

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010081.D
 Acq On : 25 Apr 2022 13:36
 Operator : CG/JU
 Sample : N2475-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BE9

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_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010081.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.205	17	20	26	rVB2	15534	20884	0.36%	0.017%
2	3.369	44	48	52	rBV	46887	60155	1.05%	0.050%
3	3.781	114	118	141	rBV	533210	872887	15.21%	0.726%
4	3.934	141	144	155	rVB2	32566	48083	0.84%	0.040%
5	4.116	171	175	181	rVB	16039	21798	0.38%	0.018%
6	4.381	216	220	228	rBV	81276	123045	2.14%	0.102%
7	4.475	233	236	245	rVB6	16381	25970	0.45%	0.022%
8	5.046	328	333	338	rBV	32744	50182	0.87%	0.042%
9	5.193	353	358	364	rBV3	11318	19257	0.34%	0.016%
10	5.363	382	387	392	rBV5	14796	21277	0.37%	0.018%
11	5.999	490	495	501	rVB	27062	42301	0.74%	0.035%
12	6.340	549	553	573	rBV	313824	617410	10.75%	0.513%
13	6.940	649	655	660	rBV	42925	61894	1.08%	0.051%
14	7.099	675	682	685	rBV	831483	1479474	25.77%	1.230%
15	7.210	697	701	704	rBV2	14215	19938	0.35%	0.017%
16	7.263	704	710	721	rVV	626919	1008929	17.57%	0.839%
17	7.352	721	725	734	rVB2	15999	28778	0.50%	0.024%
18	7.469	737	745	748	rBV	805844	1431110	24.93%	1.190%
19	7.499	748	750	757	rVV	499474	646404	11.26%	0.538%
20	7.569	759	762	771	rVB7	12657	22839	0.40%	0.019%
21	7.828	797	806	811	rBV	549970	901253	15.70%	0.749%
22	7.887	811	816	819	rVV	228579	354338	6.17%	0.295%
23	7.934	819	824	827	rVV	631033	1054714	18.37%	0.877%
24	7.969	827	830	839	rVB	602745	934460	16.28%	0.777%
25	8.175	861	865	874	rVB5	10019	19038	0.33%	0.016%
26	8.287	878	884	895	rVV	667321	1107347	19.29%	0.921%
27	8.469	904	915	920	rBV	493808	1215425	21.17%	1.011%
28	8.522	920	924	933	rVB	280854	449808	7.84%	0.374%
29	8.640	936	944	962	rBV	1039866	1728216	30.10%	1.437%
30	9.022	1002	1009	1015	rBV	458275	734962	12.80%	0.611%
31	9.093	1015	1021	1033	rVV	807324	1343418	23.40%	1.117%
32	9.199	1033	1039	1049	rVB5	17212	40295	0.70%	0.034%
33	9.769	1128	1136	1138	rBV	19211	30385	0.53%	0.025%
34	9.816	1138	1144	1154	rVB	645569	1104736	19.24%	0.919%
35	10.075	1182	1188	1198	rBV	10984	22120	0.39%	0.018%
36	10.363	1229	1237	1239	rBV	962678	1726750	30.08%	1.436%

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37	10.604	1270	1278	1289	rVB	684853	1132935	19.73%	0.942%
38	10.740	1293	1301	1305	rBV	882153	1504748	26.21%	1.251%
39	10.793	1305	1310	1317	rVV	828673	1425219	24.83%	1.185%
40	10.875	1317	1324	1343	rVB	900380	1676085	29.20%	1.394%

41	11.087	1349	1360	1372	rVB	965077	1570727	27.36%	1.306%
42	11.228	1379	1384	1389	rVB6	8500	15318	0.27%	0.013%
43	12.216	1548	1552	1556	rBV2	14118	24411	0.43%	0.020%
44	12.269	1556	1561	1566	rVV3	11368	23442	0.41%	0.019%
45	12.328	1566	1571	1580	rVB3	17792	36407	0.63%	0.030%

46	12.687	1626	1632	1635	rBV	260820	440267	7.67%	0.366%
47	12.716	1635	1637	1644	rVB	181005	250745	4.37%	0.209%
48	12.792	1644	1650	1654	rBV	60265	84435	1.47%	0.070%
49	12.910	1661	1670	1675	rBV	647556	1851061	32.24%	1.539%
50	12.957	1675	1678	1680	rVV2	100312	167024	2.91%	0.139%

51	13.163	1693	1713	1718	rBV3	427527	1542041	26.86%	1.282%
52	13.428	1751	1758	1767	rVB	75610	161561	2.81%	0.134%
53	13.504	1767	1771	1775	rBV7	10494	17997	0.31%	0.015%
54	13.640	1791	1794	1801	rVB2	11646	16900	0.29%	0.014%
55	13.975	1843	1851	1862	rBV	1990984	2767849	48.21%	2.302%

56	14.139	1872	1879	1885	rBV	1048367	1500713	26.14%	1.248%
57	14.198	1885	1889	1893	rVV2	24597	44024	0.77%	0.037%
58	14.257	1893	1899	1912	rVB	1985180	3095043	53.91%	2.574%
59	14.569	1945	1952	1957	rBV	1349942	2054474	35.79%	1.708%
60	14.634	1957	1963	1977	rVV	1325589	2025026	35.27%	1.684%

61	14.757	1978	1984	2003	rVV	730437	1323926	23.06%	1.101%
62	14.922	2006	2012	2033	rVB	1074074	1646226	28.68%	1.369%
63	15.386	2085	2091	2102	rVB	1689091	2268343	39.51%	1.886%
64	15.563	2112	2121	2126	rBV	2641760	3961922	69.01%	3.295%
65	15.616	2126	2130	2135	rVV	1511799	2119695	36.92%	1.763%

66	15.675	2135	2140	2154	rVV	1013309	1566611	27.29%	1.303%
67	15.798	2157	2161	2168	rVV	35261	63228	1.10%	0.053%
68	15.869	2168	2173	2180	rVV	33449	54429	0.95%	0.045%
69	15.939	2180	2185	2192	rVV	122864	152589	2.66%	0.127%
70	16.322	2245	2250	2261	rBV2	63916	102990	1.79%	0.086%

71	16.463	2269	2274	2280	rVB	51487	65770	1.15%	0.055%
72	16.539	2282	2287	2295	rVB	9755	18009	0.31%	0.015%
73	16.628	2295	2302	2309	rBV	1472255	2109141	36.74%	1.754%
74	16.975	2354	2361	2370	rBV	1793481	2634770	45.90%	2.191%
75	17.069	2370	2377	2391	rVV	2428175	3397762	59.19%	2.825%

