

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
 Data File : BG025077.D
 Acq On : 10 Dec 2016 20:34
 Operator : UM/SJ
 Sample : H5905-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DA805

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-BG113016.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	4.353	252	258	265	rBV9	9253	23563	0.62%	0.072%
2	4.700	309	317	319	rBV6	12274	25378	0.66%	0.077%
3	5.352	426	428	436	rBV5	8404	17285	0.45%	0.053%
4	7.379	764	773	788	rBV3	251242	565941	14.77%	1.723%
5	7.550	795	802	810	rBV2	184162	316297	8.26%	0.963%
6	7.761	827	838	847	rBV	242704	450905	11.77%	1.373%
7	7.873	851	857	862	rVB7	9175	22442	0.59%	0.068%
8	8.237	911	919	929	rVB2	241776	436064	11.38%	1.328%
9	8.936	1030	1038	1045	rVV	305648	573747	14.98%	1.747%
10	9.007	1047	1050	1057	rVB7	18333	33215	0.87%	0.101%
11	9.412	1112	1119	1131	rVB2	261302	491828	12.84%	1.498%
12	9.671	1155	1163	1167	rBV8	12945	27622	0.72%	0.084%
13	9.782	1179	1182	1187	rVB5	12062	22068	0.58%	0.067%
14	10.011	1215	1221	1230	rBV6	11794	23312	0.61%	0.071%
15	10.135	1232	1242	1251	rBV2	164366	328682	8.58%	1.001%
16	10.211	1251	1255	1259	rBV4	29779	57179	1.49%	0.174%
17	10.487	1292	1302	1304	rBV4	45883	98140	2.56%	0.299%
18	10.517	1305	1307	1320	rVB5	38880	70968	1.85%	0.216%
