

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
 Data File : BM000416.D
 Acq On : 03 Feb 2015 14:29
 Operator : TP/IZ
 Sample : G1147-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC020

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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.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	2.751	5	11	30	rBV2	49723144	90982536	100.00%	32.687%
2	2.881	30	33	42	rVV	397695	704924	0.77%	0.253%
3	2.957	42	46	55	rVB	760807	975093	1.07%	0.350%
4	3.216	86	90	96	rBV	77237	101253	0.11%	0.036%
5	3.775	180	185	191	rBV	77176	104845	0.12%	0.038%
6	4.216	256	260	268	rBV2	87190	133742	0.15%	0.048%
7	6.163	586	591	606	rBV	268085	621874	0.68%	0.223%
8	6.751	686	691	700	rVB3	58219	108714	0.12%	0.039%
9	6.922	713	720	736	rBV2	1768240	3771189	4.14%	1.355%
10	7.087	742	748	760	rVB	1199817	1826164	2.01%	0.656%
11	7.286	774	782	786	rBV	1770179	3146519	3.46%	1.130%
12	7.639	835	842	848	rBV	633783	992503	1.09%	0.357%
13	7.704	848	853	856	rVV	213645	325782	0.36%	0.117%
14	7.751	856	861	864	rVV	726129	1200664	1.32%	0.431%
15	7.786	864	867	874	rVB	677325	996837	1.10%	0.358%
16	8.104	914	921	929	rBV	764810	1240421	1.36%	0.446%
17	8.286	940	952	956	rBV2	588591	1470195	1.62%	0.528%
18	8.333	956	960	969	rVB	341294	529869	0.58%	0.190%
19	8.457	969	981	992	rBV	2319821	3778146	4.15%	1.357%
20	8.828	1036	1044	1050	rVB2	510495	868218	0.95%	0.312%
21	8.910	1050	1058	1065	rBV	1933559	3128990	3.44%	1.124%
22	9.627	1172	1180	1189	rVB	1740748	2871772	3.16%	1.032%
23	10.163	1263	1271	1285	rBV2	2539311	5611512	6.17%	2.016%
24	10.410	1305	1313	1320	rBV	801220	1341581	1.47%	0.482%
25	10.545	1328	1336	1340	rBV	974808	1688911	1.86%	0.607%
26	10.592	1340	1344	1352	rVB	961679	1564239	1.72%	0.562%
27	10.680	1352	1359	1370	rBV	2301173	3997505	4.39%	1.436%
28	10.880	1385	1393	1400	rBV	1159629	1903990	2.09%	0.684%
29	11.251	1451	1456	1460	rBV	85985	133842	0.15%	0.048%
30	12.069	1588	1595	1602	rVB3	143695	259816	0.29%	0.093%
31	12.504	1664	1669	1673	rBV2	219918	348392	0.38%	0.125%
32	12.551	1673	1677	1682	rVB	237844	330039	0.36%	0.119%
33	12.627	1684	1690	1694	rBV2	64057	111094	0.12%	0.040%
34	12.786	1706	1717	1718	rBV2	487947	1087985	1.20%	0.391%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
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 Stop Thrs : 0

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35	13.027	1737	1758	1774	rVB3	771659	3927535	4.32%	1.411%
36	13.257	1792	1797	1803	rVB	117877	161188	0.18%	0.058%
37	13.810	1885	1891	1900	rVB	4542746	6509346	7.15%	2.339%
38	13.980	1913	1920	1925	rBV	1409039	1899272	2.09%	0.682%
39	14.092	1932	1939	1949	rBV	4935820	7235846	7.95%	2.600%
40	14.398	1985	1991	1997	rBV2	1507206	2359849	2.59%	0.848%
41	14.462	1997	2002	2009	rVB	1651948	2284936	2.51%	0.821%
42	14.604	2019	2026	2039	rBV	2426666	3820933	4.20%	1.373%
43	14.768	2047	2054	2061	rVB	1459336	2016051	2.22%	0.724%
44	15.221	2125	2131	2138	rVB	1989707	2642666	2.90%	0.949%
45	15.398	2153	2161	2166	rBV	6383650	9161058	10.07%	3.291%
46	15.451	2166	2170	2175	rVV	1937923	2533537	2.78%	0.910%
47	15.515	2175	2181	2192	rVV	2699816	4027003	4.43%	1.447%
48	15.633	2194	2201	2208	rVB3	63499	113913	0.13%	0.041%
49	15.768	2219	2224	2230	rVB	881092	1047648	1.15%	0.376%
50	16.451	2334	2340	2349	rVB	1709873	2504982	2.75%	0.900%
51	16.804	2388	2400	2410	rBV	2078354	3131099	3.44%	1.125%
52	16.898	2410	2416	2422	rVB	3247381	4391735	4.83%	1.578%
53	17.145	2452	2458	2465	rBV	2195290	3082751	3.39%	1.108%
54	17.245	2468	2475	2479	rVV	6955467	10878647	11.96%	3.908%
55	17.280	2479	2481	2491	rVB	2031861	2254431	2.48%	0.810%
56	17.774	2560	2565	2569	rVB3	76755	104304	0.11%	0.037%
57	18.150	2624	2629	2637	rVB	415244	502214	0.55%	0.180%
58	19.427	2841	2846	2855	rVB3	69279	135646	0.15%	0.049%
59	19.544	2860	2866	2870	rBV	8256825	13166480	14.47%	4.730%
60	19.574	2870	2871	2877	rVB	2941124	2122613	2.33%	0.763%
61	20.474	3018	3024	3029	rBV	4707275	5378939	5.91%	1.932%
62	21.250	3151	3156	3163	rBV	4465629	4927659	5.42%	1.770%
63	21.327	3163	3169	3174	rVV2	3189874	6573772	7.23%	2.362%
64	21.374	3174	3177	3183	rVB	2834535	3260081	3.58%	1.171%
65	22.915	3432	3439	3443	rBV	2814413	4925135	5.41%	1.769%
66	22.956	3443	3446	3456	rVB	1772015	2710564	2.98%	0.974%
67	23.456	3522	3531	3536	rBV2	4626802	9787804	10.76%	3.516%
68	23.497	3536	3538	3546	rVV	1516684	2126969	2.34%	0.764%
69	23.591	3547	3554	3565	rVB	1320101	2603687	2.86%	0.935%
70	25.879	3931	3943	3957	rVB2	1946372	5779510	6.35%	2.076%

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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