76	17.322	2413	2420	2430	rBV	1682696	2461738	42.88%	2.047%
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77	17.422	2430	2437	2440	rBV2	3063234	4572389	79.65%	3.802%
78	17.457	2440	2443	2451	rVB	1547272	2101453	36.61%	1.747%
79	17.892	2512	2517	2520	rBV	33560	42146	0.73%	0.035%
80	17.945	2520	2526	2533	rVV3	60922	108347	1.89%	0.090%
81	18.175	2561	2565	2568	rBV6	12959	16928	0.29%	0.014%
82	18.275	2577	2582	2584	rBV2	9409	13313	0.23%	0.011%
83	18.310	2584	2588	2593	rVB	172707	215952	3.76%	0.180%
84	18.675	2644	2650	2653	rBV3	23947	34075	0.59%	0.028%
85	18.716	2653	2657	2661	rVB7	13486	20710	0.36%	0.017%
86	19.227	2739	2744	2749	rBV	39433	53852	0.94%	0.045%
87	19.298	2752	2756	2762	rVV	46536	61351	1.07%	0.051%
88	19.563	2797	2801	2806	rBV	25695	39837	0.69%	0.033%
89	19.651	2809	2816	2819	rBV	3802341	5740753	100.00%	4.774%
90	19.680	2819	2821	2834	rVB	2397980	2592957	45.17%	2.156%
91	20.551	2963	2969	2979	rBV	4250083	4827925	84.10%	4.015%
92	21.310	3094	3098	3103	rBV	4406848	4630549	80.66%	3.851%
93	21.392	3106	3112	3117	rVV2	2618062	5201421	90.61%	4.325%
94	21.445	3117	3121	3131	rVB	2213428	3044713	53.04%	2.532%
95	23.110	3397	3404	3408	rBV	2306919	4296968	74.85%	3.573%
96	23.157	3408	3412	3426	rVB	1418931	2497637	43.51%	2.077%
97	23.698	3496	3504	3509	rBV2	1912594	4038495	70.35%	3.358%
98	23.751	3509	3513	3522	rVV	1174093	2268348	39.51%	1.886%
99	23.857	3524	3531	3543	rVB2	934527	1972445	34.36%	1.640%
100	26.427	3956	3968	3990	rBV2	1436820	5099892	88.84%	4.241%

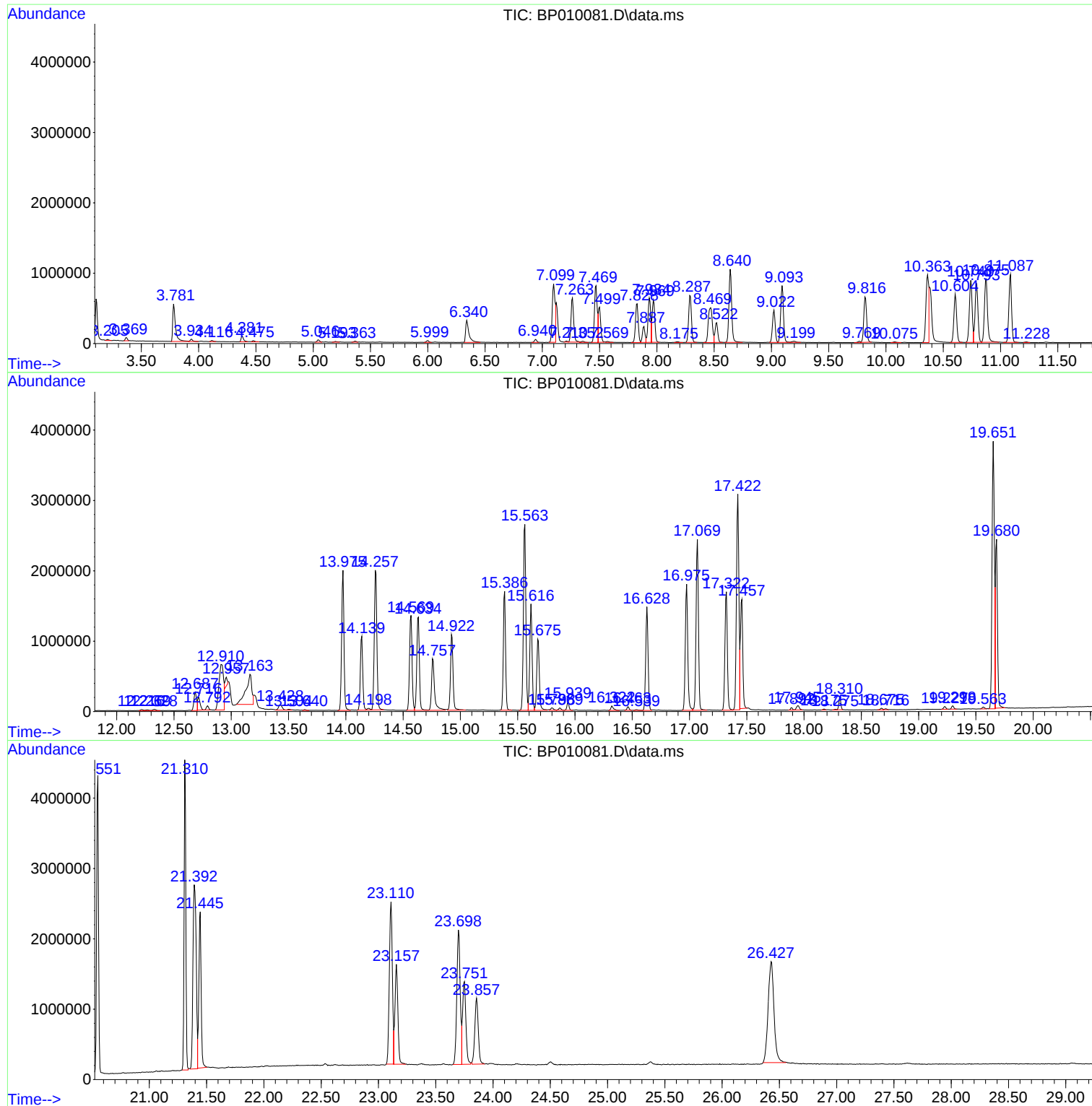
Sum of corrected areas: 120257937

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
BNA_P
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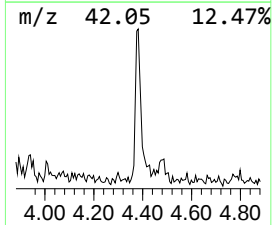
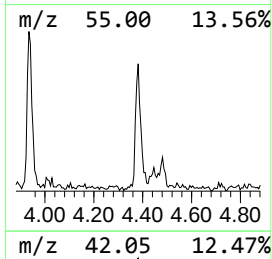
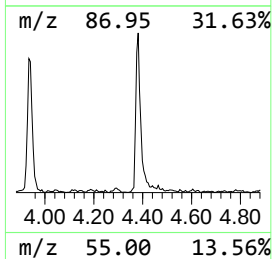
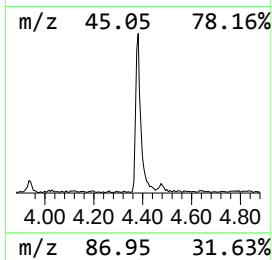
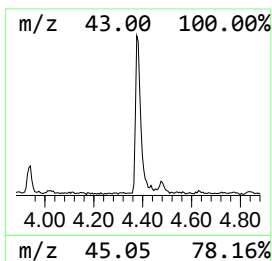
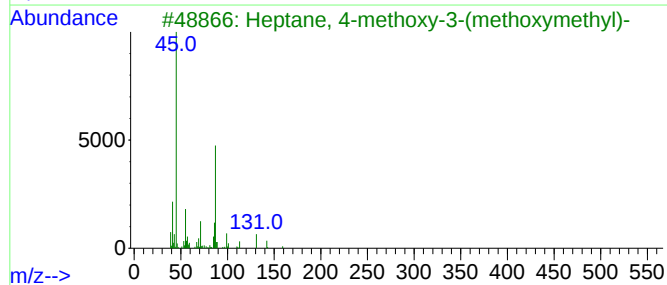
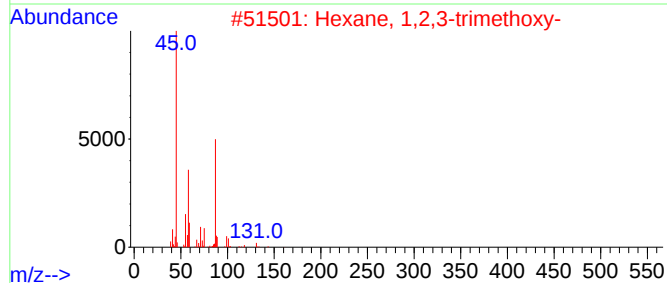
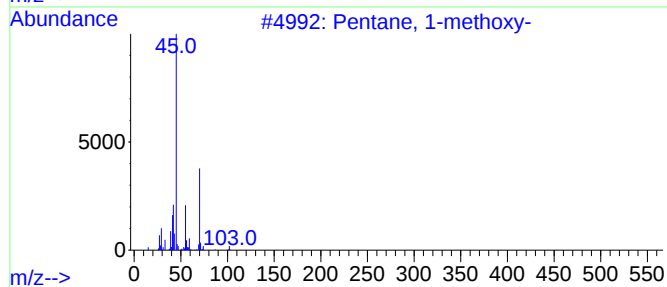
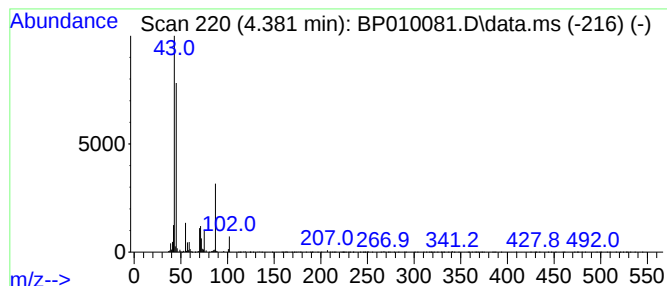
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Pentane, 1-methoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.381	2.33 ng/ul	123045	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 1-methoxy-	102	C6H14O	000628-80-8	50
2			Hexane, 1,2,3-trimethoxy-	176	C9H20O3	020637-29-0	50
3			Heptane, 4-methoxy-3-(methoxymet...	174	C10H22O2	020637-47-2	45
4			Heptane, 4-methoxy-3-(methoxymet...	174	C10H22O2	020637-47-2	45
5			Propane, 2,2'-[ethylidenebis(oxy...	146	C8H18O2	004285-59-0	42