19	10.675	1326	1334	1345	rBV2	330547	636573	16.62%	1.938%
20	10.875	1362	1368	1370	rBV6	9647	17981	0.47%	0.055%
21	10.922	1370	1376	1380	rVV5	12933	25347	0.66%	0.077%
22	11.063	1392	1400	1415	rVB	329345	662377	17.29%	2.017%
23	11.198	1415	1423	1434	rBV2	321593	641510	16.75%	1.953%
24	11.298	1435	1440	1445	rVB7	11856	23756	0.62%	0.072%
25	11.410	1453	1459	1463	rBV7	27337	57232	1.49%	0.174%
26	11.463	1464	1468	1476	rVV5	42869	97240	2.54%	0.296%
27	11.521	1476	1478	1484	rVV5	12517	19341	0.50%	0.059%
28	11.586	1484	1489	1494	rVB5	23866	40844	1.07%	0.124%
29	11.639	1494	1498	1507	rVB9	16011	41637	1.09%	0.127%
30	11.862	1525	1536	1549	rBV3	69426	173253	4.52%	0.528%
31	12.062	1564	1570	1573	rBV4	18776	33045	0.86%	0.101%
32	12.115	1574	1579	1583	rVV8	13698	21072	0.55%	0.064%
33	12.191	1588	1592	1599	rBV9	17260	34487	0.90%	0.105%
34	12.367	1618	1622	1626	rVB5	10579	16465	0.43%	0.050%

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35	12.432	1626	1633	1643	rBV9	16627	44420	1.16%	0.135%
36	12.579	1654	1658	1660	rBV4	13788	23050	0.60%	0.070%
37	12.702	1673	1679	1682	rBV	60742	105359	2.75%	0.321%
38	12.732	1683	1684	1689	rVB4	28397	42083	1.10%	0.128%
39	12.826	1696	1700	1701	rBV3	17738	19495	0.51%	0.059%
40	12.920	1711	1716	1723	rVB3	48534	83060	2.17%	0.253%
