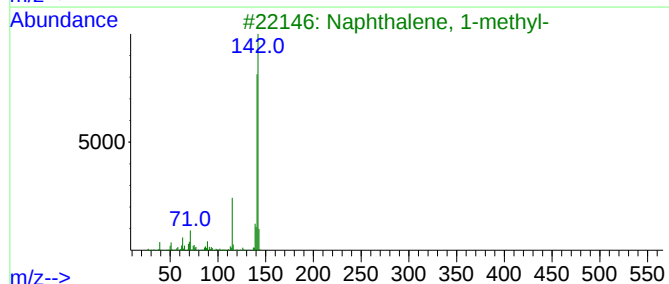
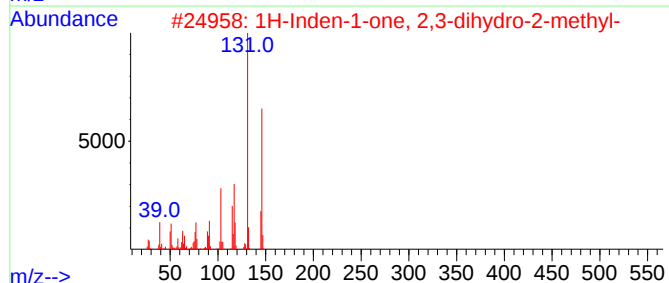
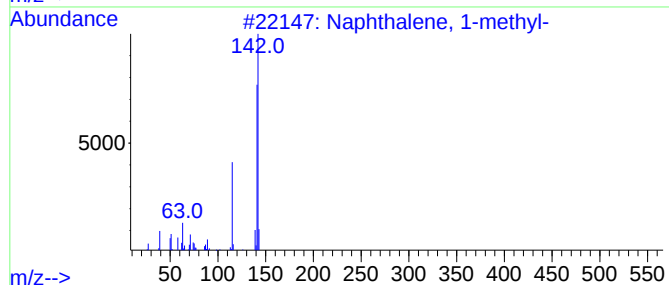
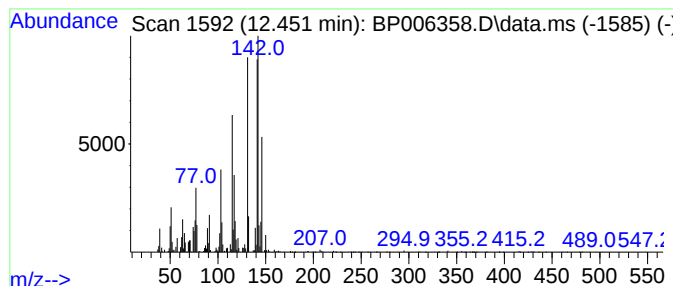
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.450	2.43 ng/ul	233922	Naphthalene-d8	10.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	83
2			1H-Inden-1-one, 2,3-dihydro-2-methyl-	146	C10H10O	017496-14-9	50
3			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	50
4			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	50
5			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
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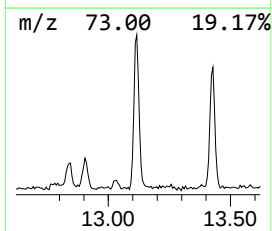
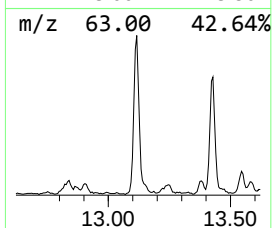
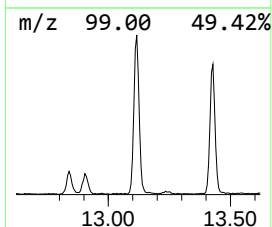
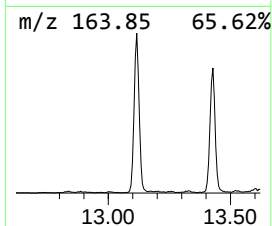
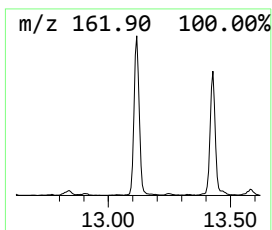
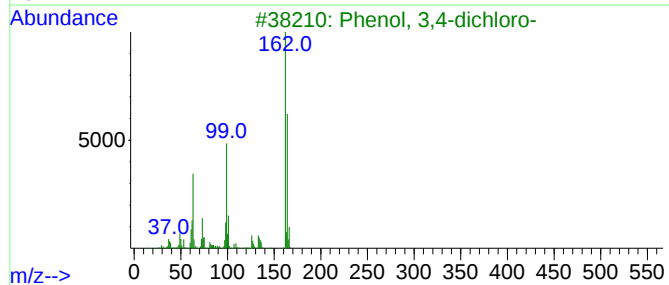
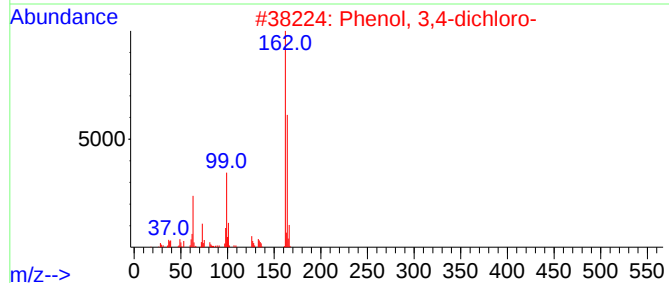
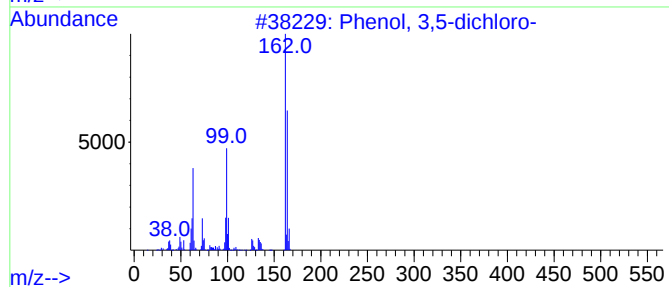
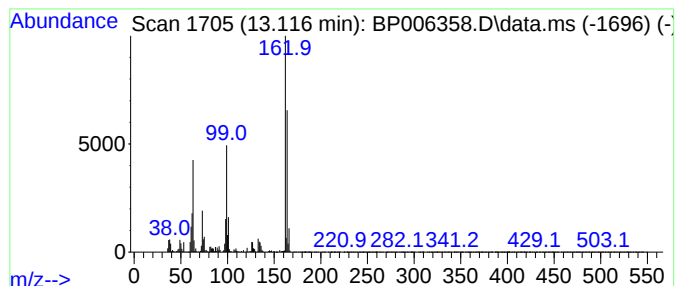
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 3,5-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.	
13.120	6.05 ng/ul	801526	Acenaphthene-d10			14.410	
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol,	3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98	
2	Phenol,	3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97	
3	Phenol,	3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97	
4	Phenol,	3,5-dichloro-	162	C6H4Cl2O	000591-35-5	96	
5	Phenol,	3,4-dichloro-	162	C6H4Cl2O	000095-77-2	96	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
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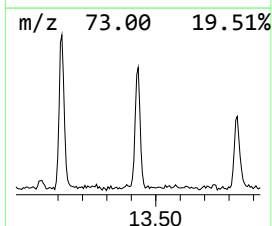
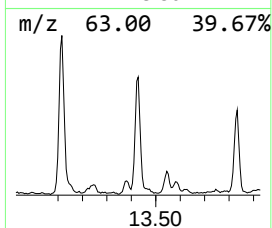
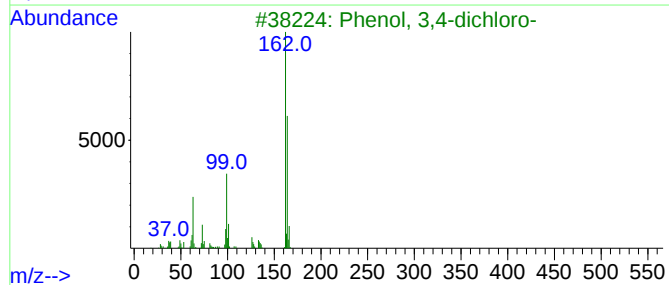
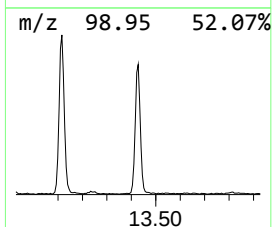
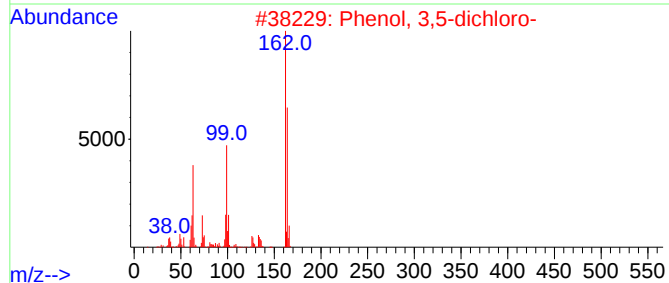
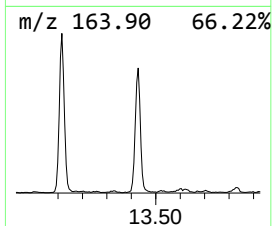
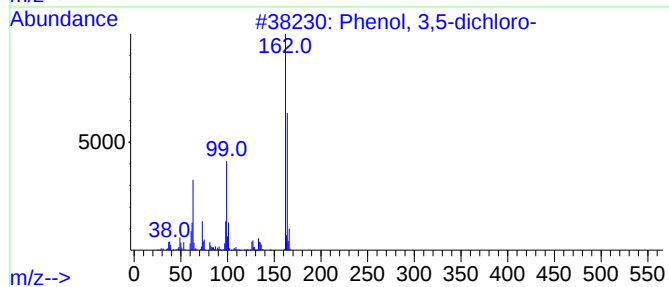
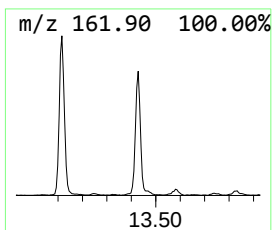
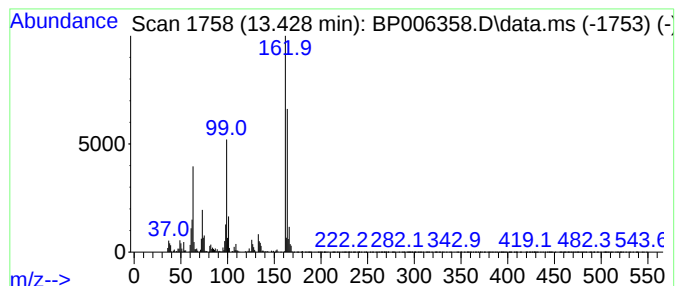
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phenol, 3,4-dichloro- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.430	4.04 ng/ul	535239	Acenaphthene-d10	14.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
5			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

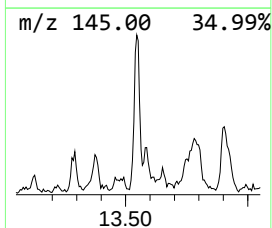
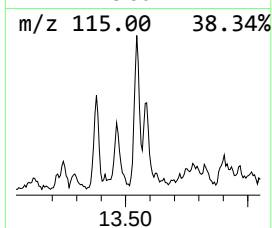
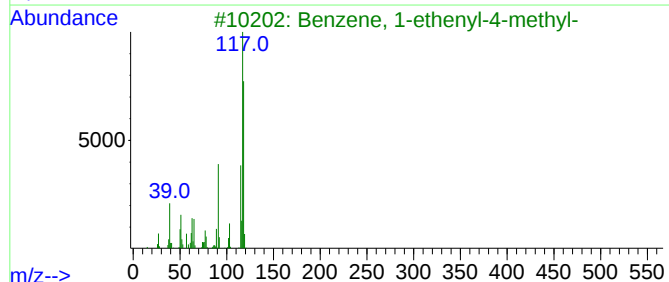
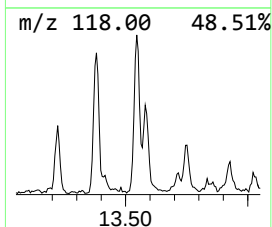
