

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
 Data File : BP004391.D
 Acq On : 19 Dec 2020 13:22
 Operator : CG/JU
 Sample : L5128-05
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8C6

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : 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\SOM-EPA-BP121820MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.481	216	220	235	rVB	110001	183677	2.21%	0.215%
2	5.158	331	335	343	rVB	50065	77630	0.93%	0.091%
3	6.316	529	532	542	rVB3	23284	42848	0.52%	0.050%
4	6.993	638	647	660	rVB	730238	1425707	17.15%	1.666%
5	7.152	668	674	684	rBV	885310	1433389	17.24%	1.675%
6	7.357	699	709	717	rBV	1103508	1810458	21.78%	2.116%
7	7.428	717	721	727	rVV4	25365	53358	0.64%	0.062%
8	7.481	727	730	736	rVB5	17911	27960	0.34%	0.033%
9	7.681	759	764	768	rBV6	9436	18403	0.22%	0.022%
10	7.822	781	788	795	rBV	804780	1270173	15.28%	1.484%
11	7.887	795	799	807	rVV4	37444	85600	1.03%	0.100%
12	7.963	808	812	821	rVB	26806	53510	0.64%	0.063%
13	8.199	846	852	858	rVV	58269	106032	1.28%	0.124%
14	8.281	862	866	872	rVV8	7810	19641	0.24%	0.023%
15	8.357	875	879	885	rVB6	8166	17270	0.21%	0.020%
16	8.469	893	898	901	rBV3	26025	42328	0.51%	0.049%
17	8.528	901	908	918	rVB	1273955	2029344	24.41%	2.372%
18	8.716	934	940	948	rVB5	20343	47509	0.57%	0.056%
19	8.804	948	955	965	rBV	325589	564921	6.80%	0.660%
20	8.922	970	975	978	rBV2	45165	71723	0.86%	0.084%
21	8.975	978	984	995	rVV	1073638	1845109	22.19%	2.156%
22	9.181	1011	1019	1026	rVB10	13190	41749	0.50%	0.049%
23	9.334	1039	1045	1051	rBV3	57775	100185	1.21%	0.117%
24	9.416	1051	1059	1072	rVB3	102217	220142	2.65%	0.257%
25	9.563	1079	1084	1089	rVB3	26524	38335	0.46%	0.045%
26	9.699	1100	1107	1121	rBV	944058	1595701	19.19%	1.865%
27	9.828	1121	1129	1138	rBV	23296	59087	0.71%	0.069%
28	10.028	1158	1163	1171	rVB7	22803	54169	0.65%	0.063%
29	10.110	1171	1177	1179	rBV4	18220	33155	0.40%	0.039%
30	10.151	1179	1184	1192	rVB4	41837	102047	1.23%	0.119%
31	10.240	1192	1199	1218	rVB	1667176	2899875	34.88%	3.389%
32	10.399	1218	1226	1227	rBV7	15145	29551	0.36%	0.035%
33	10.528	1243	1248	1256	rVB3	48221	84852	1.02%	0.099%