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Title : SVOA CALIBRATION

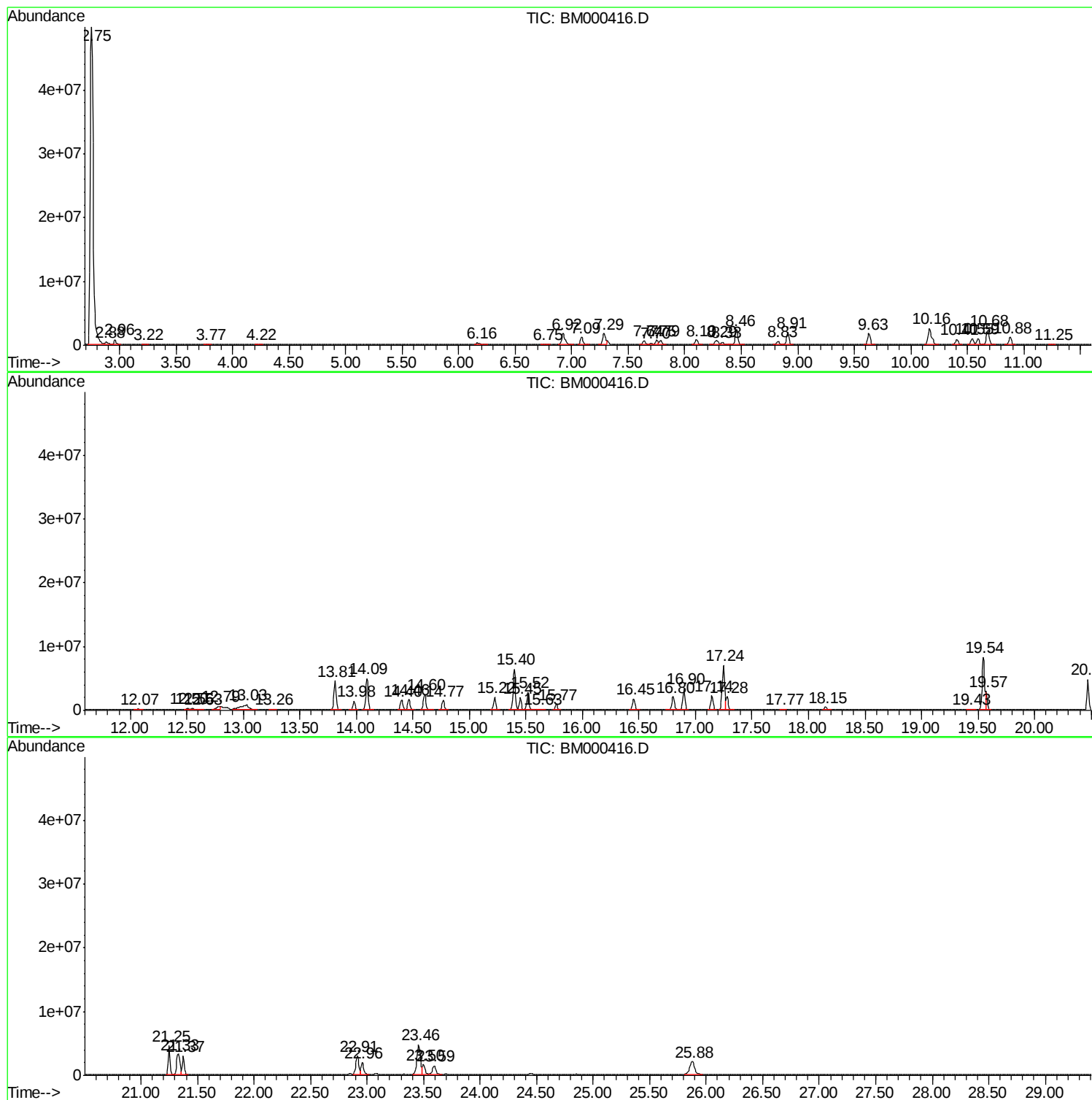
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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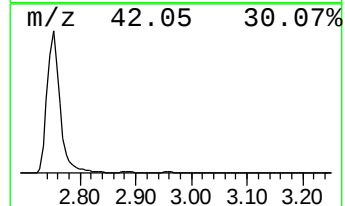
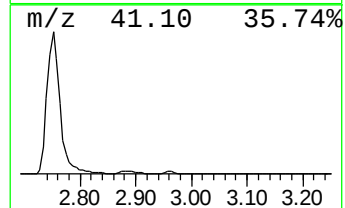
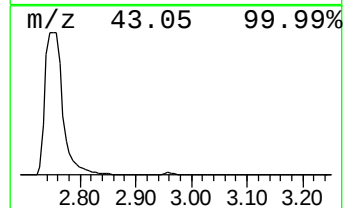
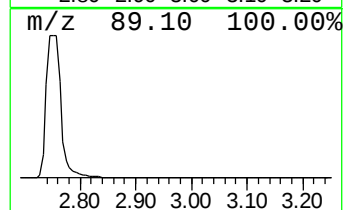
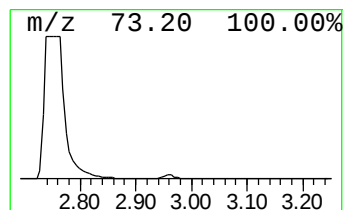
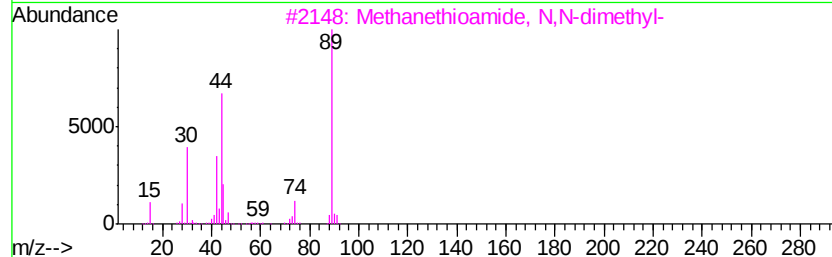
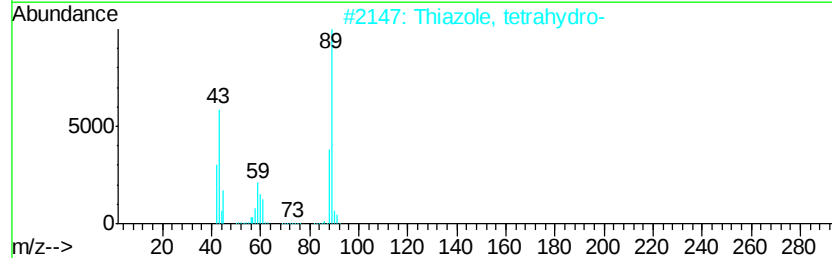
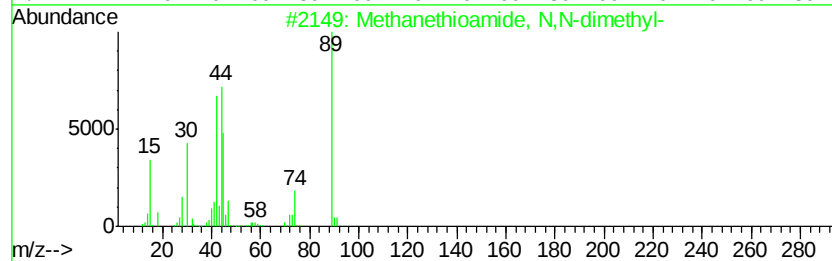
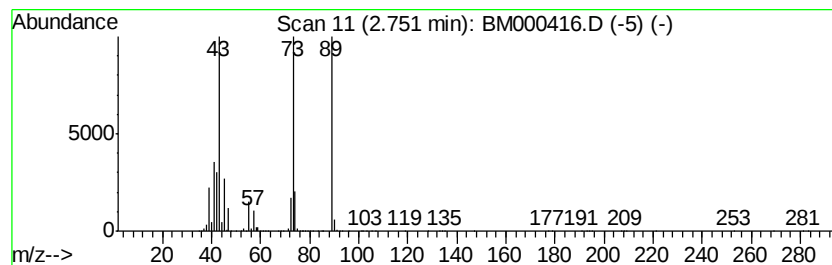
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Methanethioamide, N,N-dimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.75	1515.54 ng/ul	90982500	1,4-Dichlorobenzene-d4	7.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	56
2			Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	50
3			Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	9
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
5			Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	9