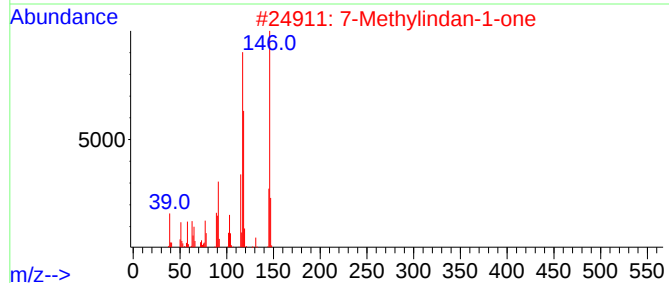
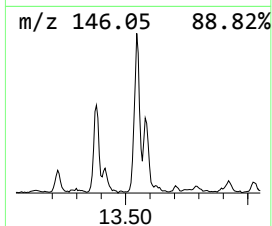
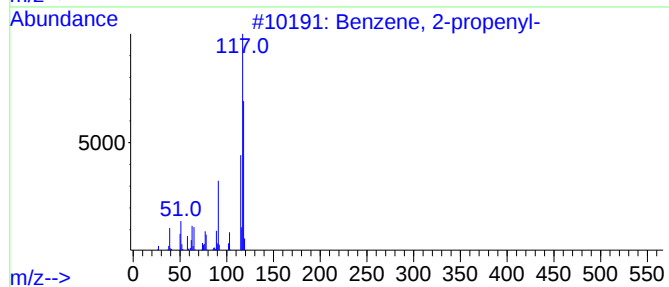
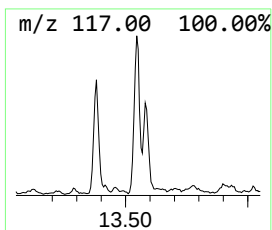
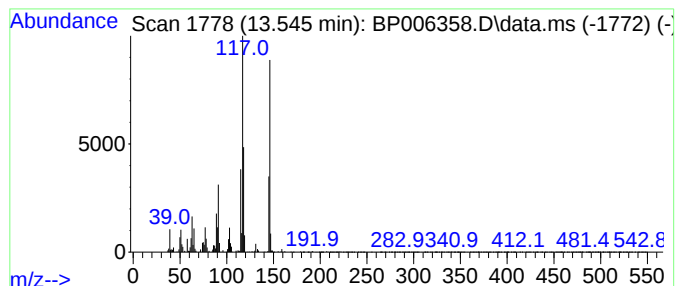
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 2-propenyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.550	2.51 ng/ul	332454	Acenaphthene-d10	14.410	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
2	7-Methylindan-1-one	146	C10H10O	039627-61-7	76
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	64
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	52
5	Indane	118	C9H10	000496-11-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
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Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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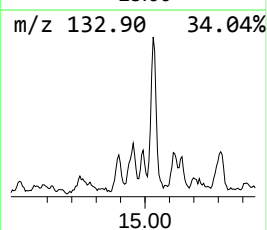
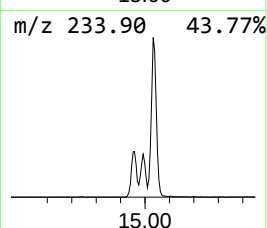
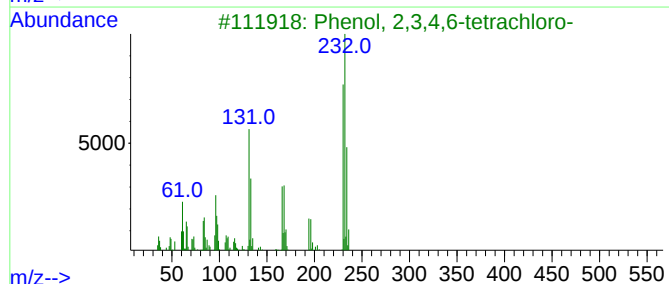
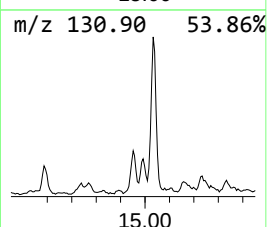
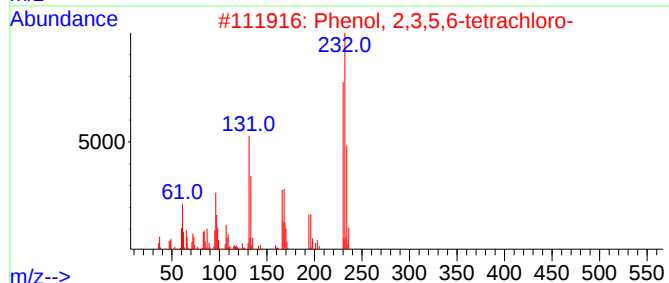
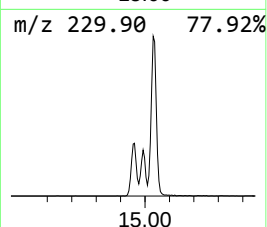
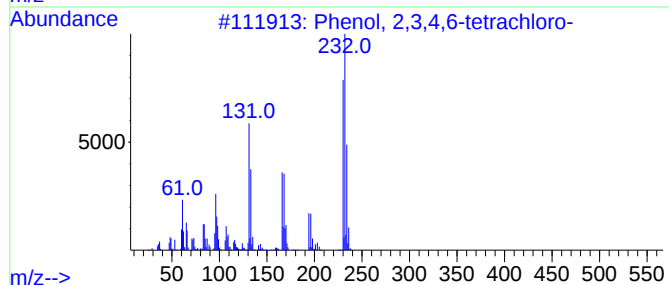
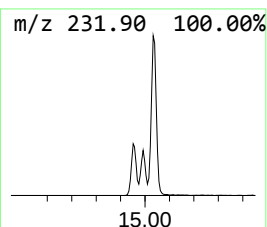
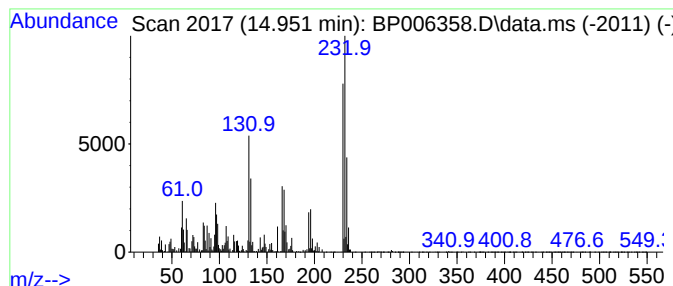
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Phenol, 2,3,4,6-tetrachloro- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.950	2.76 ng/ul	365126	Acenaphthene-d10	14.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
2			Phenol, 2,3,5,6-tetrachloro-	230	C6H2Cl4O	000935-95-5	99
3			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	99
4			Phenol, 2,3,4,6-tetrachloro-	230	C6H2Cl4O	000058-90-2	98
5			Phenol, 2,3,4,5-tetrachloro-	230	C6H2Cl4O	004901-51-3	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
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Acq On : 27 Jul 2021 10:51
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Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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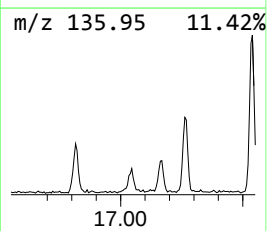
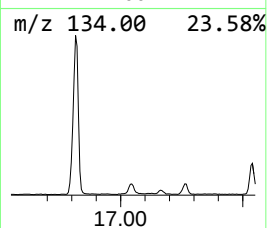
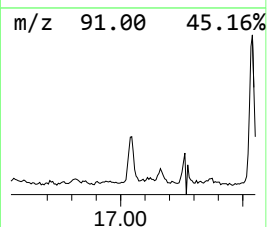
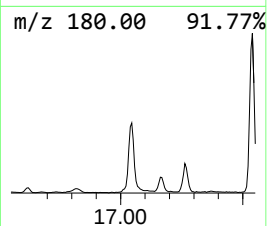
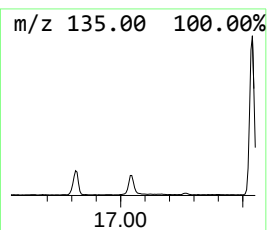
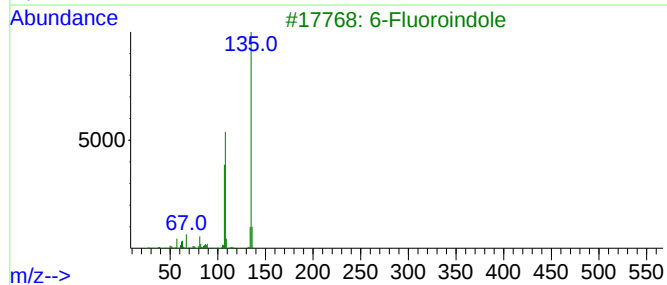
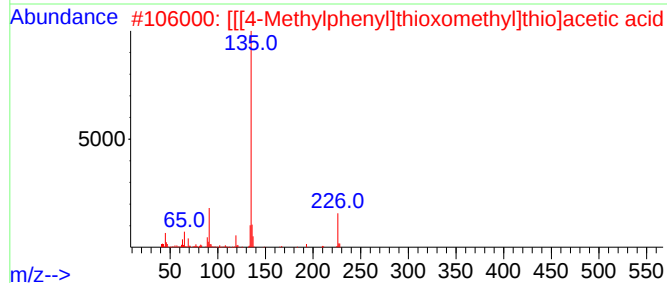
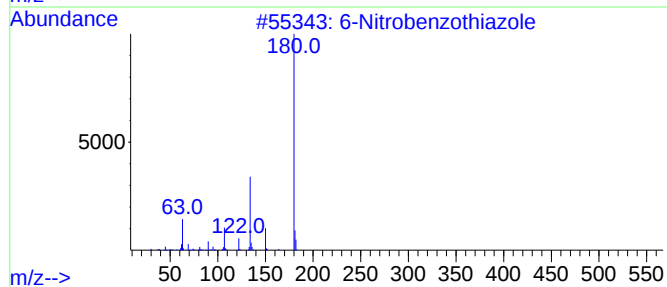
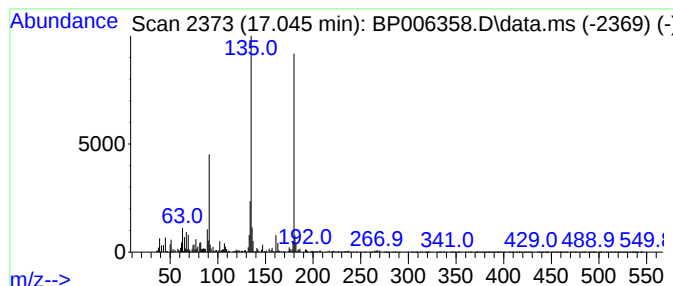