41	13.008	1724	1731	1736	rBV6	24615	62571	1.63%	0.191%
42	13.084	1742	1744	1753	rVB8	20723	35473	0.93%	0.108%
43	13.202	1760	1764	1770	rBV9	17127	37055	0.97%	0.113%
44	13.278	1772	1777	1780	rBV6	13779	22731	0.59%	0.069%
45	13.343	1785	1788	1795	rVB8	19762	32706	0.85%	0.100%
46	13.548	1816	1823	1826	rBV6	17422	36650	0.96%	0.112%
47	13.642	1833	1839	1844	rBV7	24721	46977	1.23%	0.143%
48	13.689	1844	1847	1853	rVB6	12265	20317	0.53%	0.062%
49	13.883	1878	1880	1891	rBV6	21847	56128	1.47%	0.171%
50	14.042	1897	1907	1916	rVB9	33074	102073	2.66%	0.311%
51	14.183	1925	1931	1935	rBV5	29837	56374	1.47%	0.172%
52	14.248	1935	1942	1947	rVV	740206	1166304	30.44%	3.551%
53	14.295	1947	1950	1956	rVB	209311	276998	7.23%	0.843%
54	14.412	1966	1970	1973	rBV6	17254	19605	0.51%	0.060%
55	14.553	1987	1994	2001	rBV	706840	1170483	30.55%	3.564%
56	14.706	2015	2020	2025	rBV5	48735	63895	1.67%	0.195%
57	14.812	2034	2038	2040	rBV2	40274	54344	1.42%	0.165%
58	14.859	2040	2046	2055	rVB	624149	1047043	27.33%	3.188%
59	15.053	2072	2079	2090	rBV2	222528	478114	12.48%	1.456%
60	15.147	2092	2095	2101	rVB6	23491	39964	1.04%	0.122%
61	15.317	2121	2124	2129	rVB7	18559	22593	0.59%	0.069%
62	15.458	2142	2148	2154	rBV10	19844	44171	1.15%	0.134%
