

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
 Data File : BG024784.D
 Acq On : 15 Nov 2016 4:30
 Operator : UM/SJ
 Sample : H5626-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 F2J03

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.1
 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: 1

Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.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	3.622	129	133	138	rVV7	9247	14697	0.30%	0.028%
2	4.386	257	263	271	rVB6	21180	44573	0.90%	0.084%
3	5.320	418	422	432	rVB5	17344	38840	0.78%	0.073%
4	5.737	487	493	498	rVB2	25718	36446	0.74%	0.068%
5	5.867	509	515	523	rBV4	24750	60029	1.21%	0.113%
6	5.978	530	534	542	rVB6	7473	20750	0.42%	0.039%
7	6.243	570	579	586	rVV5	23130	48948	0.99%	0.092%
8	6.313	586	591	595	rVB5	6834	15955	0.32%	0.030%
9	7.330	760	764	768	rVB5	13800	26264	0.53%	0.049%
10	7.400	768	776	782	rBV2	122463	245936	4.97%	0.462%
11	7.465	782	787	793	rVB6	10177	17076	0.35%	0.032%
12	7.576	797	806	814	rBV	364350	624667	12.62%	1.173%
13	7.782	834	841	849	rVB	397177	716200	14.47%	1.345%
14	7.888	853	859	867	rVB5	24084	50157	1.01%	0.094%
15	7.976	869	874	882	rBV2	50050	91545	1.85%	0.172%
16	8.258	915	922	935	rVV	376072	704953	14.25%	1.323%
17	8.375	935	942	946	rVB5	17225	36978	0.75%	0.069%
18	8.634	978	986	995	rVB	233352	408655	8.26%	0.767%
19	8.804	1005	1015	1023	rVB	202722	380029	7.68%	0.713%
20	8.881	1023	1028	1033	rBV8	8372	18917	0.38%	0.036%
21	8.951	1033	1040	1053	rVB	345076	630157	12.74%	1.183%
22	9.116	1059	1068	1074	rVB9	19032	49209	0.99%	0.092%
23	9.362	1101	1110	1115	rBV6	10994	26488	0.54%	0.050%
24	9.433	1115	1122	1137	rVV	506953	950531	19.21%	1.784%
25	9.621	1147	1154	1161	rBV4	31083	59813	1.21%	0.112%
26	9.744	1166	1175	1182	rBV4	73960	176649	3.57%	0.332%
27	9.809	1182	1186	1194	rVB5	24796	43598	0.88%	0.082%
28	10.162	1238	1246	1258	rBV2	443875	840385	16.98%	1.578%
29	10.338	1269	1276	1278	rBV6	7576	15318	0.31%	0.029%
30	10.467	1291	1298	1309	rBV8	16107	45491	0.92%	0.085%
31	10.561	1311	1314	1322	rBV7	9923	18694	0.38%	0.035%
32	10.696	1326	1337	1346	rBV	628060	1133732	22.91%	2.128%
33	10.831	1354	1360	1365	rBV7	6058	18102	0.37%	0.034%
34	11.090	1395	1404	1407	rBV	539533	960910	19.42%	1.804%

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35	11.143	1407	1413	1421	rVV	2591438	4869093	98.41%	9.141%
36	11.260	1421	1433	1441	rVB2	260652	527246	10.66%	0.990%
37	11.848	1527	1533	1539	rVB7	9332	21830	0.44%	0.041%
38	12.189	1586	1591	1600	rVB7	15293	33209	0.67%	0.062%
39	12.617	1660	1664	1666	rBV5	12357	15700	0.32%	0.029%
40	12.653	1666	1670	1674	rVB4	16229	26578	0.54%	0.050%
41	12.723	1674	1682	1694	rVB	146331	239772	4.85%	0.450%
42	12.841	1695	1702	1709	rVV5	17289	29098	0.59%	0.055%
43	12.935	1709	1718	1727	rVB2	115924	229920	4.65%	0.432%
44	13.710	1845	1850	1860	rVB5	36592	90148	1.82%	0.169%
45	13.804	1862	1866	1873	rVB9	9820	18286	0.37%	0.034%
46	14.039	1896	1906	1907	rBV7	16151	37842	0.76%	0.071%
47	14.268	1938	1945	1952	rBV	1442447	2131058	43.07%	4.001%
48	14.374	1957	1963	1974	rBV	376968	562128	11.36%	1.055%
49	14.503	1980	1985	1990	rVB6	24565	39854	0.81%	0.075%
50	14.574	1990	1997	2007	rBV	1383802	2267746	45.83%	4.257%
51	14.733	2020	2024	2028	rBV6	13667	17931	0.36%	0.034%
52	14.880	2033	2049	2055	rBV	1038963	1848164	37.35%	3.470%
53	14.938	2055	2059	2065	rVV	82020	134628	2.72%	0.253%
54	15.009	2065	2071	2075	rVV	93123	155550	3.14%	0.292%
55	15.056	2075	2079	2091	rVV2	188619	367213	7.42%	0.689%
56	15.138	2091	2093	2104	rVB9	12775	26875	0.54%	0.050%
57	15.273	2111	2116	2126	rVB2	88314	160234	3.24%	0.301%
58	15.373	2126	2133	2141	rBV9	20656	57890	1.17%	0.109%
59	15.508	2151	2156	2161	rBV6	39079	58638	1.19%	0.110%
60	15.590	2164	2170	2178	rVB	139671	211120	4.27%	0.396%
61	15.667	2178	2183	2190	rVB3	28514	42935	0.87%	0.081%
62	15.796	2195	2205	2209	rBV7	78859	150860	3.05%	0.283%
63	15.867	2209	2217	2223	rVV	1945879	3168230	64.03%	5.948%
64	15.925	2223	2227	2229	rVV3	46394	69659	1.41%	0.131%
65	15.972	2229	2235	2246	rVB	865590	1338637	27.05%	2.513%
66	16.102	2249	2257	2264	rVB9	32614	60954	1.23%	0.114%
67	16.219	2275	2277	2283	rVB5	18995	25100	0.51%	0.047%
68	16.313	2284	2293	2298	rBV8	31010	57499	1.16%	0.108%
69	16.366	2300	2302	2306	rVB5	19359	25961	0.52%	0.049%
70	16.425	2306	2312	2320	rBV5	28039	79344	1.60%	0.149%
71	16.589	2336	2340	2348	rVB9	18627	40396	0.82%	0.076%