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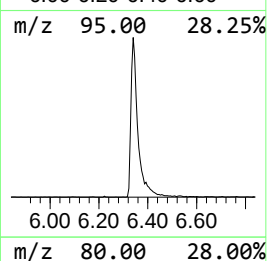
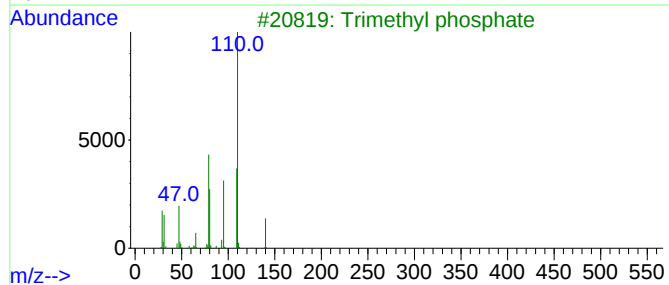
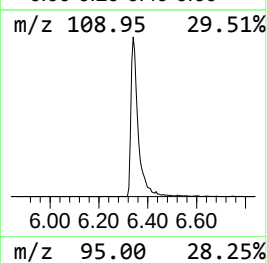
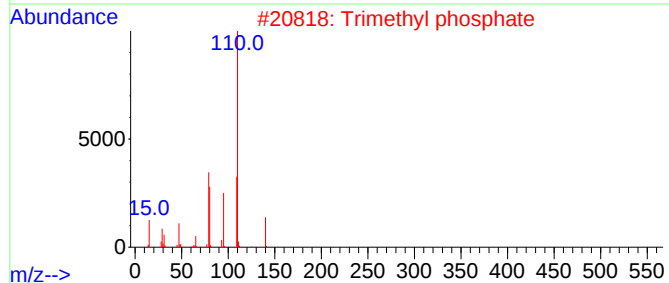
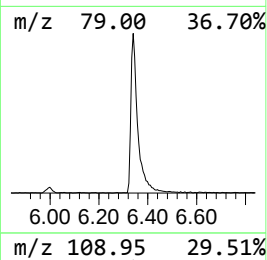
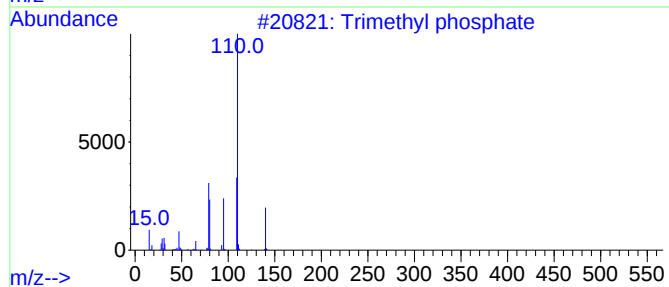
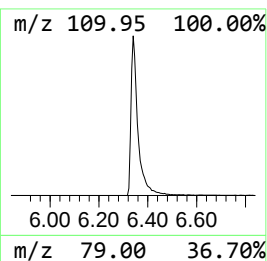
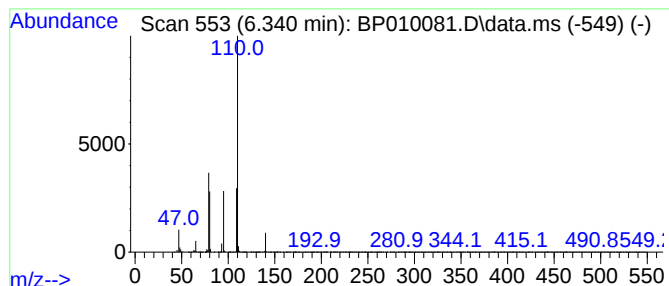
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Trimethyl phosphate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.340	11.71 ng/ul	617410	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trimethyl phosphate	140	C3H9O4P	000512-56-1	91
2			Trimethyl phosphate	140	C3H9O4P	000512-56-1	91
3			Trimethyl phosphate	140	C3H9O4P	000512-56-1	91
4			Dimethyl isopropylphosphonate	152	C5H13O3P	054552-77-1	56
5			Butanol, 1-dimethylphosphonato)-	182	C6H15O4P	023054-81-1	50