63	15.517	2154	2158	2162	rBV6	19745	33737	0.88%	0.103%
64	15.611	2167	2174	2177	rBV6	45528	87924	2.30%	0.268%
65	15.646	2177	2180	2191	rVB2	89682	144992	3.78%	0.441%
66	15.846	2206	2214	2228	rVB	1009639	1673854	43.69%	5.097%
67	15.963	2228	2234	2239	rBV10	43735	83072	2.17%	0.253%
68	16.045	2246	2248	2252	rBV4	12648	21054	0.55%	0.064%
69	16.145	2263	2265	2270	rBV5	11879	20988	0.55%	0.064%
70	16.451	2311	2317	2322	rVV9	26361	54903	1.43%	0.167%
71	16.504	2322	2326	2338	rVB2	123080	229318	5.99%	0.698%

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72	16.727	2361	2364	2369	rVB7	33784	51031	1.33%	0.155%
73	16.809	2376	2378	2382	rVB5	19112	21127	0.55%	0.064%
74	17.003	2409	2411	2416	rBV6	13447	18357	0.48%	0.056%
75	17.056	2416	2420	2427	rBV2	49573	92194	2.41%	0.281%
76	17.121	2428	2431	2439	rVB9	16204	30928	0.81%	0.094%
77	17.297	2457	2461	2466	rBV	110613	122065	3.19%	0.372%
78	17.438	2482	2485	2492	rVB9	20336	40619	1.06%	0.124%
79	17.597	2505	2512	2523	rBV	933324	1470581	38.39%	4.478%
80	17.697	2523	2529	2538	rVB2	1214080	1896640	49.51%	5.775%
81	17.773	2539	2542	2548	rVB8	23022	44032	1.15%	0.134%
82	17.855	2549	2556	2560	rBV10	9606	24015	0.63%	0.073%
83	18.037	2583	2587	2593	rVV	96038	124549	3.25%	0.379%
84	18.431	2650	2654	2659	rBV7	16571	36355	0.95%	0.111%