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72	16.713	2357	2361	2367	rVB9	16277	27161	0.55%	0.051%
73	16.789	2370	2374	2384	rVB9	11645	33398	0.67%	0.063%
74	16.901	2390	2393	2397	rVB4	9020	14127	0.29%	0.027%
75	17.053	2414	2419	2423	rBV6	9225	17023	0.34%	0.032%
76	17.124	2423	2431	2438	rVV5	33971	88277	1.78%	0.166%
77	17.189	2438	2442	2448	rBV2	46577	63794	1.29%	0.120%
78	17.265	2450	2455	2464	rVB4	40133	78101	1.58%	0.147%
79	17.406	2473	2479	2483	rBV7	22266	36110	0.73%	0.068%
80	17.553	2499	2504	2508	rVB6	14741	27905	0.56%	0.052%
81	17.617	2508	2515	2526	rBV	1350896	2213161	44.73%	4.155%
82	17.717	2526	2532	2544	rVB2	2406063	3886495	78.55%	7.296%
83	17.870	2553	2558	2563	rVB	127520	176800	3.57%	0.332%
84	18.023	2576	2584	2588	rVV	361752	552882	11.17%	1.038%
85	18.082	2588	2594	2613	rVB	626899	1006857	20.35%	1.890%
86	18.205	2613	2615	2618	rBV4	14849	16549	0.33%	0.031%
87	18.258	2619	2624	2632	rVB6	119318	179893	3.64%	0.338%
88	18.546	2668	2673	2679	rVB	415834	547021	11.06%	1.027%
89	18.687	2692	2697	2705	rBV2	202587	301036	6.08%	0.565%
90	18.769	2709	2711	2718	rBV8	14502	18438	0.37%	0.035%
91	19.063	2757	2761	2766	rVB6	44423	63577	1.28%	0.119%
92	19.239	2787	2791	2800	rVB3	84884	130274	2.63%	0.245%
93	19.627	2852	2857	2864	rBV	37534	65673	1.33%	0.123%
94	19.880	2897	2900	2905	rVB2	30574	41826	0.85%	0.079%
95	19.991	2911	2919	2932	rBV	3353737	4910538	99.24%	9.219%
96	20.861	3064	3067	3073	rVB7	54097	64788	1.31%	0.122%
97	21.918	3239	3247	3256	rVB	1573178	2860183	57.81%	5.370%
98	25.109	3779	3790	3802	rVB3	1636402	4947949	100.00%	9.289%
99	25.344	3821	3830	3849	rVB2	970586	3029457	61.23%	5.687%
100	27.905	4263	4266	4268	rBV3	50667	59792	1.21%	0.112%

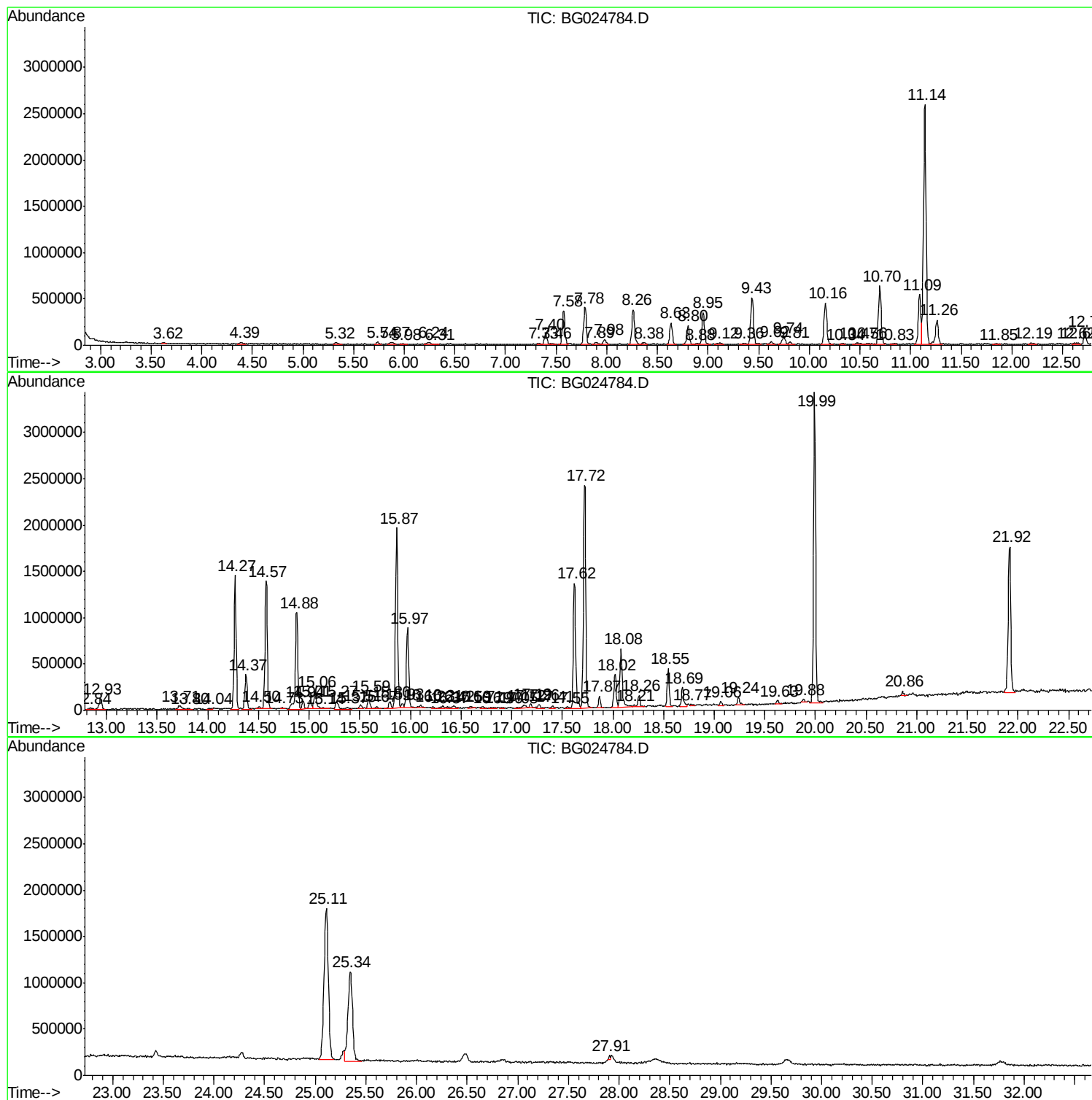
Sum of corrected areas: 53267103

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
Quant Title : SVOA CALIBRATION