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-02 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.050	2.11 ng/ul	284169	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Nitrobenzothiazole	180	C7H4N2O2S	002942-06-5	43
2			[[[4-Methylphenyl]thioxomethyl]t...	226	C10H10O2S2	057149-19-6	38
3			6-Fluoroindole	135	C8H6FN	000399-51-9	35
4			Benzothiazole	135	C7H5NS	000095-16-9	30
5			1-Adamantanecarboxylic acid, 2-m...	270	C18H22O2	1010293-75-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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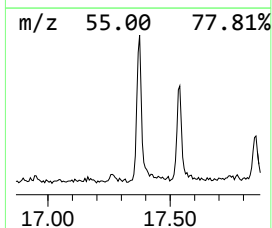
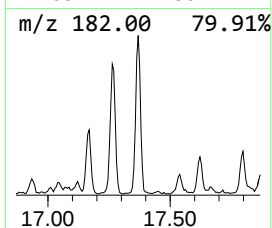
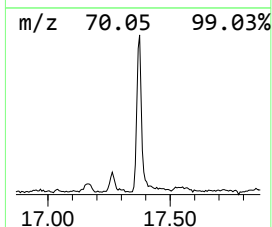
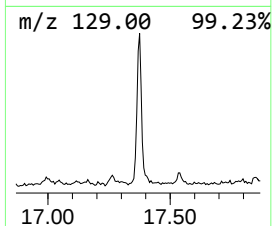
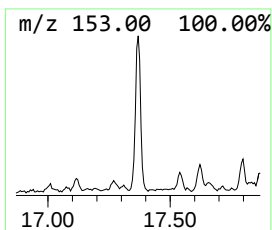
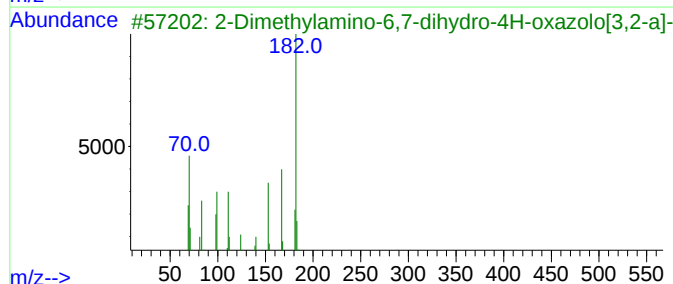
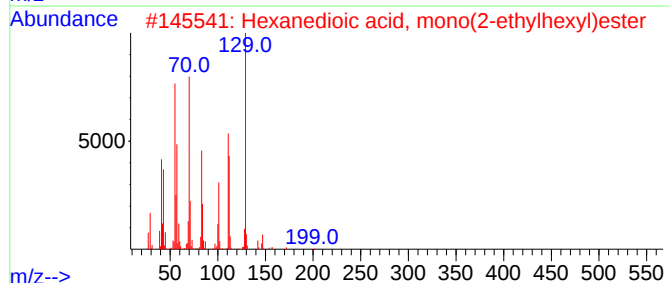
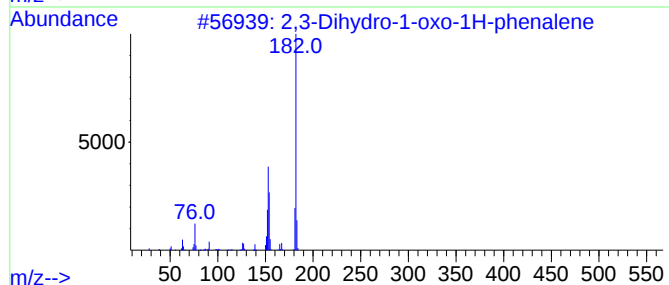
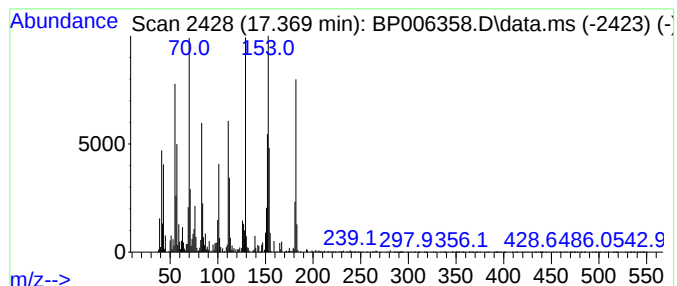
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-03 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.370	5.72 ng/ul	770560	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	46		
2	Hexanedioic acid, mono(2-ethylhe...	258	C14H26O4	004337-65-9	46		
3	2-Dimethylamino-6,7-dihydro-4H-o...	182	C7H10N4O2	062626-99-7	38		
4	5-Acenaphthylenecarboxaldehyde, ...	182	C13H10O	1000362-64-3	35		
5	5,10-Methanobenzocycloocten-11-o...	216	C13H9ClO	033655-73-1	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
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Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

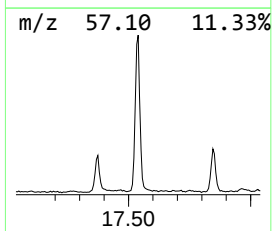
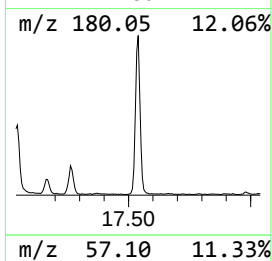
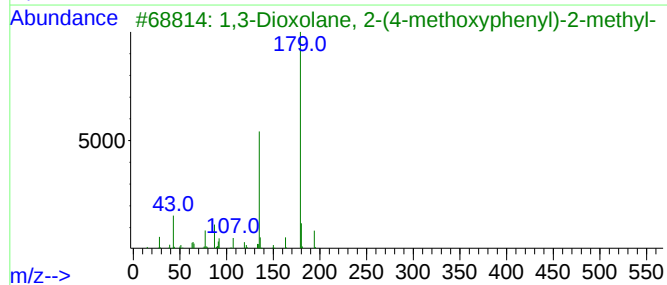
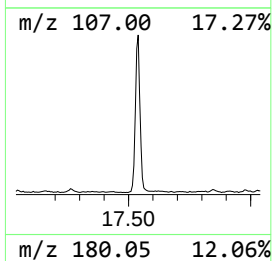
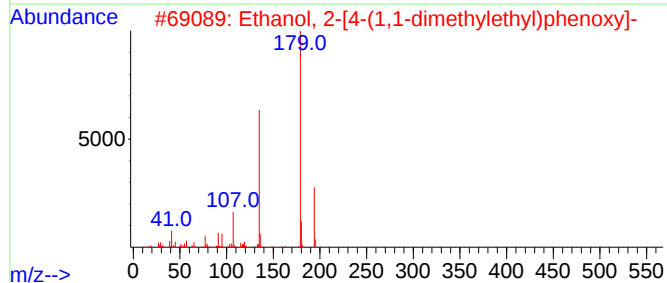
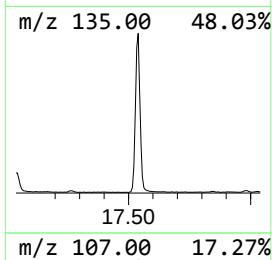
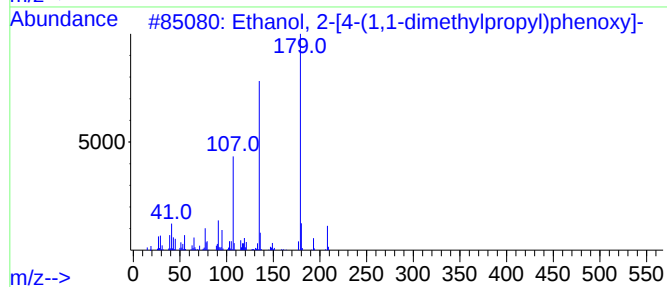
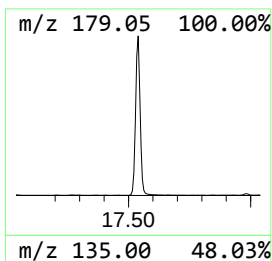
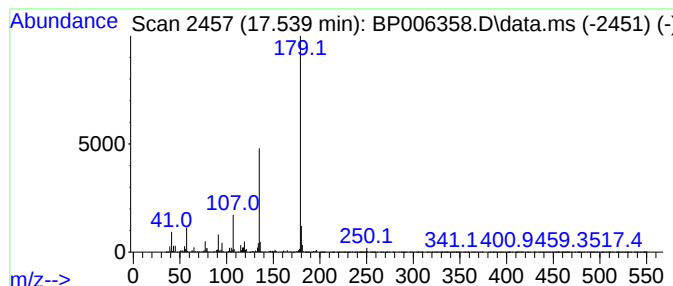
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Ethanol, 2-[4-(1,1-dimethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.540	21.18 ng/ul	2852950	Phenanthrene-d10	17.163	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol, 2-[4-(1,1-dimethylpropyl)phenoxy]-	208	C13H20O2	006382-07-6	78
2	Ethanol, 2-[4-(1,1-dimethylethyl)phenoxy]-	194	C12H18O2	000713-46-2	78
3	1,3-Dioxolane, 2-(4-methoxyphenyl)-2-methyl-	194	C11H14O3	036881-00-2	64
4	2-Mercapto-3-(trifluoromethyl)pyridine	179	C6H4F3NS	104040-74-6	38