85	18.595	2679	2682	2687	rBV6	19962	26043	0.68%	0.079%
86	18.637	2687	2689	2695	rBV7	10870	23414	0.61%	0.071%
87	18.719	2699	2703	2707	rBV2	53549	61331	1.60%	0.187%
88	19.342	2806	2809	2816	rVB	50544	74915	1.96%	0.228%
89	19.606	2851	2854	2859	rBV6	26593	42587	1.11%	0.130%
90	19.765	2879	2881	2889	rVB9	19924	29510	0.77%	0.090%
91	19.859	2892	2897	2901	rBV	230603	303899	7.93%	0.925%
92	19.970	2910	2916	2923	rBV	1724180	2465658	64.36%	7.507%
93	20.417	2990	2992	2993	rBV2	26432	16908	0.44%	0.051%
94	20.840	3055	3064	3075	rBV	2778067	3831005	100.00%	11.665%
95	20.946	3078	3082	3090	rVB	177514	282087	7.36%	0.859%
96	21.463	3166	3170	3177	rBV2	127762	239316	6.25%	0.729%
97	21.886	3235	3242	3253	rVB	1093608	1969170	51.40%	5.996%
98	23.384	3489	3497	3509	rBV	759307	1600217	41.77%	4.872%
99	25.064	3772	3783	3794	rVB	793029	2309784	60.29%	7.033%
100	25.299	3814	3823	3837	rVB2	624357	1908181	49.81%	5.810%

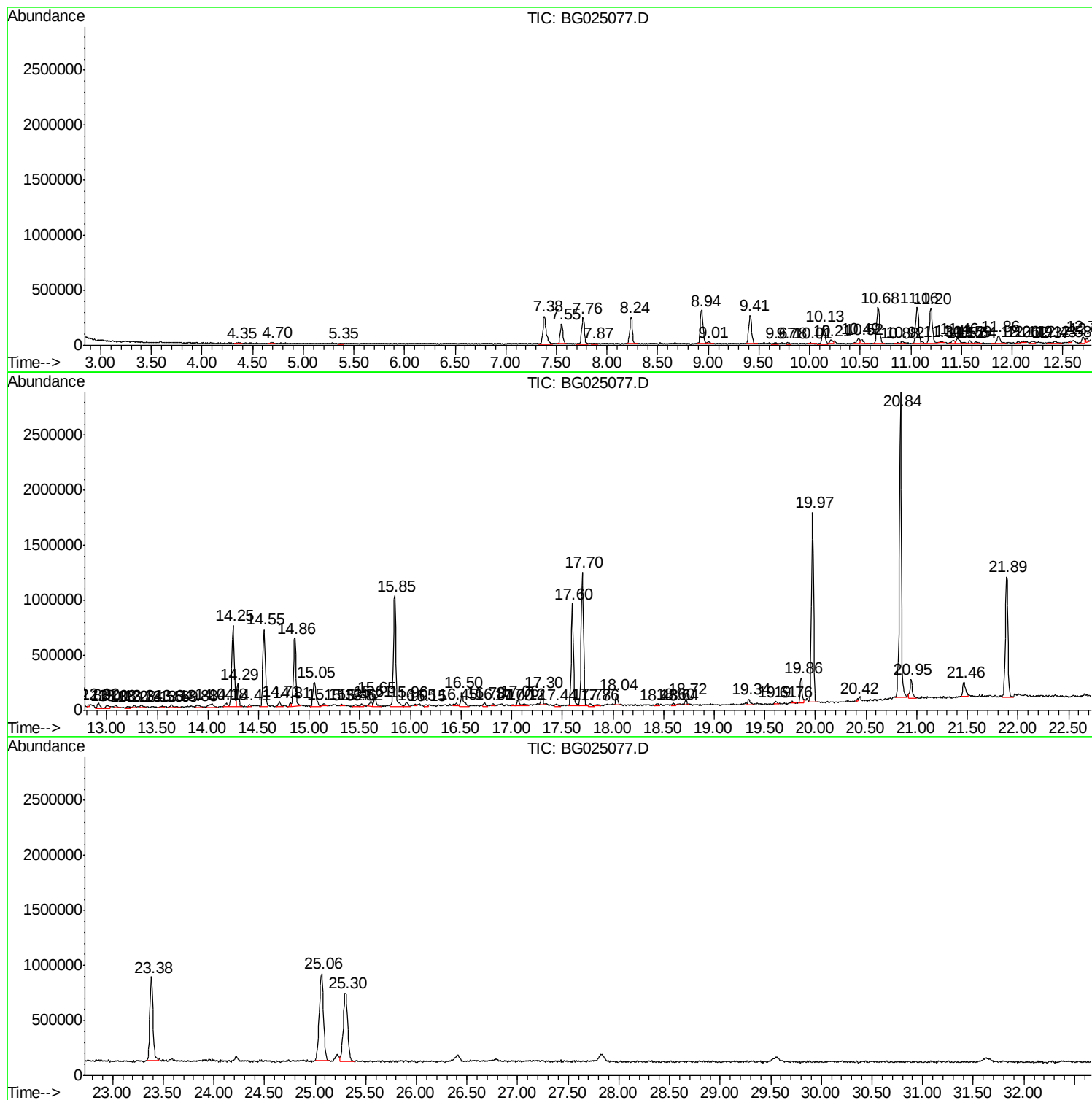
Sum of corrected areas: 32843192

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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



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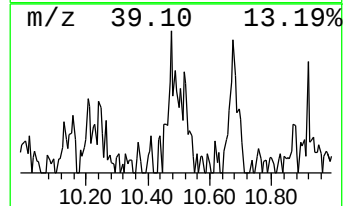
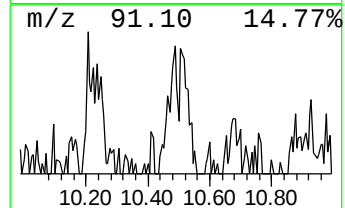
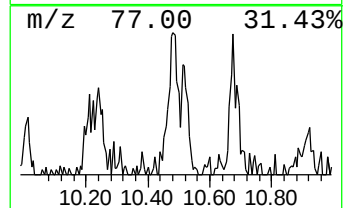