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



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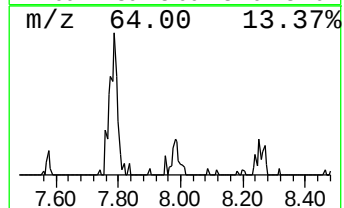
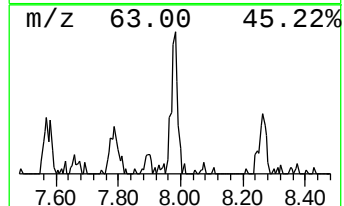
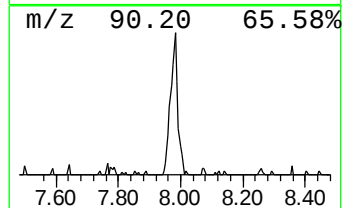
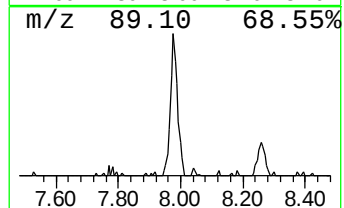
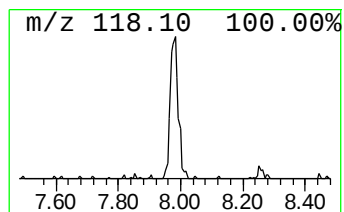
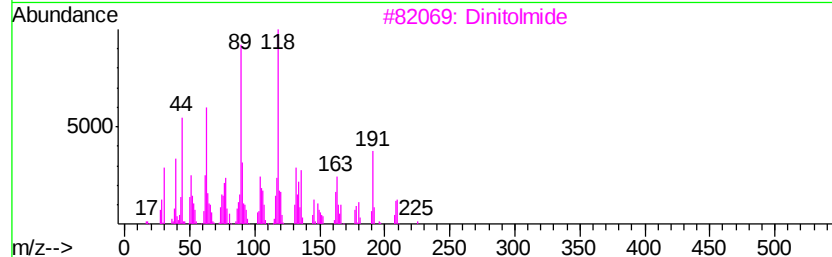
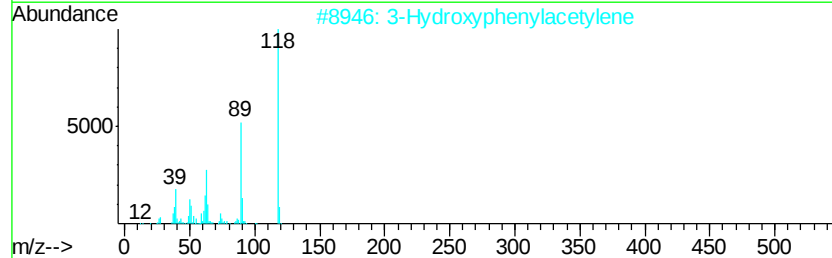
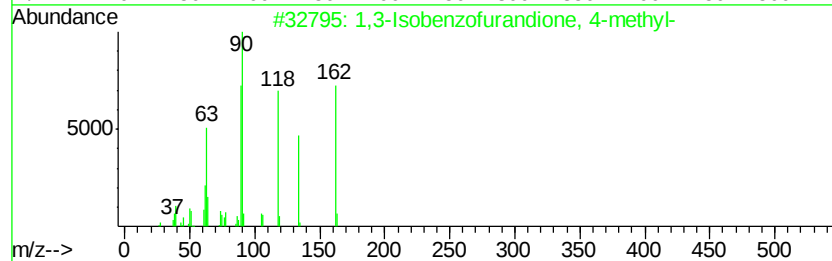
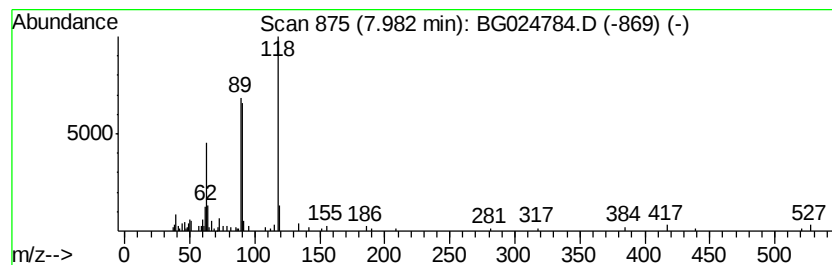
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1,3-Isobenzofurandione, 4-m... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.98	2.60 ng/ul	91545	1,4-Dichlorobenzene-d4	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Isobenzofurandione, 4-methyl-	162	C9H6O3	004792-30-7	86
2		3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	83
3		Dinitolmide	225	C8H7N3O5	000148-01-6	83
4		Benzofuran	118	C8H6O	000271-89-6	70
5		6-Nitro-o-tolunitrile	162	C8H6N2O2	001885-76-3	64



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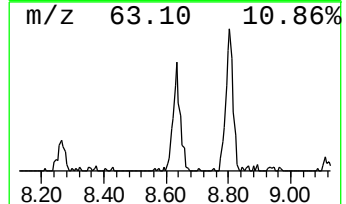
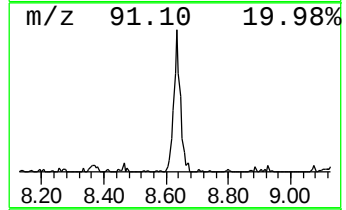
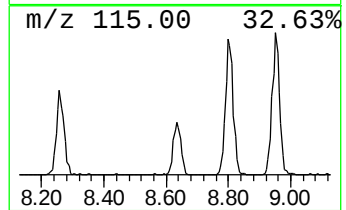
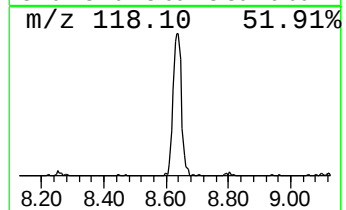
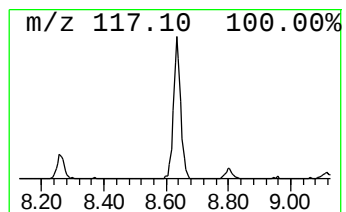
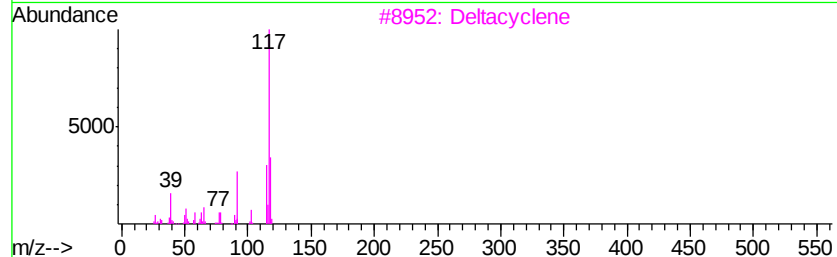
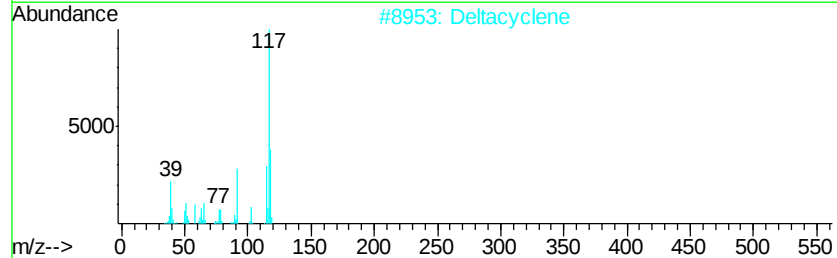
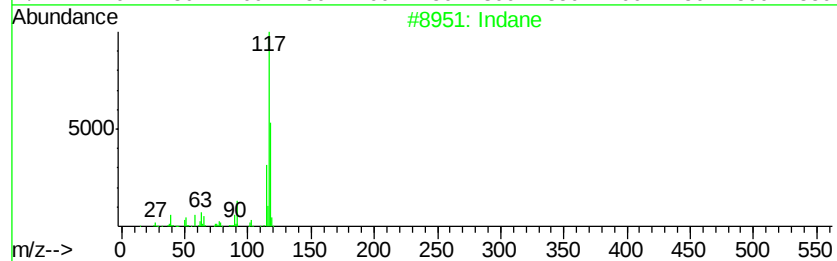
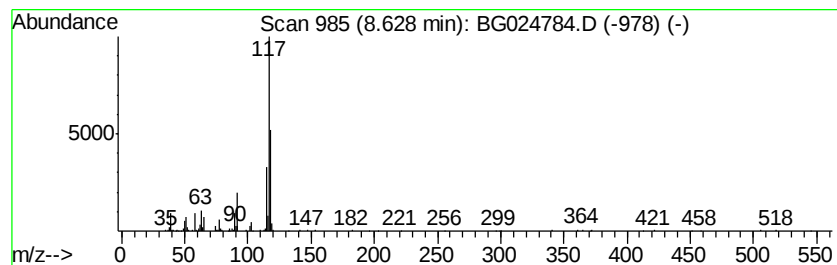
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.63	11.59 ng/ul	408655	1,4-Dichlorobenzene-d4	8.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	74
2	Deltacyclene		118	C9H10	007785-10-6	72
3	Deltacyclene		118	C9H10	007785-10-6	72
4	Benzene, cyclopropyl-		118	C9H10	000873-49-4	64
5	Indane		118	C9H10	000496-11-7	55