5	Hydrobenzoin, 2TMS derivative	358	C20H30O2Si2	035479-44-8	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
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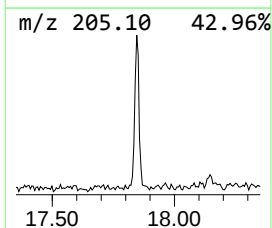
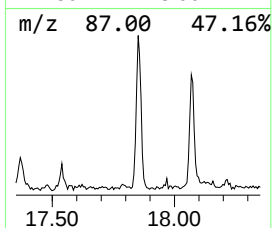
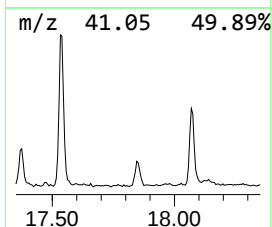
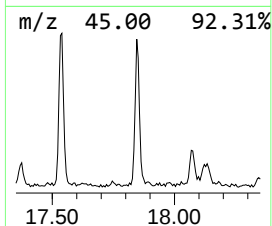
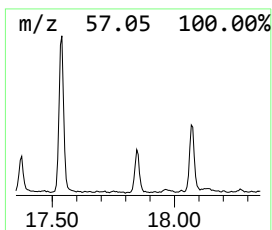
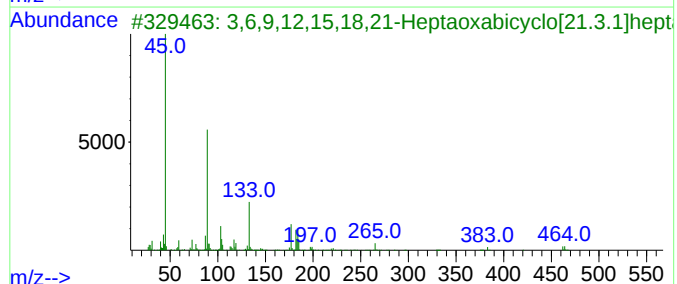
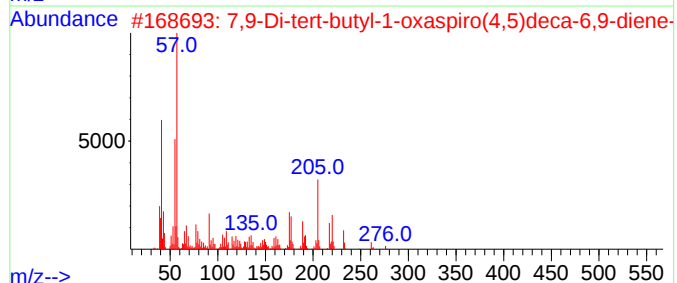
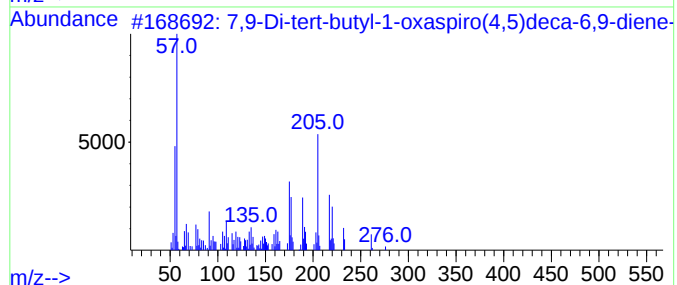
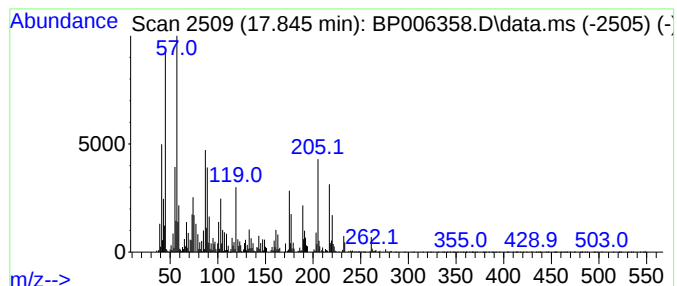
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.850	2.82 ng/ul	379569	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	95		
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	45		
3	3,6,9,12,15,18,21-Heptaoxabicycl...	462	C20H31BrO7	111982-23-1	27		
4	Heptaethylene glycol, TMS deriva...	398	C17H38O8Si	1000352-06-3	22		
5	Ethanol, 2-butoxy-	118	C6H14O2	000111-76-2	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT8

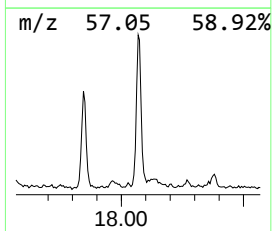
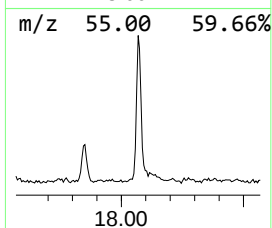
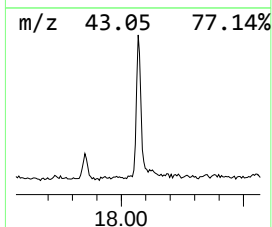
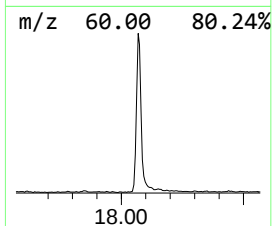
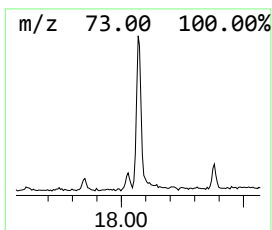
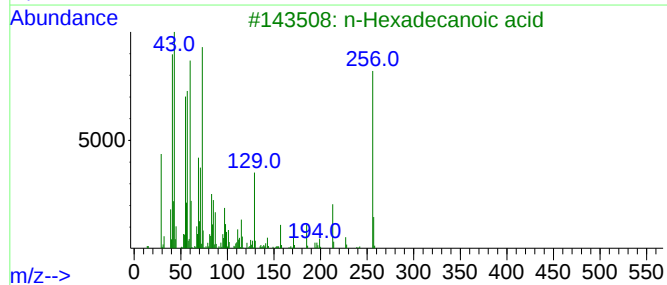
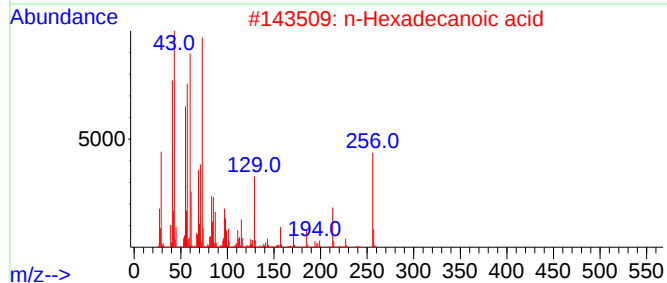
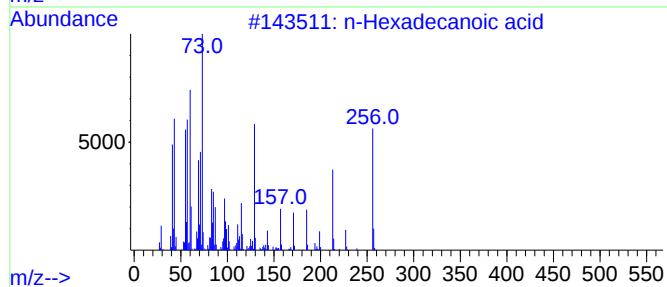
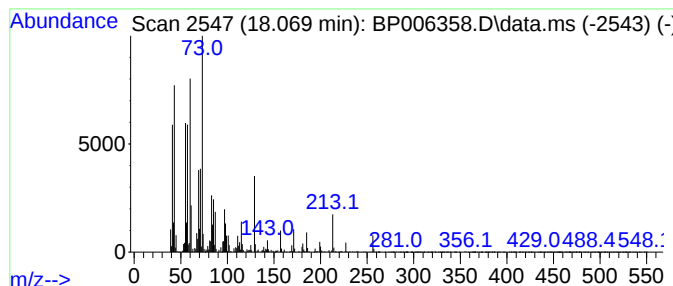
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.070	6.29 ng/ul	847965	Phenanthrene-d10	17.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	99
4	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid			256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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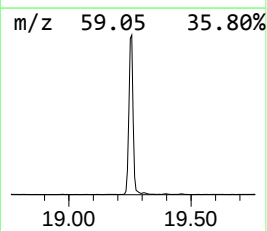
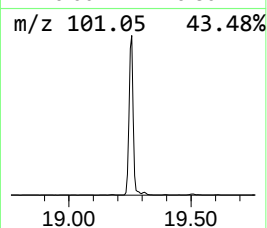
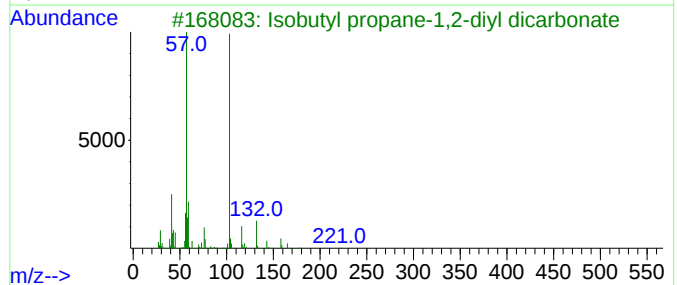
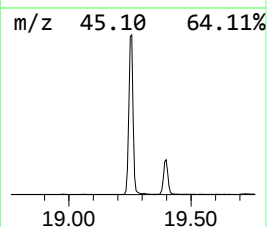
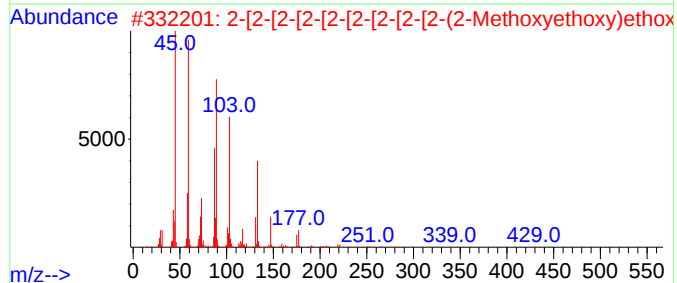
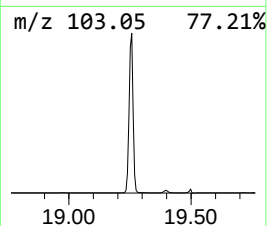
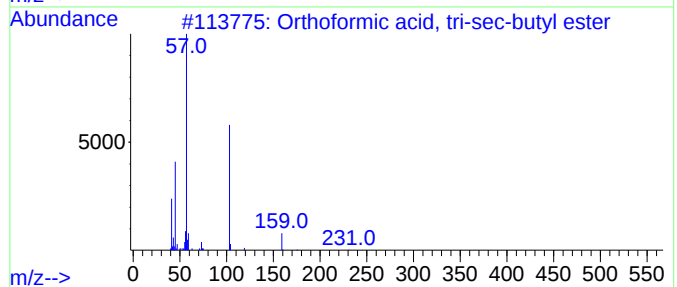
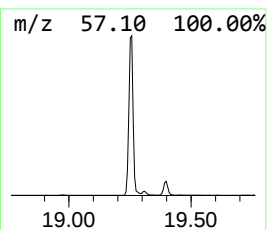
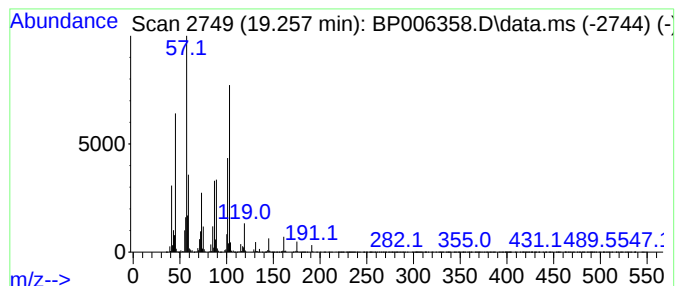
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.260	82.45 ng/ul	13646700	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Orthoformic acid, tri-sec-butyl ...	232	C13H28O3	016754-48-6	38
2			2-[2-[2-[2-[2-[2-[2-(2-Met...	472	C21H44O11	027425-92-9	30
3			Isobutyl propane-1,2-diyl dicarb...	276	C13H24O6	1000372-79-7	27
4			3,6,9,12,15-Pentaoxanonadecan-1-ol	294	C14H30O6	001786-94-3	27