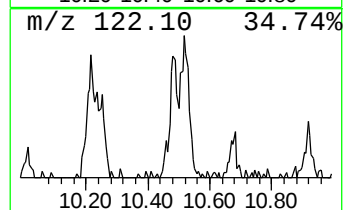
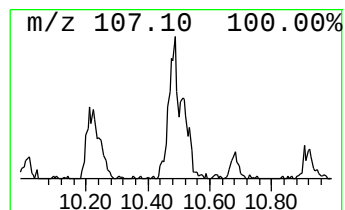
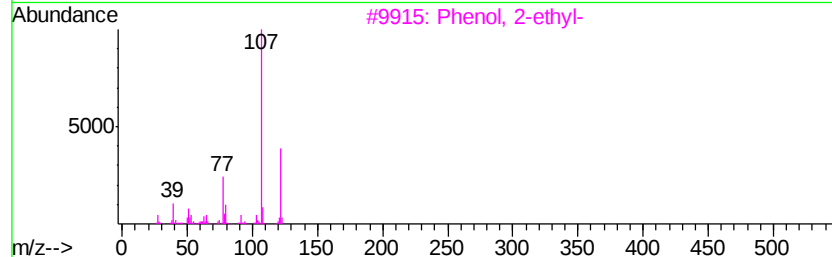
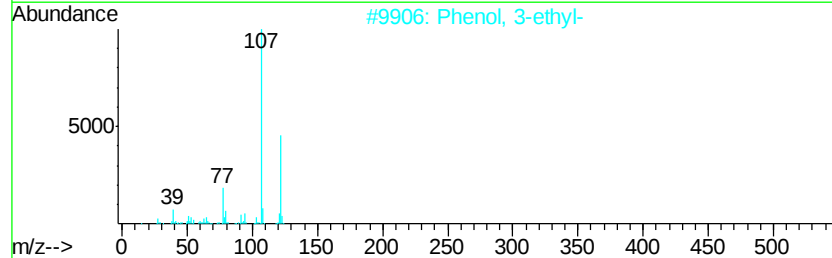
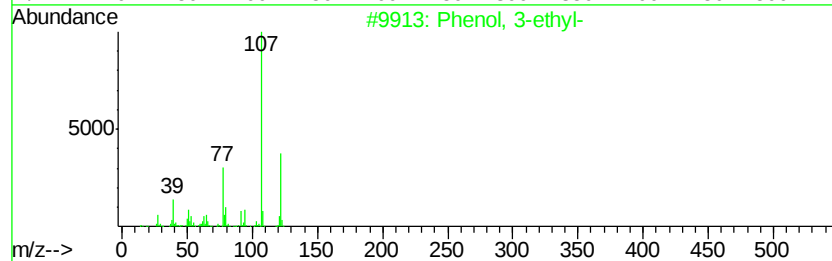
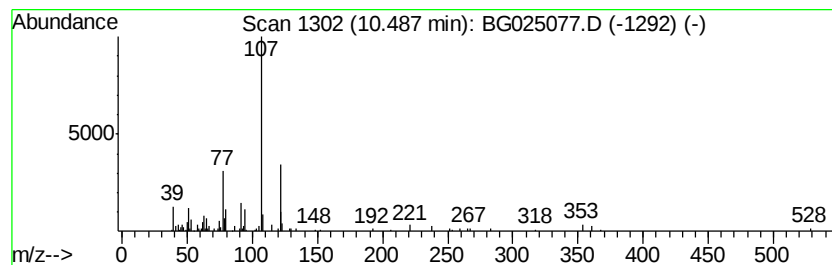
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	2.96 ng/ul	98140	Napthalene-d8	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	90	
2	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	86	
3	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	81	
4	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	81	
5	Phenol, 3-ethyl-	122	C8H10O	000620-17-7	80	



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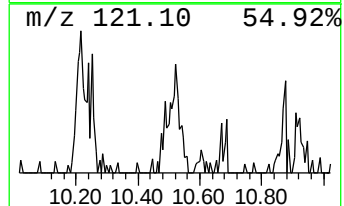
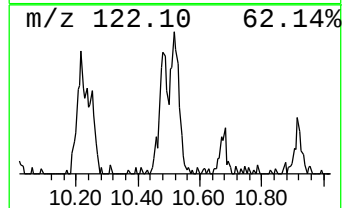
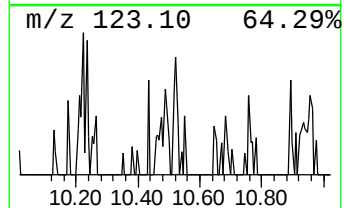
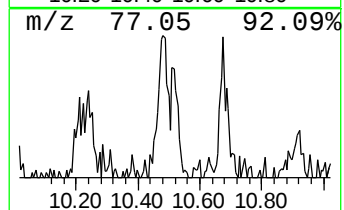
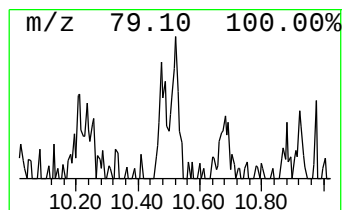
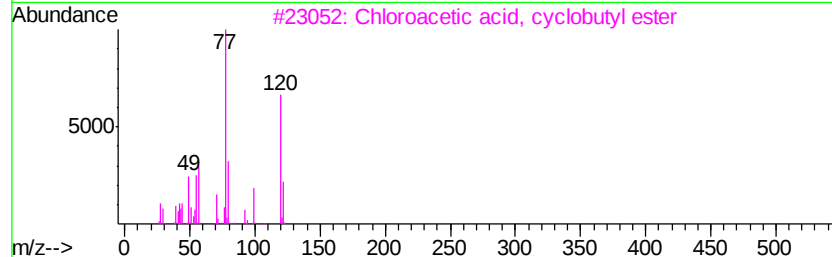
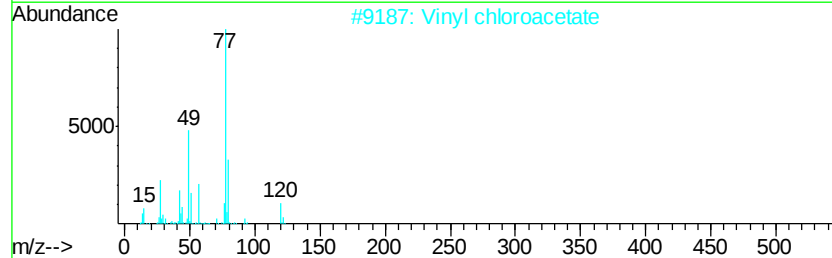
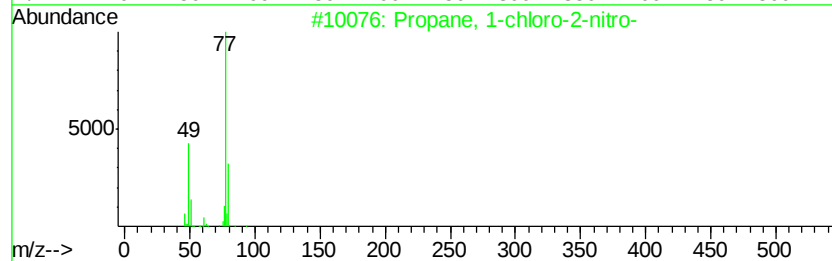
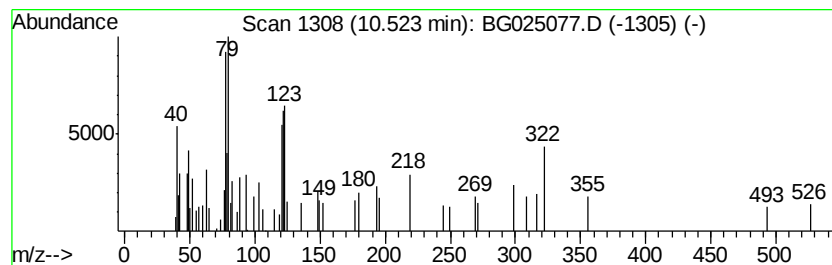
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.52	2.14 ng/ul	70968	Naphthalene-d8	11.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1-chloro-2-nitro-	123	C3H6ClNO2	002425-66-3	27
2		Vinyl chloroacetate	120	C4H5ClO2	002549-51-1	27