34	10.616	1256	1263	1270	rBV	1125844	1897824	22.83%	2.218%

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35	10.669	1270	1272	1278	rVB7	11677	17763	0.21%	0.021%
36	10.751	1278	1286	1301	rBV	1412900	2468974	29.70%	2.885%
37	10.875	1301	1307	1314	rBV5	31772	66960	0.81%	0.078%
38	11.040	1329	1335	1342	rVB9	14049	33089	0.40%	0.039%
39	11.181	1352	1359	1361	rBV8	15605	31176	0.38%	0.036%
40	11.328	1380	1384	1389	rBV3	17965	27300	0.33%	0.032%
41	11.463	1395	1407	1412	rBV5	31328	95013	1.14%	0.111%
42	11.581	1423	1427	1435	rVB10	18666	38015	0.46%	0.044%
43	11.681	1440	1444	1450	rVB7	13046	28150	0.34%	0.033%
44	11.822	1463	1468	1473	rVB3	25598	38409	0.46%	0.045%
45	11.875	1473	1477	1480	rBV4	10862	17636	0.21%	0.021%
46	11.963	1484	1492	1498	rVV	245203	436046	5.25%	0.510%
47	12.010	1498	1500	1505	rVB2	24429	31495	0.38%	0.037%
48	12.122	1513	1519	1525	rBV2	94181	174773	2.10%	0.204%
49	12.228	1525	1537	1546	rVV10	56256	180770	2.17%	0.211%
50	12.328	1550	1554	1560	rVB5	28023	53600	0.64%	0.063%
51	12.387	1561	1564	1569	rVV3	18645	28845	0.35%	0.034%
52	12.498	1578	1583	1588	rBV6	32721	51560	0.62%	0.060%
53	12.569	1589	1595	1598	rVV7	15211	32388	0.39%	0.038%
54	12.616	1599	1603	1607	rVV3	35349	54467	0.66%	0.064%
55	12.657	1607	1610	1615	rVV6	18863	31150	0.37%	0.036%
56	12.722	1616	1621	1625	rVB4	26972	51599	0.62%	0.060%
57	12.804	1631	1635	1639	rBV6	18198	33427	0.40%	0.039%
58	13.175	1694	1698	1708	rVV6	32459	94337	1.13%	0.110%
59	13.269	1708	1714	1717	rVV4	43476	91156	1.10%	0.107%
60	13.310	1718	1721	1728	rVV3	40674	82464	0.99%	0.096%
61	13.369	1728	1731	1735	rVV5	11748	20919	0.25%	0.024%
62	13.422	1736	1740	1747	rVB7	19459	40225	0.48%	0.047%
63	13.522	1751	1757	1761	rBV5	39016	70780	0.85%	0.083%
64	13.651	1774	1779	1781	rBV5	23865	36202	0.44%	0.042%
65	13.692	1782	1786	1791	rBV6	30746	60324	0.73%	0.070%
66	13.875	1812	1817	1822	rBV	2771524	3860310	46.43%	4.511%
67	13.922	1822	1825	1830	rVB	142199	175739	2.11%	0.205%
68	14.034	1837	1844	1847	rBV6	45693	99566	1.20%	0.116%
69	14.163	1856	1866	1879	rVB	3209447	4945070	59.48%	5.779%
70	14.281	1881	1886	1893	rBV2	97181	181430	2.18%	0.212%
71	14.392	1899	1905	1909	rBV3	114158	180433	2.17%	0.211%