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Instrument :
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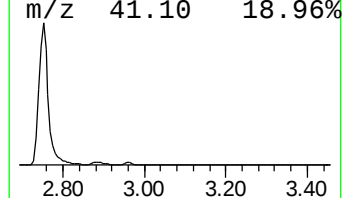
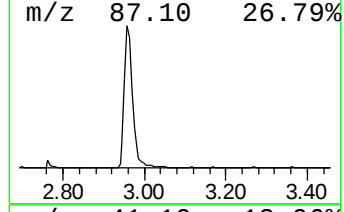
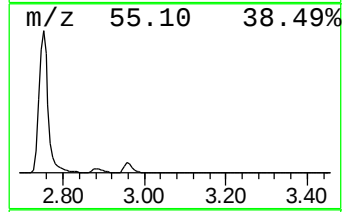
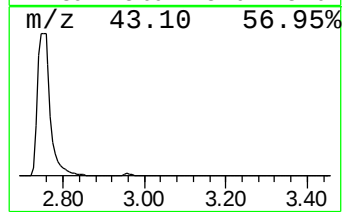
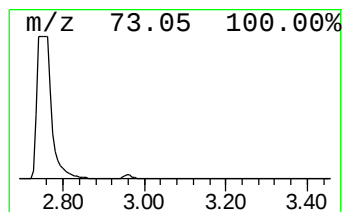
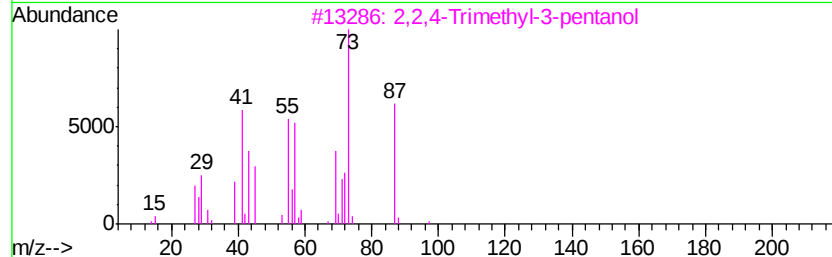
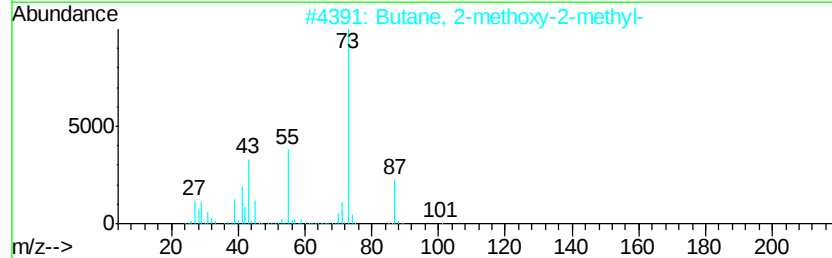
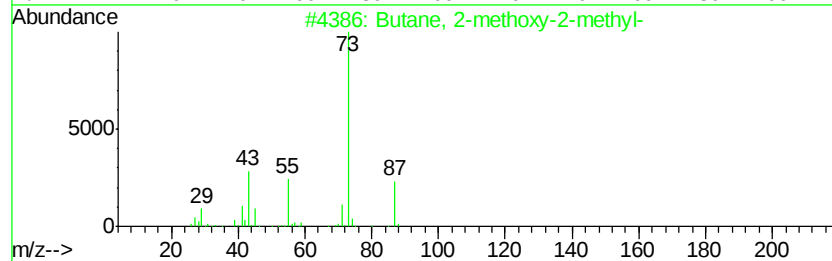
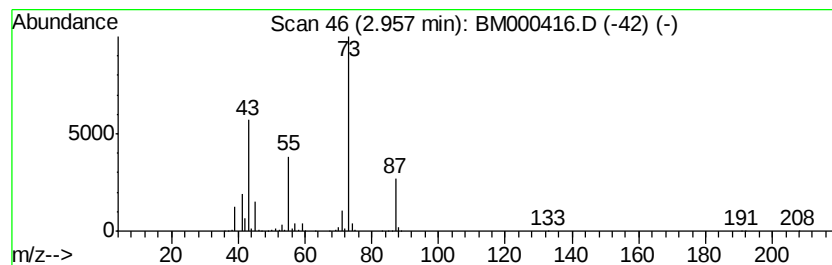
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.96	16.24 ng/ul	975093	1,4-Dichlorobenzene-d4	7.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4		2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	25
5		2,5-Dimethyl-4-hydroxy-3-hexanone	144	C8H16O2	000815-77-0	25



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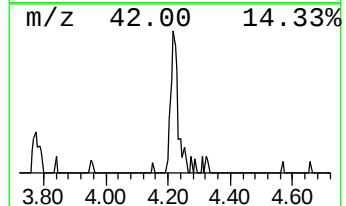
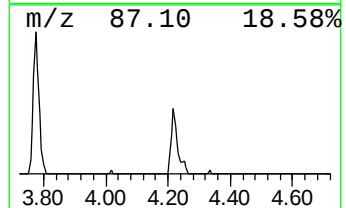
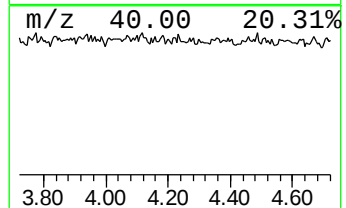
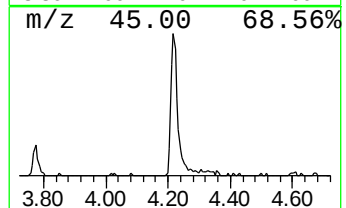
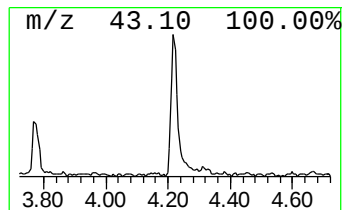
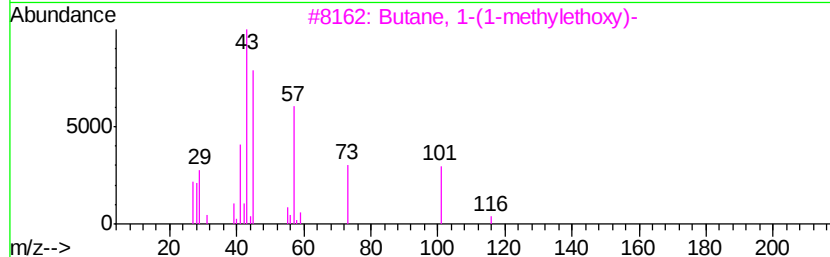
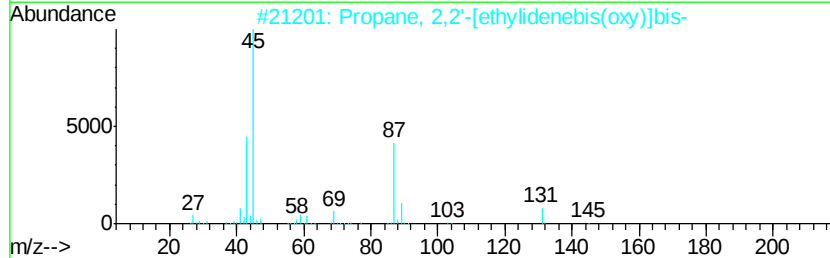
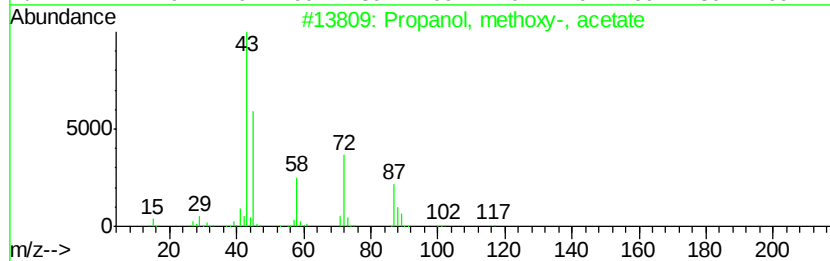
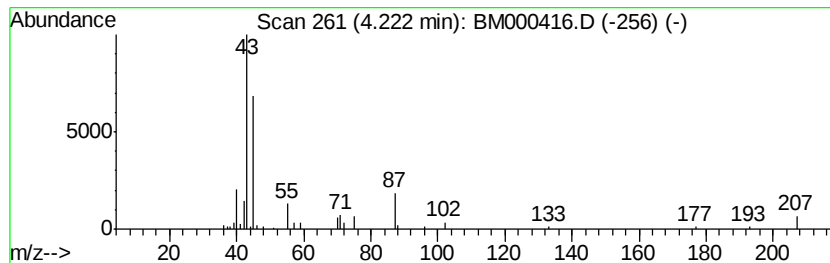
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Propanol, methoxy-, acetate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.22	2.23 ng/ul	133742	1,4-Dichlorobenzene-d4	7.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propanol, methoxy-, acetate	132	C6H12O3	084540-57-8	59
2		Propane, 2,2'-[ethylidenebis(oxy...]	146	C8H18O2	004285-59-0	50
3		Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	40
4		Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	40
5		2-Ethyl-2-methyl-1,3-propanediol	118	C6H14O2	000077-84-9	40