3		Chloroacetic acid, cyclobutyl ester	148	C6H9ClO2	1000282-92-5	17
4		2-Propanone, 1,1,3-trichloro-	160	C3H3Cl3O	000921-03-9	16
5		.alpha.-d-Glucopyranose, 2-azido...	399	C16H18ClN3O5S	146934-31-8	16



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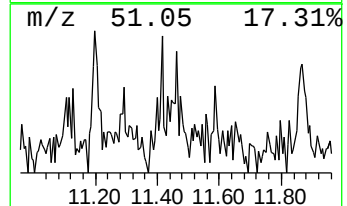
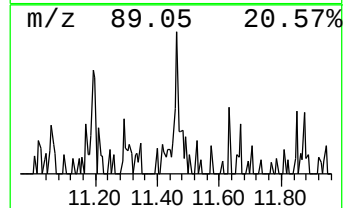
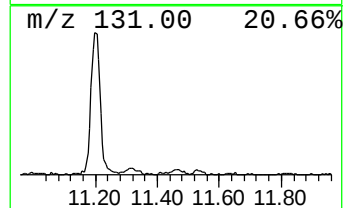
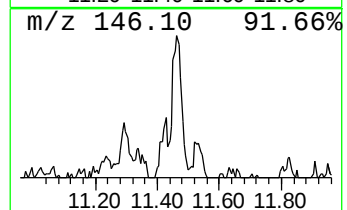
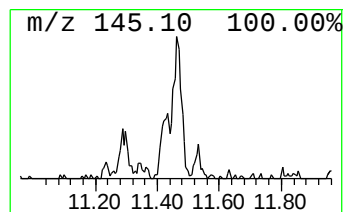
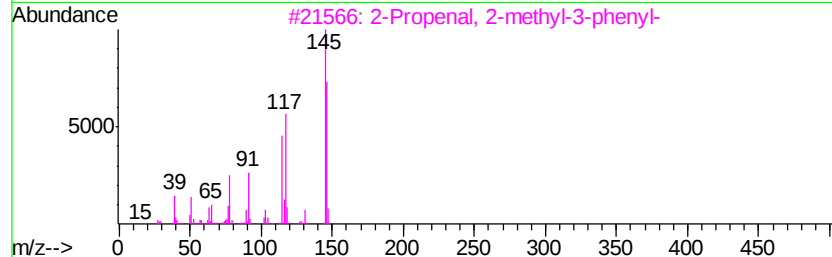
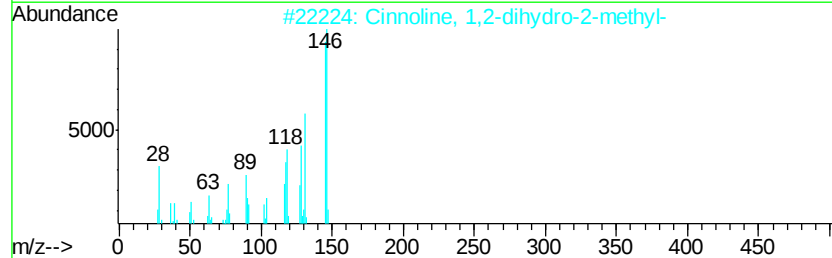
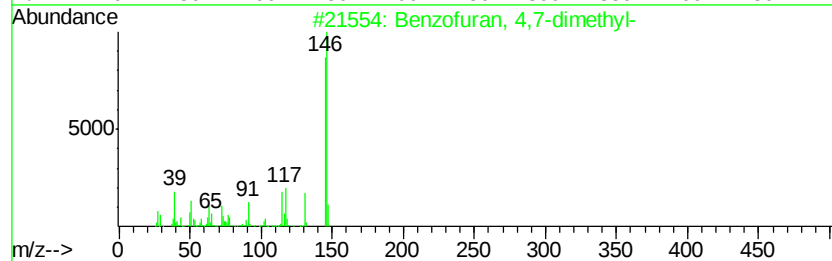
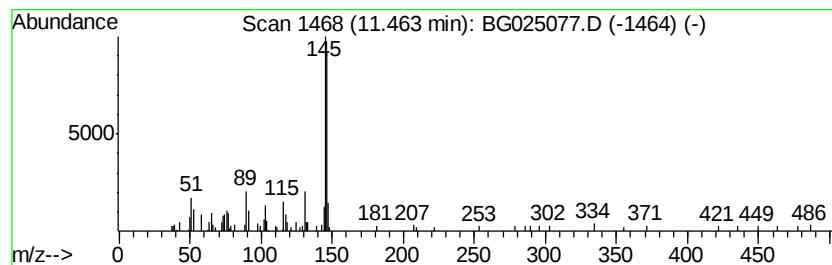
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Peak Number 3 Benzofuran, 4,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.46	2.94 ng/ul	97240	Napthalene-d8	11.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzofuran, 4,7-dimethyl-	146 C10H10O	028715-26-6 80
2		Cinnoline, 1,2-dihydro-2-methyl-	146 C9H10N2	054063-10-4 72
3		2-Propenal, 2-methyl-3-phenyl-	146 C10H10O	000101-39-3 64
4		Benzonitrile, 4-(dimethylamino)-	146 C9H10N2	001197-19-9 64
5		1H-Benzimidazole, 2-ethyl-	146 C9H10N2	001848-84-6 64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
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Operator : UM/SJ