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72	14.469	1910	1918	1926	rVB	1739206	2656445	31.95%	3.104%
73	14.681	1948	1954	1962	rBV	673028	1076327	12.95%	1.258%
74	15.216	2040	2045	2052	rBV	1331414	1857549	22.34%	2.171%
75	15.334	2062	2065	2067	rBV3	23691	29292	0.35%	0.034%
76	15.469	2081	2088	2097	rVB	3938959	5805231	69.83%	6.784%
77	15.592	2103	2109	2122	rBV	1083428	1585221	19.07%	1.853%
78	15.698	2122	2127	2142	rVB2	140735	298434	3.59%	0.349%
79	15.986	2169	2176	2187	rBV	1373007	1957405	23.55%	2.287%
80	16.157	2198	2205	2210	rBV2	315205	514752	6.19%	0.602%
81	16.316	2228	2232	2237	rBV	70561	96952	1.17%	0.113%
82	16.498	2257	2263	2267	rBV	768925	1049623	12.63%	1.227%
83	16.545	2267	2271	2277	rVB	153547	220988	2.66%	0.258%
84	16.763	2305	2308	2318	rVB	89973	141471	1.70%	0.165%
85	17.028	2350	2353	2359	rVB	108974	150430	1.81%	0.176%
86	17.228	2382	2387	2397	rVB	2172987	3111465	37.43%	3.636%
87	17.328	2398	2404	2414	rVB2	4512426	6493987	78.11%	7.589%
88	17.769	2477	2479	2483	rVB3	32155	29479	0.35%	0.034%
89	18.122	2536	2539	2545	rBV3	83372	135502	1.63%	0.158%
90	18.516	2603	2606	2610	rVB	38557	48103	0.58%	0.056%
91	18.875	2664	2667	2676	rVB2	89027	123571	1.49%	0.144%
92	19.569	2779	2785	2790	rBV	5963525	8019940	96.47%	9.372%
93	19.610	2790	2792	2804	rVB	633986	850913	10.24%	0.994%
94	20.563	2952	2954	2963	rVB3	79399	120509	1.45%	0.141%
95	21.033	3031	3034	3042	rBV2	393084	621766	7.48%	0.727%
96	21.098	3043	3045	3051	rVB	128079	167119	2.01%	0.195%
97	21.321	3078	3083	3090	rVB	2894397	3624332	43.60%	4.235%
98	22.551	3288	3292	3301	rVB	256636	373796	4.50%	0.437%
99	23.527	3451	3458	3471	rVB2	4088755	8313404	100.00%	9.715%
100	23.674	3476	3483	3495	rVB	1920532	3751715	45.13%	4.384%

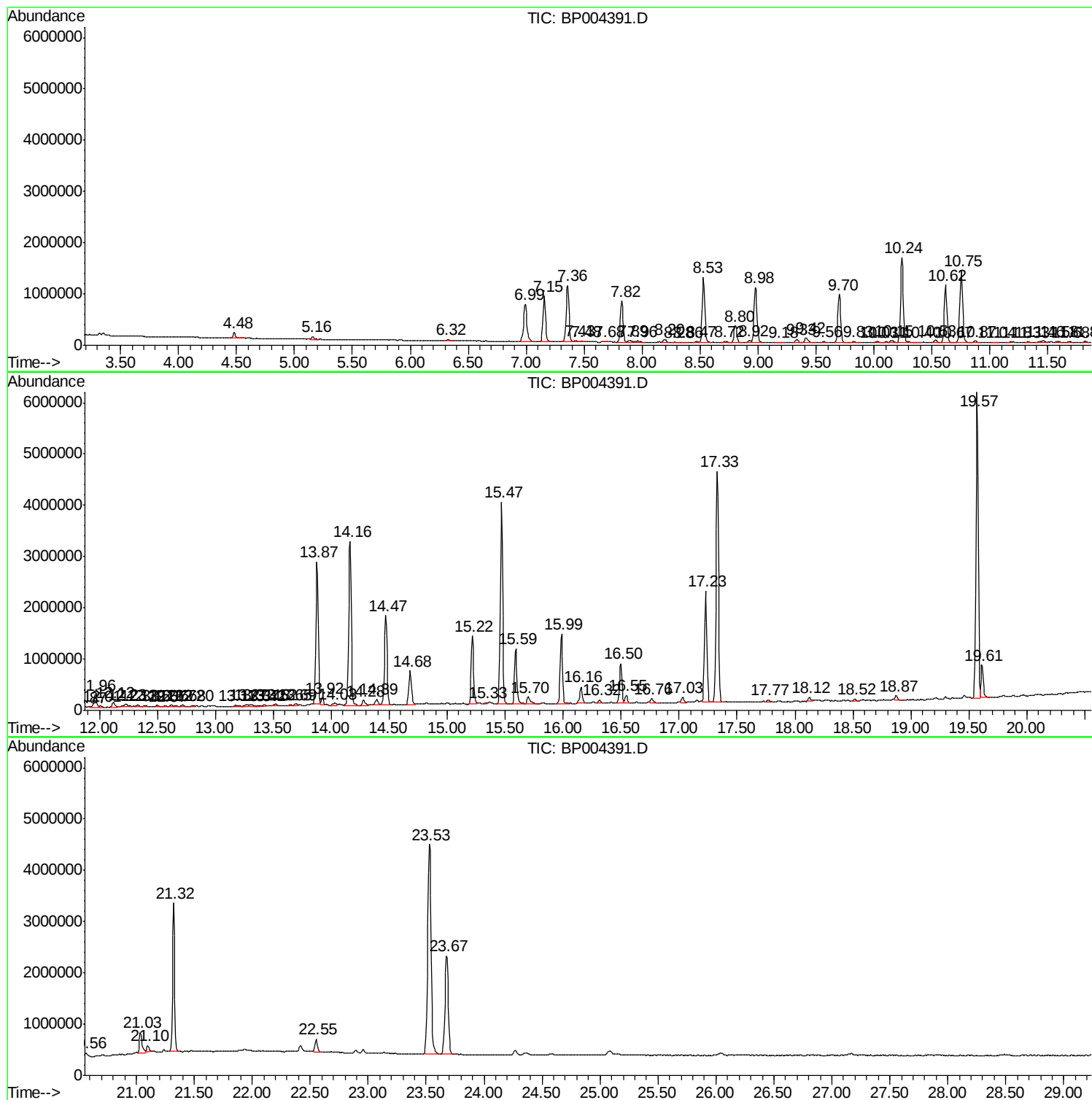
Sum of corrected areas: 85570543

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Instrument :
BNA_P
ClientSampled :
CB8C6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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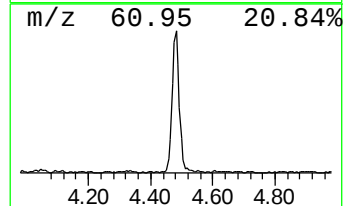
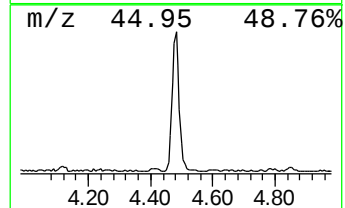
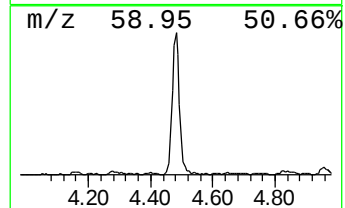