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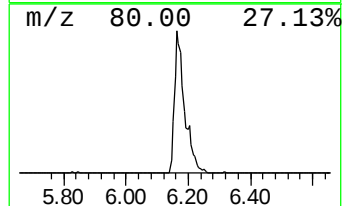
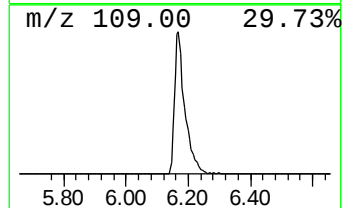
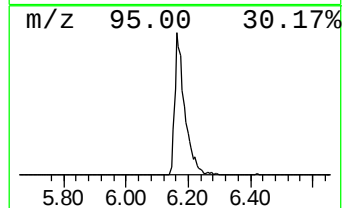
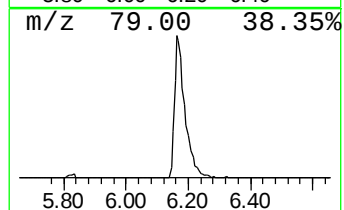
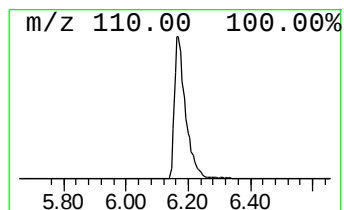
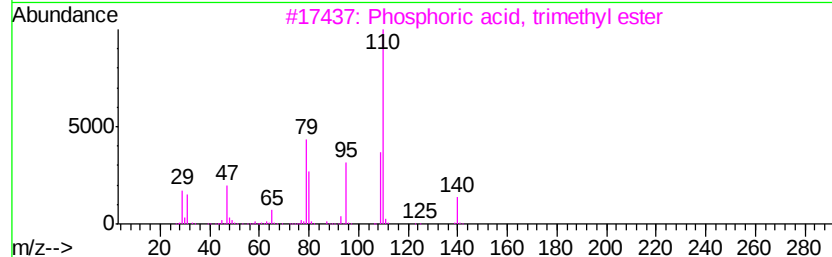
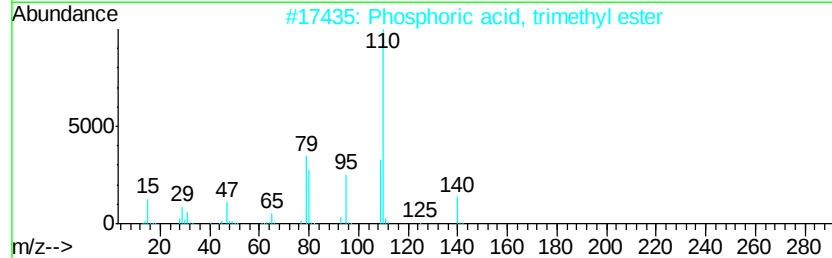
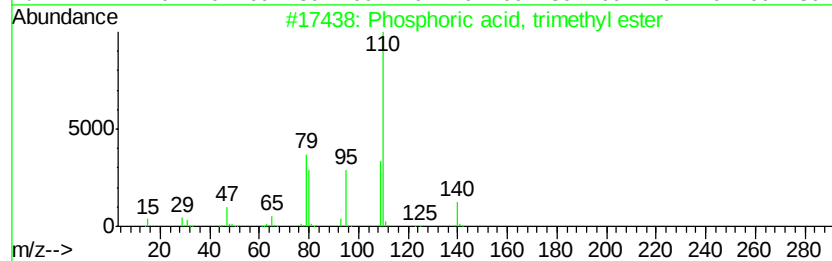
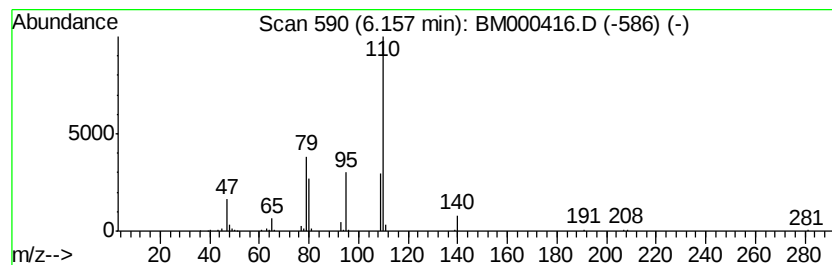
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phosphoric acid, trimethyl ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.16	10.36 ng/ul	621874	1,4-Dichlorobenzene-d4	7.75

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phosphoric acid, trimethyl ester	140	C3H9O4P	000512-56-1	95
2		Phosphoric acid, trimethyl ester	140	C3H9O4P	000512-56-1	91
3		Phosphoric acid, trimethyl ester	140	C3H9O4P	000512-56-1	91
4		Phosphoric acid, trimethyl ester	140	C3H9O4P	000512-56-1	90
5		Phosphonic acid, ethyl-, dimethy...	138	C4H11O3P	006163-75-3	64



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

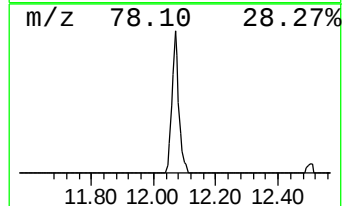
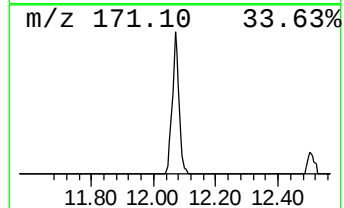
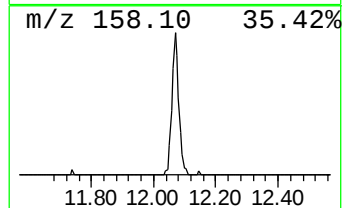
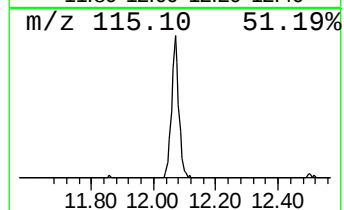
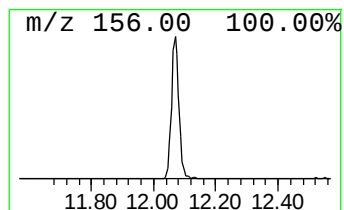
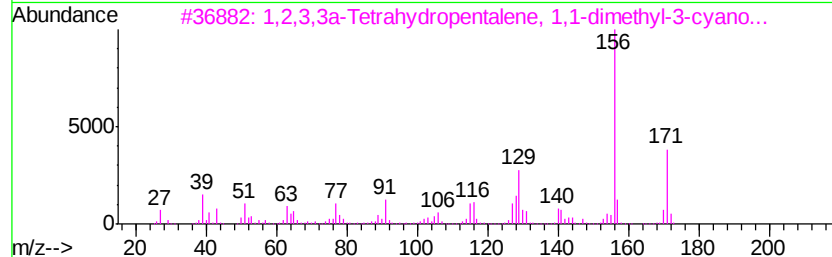
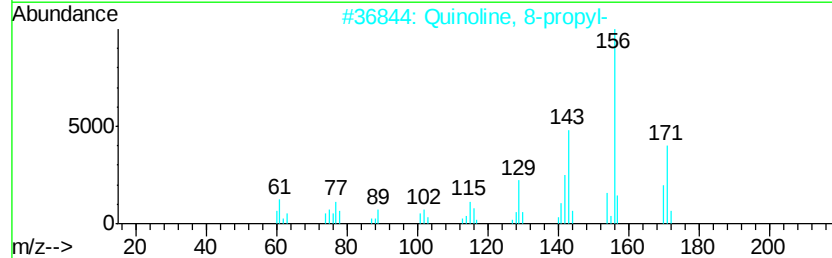
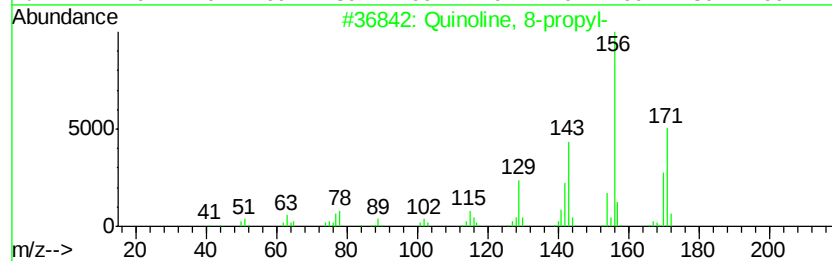
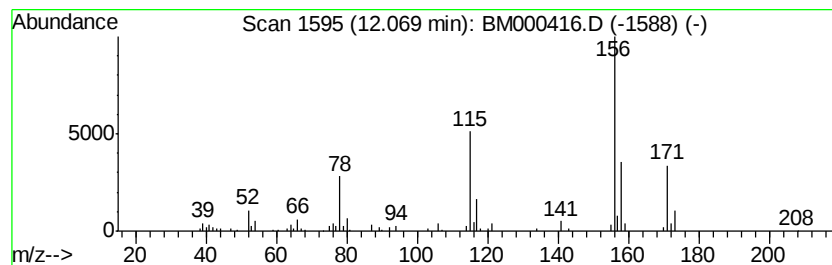
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	3.08 ng/ul	259816	Naphthalene-d8	10.55