5			2-[2-[2-[2-[2-[2-(2-Methoxyet...	384	C17H36O9	025990-96-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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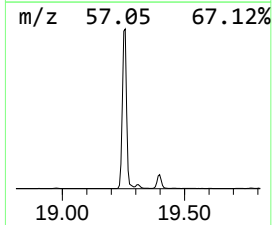
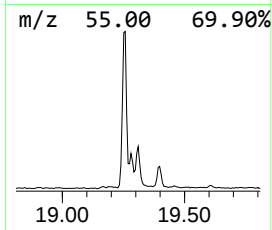
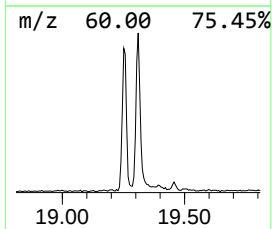
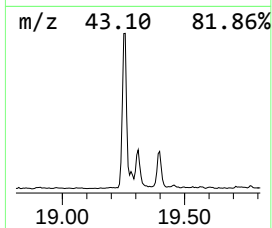
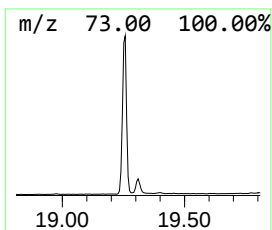
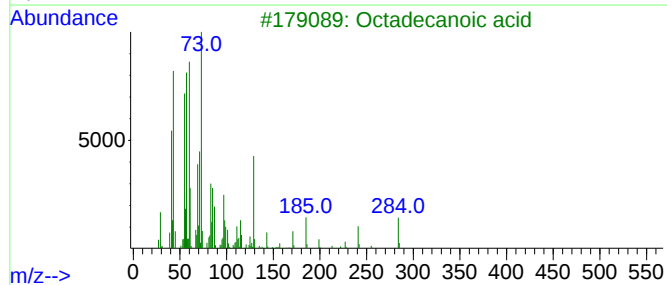
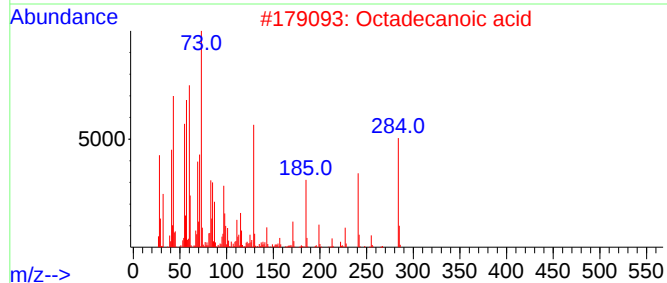
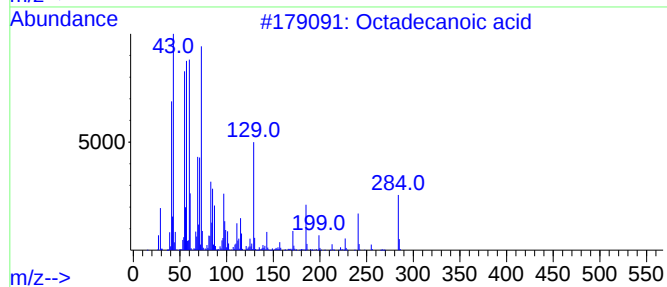
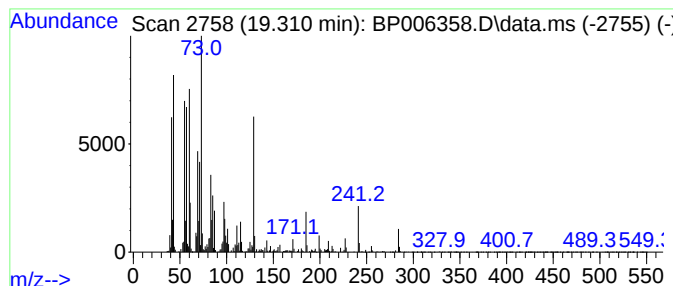
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Octadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	5.35 ng/ul	884743	Chrysene-d12	21.263

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Octadecanoic acid	284	C18H36O2	000057-11-4	98
3		Octadecanoic acid	284	C18H36O2	000057-11-4	96
4		Octadecanoic acid	284	C18H36O2	000057-11-4	95
5		Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

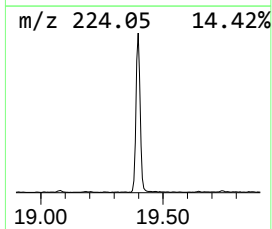
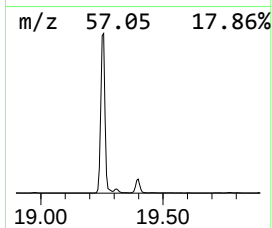
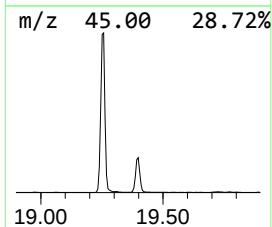
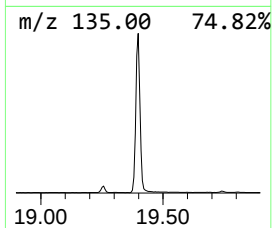
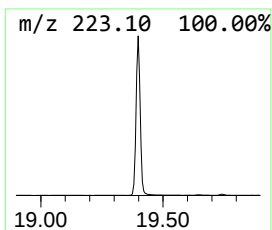
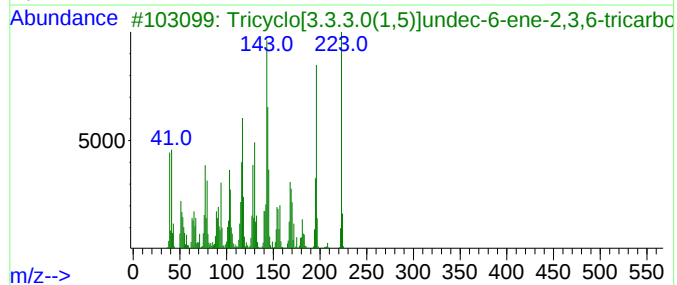
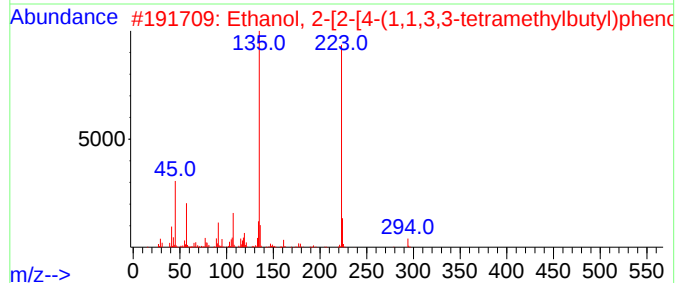
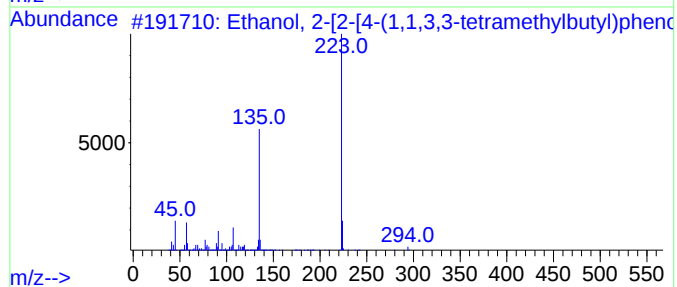
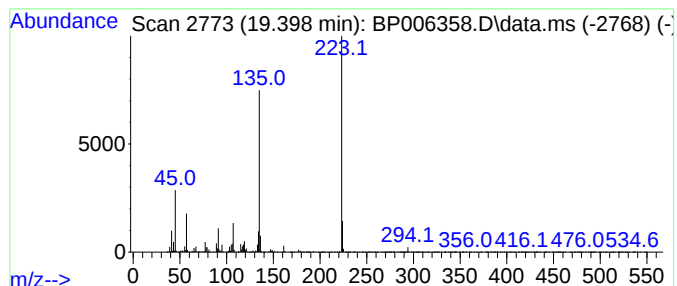
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.400	24.86 ng/ul	4114210	Chrysene-d12		21.263	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Ethanol, 2-[2-[4-(1,1,3,3-tetram...		294	C18H30O3	002315-61-9	96
2	Ethanol, 2-[2-[4-(1,1,3,3-tetram...		294	C18H30O3	002315-61-9	93
3	Tricyclo[3.3.3.0(1,5)]undec-6-en...		223	C14H13N3	118894-71-6	93
4	Pyrrolo[1,2-a]-1,3,5-triazine-7-...		223	C9H9N3O4	054449-89-7	90
5	Ethanol, 2-[2-[4-(1,1,3,3-tetram...		294	C18H30O3	002315-61-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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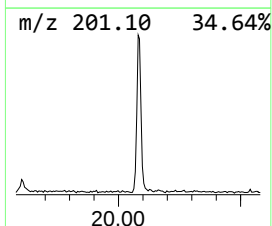
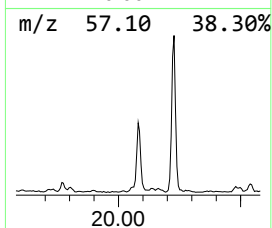
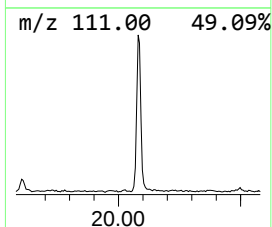
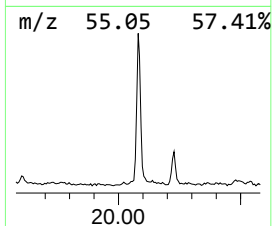
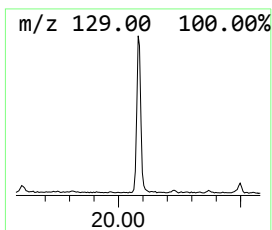
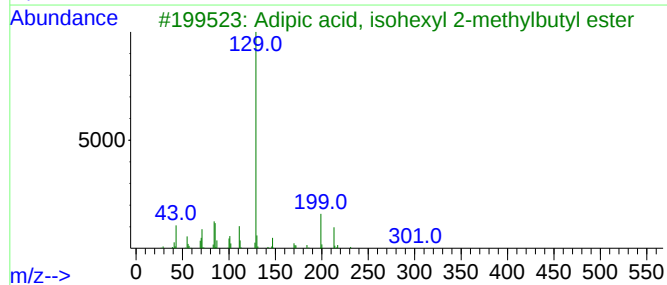
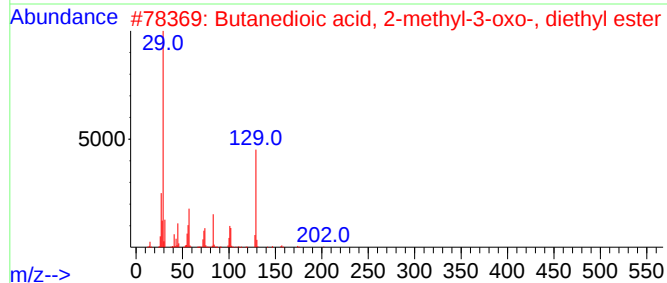
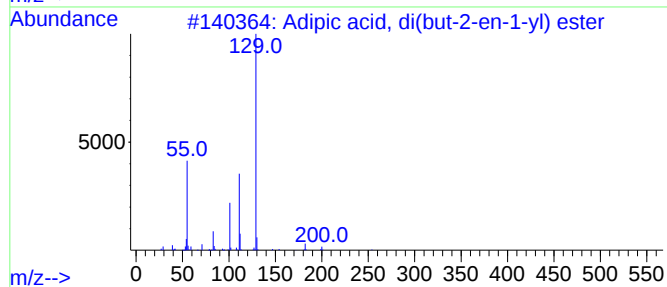
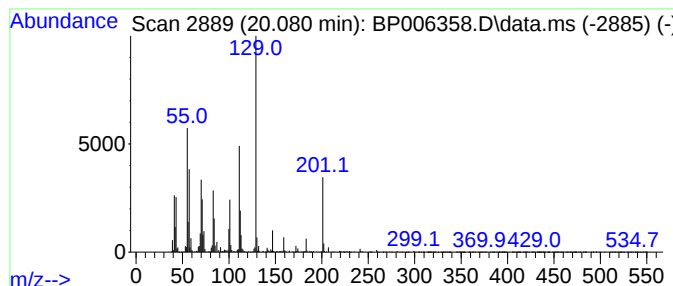
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-05 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.080	4.95 ng/ul	819366	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Adipic acid, di(but-2-en-1-yl) e...	254	C14H22O4	1000324-71-1	47