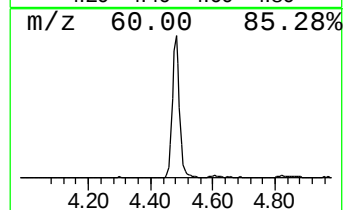
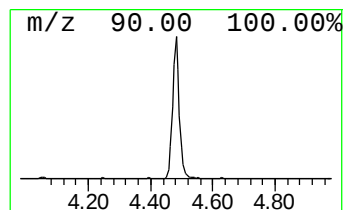
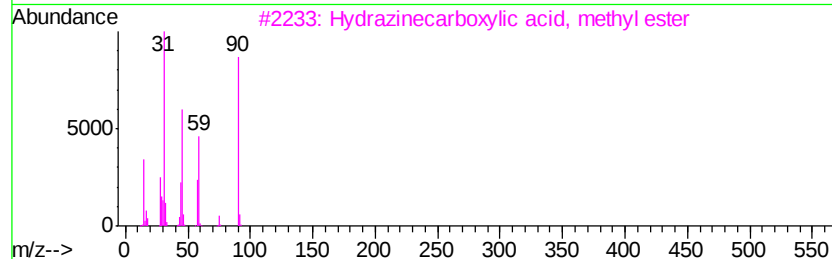
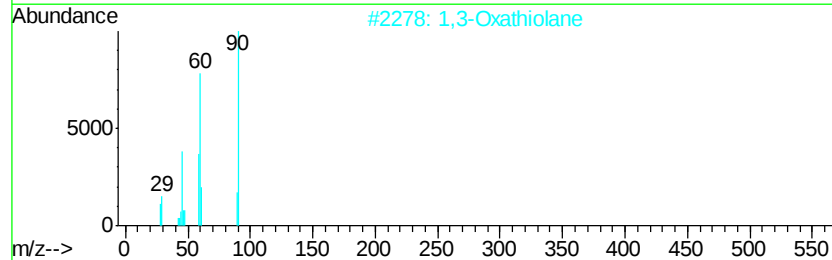
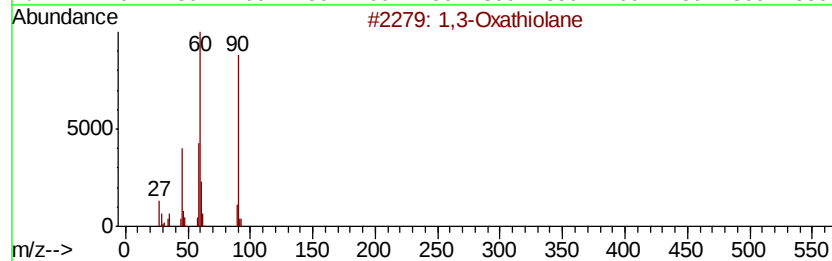
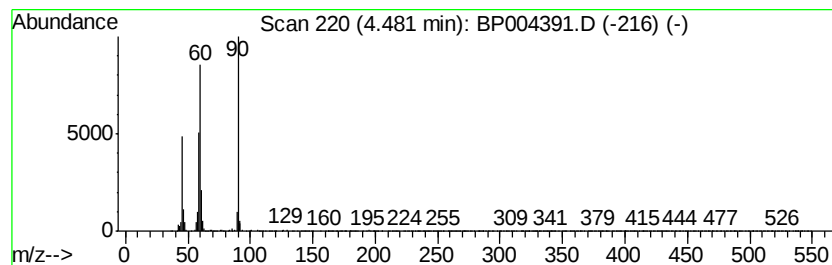
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 1,3-Oxathiolane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.48	2.89 ng/ul	183677	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Oxathiolane	90	C3H6OS	002094-97-5	90
2		1,3-Oxathiolane	90	C3H6OS	002094-97-5	74
3		Hvdrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	59
4		Heptanoic acid, 2-hydroxy-, methyl...	160	C8H16O3	054340-91-9	53
5		Thiourea, methyl-	90	C2H6N2S	000598-52-7	50



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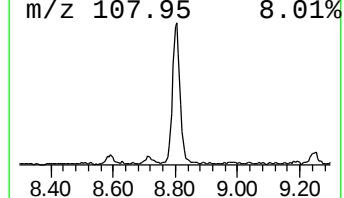
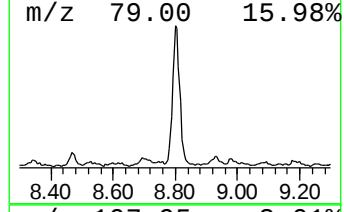
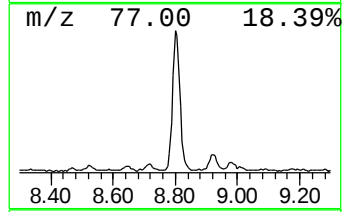
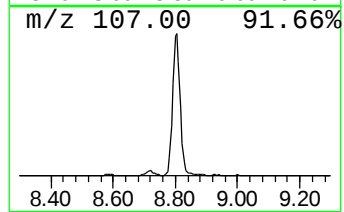
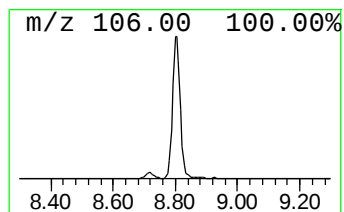
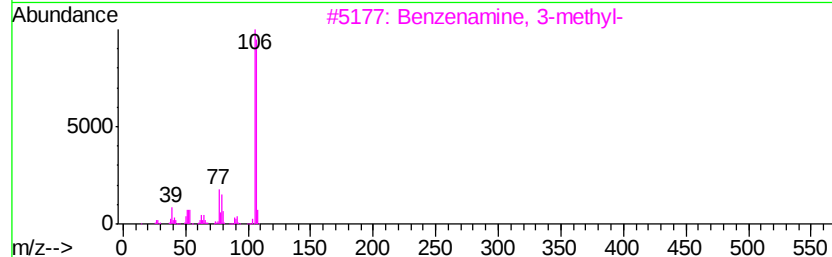
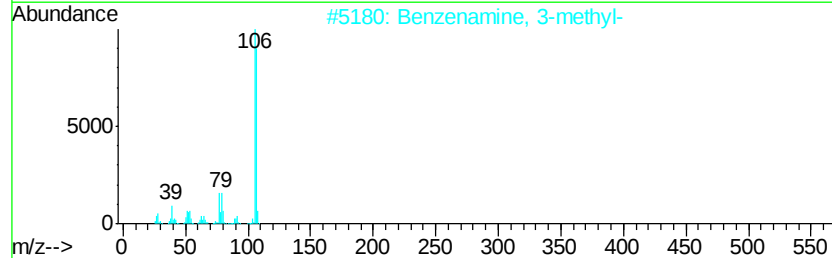
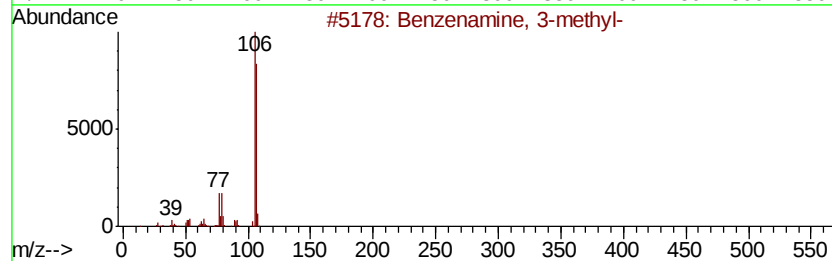
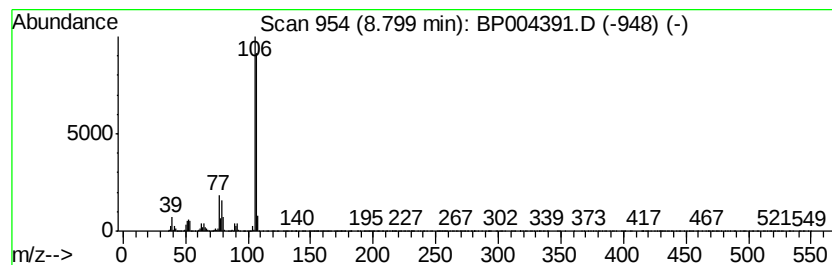
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzenamine, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	8.90 ng/ul	564921	1,4-Dichlorobenzene-d4	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
2		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
3		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	97
4		Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	95
5		o-Toluidine	107	C7H9N	000095-53-4	94



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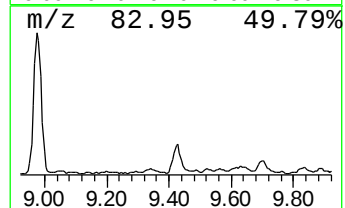
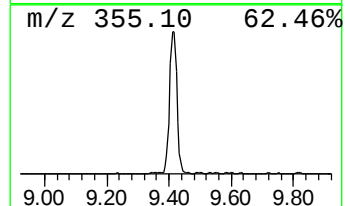
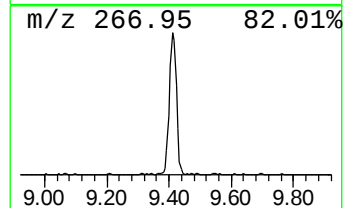
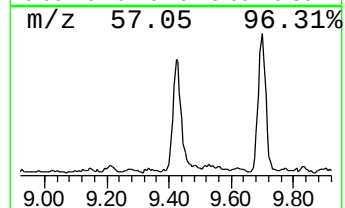
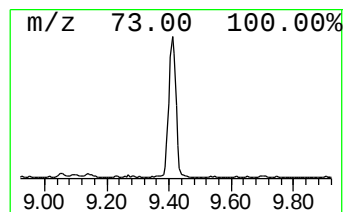
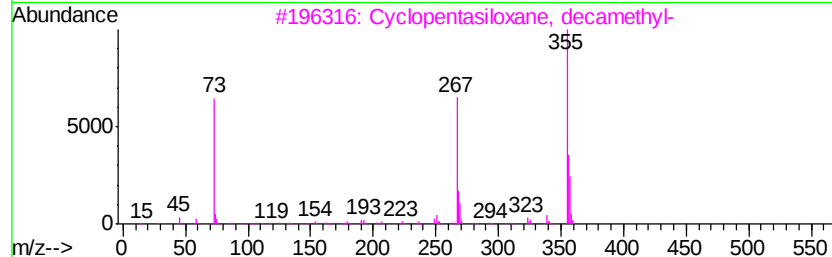
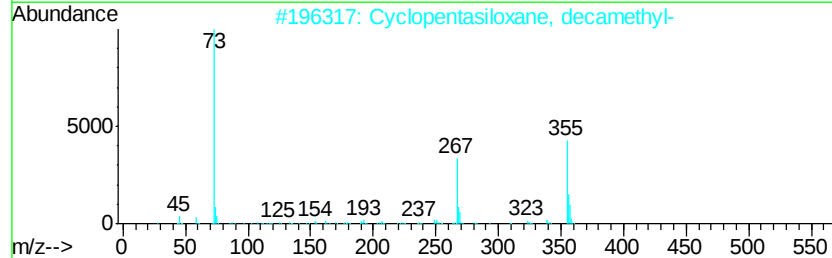
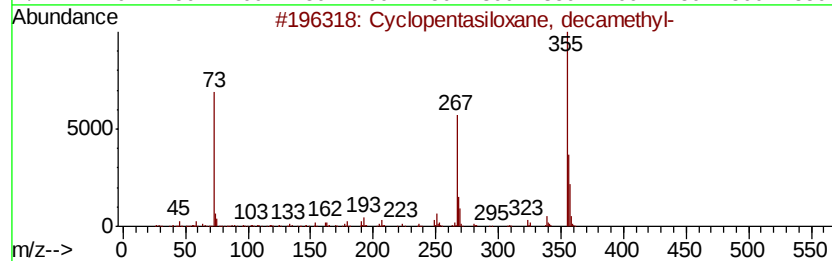
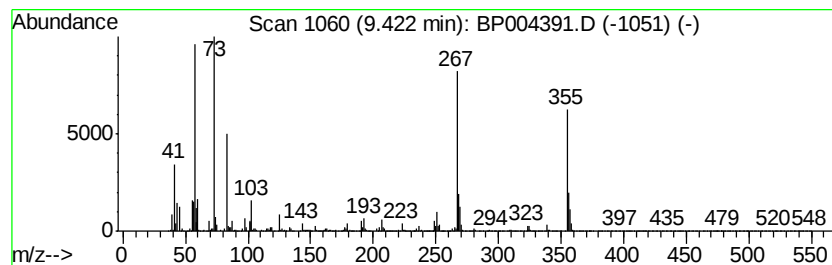
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Cyclopentasiloxane, decamet... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.42	2.32 ng/ul	220142	Naphthalene-d8	10.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