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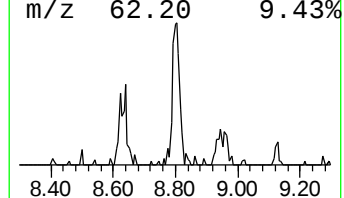
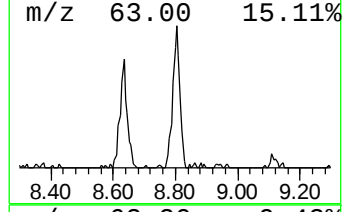
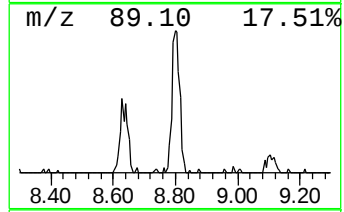
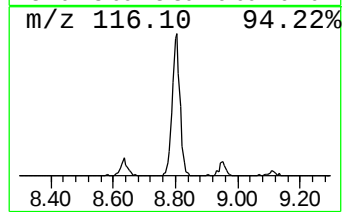
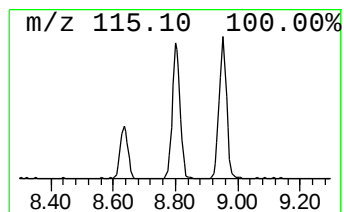
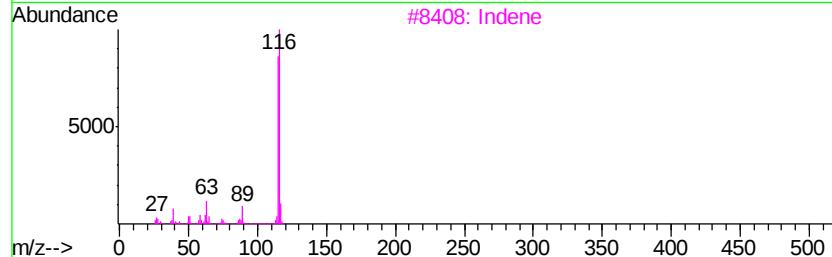
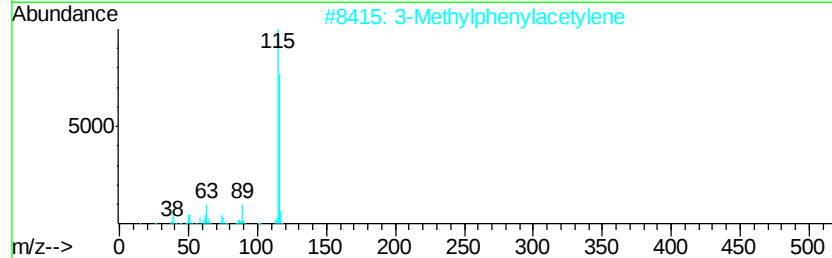
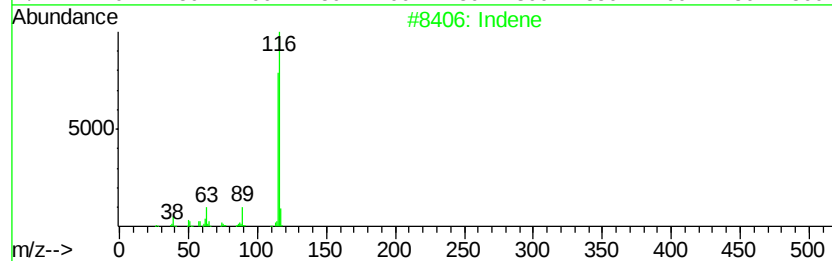
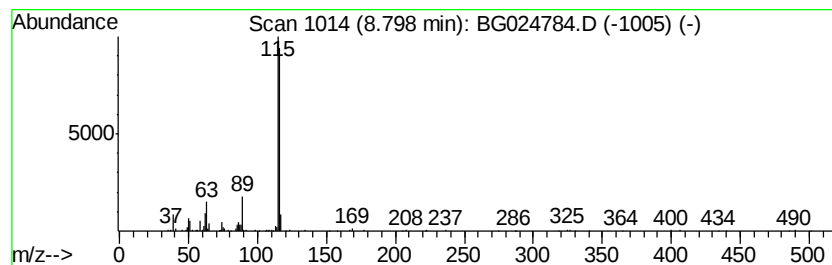
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Peak Number 3 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	10.78 ng/ul	380029	1,4-Dichlorobenzene-d4	8.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	96
2		3-Methylphenylacetylene	116	C9H8	000766-82-5	91
3		Indene	116	C9H8	000095-13-6	91
4		Indene	116	C9H8	000095-13-6	91
5		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
Acq On : 15 Nov 2016 4:30
Operator : UM/SJ
Sample : H5626-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
F2J03