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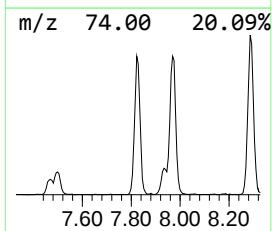
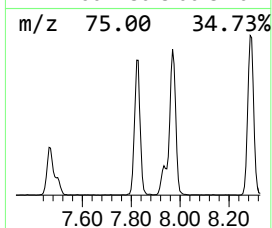
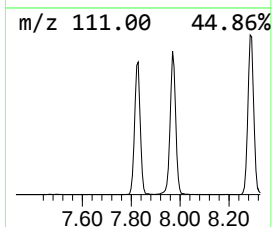
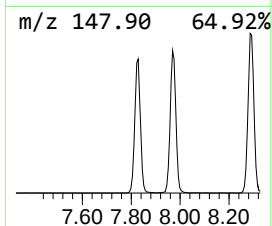
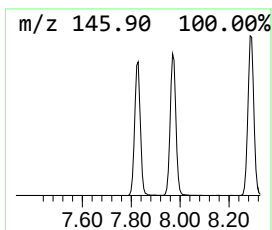
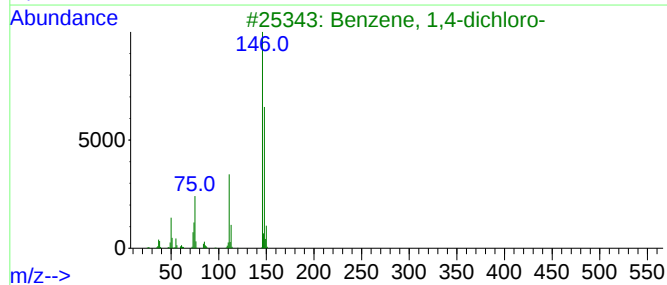
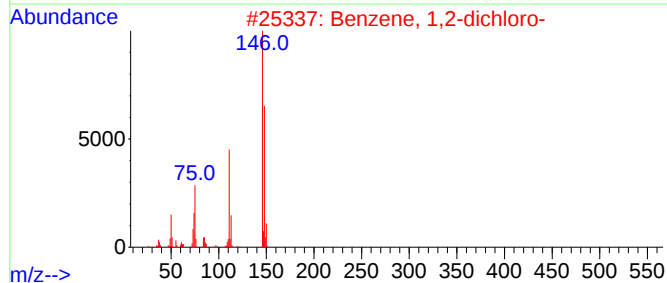
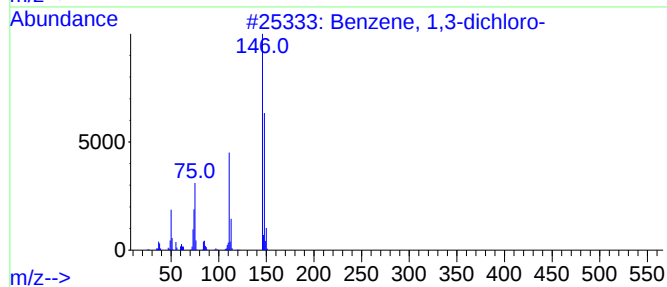
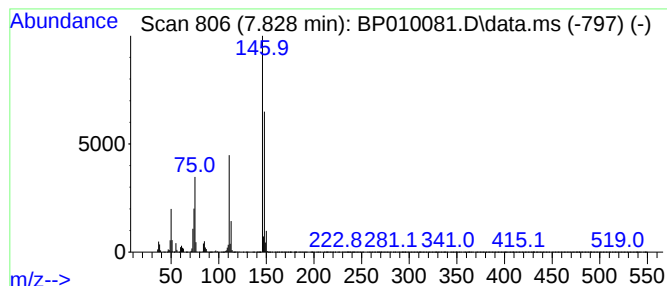
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.828	17.09 ng/ul	901253	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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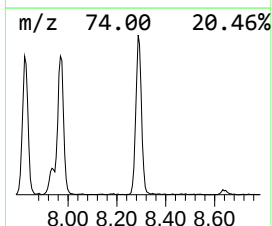
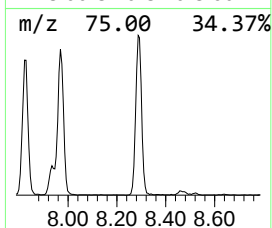
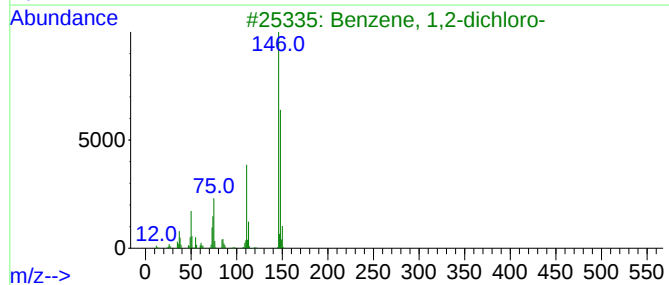
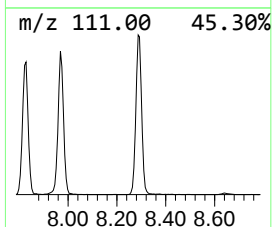
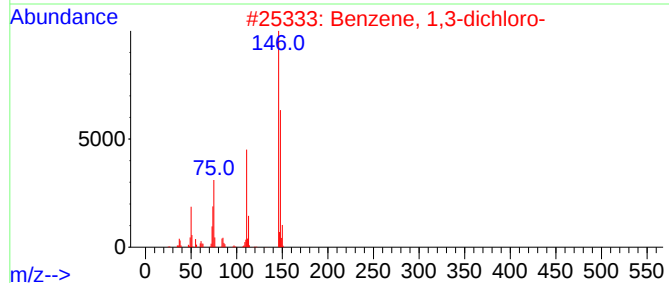
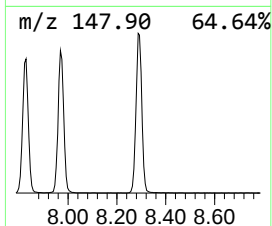
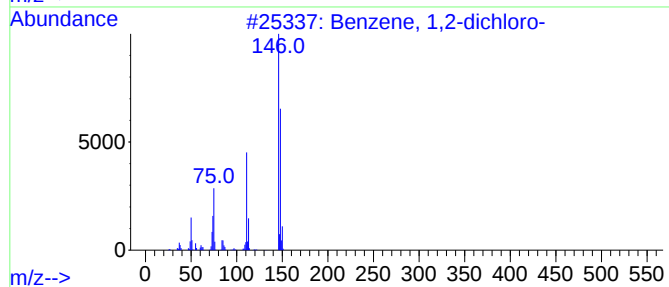
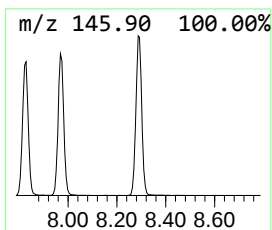
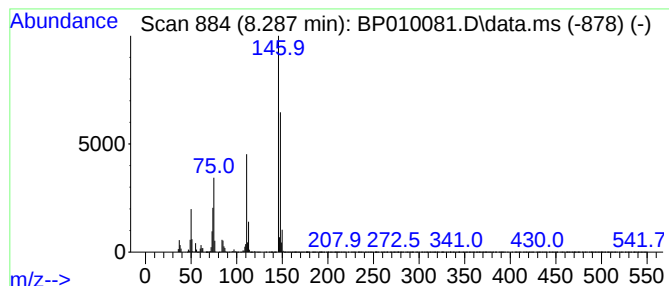
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,2-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.287	21.00 ng/ul	1107350	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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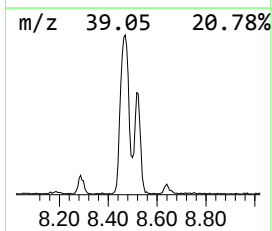
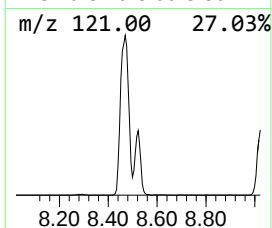
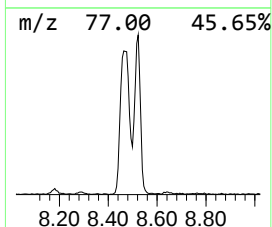
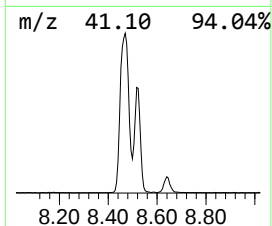