2		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
3		Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	91
4		3-Amino-2-phenazinol ditms	355	C18H25N3OSi2	1000332-15-5	47
5		Benzaldehyde, 2,4-bis(trimethyls...	282	C13H22O3Si2	033617-38-8	38



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Data File : BP004391.D
Acq On : 19 Dec 2020 13:22
Operator : CG/JU
Sample : L5128-05
Misc :
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ClientSampleId :
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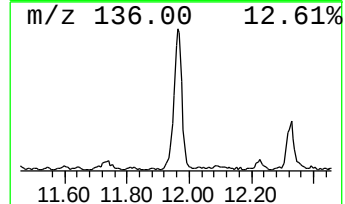
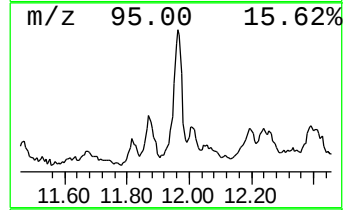
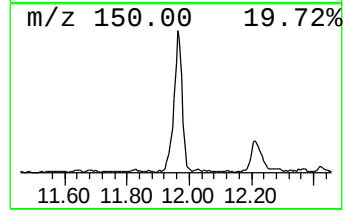
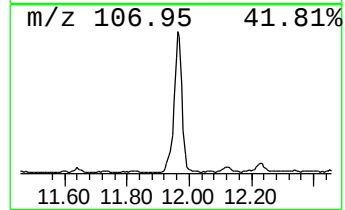
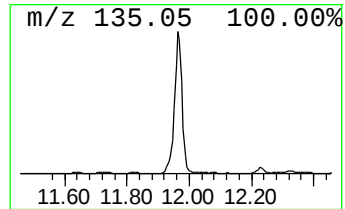
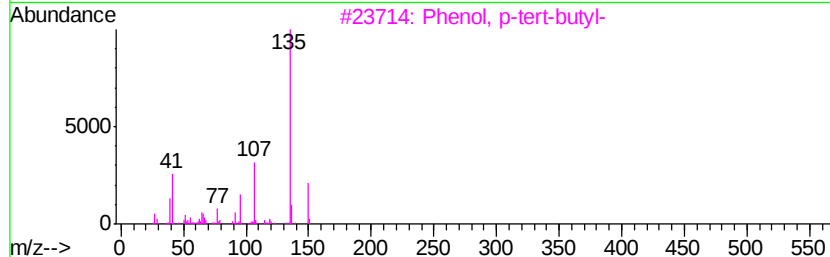
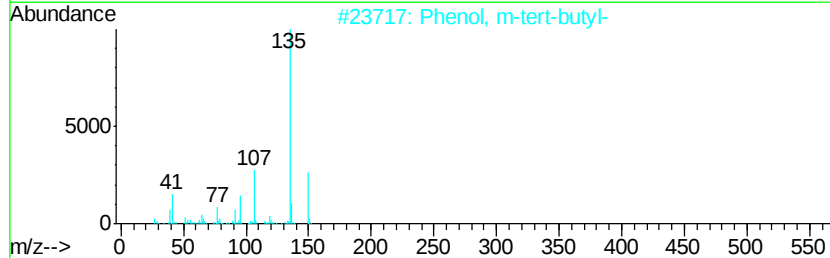
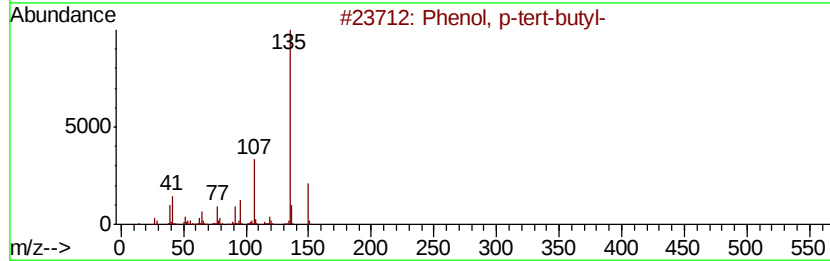
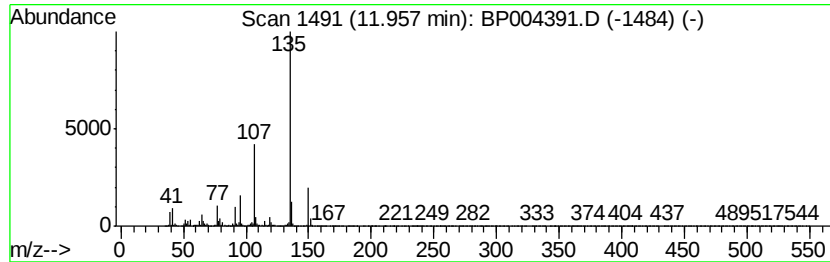
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenol, p-tert-butyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	4.60 ng/ul	436046	Napthalene-d8	10.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	96
2		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
3		Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	94
4		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91
5		Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	91



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Data File : BP004391.D
Acq On : 19 Dec 2020 13:22
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Misc :
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ClientSampleId :
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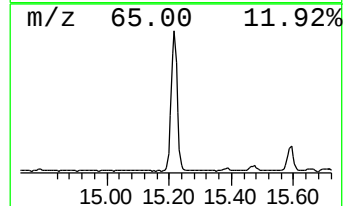
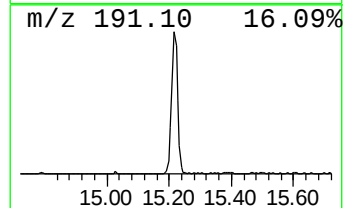
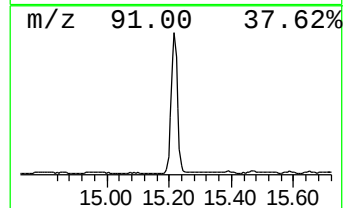
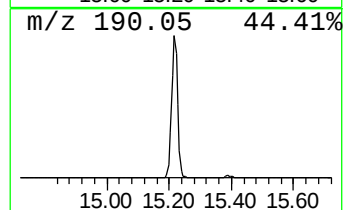
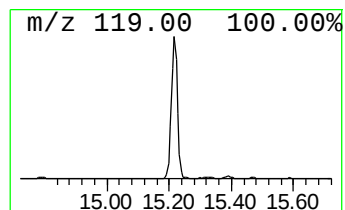
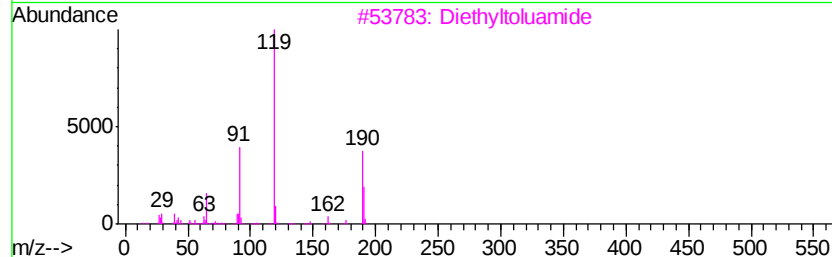
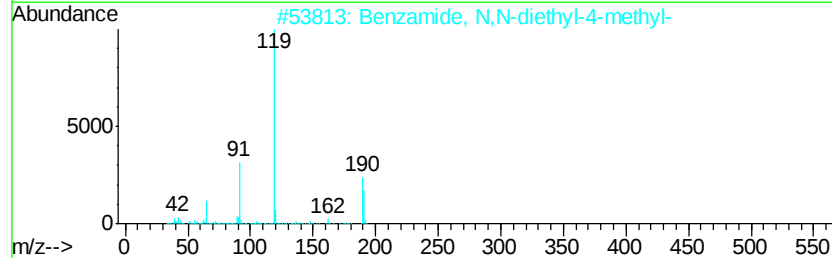
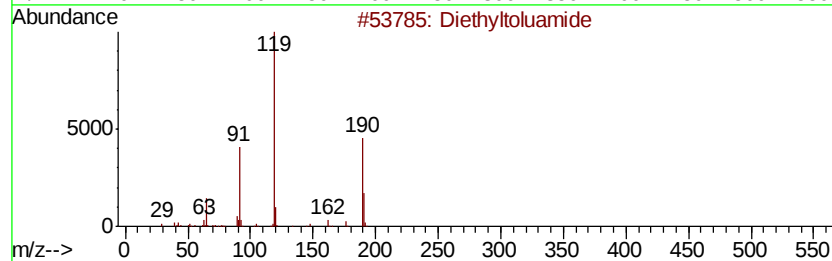
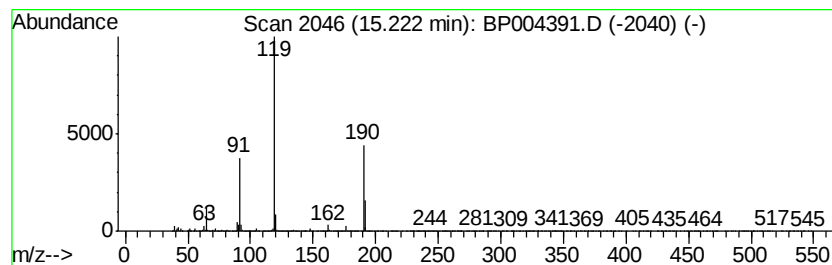