2			Butanedioic acid, 2-methyl-3-oxo...	202	C9H14O5	000759-65-9	43
3			Adipic acid, isohexyl 2-methylbu...	300	C17H32O4	1000324-67-4	38
4			2-Heptenoic acid, nonyl ester	254	C16H30O2	1000406-09-5	38
5			Adipic acid, di(3-hexyl) ester	314	C18H34O4	1000353-63-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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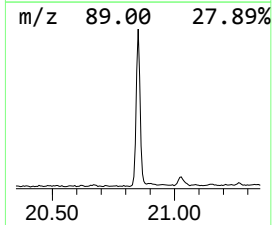
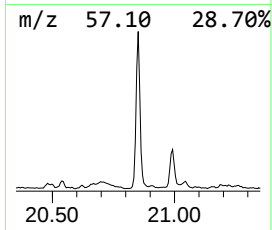
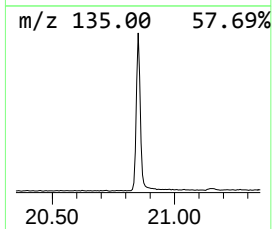
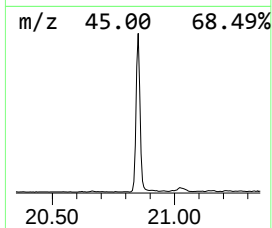
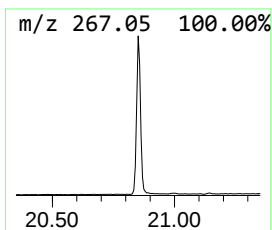
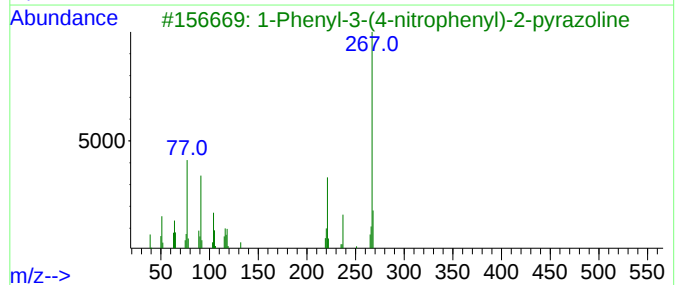
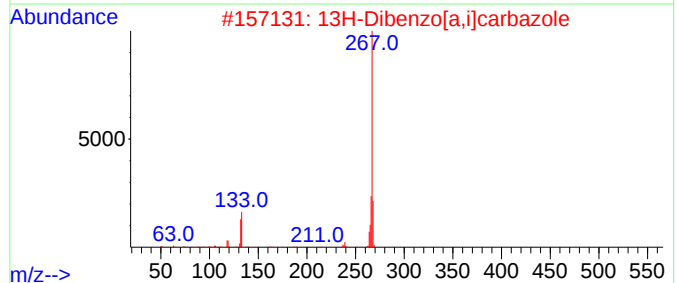
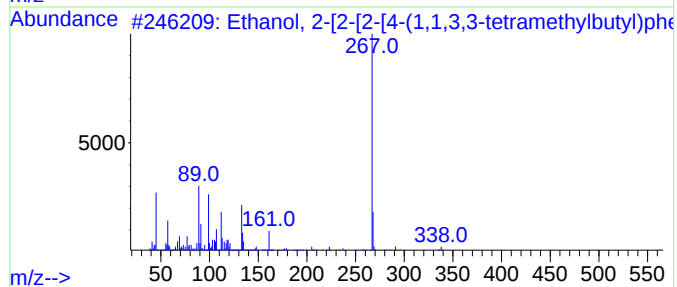
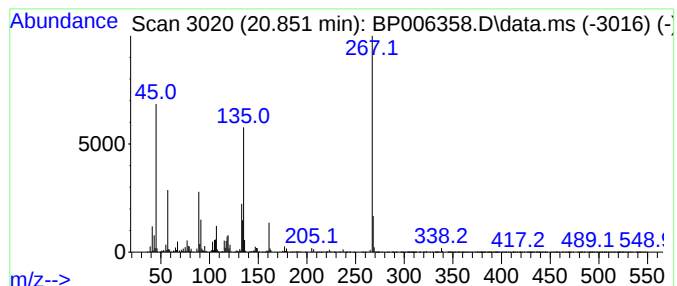
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 Ethanol, 2-[2-[2-[4-(1,1,3,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.850	13.73 ng/ul	2273180	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[2-[4-(1,1,3,3-tet...	338	C20H34O4	002315-62-0	96
2			13H-Dibenzo[a,i]carbazole	267	C20H13N	000239-64-5	38
3			1-Phenyl-3-(4-nitrophenyl)-2-pyr...	267	C15H13N3O2	003314-41-8	22
4			2-Methyl-5-(3,4-methylenedioxy...	267	C16H13NO3	094298-70-1	22
5			2-Benzylidenehydrazono-3-methyl...	267	C15H13N3S	1000146-59-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
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Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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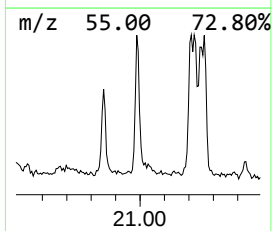
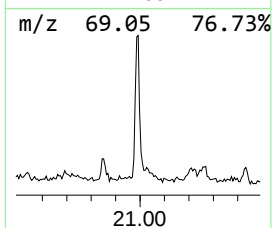
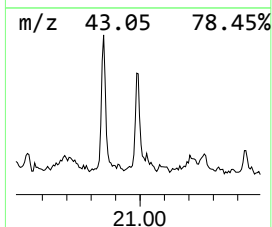
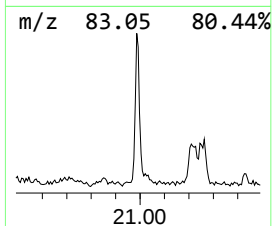
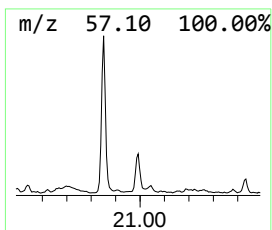
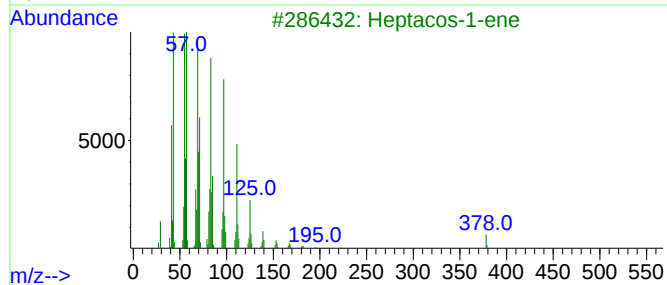
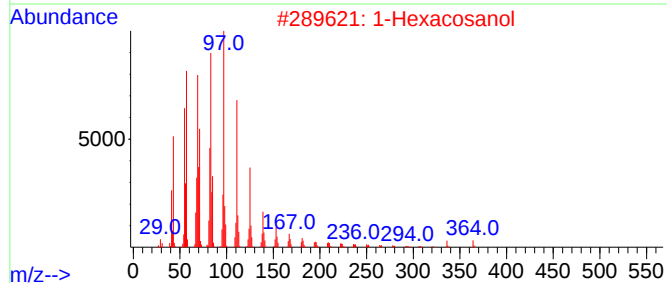
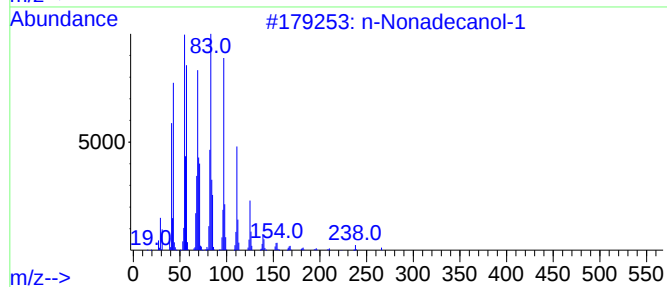
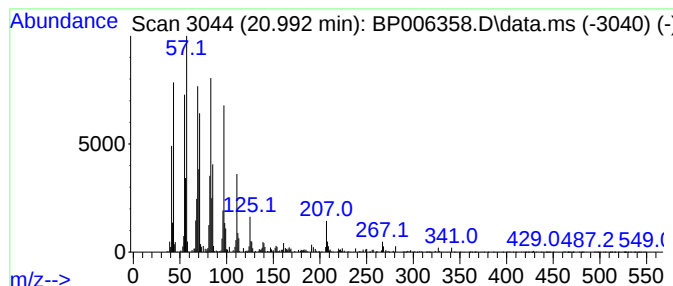
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 n-Nonadecanol-1 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.990	2.03 ng/ul	336772	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Nonadecanol-1	284	C19H40O	001454-84-8	93
2			1-Hexacosanol	382	C26H54O	000506-52-5	90
3			Heptacos-1-ene	378	C27H54	015306-27-1	90
4			Octacosanol	410	C28H58O	000557-61-9	90
5			Cyclotetradecane	196	C14H28	000295-17-0	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT8

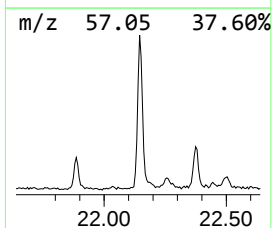