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Quinoline, 8-propyl-	171	C12H13N	007661-53-2	18
2		Quinoline, 8-propyl-	171	C12H13N	007661-53-2	17
3		1,2,3,3a-Tetrahydropentalene, 1,...	171	C12H13N	1000217-17-0	14
4		3H-1,2,4-Triazol-3-one, 2,4-diet...	171	C7H13N3O2	039636-06-1	12
5		9-Thiabicyclo[3.3.1]nonan-3-one	156	C8H12OS	067194-70-1	12



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

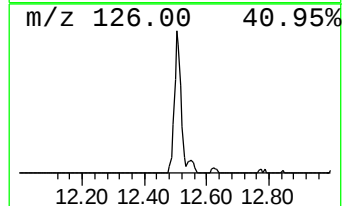
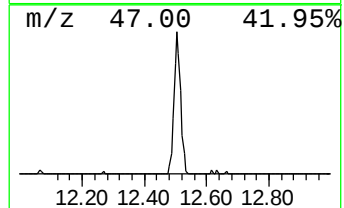
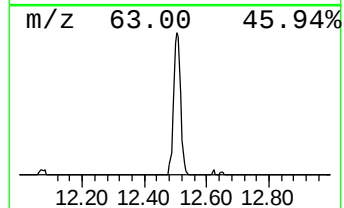
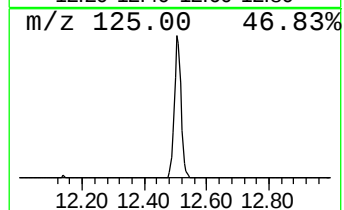
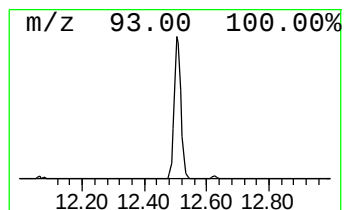
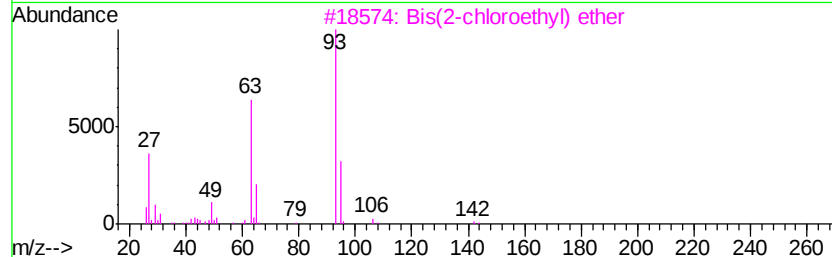
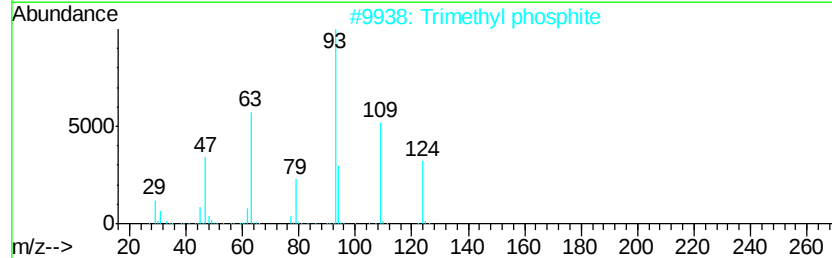
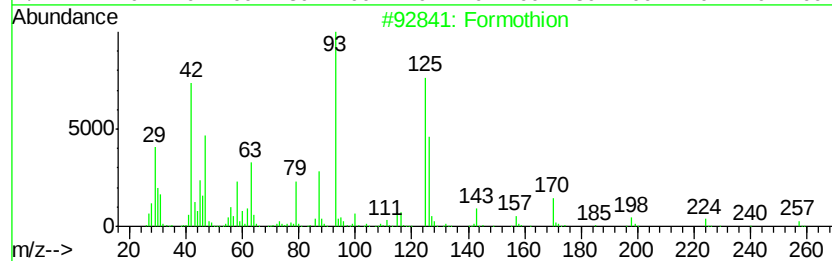
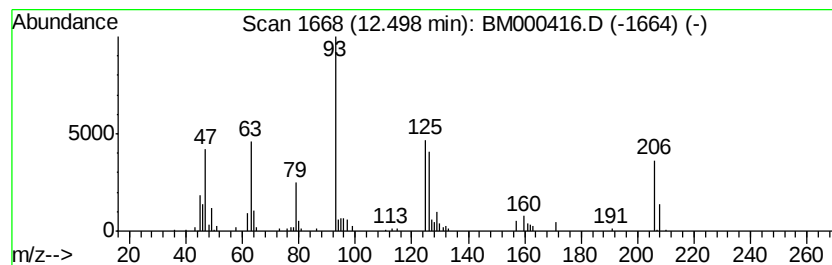
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Formothion Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.50	2.95 ng/ul	348392	Acenaphthene-d10	14.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Formothion	257	C6H12N04PS2	002540-82-1	59
2		Trimethyl phosphite	124	C3H9O3P	000121-45-9	33
3		Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	9
4		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	9
5		Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