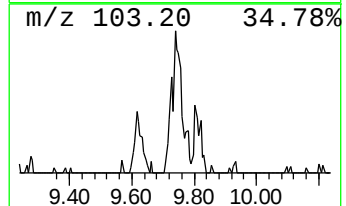
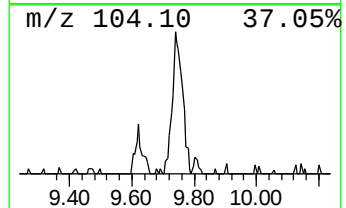
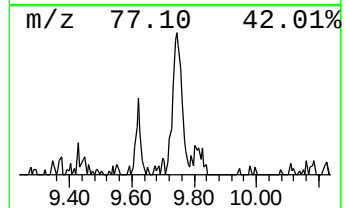
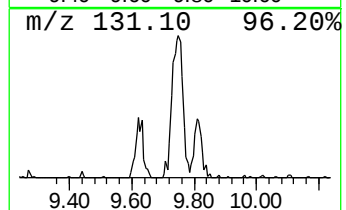
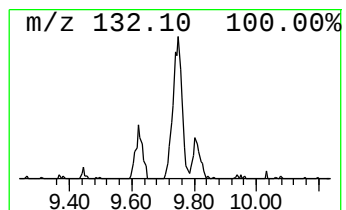
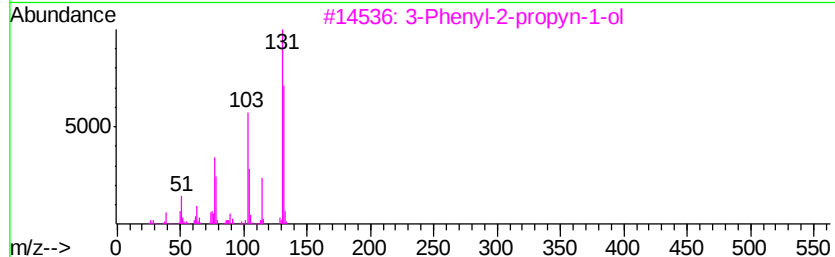
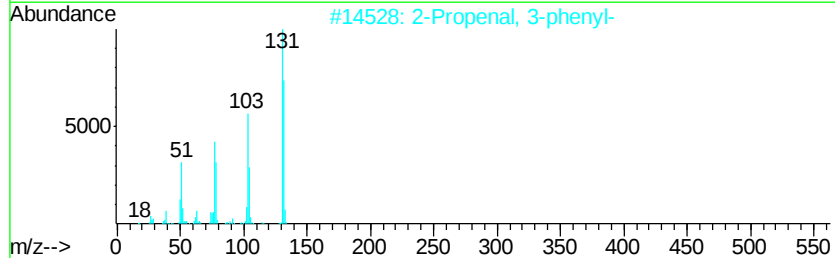
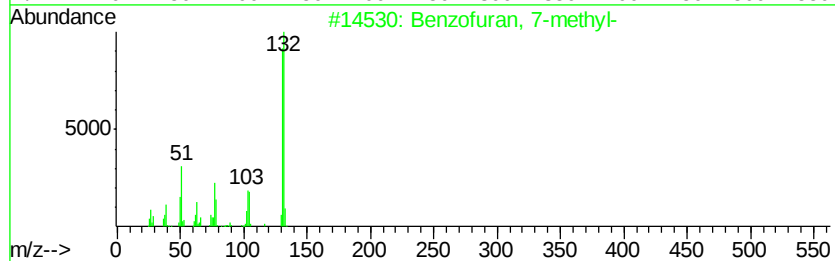
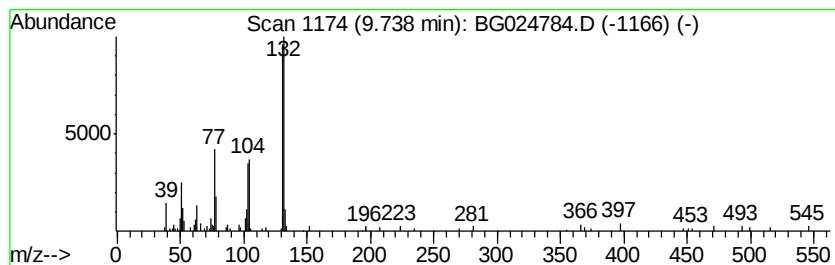
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzofuran, 7-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.74	3.68 ng/ul	176649	Naphthalene-d8	11.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	90
3		3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	87
4		2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	86
5		1H-Indazole, 3-methyl-	132	C8H8N2	1000316-00-2	81



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
Acq On : 15 Nov 2016 4:30
Operator : UM/SJ
Sample : H5626-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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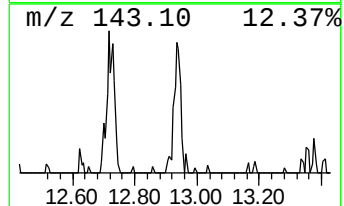
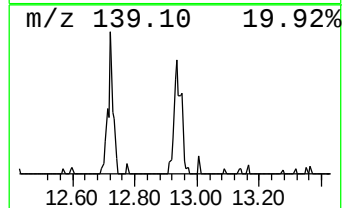
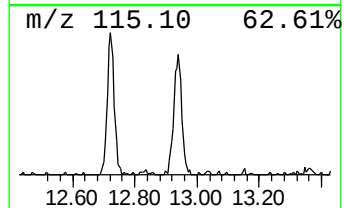
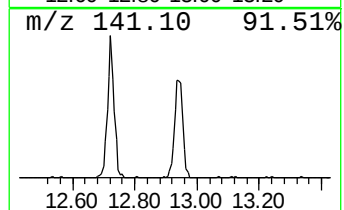
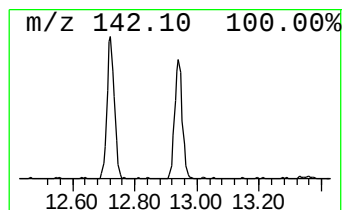
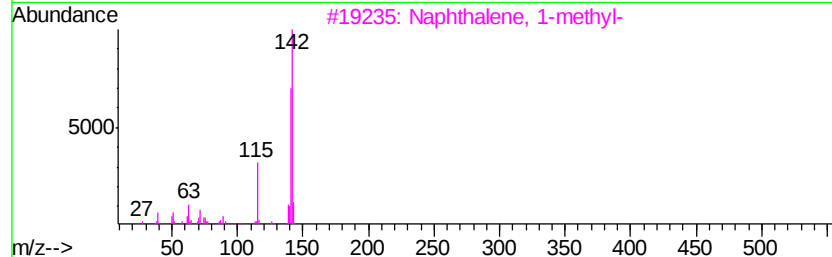
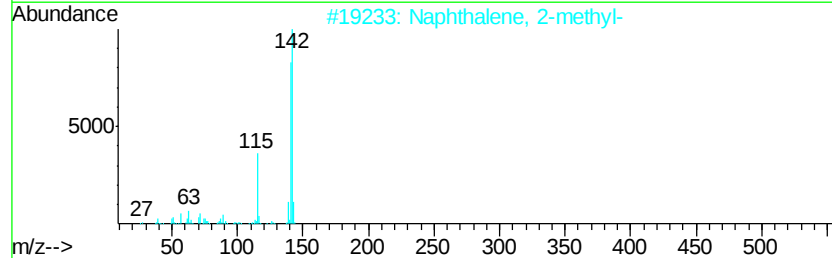
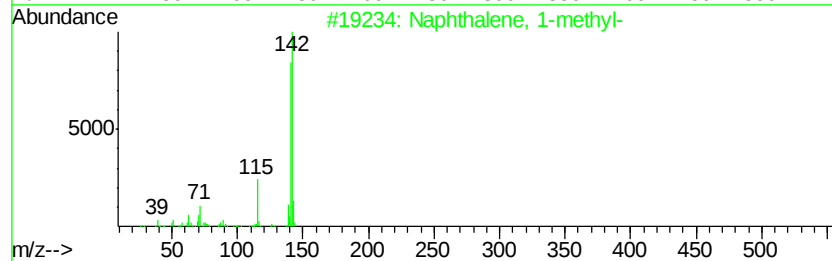
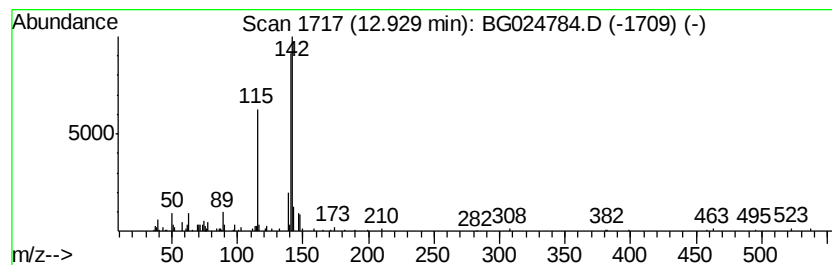
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	4.79 ng/ul	229920	Naphthalene-d8	11.09