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Diethyltoluamide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.22	13.99 ng/ul	1857550	Acenaphthene-d10	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	97
2		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
3		Diethyltoluamide	191	C12H17NO	000134-62-3	91
4		Diethyltoluamide	191	C12H17NO	000134-62-3	60
5		Benzamide, N-(1,1-dimethylethyl)...	191	C12H17NO	042498-32-8	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004391.D
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Misc :
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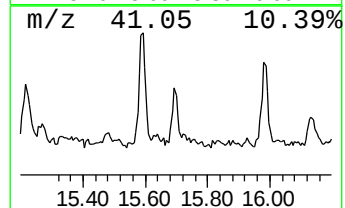
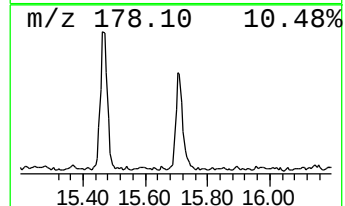
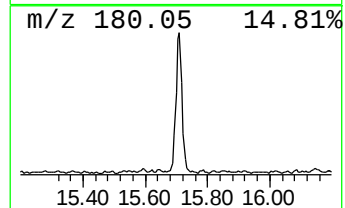
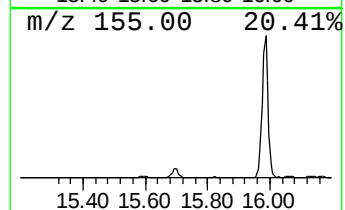
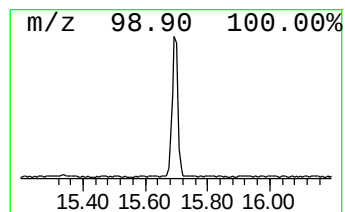
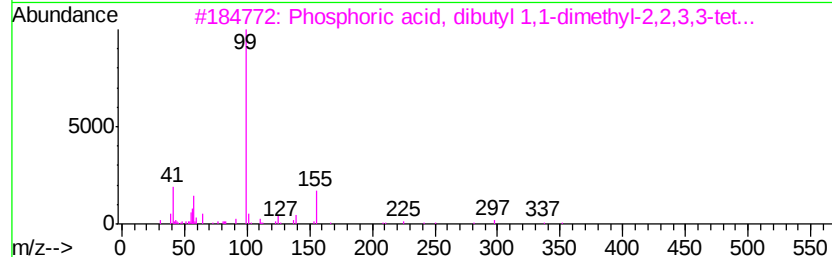
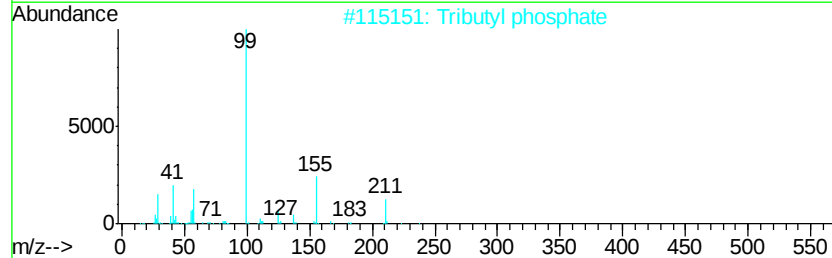
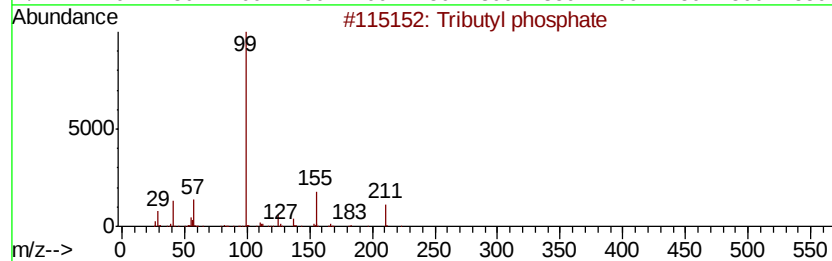
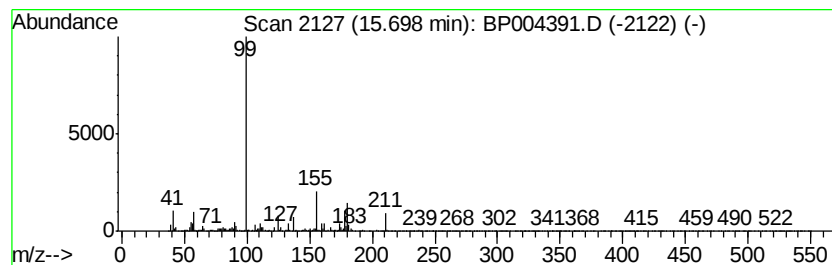
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Tributyl phosphate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	2.25 ng/ul	298434	Acenaphthene-d10	14.47

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tributyl phosphate	266	C12H27O4P	000126-73-8	64
2		Tributyl phosphate	266	C12H27O4P	000126-73-8	43
3		Phosphoric acid, dibutyl 1,1-dim...	352	C13H25F4O4P	1000298-93-9	40
4		Tributyl phosphate	266	C12H27O4P	000126-73-8	38
5		Phosphoric acid, dibutyl 3-trifl...	348	C14H28F3O4P	1000298-93-8	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004391.D
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Sample : L5128-05
Misc :
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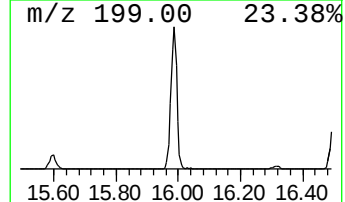
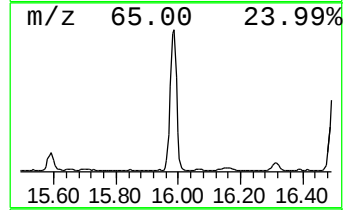
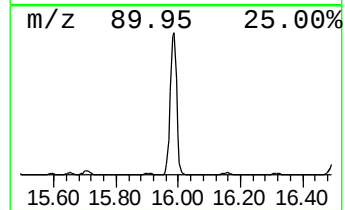
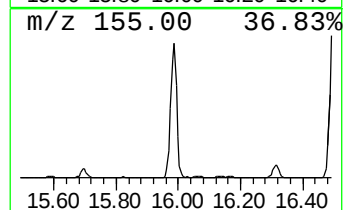
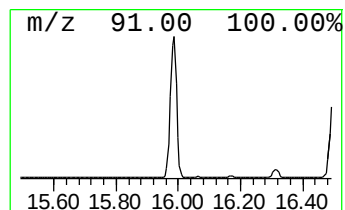
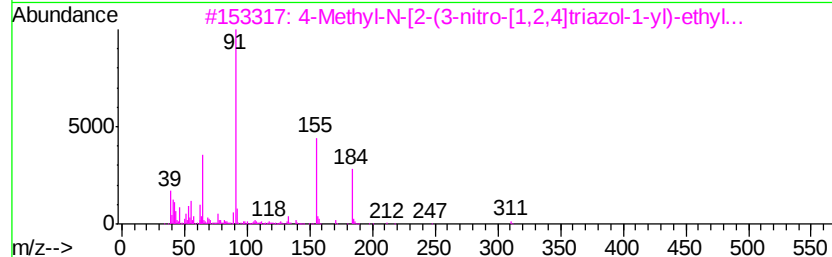
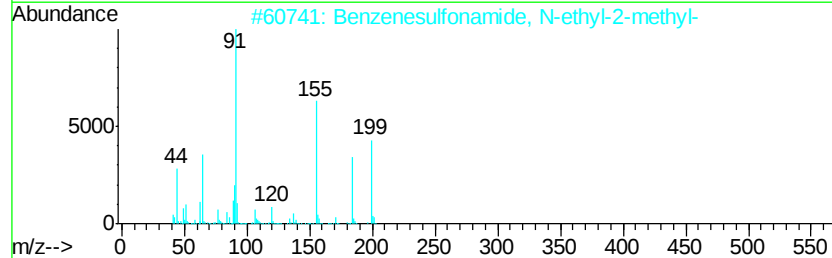
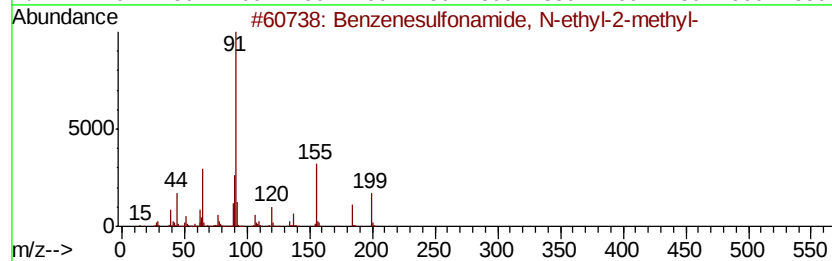
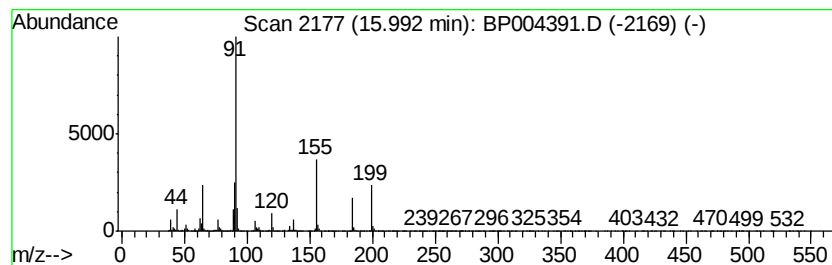
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzenesulfonamide, N-ethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	12.58 ng/ul	1957410	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	97