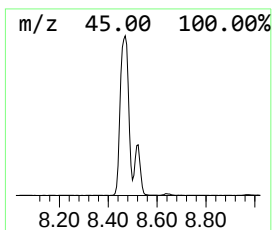
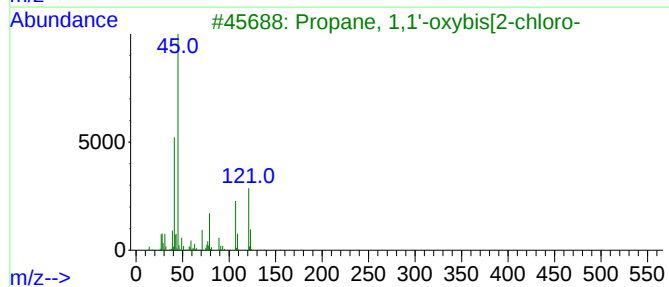
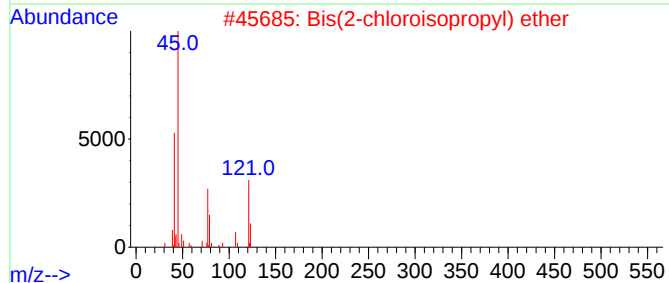
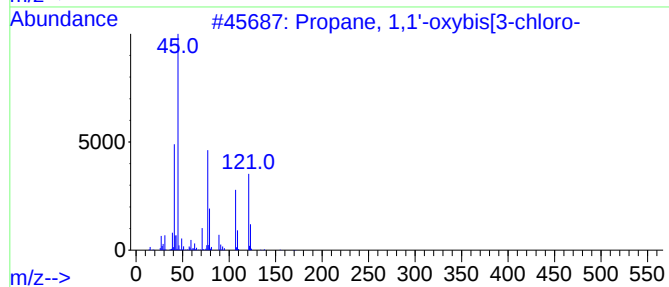
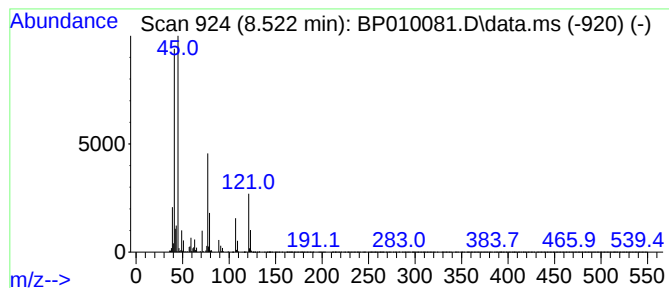
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Propane, 1,1'-oxybis[3-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.522	8.53 ng/ul	449808	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propane, 1,1'-oxybis[3-chloro-	170	C6H12Cl2O	000629-36-7	83
2			Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	59
3			Propane, 1,1'-oxybis[2-chloro-	170	C6H12Cl2O	054460-96-7	59
4			Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	50
5			Bis(2-chloroisopropyl) ether	170	C6H12Cl2O	039638-32-9	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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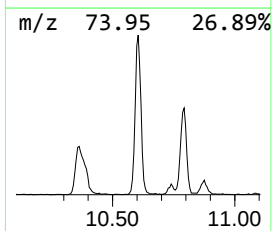
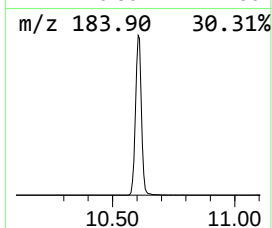
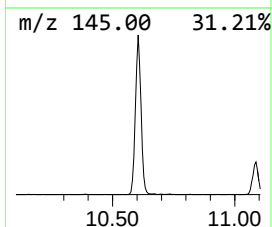
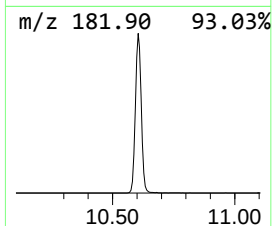
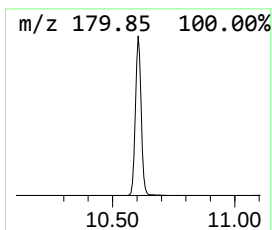
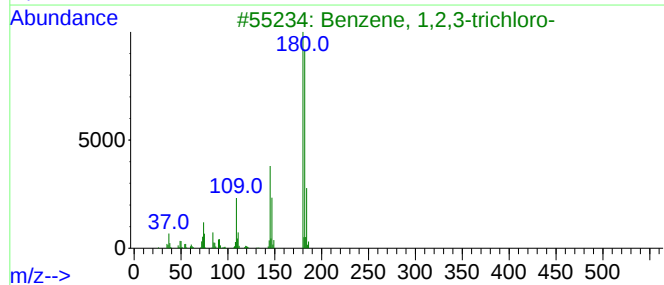
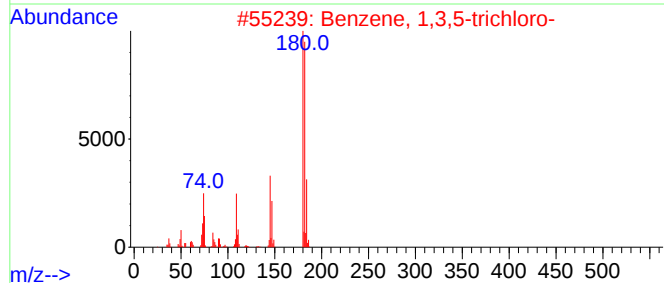
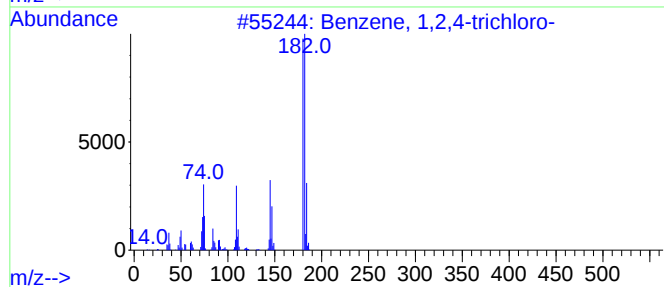
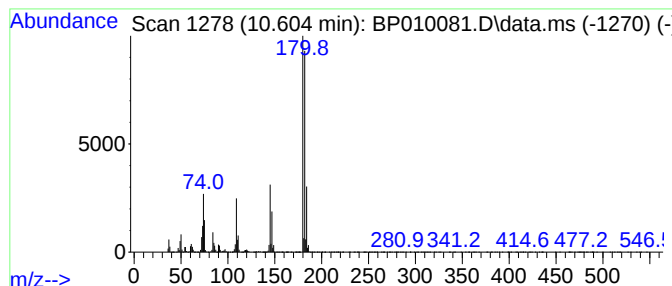
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,4-trichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.604	15.06 ng/ul	1132940	Naphthalene-d8	10.740