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
4		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
Acq On : 15 Nov 2016 4:30
Operator : UM/SJ
Sample : H5626-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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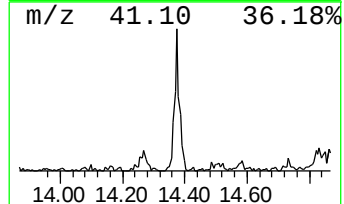
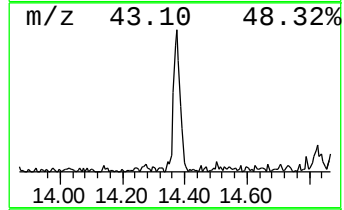
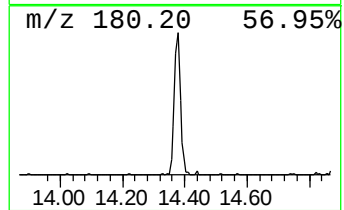
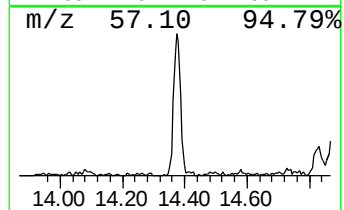
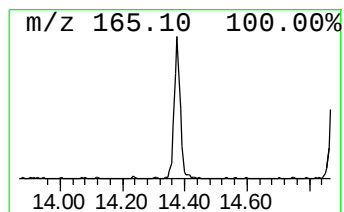
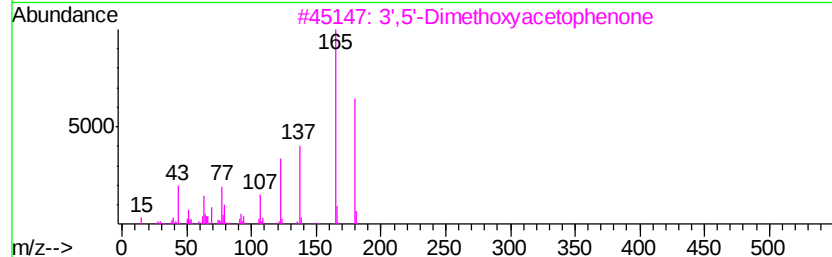
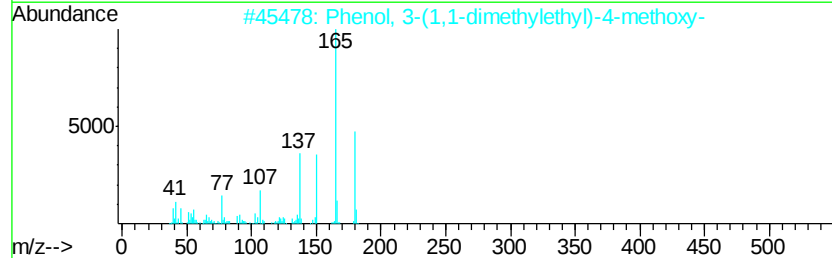
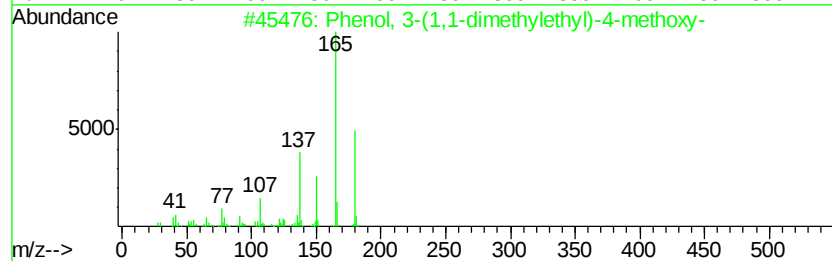
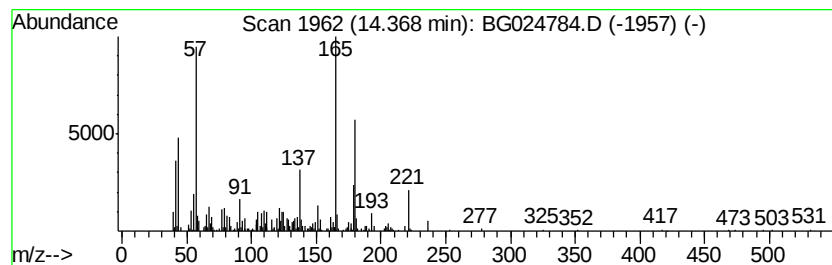
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 3-(1,1-dimethylethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.37	6.08 ng/ul	562128	Acenaphthene-d10	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 3-(1,1-dimethylethyl)-4-...	180	C11H16O2	000088-32-4	62
2		Phenol, 3-(1,1-dimethylethyl)-4-...	180	C11H16O2	000088-32-4	62
3		3',5'-Dimethoxyacetophenone	180	C10H12O3	039151-19-4	58
4		3-tert-Butyl-4-hydroxyanisole	180	C11H16O2	000121-00-6	58
5		2-Benzothiazolamine, 6-methoxy-	180	C8H8N2OS	001747-60-0	53