2		Benzenesulfonamide, N-ethyl-2-me...	199	C9H13NO2S	001077-56-1	52
3		4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	47
4		4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	47
5		4-Toluenesulfonylmethyl isocyanide	195	C9H9NO2S	036635-61-7	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121820\
Data File : BP004391.D
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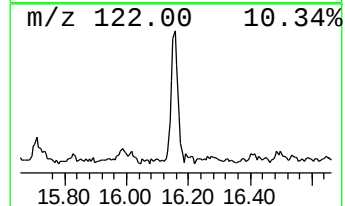
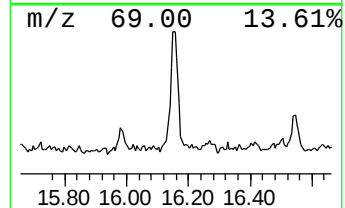
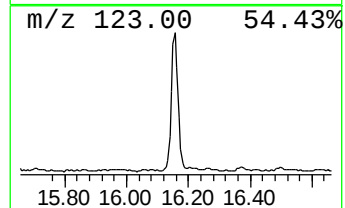
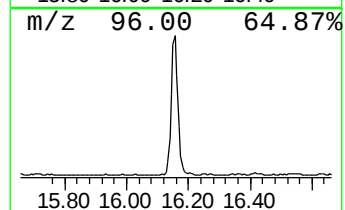
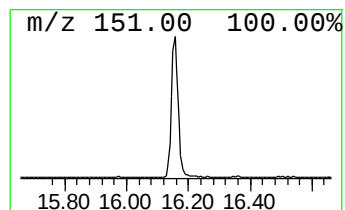
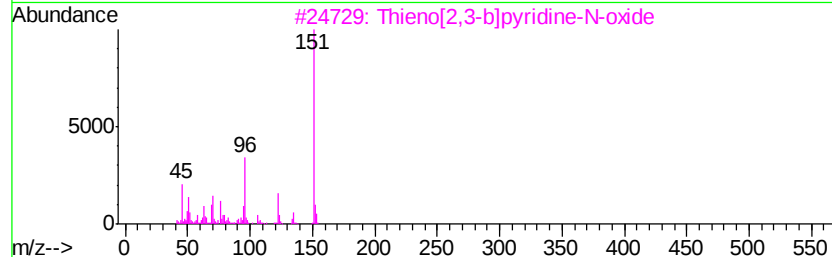
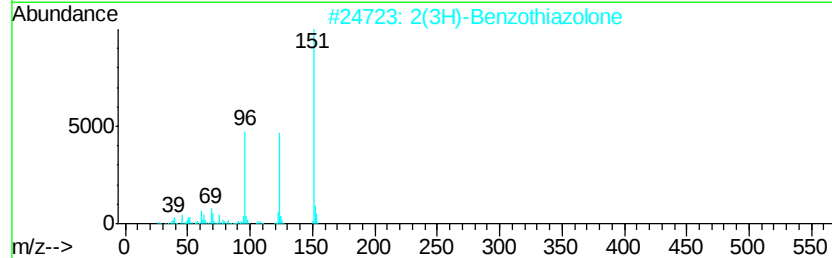
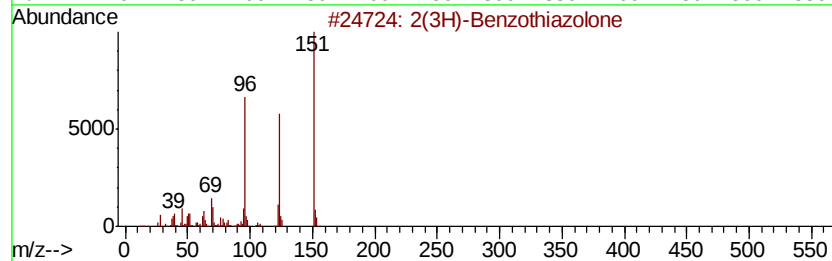
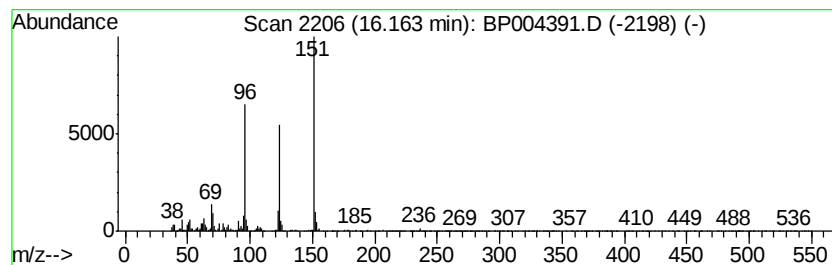
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 2(3H)-Benzothiazolone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	3.31 ng/ul	514752	Phenanthrene-d10	17.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2		2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
3		Thieno[2,3-b]pyridine-N-oxide	151	C7H5NOS	025557-50-0	47
4		4-Methoxyformanilide	151	C8H9NO2	005470-34-8	46
5		2-Aminothiazolo(5,4-b)pyridine	151	C6H5N3S	031784-70-0	38



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Data File : BP004391.D
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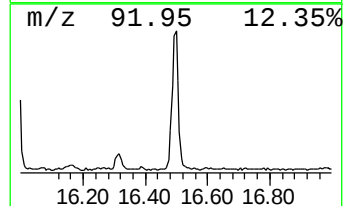
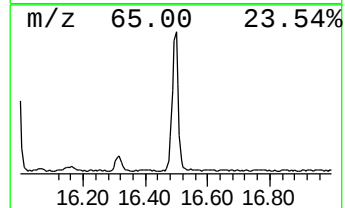
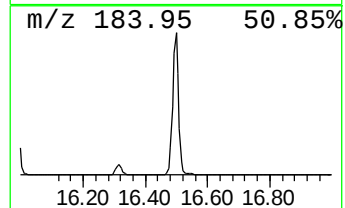
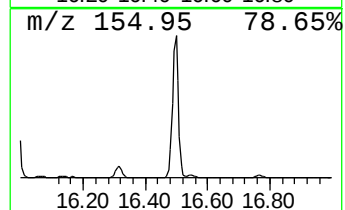
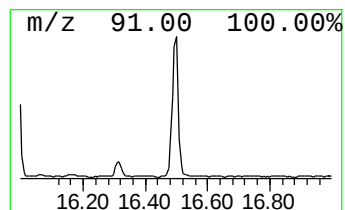
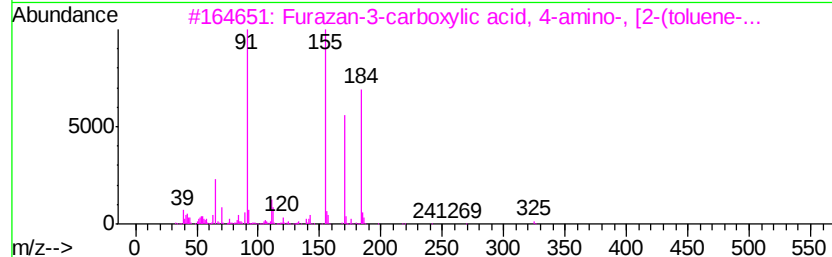
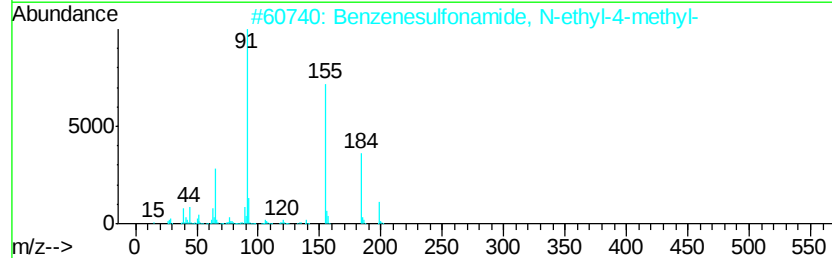
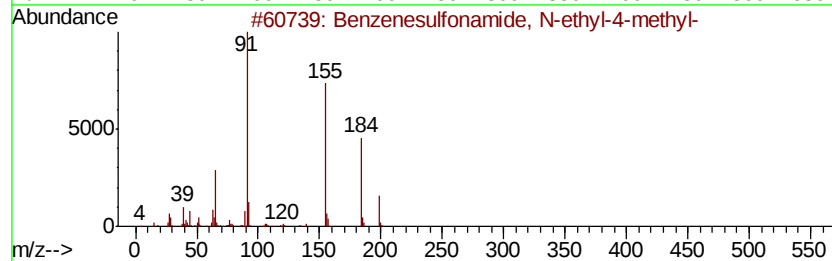
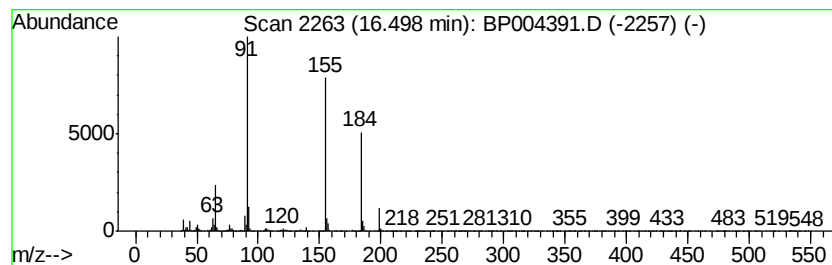
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzenesulfonamide, N-ethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	6.75 ng/ul	1049620	Phenanthrene-d10	17.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	96
3			Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	83
4			4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83