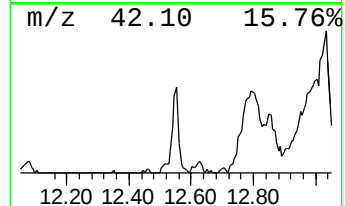
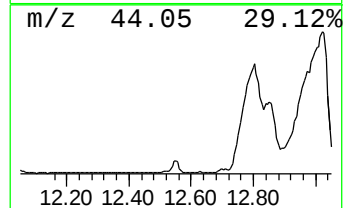
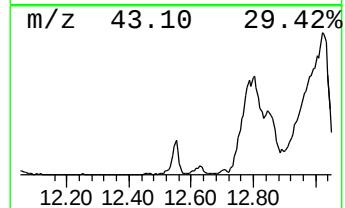
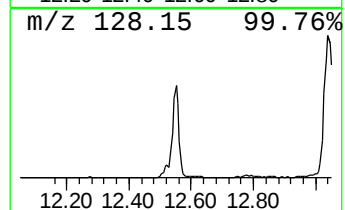
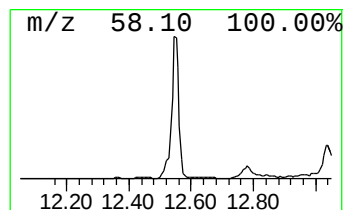
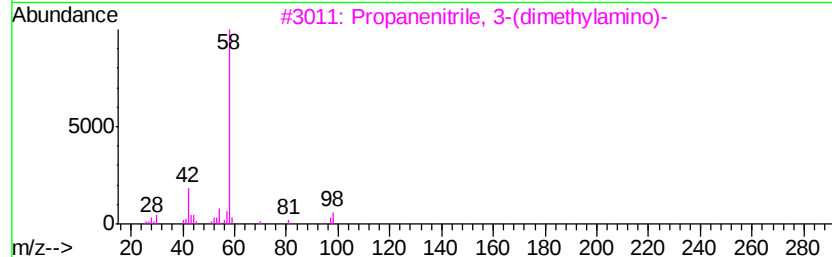
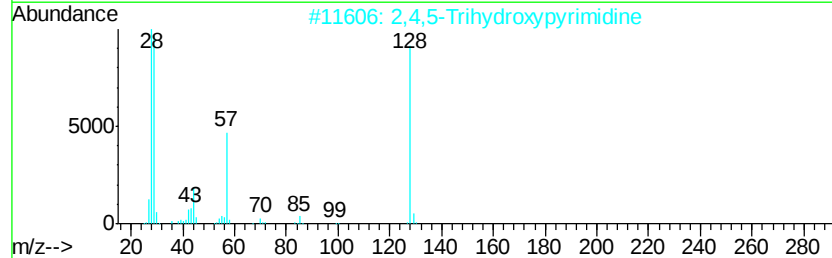
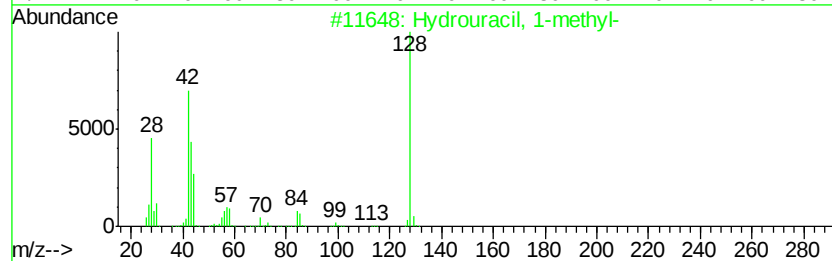
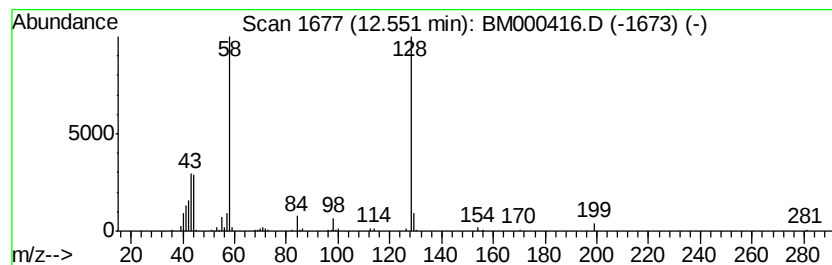
Instrument :
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LabSampleId :
SSTDCCC020

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown-02 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.55	2.80 ng/ul	330039	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Hydrouracil, 1-methyl-	128 C5H8N2O2	000696-11-7 43
2		2,4,5-Trihydroxypyrimidine	128 C4H4N2O3	000496-76-4 38
3		Propanenitrile, 3-(dimethylamino)-	98 C5H10N2	001738-25-6 38
4		d-Ribitol, 1-deoxy-1-heptylamino-	249 C12H27NO4	1000158-66-6 32
5		2-Pyrimidinone, 1,2,5,6-tetrahyd...	128 C5H8N2O2	1000261-72-3 32



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

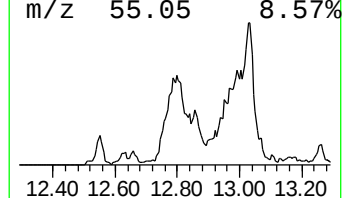
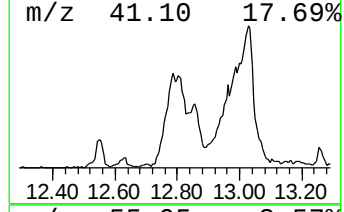
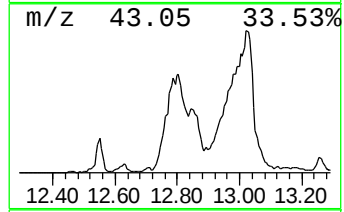
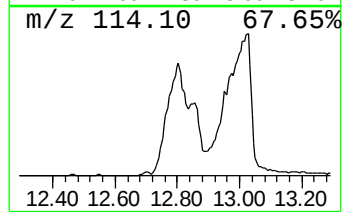
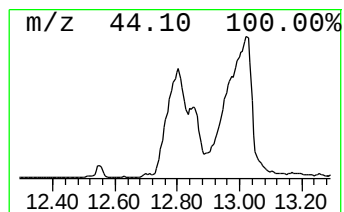
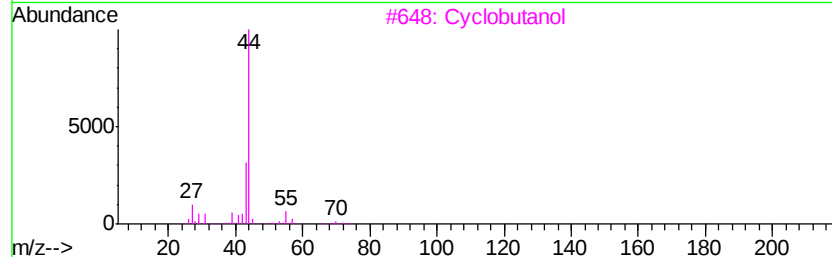
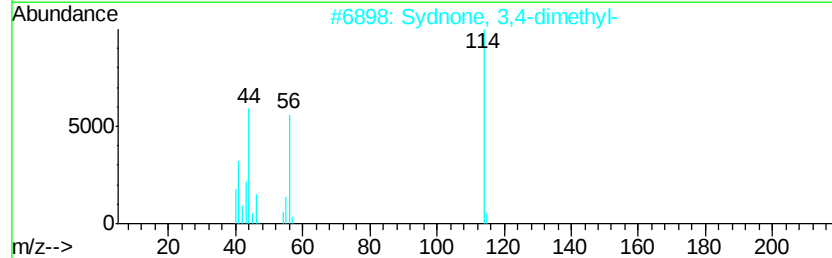
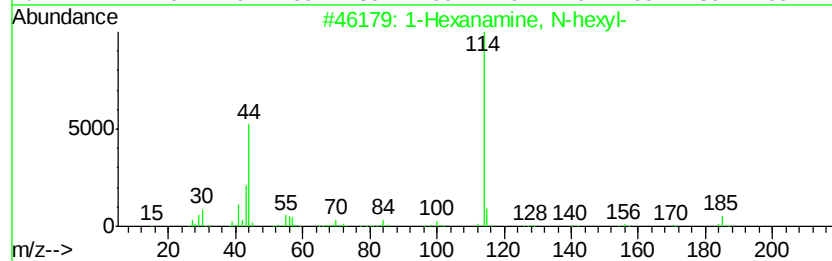
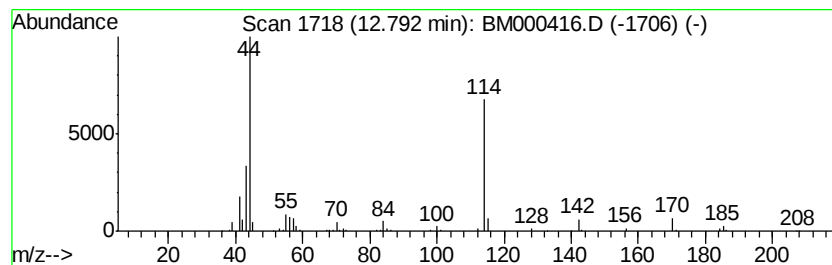
Instrument :
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LabSampleId :
SSTDCCC020