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
Acq On : 15 Nov 2016 4:30
Operator : UM/SJ
Sample : H5626-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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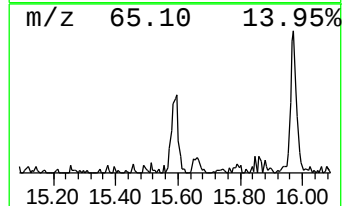
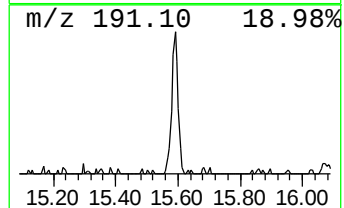
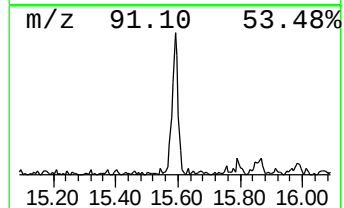
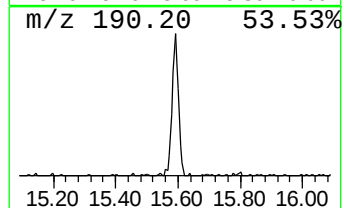
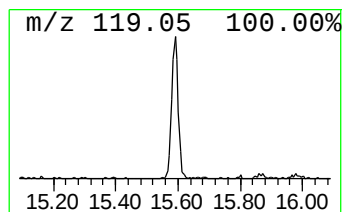
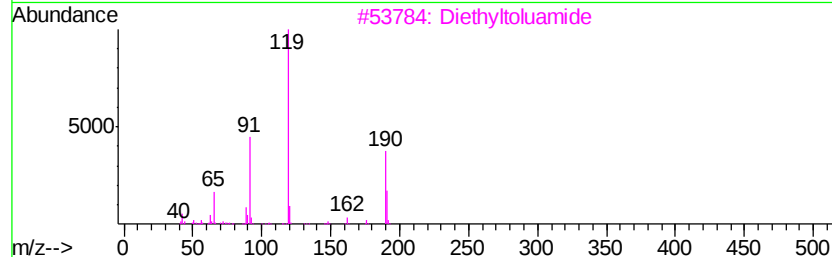
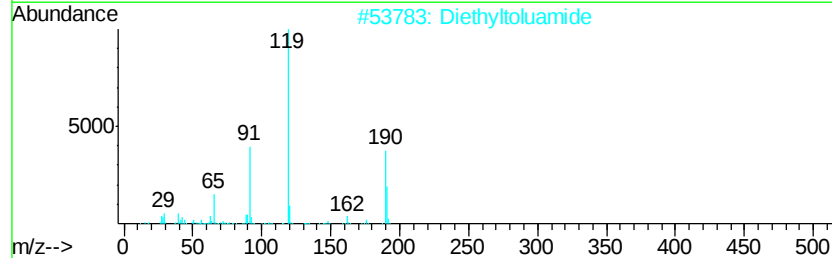
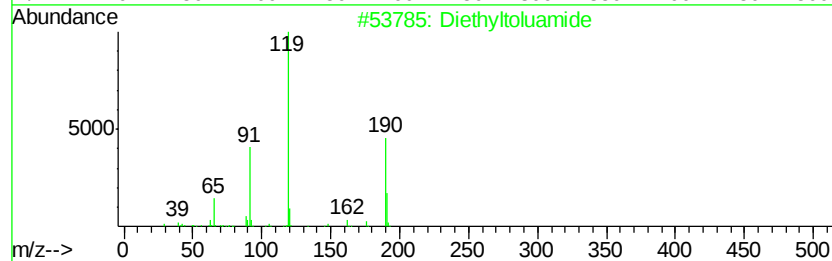
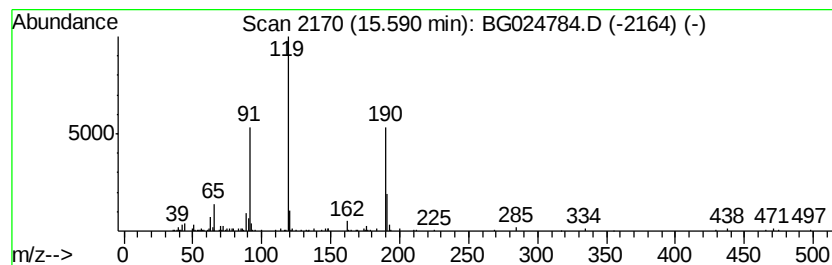
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Diethyltoluamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.59	2.28 ng/ul	211120	Acenaphthene-d10	14.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	95
2		Diethyltoluamide	191	C12H17NO	000134-62-3	94
3		Diethyltoluamide	191	C12H17NO	000134-62-3	93
4		Diethyltoluamide	191	C12H17NO	000134-62-3	76
5		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
Acq On : 15 Nov 2016 4:30
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Sample : H5626-04
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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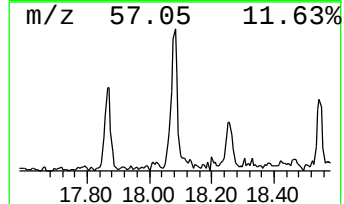
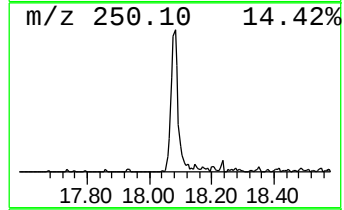
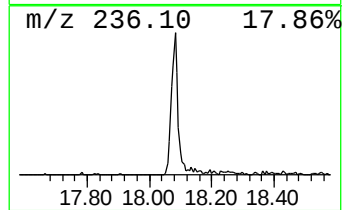
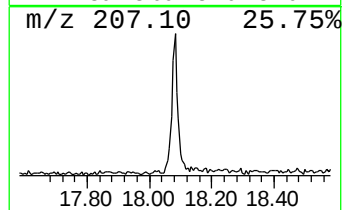
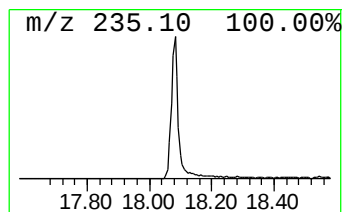
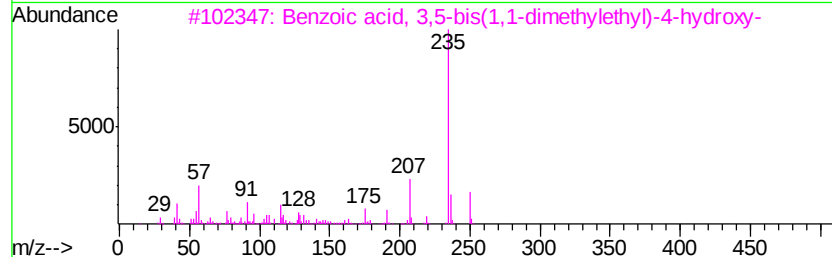
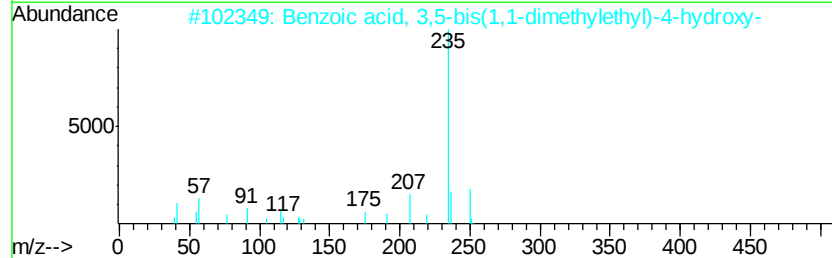
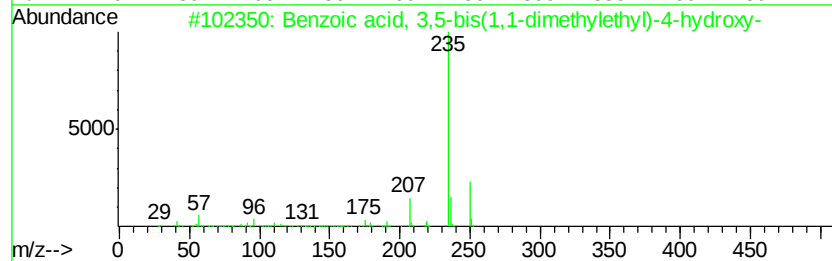
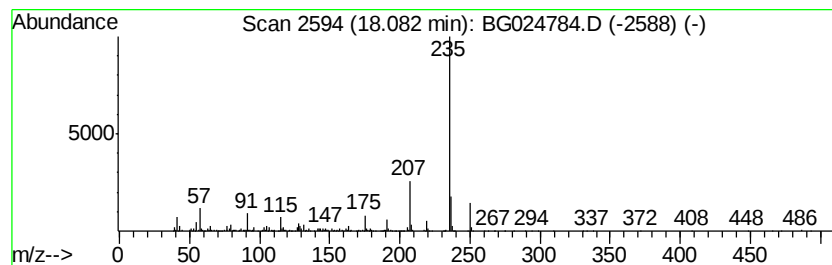
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.08	9.10 ng/ul	1006860	Phenanthrene-d10	17.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	94
2		Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	90
3		Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	90
4		-4'-Dimethylamino-2'-(trimethyls...	250	C13H22N2OSi	018410-40-7	64
5		N-(4-Hydroxyphenyl)-2-naphthylamine	235	C16H13NO	000093-45-8	59