5			(Toluene-4-sulfonylamino)-acetic...	283	C13H17N04S	1000190-48-0	72



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Data File : BP004391.D
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Sample : L5128-05
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ALS Vial : 38 Sample Multiplier: 1

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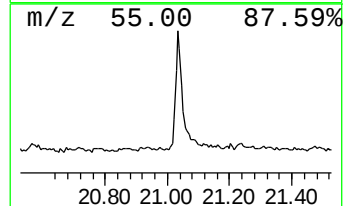
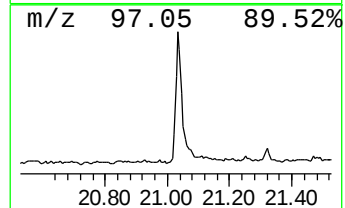
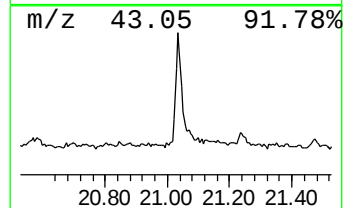
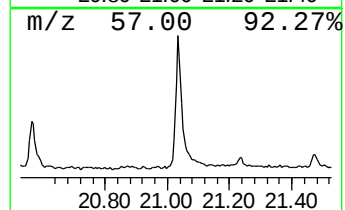
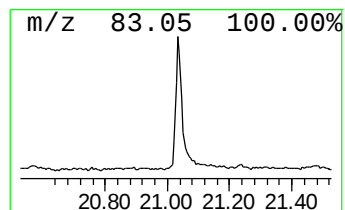
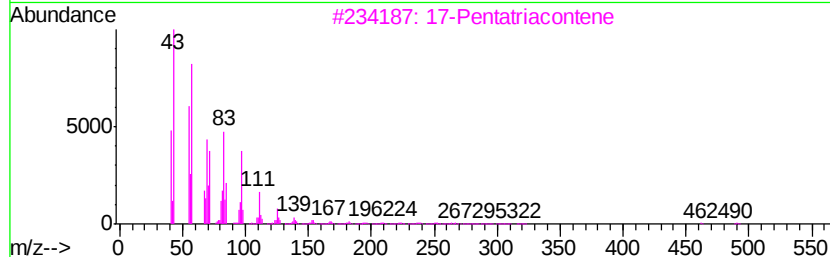
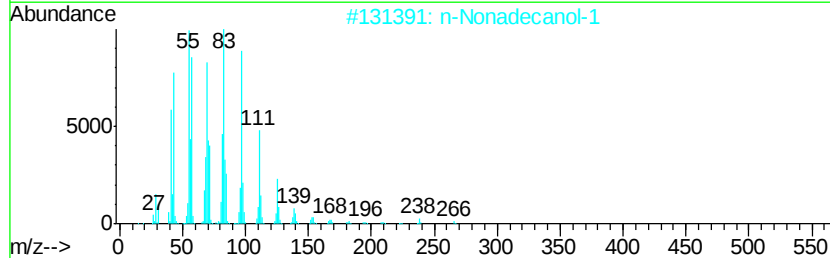
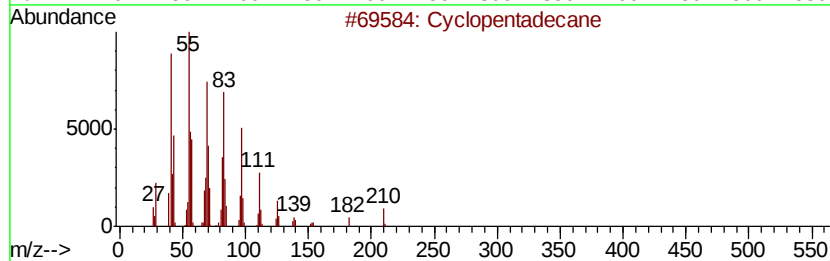
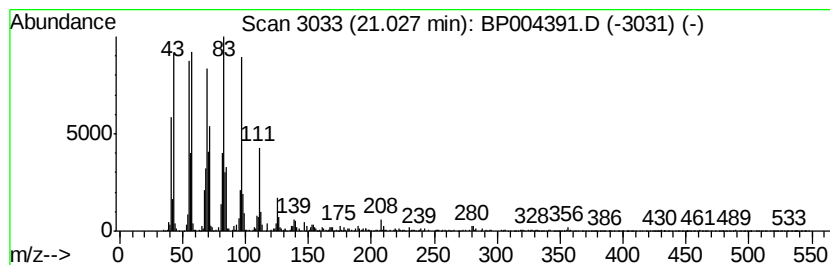
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Cyclic21.03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.03	3.43 ng/ul	621766	Chrysene-d12	21.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentadecane	210	C15H30	000295-48-7	94
2		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
3		17-Pentatriacontene	491	C35H70	006971-40-0	94
4		Behenic alcohol	326	C22H46O	000661-19-8	94
5		1-Heneicosanol	312	C21H44O	015594-90-8	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121820\
 Data File : BP004391.D
 Acq On : 19 Dec 2020 13:22
 Operator : CG/JU
 Sample : L5128-05
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB8C6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121820MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1,3-Oxathiolane	4.48	2.9	ng/ul	183677	1	7.82	1270170	20.0
Benzenamine, 3-me...	8.80	8.9	ng/ul	564921	1	7.82	1270170	20.0
Cyclopentasiloxan...	9.42	2.3	ng/ul	220142	2	10.62	1897820	20.0
Phenol, p-tert-bu...	11.96	4.6	ng/ul	436046	2	10.62	1897820	20.0
Diethyltoluamide	15.22	14.0	ng/ul	1857550	3	14.47	2656450	20.0
Tributyl phosphate	15.70	2.3	ng/ul	298434	3	14.47	2656450	20.0
Benzenesulfonamid...	15.99	12.6	ng/ul	1957410	4	17.23	3111470	20.0
2(3H)-Benzothiazo...	16.16	3.3	ng/ul	514752	4	17.23	3111470	20.0
Benzenesulfonamid...	16.50	6.8	ng/ul	1049620	4	17.23	3111470	20.0
(DEL) Alkane: Cyc...	21.03	3.4	ng/ul	621766	5	21.32	3624330	20.0