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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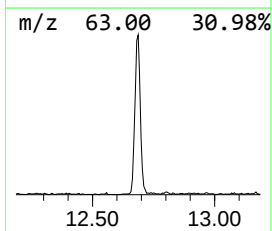
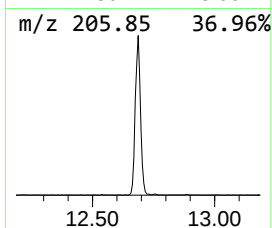
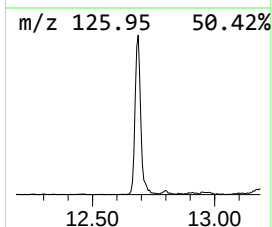
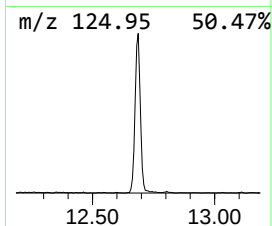
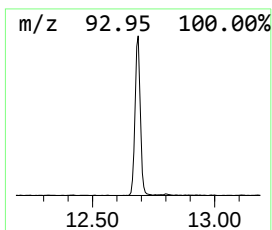
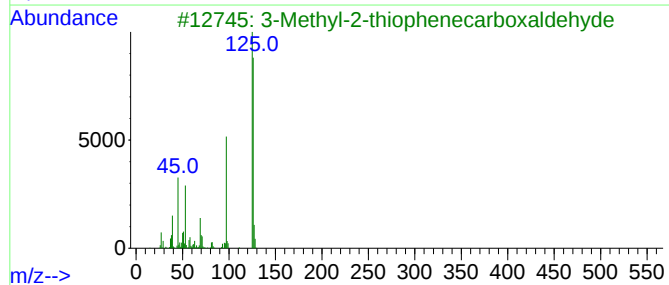
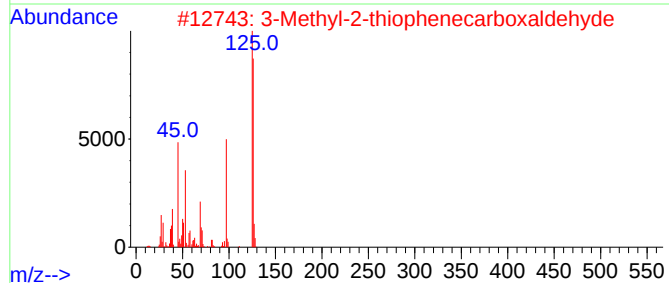
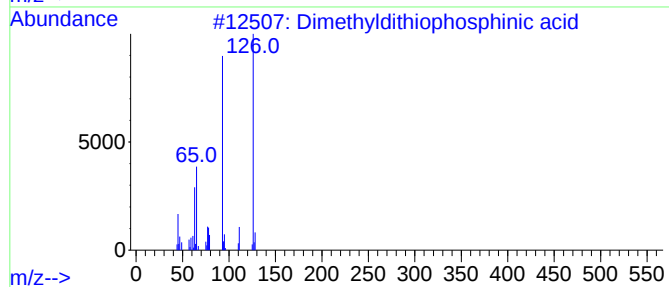
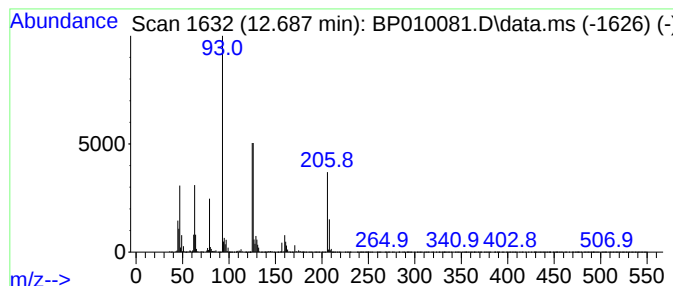
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	4.29 ng/ul	440267	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyldithiophosphinic acid	126	C2H7PS2	016367-68-3	35
2			3-Methyl-2-thiophenecarboxaldehyde	126	C6H6OS	005834-16-2	18
3			3-Methyl-2-thiophenecarboxaldehyde	126	C6H6OS	005834-16-2	18
4			5-Methyl-2-thiophenecarboxaldehyde	126	C6H6OS	013679-70-4	18
5			5-Methyl-2-thiophenecarboxaldehyde	126	C6H6OS	013679-70-4	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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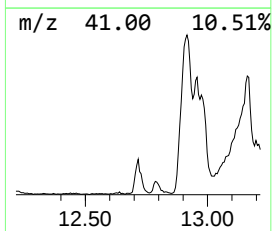
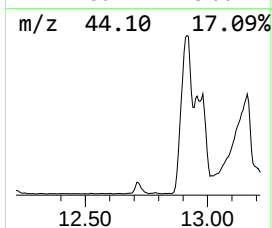
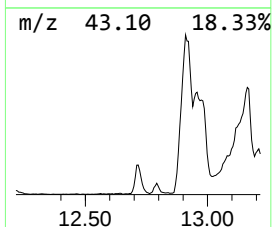
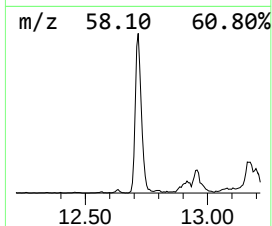
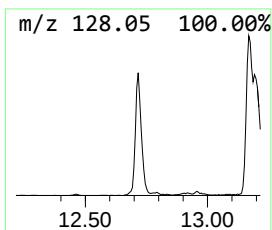
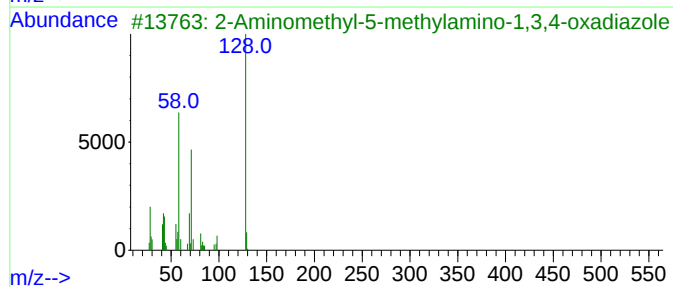
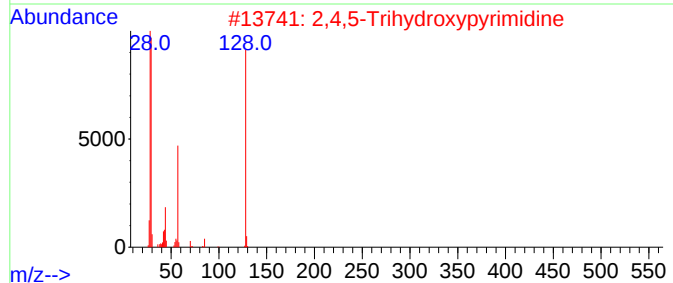
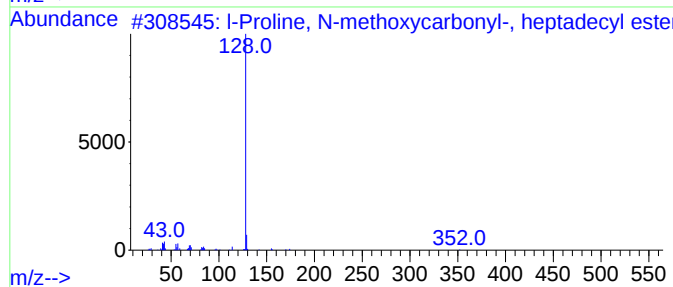
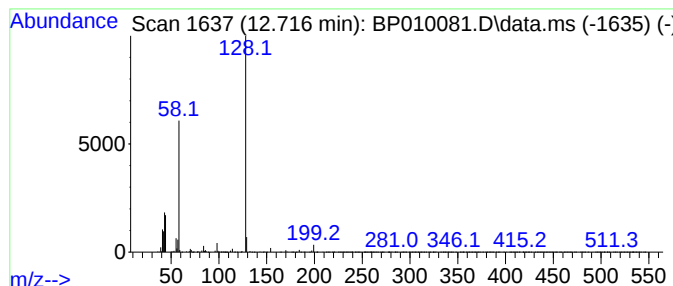
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.716	2.44 ng/ul	250745	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Proline, N-methoxycarbonyl-, h...	411	C24H45NO4	1000313-77-4	47
2			2,4,5-Trihydroxypyrimidine	128	C4H4N2O3	000496-76-4	40
3			2-Aminomethyl-5-methylamino-1,3,...	128	C4H8N4O	002937-92-0	38
4			1-Prolylglycine, N-methoxycarbon...	244	C10H16N2O5	1000320-27-5	38
5			1-Prolylglycine, N-methoxycarbon...	362	C12H15F5N2O5	1000322-64-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
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Misc :