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 1-Hexanamine, N-hexyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	9.22 ng/ul	1087990	Acenaphthene-d10	14.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1-Hexanamine, N-hexyl-	185 C12H27N	000143-16-8 64
2		Sydnone, 3,4-dimethyl-	114 C4H6N2O2	004007-18-5 45
3		Cyclobutanol	72 C4H8O	002919-23-5 43
4		Butanoic acid, 2-hexylimino-	185 C10H19NO2	1000258-67-1 32
5		1-Hexanamine, N-hexyl-	185 C12H27N	000143-16-8 27



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

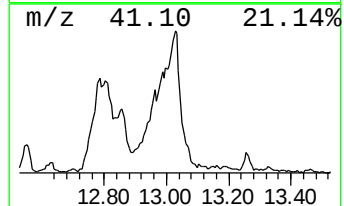
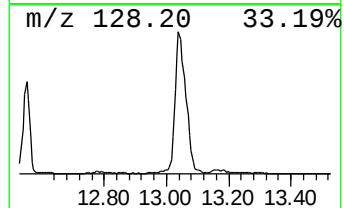
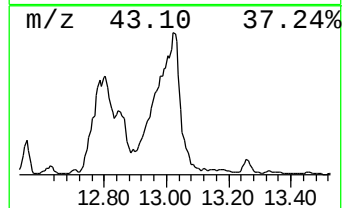
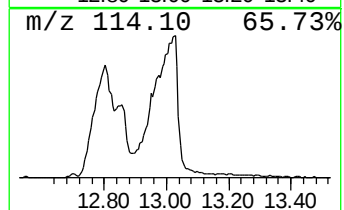
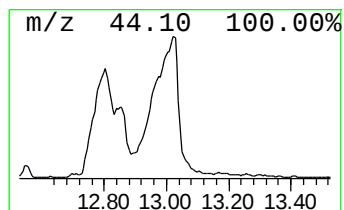
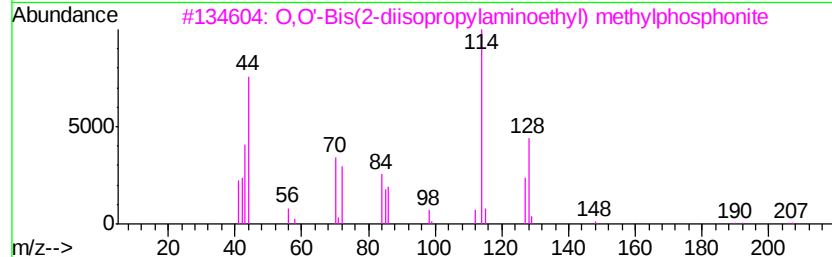
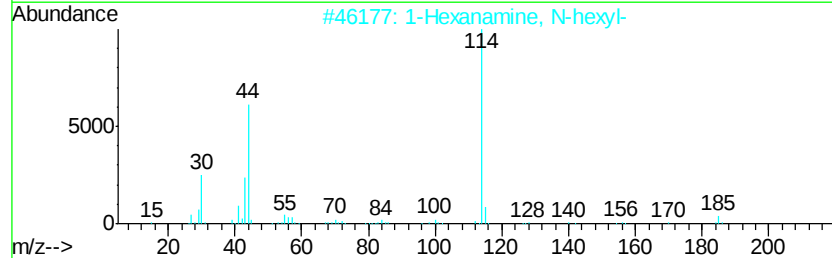
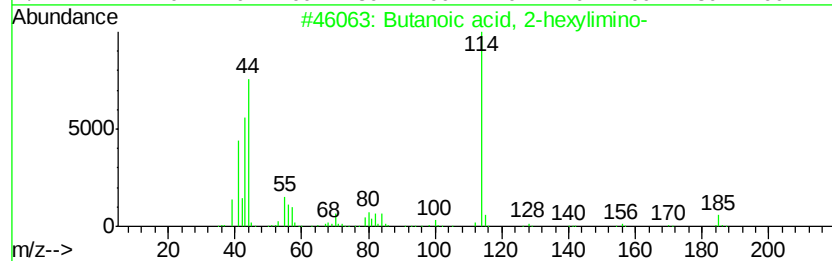
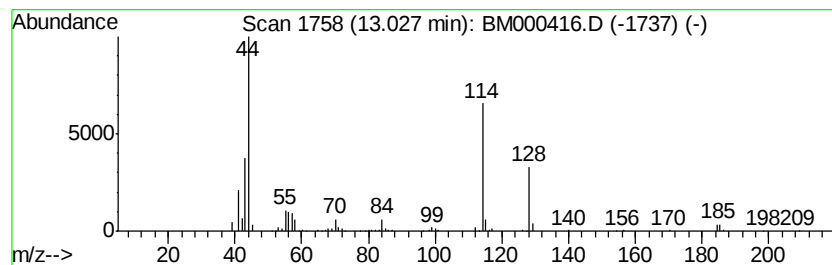
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	33.29 ng/ul	3927540	Acenaphthene-d10	14.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 2-hexylimino-	185	C10H19NO2	1000258-67-1	43
2		1-Hexanamine, N-hexyl-	185	C12H27N	000143-16-8	43
3		O,O'-Bis(2-diisopropylaminoethyl)...	334	C17H39N2O2P	057856-12-9	38
4		Cyclobutanol	72	C4H8O	002919-23-5	35
5		1-Octadecanamine, N-methyl-	283	C19H41N	002439-55-6	32



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

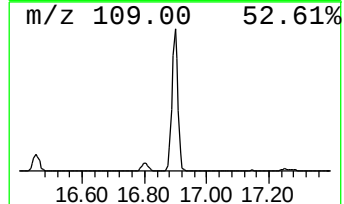
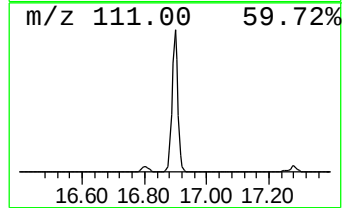
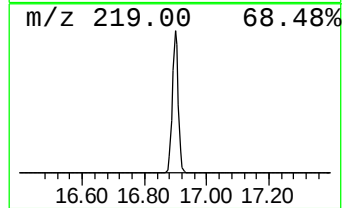
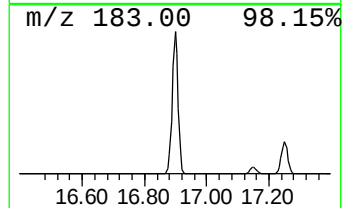
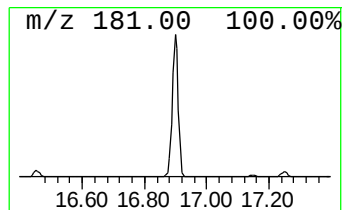
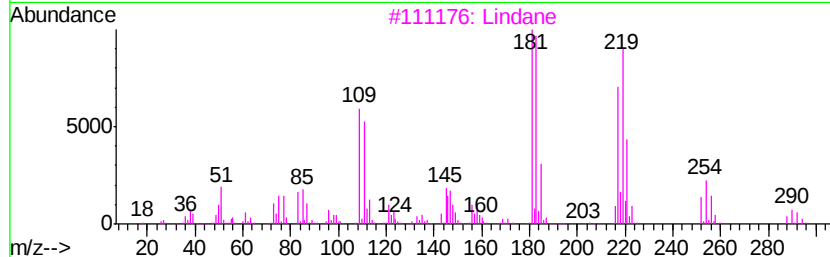
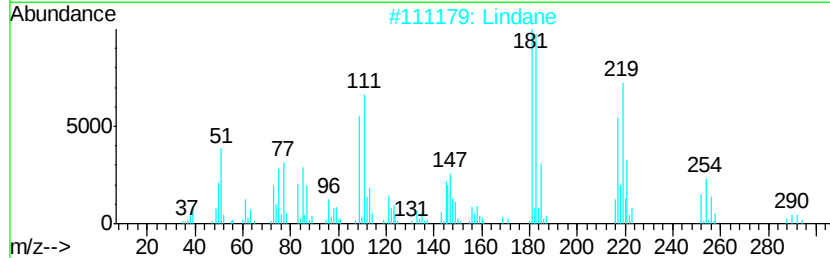
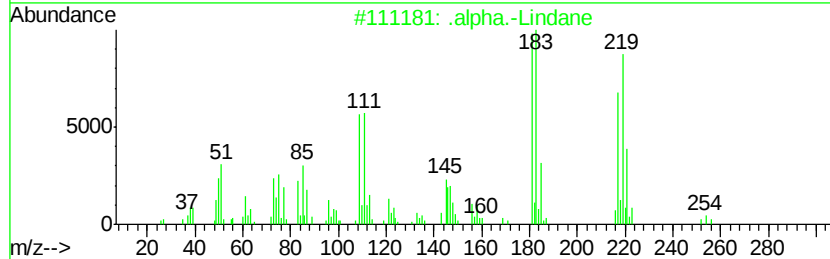
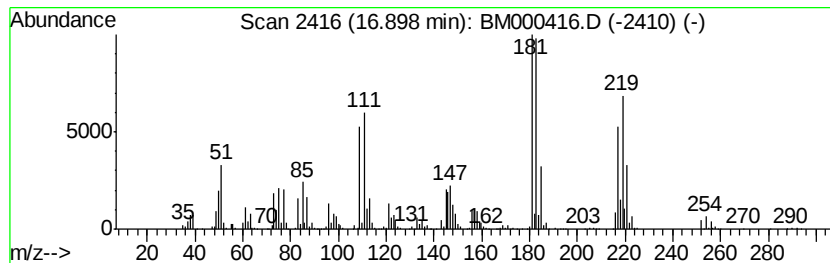
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 .alpha.-Lindane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.90	28.49 ng/ul	4391740	Phenanthrene-d10	17.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.alpha.-Lindane	288	C6H6Cl6	000319-84-6	95
2		Lindane	288	C6H6Cl6	000058-89-9	94
3		Lindane	288	C6H6Cl6	000058-89-9	91
4		Cyclohexane, 1,2,3,4,5,6-hexachl...	288	C6H6Cl6	000319-85-7	91
5		.alpha.-Lindane	288	C6H6Cl6	000319-84-6	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
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Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