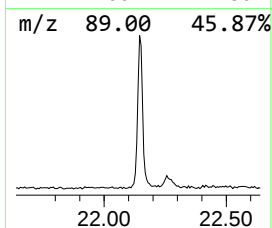
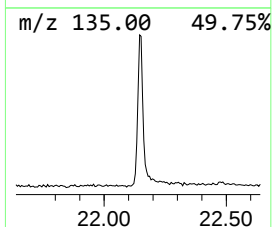
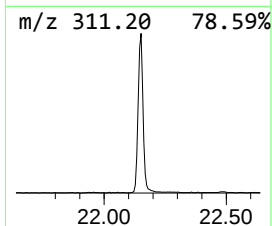
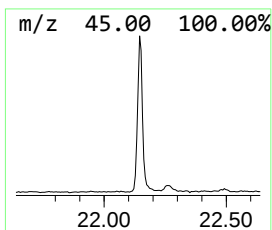
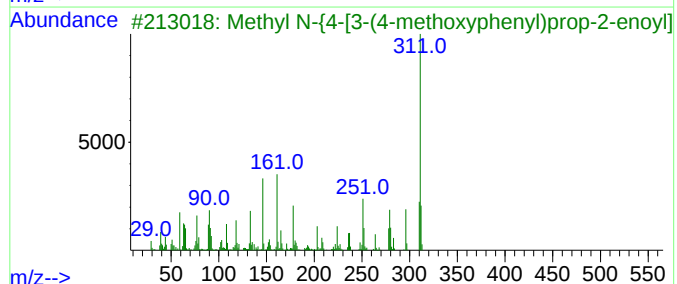
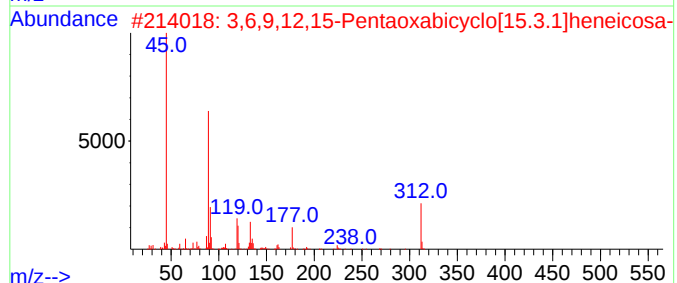
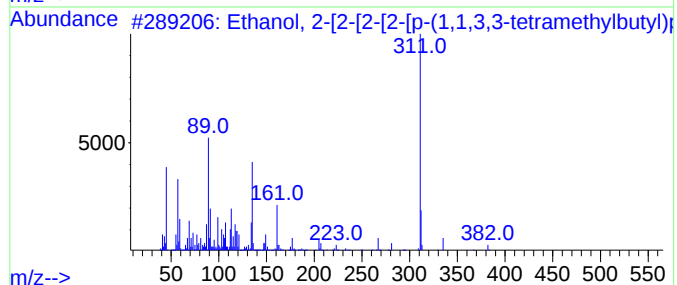
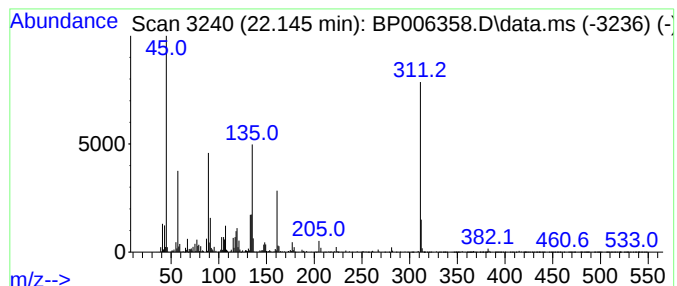
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-07 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.140	7.00 ng/ul	115880	Chrysene-d12	21.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-[2-[2-[p-(1,1,3,3-...	382	C22H38O5	002315-63-1	46
2			3,6,9,12,15-Pentaoxabicyclo[15.3...	312	C16H24O6	065112-36-9	43
3			Methyl N-{4-[3-(4-methoxyphenyl)...}	311	C18H17NO4	1000444-07-2	30
4			3,6,9,12,15-Pentaoxabicyclo[15.3...	374	C16H23BrO5	055440-88-5	27
5			Isoxazole, 4-(3,5-dimethoxypheny...	311	C18H17NO4	1000460-88-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AT8

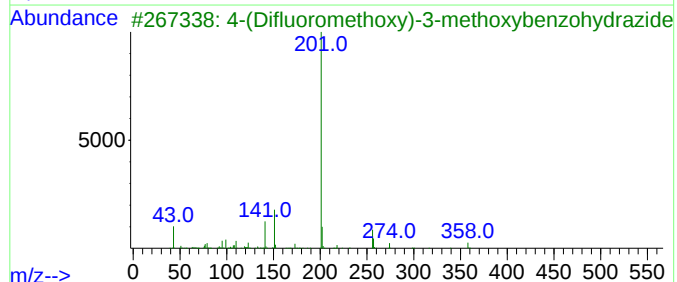
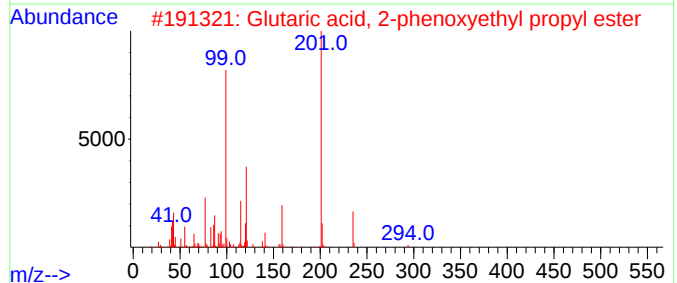
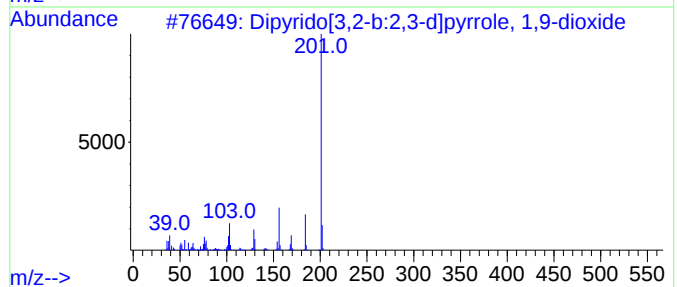
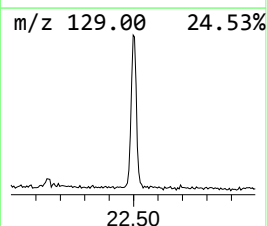
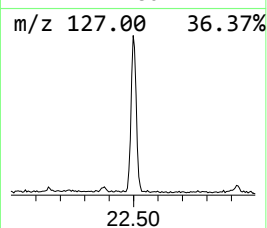
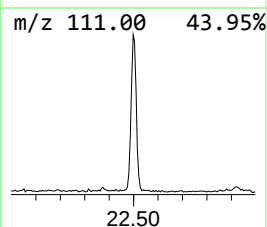
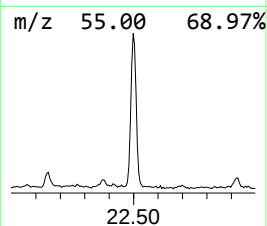
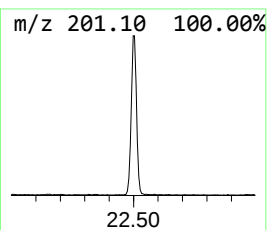
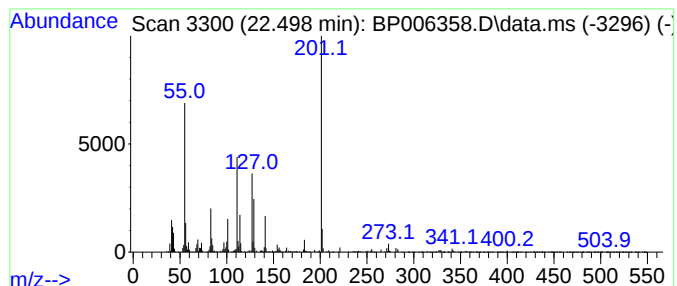
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-08 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.500	4.57 ng/ul	726863	Perylene-d12	23.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dipyrido[3,2-b:2,3-d]pyrrole, 1,...	201	C10H7N3O2	108349-61-7	27
2			Glutaric acid, 2-phenoxyethyl pr...	294	C16H22O5	1000376-91-3	17
3			4-(Difluoromethoxy)-3-methoxyben...	358	C15H16F2N2O6	1010512-12-8	17
4			1-Dimethylisopropylsilyloxynonane	244	C14H32OSi	1000283-07-0	17
5			Succinic acid, butyl 2-phenoxyet...	294	C16H22O5	1000381-19-4	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
Data File : BP006358.D
Acq On : 27 Jul 2021 10:51
Operator : CG/JU
Sample : M3063-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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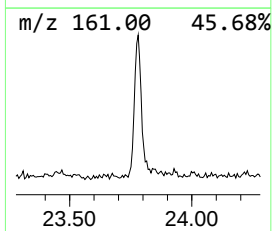
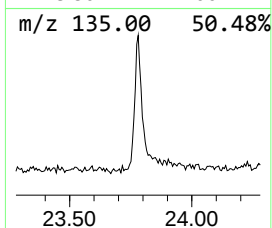
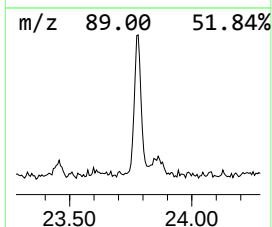
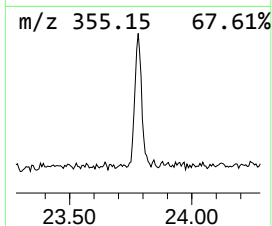
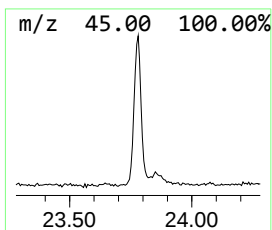
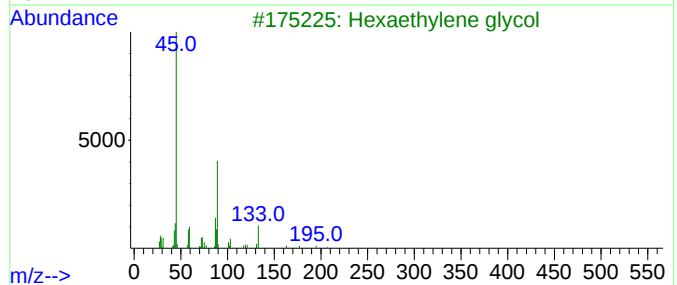
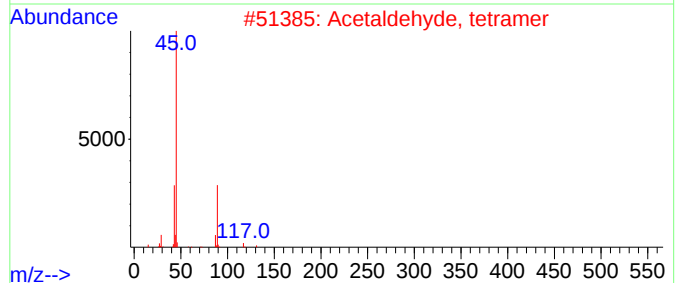
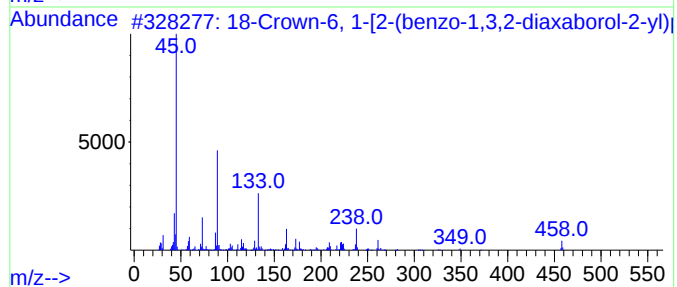
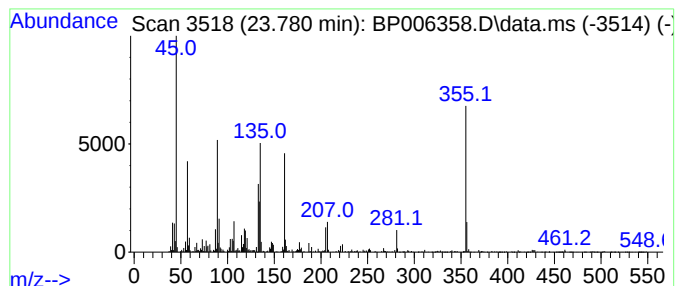