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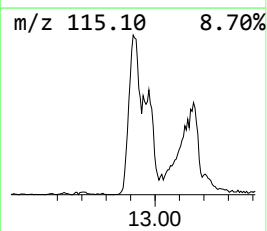
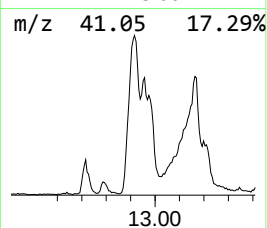
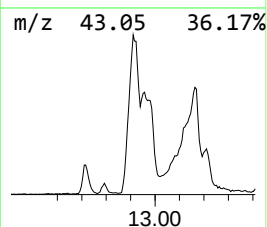
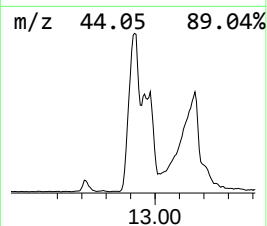
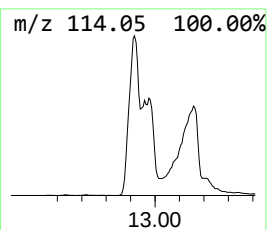
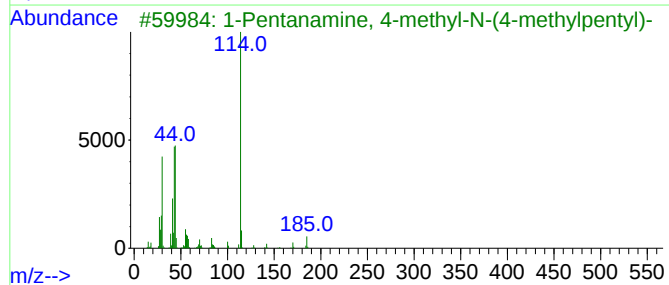
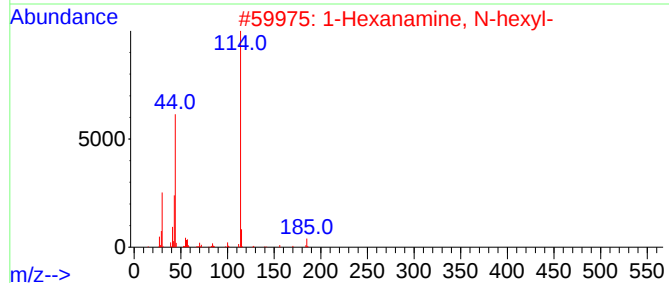
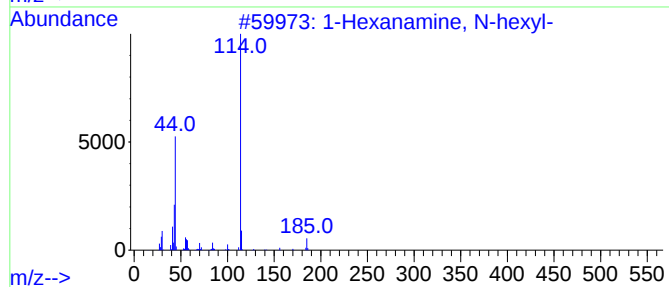
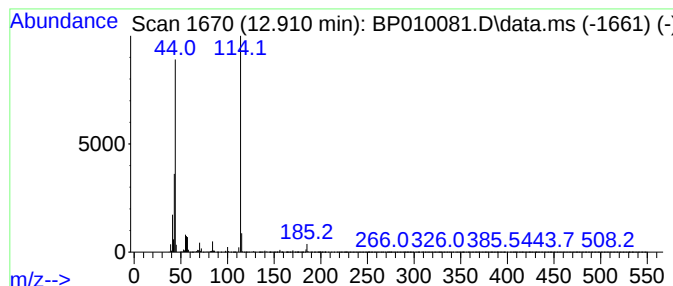
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Hexanamine, N-hexyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.910	18.02 ng/ul	1851060	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanamine, N-hexyl-	185	C12H27N	000143-16-8	78
2			1-Hexanamine, N-hexyl-	185	C12H27N	000143-16-8	68
3			1-Pentanamine, 4-methyl-N-(4-met...	185	C12H27N	054775-00-7	47
4			d-Ribitol, 1-deoxy-1-hexylamino-	235	C11H25NO4	1000158-64-0	40
5			Tetrapropylammonium iodide	313	C12H28IN	000631-40-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
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Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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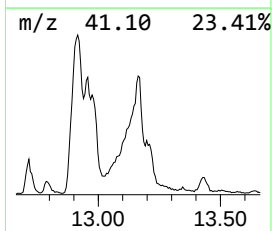
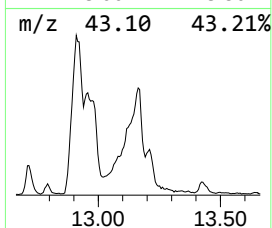
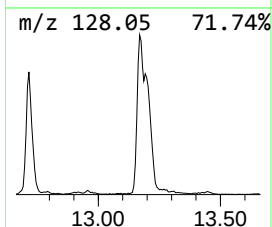
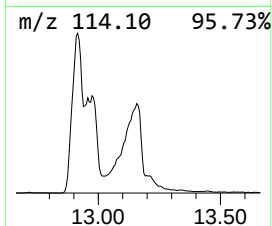
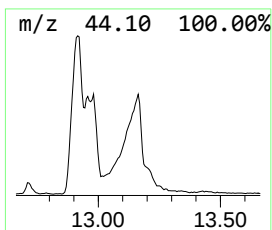
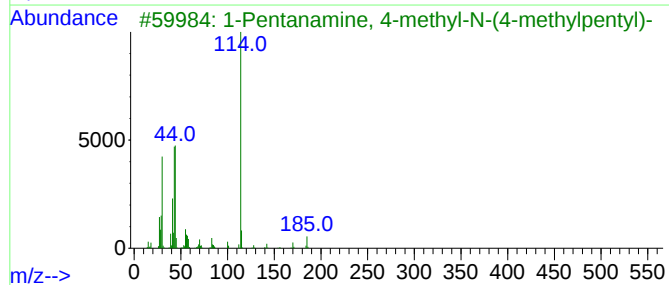
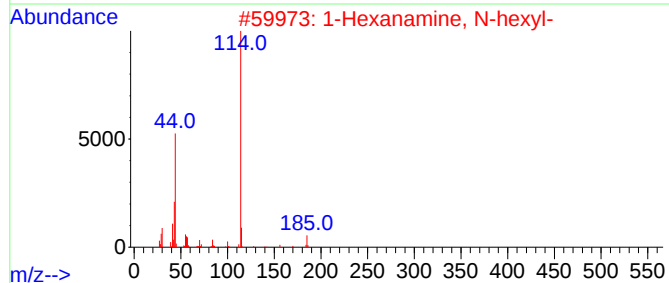
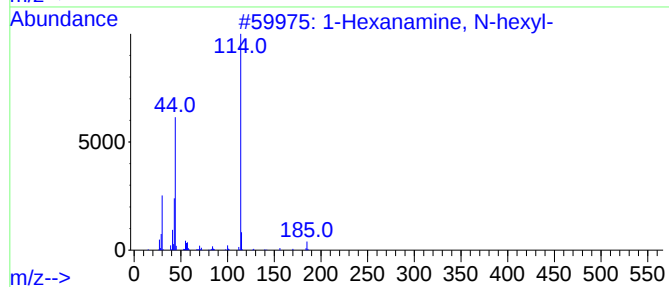
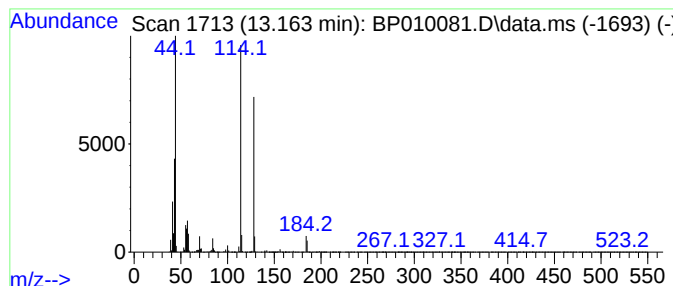
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.163	15.01 ng/ul	1542040	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexanamine, N-hexyl-	185	C12H27N	000143-16-8	46
2			1-Hexanamine, N-hexyl-	185	C12H27N	000143-16-8	46
3			1-Pentanamine, 4-methyl-N-(4-met...	185	C12H27N	054775-00-7	43
4			3-(Ethyl-hydrazono)-butan-2-one	128	C6H12N2O	1000194-94-0	38
5			Butanoic acid, 2-hexylimino-	185	C10H19NO2	1000258-67-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0BE9