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
Data File : BG024784.D
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ClientSampleId :
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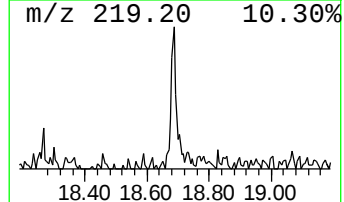
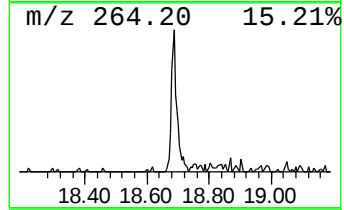
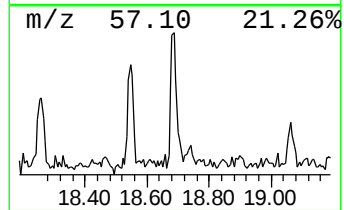
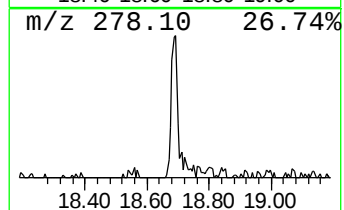
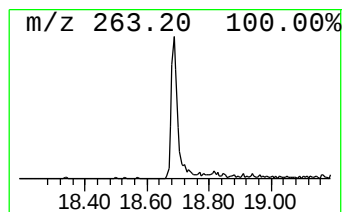
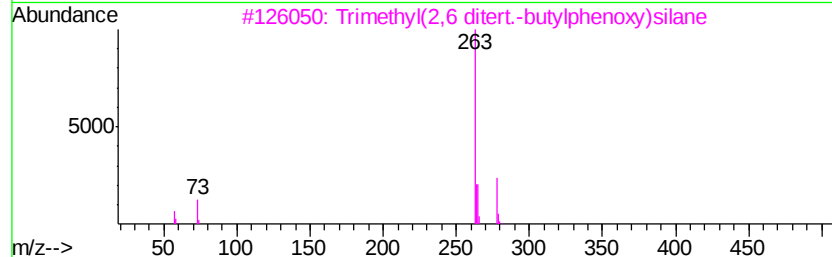
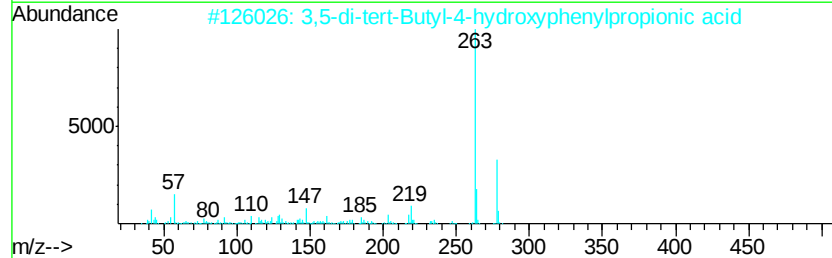
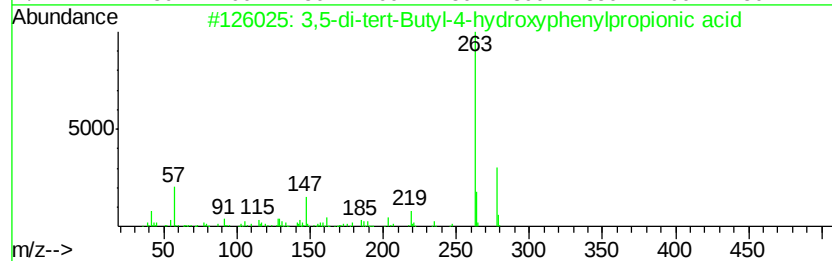
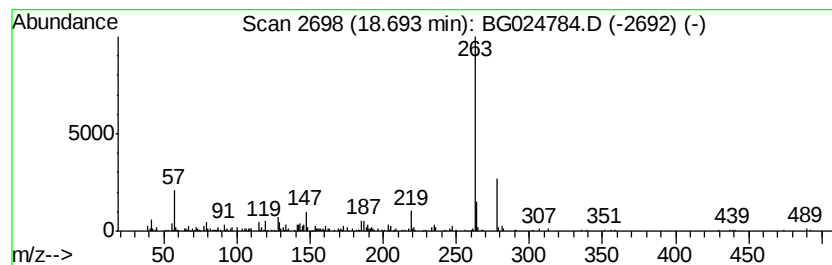
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.69	2.72 ng/ul	301036	Phenanthrene-d10	17.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	91
2		3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	87
3		Trimethyl(2,6 ditert.-butylpheno...	278	C17H30OSi	010416-73-6	72
4		Phosphenimidous amide, tris(trim...	278	C9H27N2PSi3	050732-21-3	64
5		Benzoic acid, 3,5-bis(1,1-dimeth...	278	C17H26O3	001620-64-0	45



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
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ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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-Isobenzofuran...	7.98	2.6	ng/ul	91545	1	8.26	704953	20.0
Indane	8.63	11.6	ng/ul	408655	1	8.26	704953	20.0
Indene	8.80	10.8	ng/ul	380029	1	8.26	704953	20.0
Benzofuran, 7-met...	9.74	3.7	ng/ul	176649	2	11.09	960910	20.0
Naphthalene, 1-me...	12.93	4.8	ng/ul	229920	2	11.09	960910	20.0
Phenol, 3-(1,1-di...	14.37	6.1	ng/ul	562128	3	14.88	1848160	20.0
Diethyltoluamide	15.59	2.3	ng/ul	211120	3	14.88	1848160	20.0
Benzoic acid, 3,5...	18.08	9.1	ng/ul	1006860	4	17.62	2213160	20.0
3,5-di-tert-Butyl...	18.69	2.7	ng/ul	301036	4	17.62	2213160	20.0