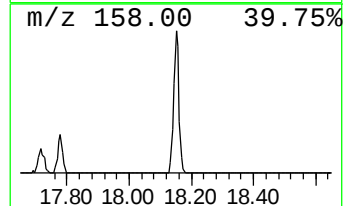
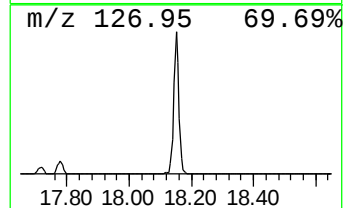
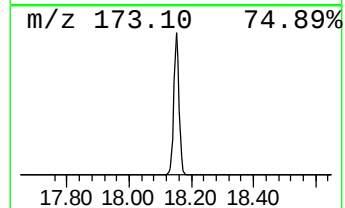
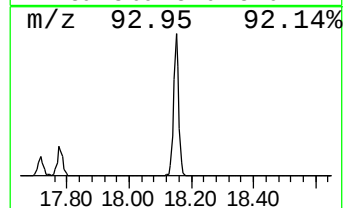
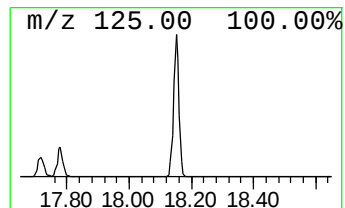
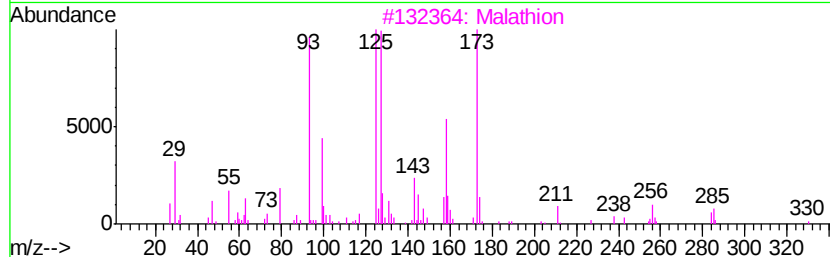
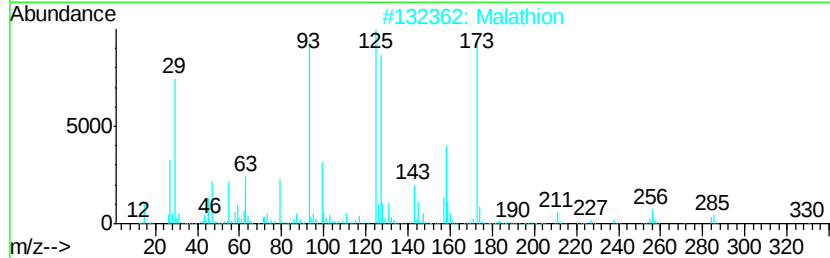
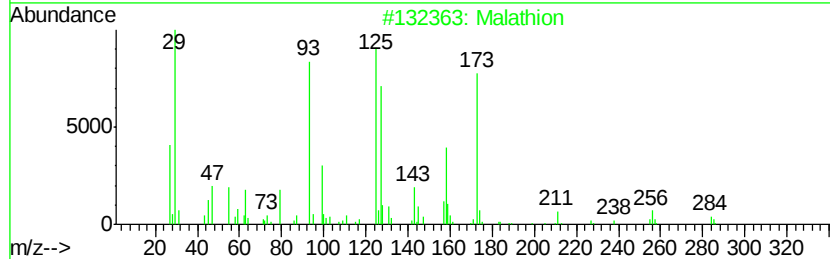
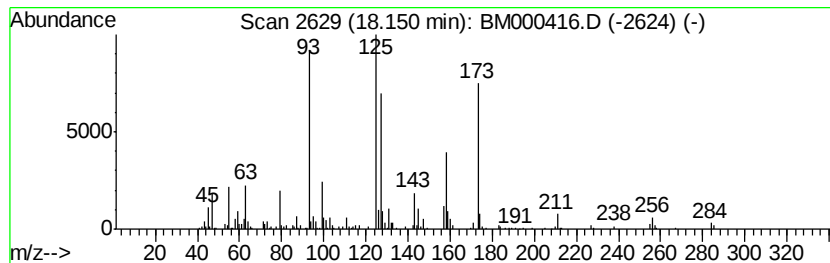
Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Malathion Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.15	3.26 ng/ul	502214	Phenanthrene-d10	17.14
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Malathion		330 C10H19O6PS2	000121-75-5 95
2	Malathion		330 C10H19O6PS2	000121-75-5 94
3	Malathion		330 C10H19O6PS2	000121-75-5 53
4	Malathion		330 C10H19O6PS2	000121-75-5 47
5	Silane, dichloroethenylmethyl-		140 C3H6Cl2Si	000124-70-9 22



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020315\
Data File : BM000416.D
Acq On : 03 Feb 2015 14:29
Operator : TP/IZ
Sample : G1147-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC020

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM020315.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Methanethioamide,...	2.75	1515.5	ng/ul	90982500	1	7.75	1200660	20.0
Butane, 2-methoxy...	2.96	16.2	ng/ul	975093	1	7.75	1200660	20.0
Propanol, methoxy...	4.22	2.2	ng/ul	133742	1	7.75	1200660	20.0
Phosphoric acid, ...	6.16	10.4	ng/ul	621874	1	7.75	1200660	20.0
unknown-01	12.07	3.1	ng/ul	259816	2	10.55	1688910	20.0
Formothion	12.50	3.0	ng/ul	348392	3	14.40	2359850	20.0
unknown-02	12.55	2.8	ng/ul	330039	3	14.40	2359850	20.0
1-Hexanamine, N-h...	12.79	9.2	ng/ul	1087990	3	14.40	2359850	20.0
unknown-03	13.03	33.3	ng/ul	3927540	3	14.40	2359850	20.0
.alpha.-Lindane	16.90	28.5	ng/ul	4391740	4	17.14	3082750	20.0
Malathion	18.15	3.3	ng/ul	502214	4	17.14	3082750	20.0