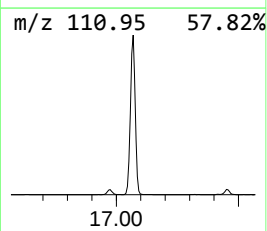
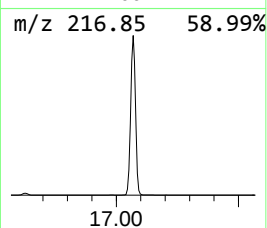
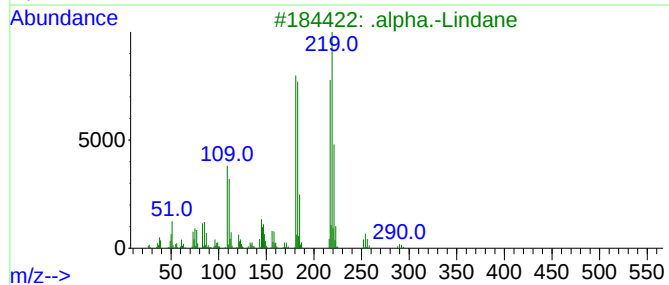
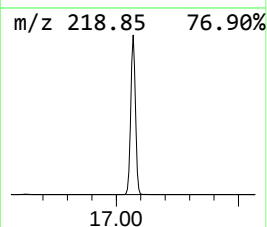
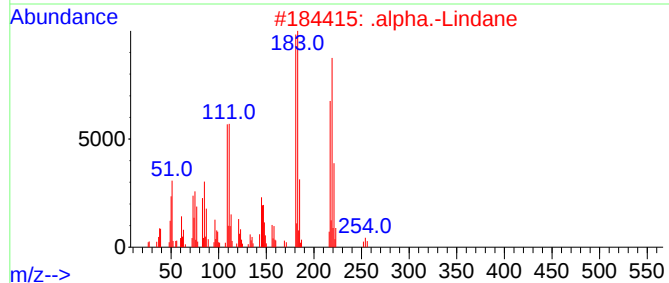
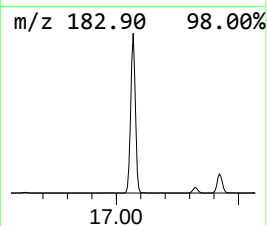
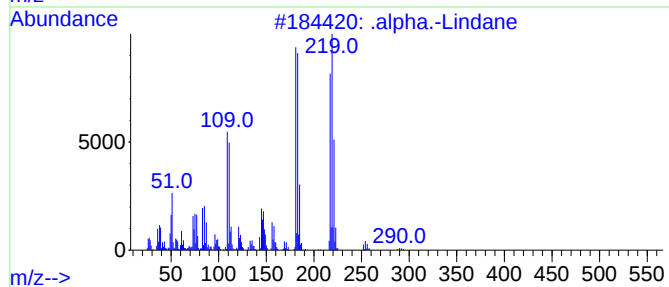
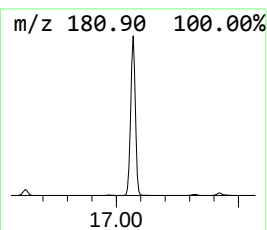
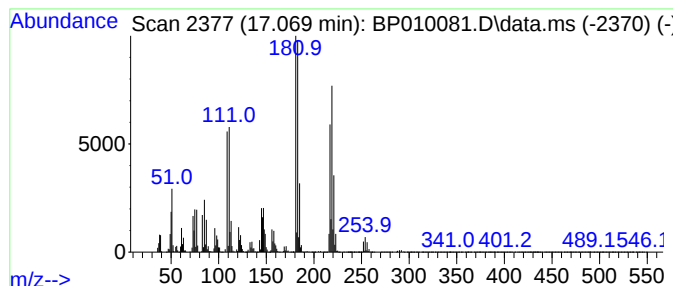
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 .alpha.-Lindane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.069	27.60 ng/ul	3397760	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	.	alpha.-Lindane	288	C6H6Cl6	000319-84-6	96
2	.	.	alpha.-Lindane	288	C6H6Cl6	000319-84-6	95
3	.	.	alpha.-Lindane	288	C6H6Cl6	000319-84-6	91
4	.	.	beta.-Hexachlorocyclohexane	288	C6H6Cl6	000319-85-7	91
5	.	.	delta.-Lindane	288	C6H6Cl6	000319-86-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010081.D
Acq On : 25 Apr 2022 13:36
Operator : CG/JU
Sample : N2475-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0BE9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentane, 1-meth...	4.381	2.3	ng/ul	123045	1	7.934	1054710	20.0
Trimethyl phosph...	6.340	11.7	ng/ul	617410	1	7.934	1054710	20.0
Benzene, 1,3-di...	7.828	17.1	ng/ul	901253	1	7.934	1054710	20.0
Benzene, 1,2-di...	8.287	21.0	ng/ul	1107350	1	7.934	1054710	20.0
Propane, 1,1'-o...	8.522	8.5	ng/ul	449808	1	7.934	1054710	20.0
Benzene, 1,2,4-...	10.604	15.1	ng/ul	1132940	2	10.740	1504750	20.0
unknown-01	12.687	4.3	ng/ul	440267	3	14.569	2054470	20.0
unknown-02	12.716	2.4	ng/ul	250745	3	14.569	2054470	20.0
1-Hexanamine, N...	12.910	18.0	ng/ul	1851060	3	14.569	2054470	20.0
unknown-03	13.163	15.0	ng/ul	1542040	3	14.569	2054470	20.0
.alpha.-Lindane	17.069	27.6	ng/ul	3397760	4	17.322	2461740	20.0