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 unknown-09 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.780	3.07 ng/ul	489019	Perylene-d12	23.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	18-Crown-6, 1-[2-(benzo-1,3,2-di...	458	C24H31B08	1000161-60-9	35		
2	Acetaldehyde, tetramer	176	C8H16O4	000108-62-3	27		
3	Hexaethylene glycol	282	C12H26O7	002615-15-8	25		
4	Pentaethylene glycol	238	C10H22O6	004792-15-8	25		
5	3,6,9,12,15,18-Hexaoxabicyclo[18...	418	C18H27BrO6	111982-22-0	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072721\
 Data File : BP006358.D
 Acq On : 27 Jul 2021 10:51
 Operator : CG/JU
 Sample : M3063-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP072421.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
Toluene	4.020	12.6	ng/ul	872981	1	7.787	1390320	20.0
1-Hexanol, 2-et...	7.850	4.2	ng/ul	290208	1	7.787	1390320	20.0
Benzene, 1,2,3-...	7.900	2.4	ng/ul	169960	1	7.787	1390320	20.0
Indane	8.150	3.6	ng/ul	253278	1	7.787	1390320	20.0
Ethanol, 2-(2-b...	10.360	9.2	ng/ul	884519	2	10.569	1926790	20.0
Benzo[b]thiophe...	11.680	2.0	ng/ul	195334	2	10.569	1926790	20.0
unknown-01	12.100	8.3	ng/ul	803877	2	10.569	1926790	20.0
Naphthalene, 1-...	12.450	2.4	ng/ul	233922	2	10.569	1926790	20.0
Phenol, 3,5-dic...	13.120	6.0	ng/ul	801526	3	14.410	2648550	20.0
Phenol, 3,4-dic...	13.430	4.0	ng/ul	535239	3	14.410	2648550	20.0
Benzene, 2-prop...	13.550	2.5	ng/ul	332454	3	14.410	2648550	20.0
Phenol, 2,3,4,6...	14.950	2.8	ng/ul	365126	3	14.410	2648550	20.0
unknown-02	17.050	2.1	ng/ul	284169	4	17.163	2694440	20.0
unknown-03	17.370	5.7	ng/ul	770560	4	17.163	2694440	20.0
Ethanol, 2-[4-(...	17.540	21.2	ng/ul	2852950	4	17.163	2694440	20.0
7,9-Di-tert-but...	17.850	2.8	ng/ul	379569	4	17.163	2694440	20.0
n-Hexadecanoic ...	18.070	6.3	ng/ul	847965	4	17.163	2694440	20.0
unknown-04	19.260	82.5	ng/ul	13646700	5	21.263	3310380	20.0
Octadecanoic acid	19.310	5.3	ng/ul	884743	5	21.263	3310380	20.0
Ethanol, 2-[2-[...	19.400	24.9	ng/ul	4114210	5	21.263	3310380	20.0
unknown-05	20.080	5.0	ng/ul	819366	5	21.263	3310380	20.0
unknown-06	20.230	4.9	ng/ul	808806	5	21.263	3310380	20.0
Ethanol, 2-[2-[...	20.850	13.7	ng/ul	2273180	5	21.263	3310380	20.0
n-Nonadecanol-1	20.990	2.0	ng/ul	336772	5	21.263	3310380	20.0
unknown-07	22.140	7.0	ng/ul	1158880	5	21.263	3310380	20.0
unknown-08	22.500	4.6	ng/ul	726863	6	23.610	3182060	20.0
unknown-09	23.780	3.1	ng/ul	489019	6	23.610	3182060	20.0