Sample : H5905-03
Misc :
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Instrument :
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ClientSampleId :
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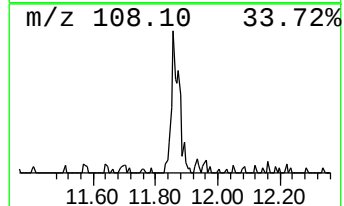
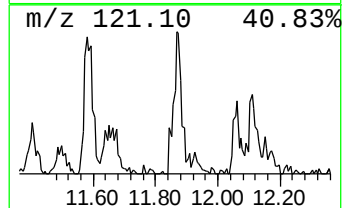
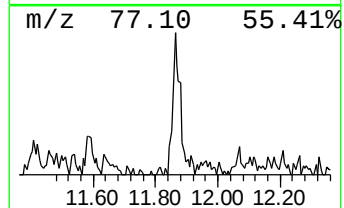
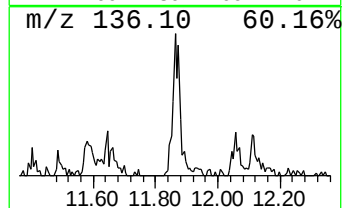
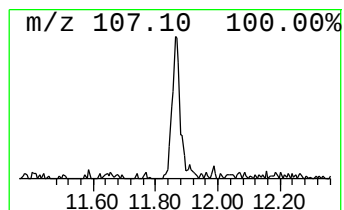
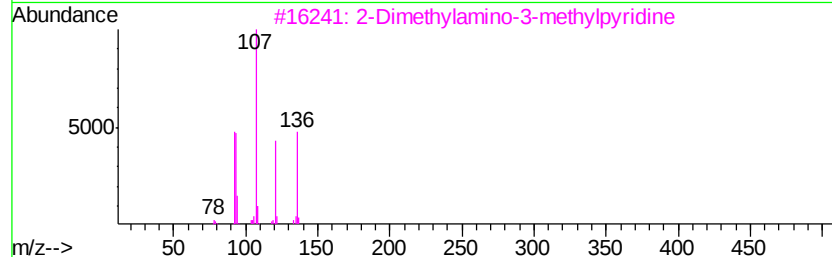
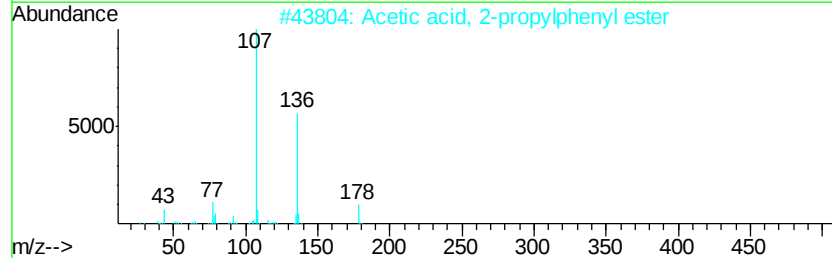
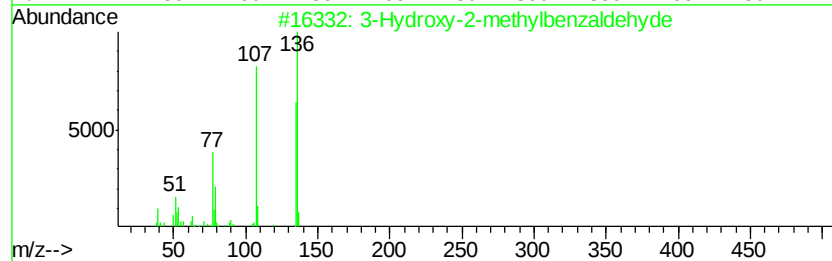
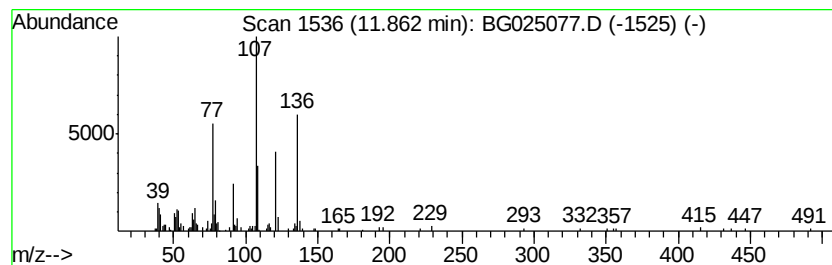
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 3-Hydroxy-2-methylbenzaldehyde Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	5.23 ng/ul	173253	Naphthalene-d8	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Hydroxy-2-methylbenzaldehyde	136	C8H8O2	090111-15-2	64
2		Acetic acid, 2-propylphenyl ester	178	C11H14O2	035922-83-9	53
3		2-Dimethylamino-3-methylpyridine	136	C8H12N2	061713-46-0	50
4		Benzaldehyde, 3-methoxy-	136	C8H8O2	000591-31-1	38
5		Phenol, 2-propyl-	136	C9H12O	000644-35-9	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
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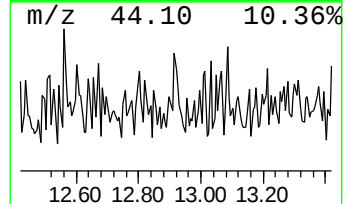
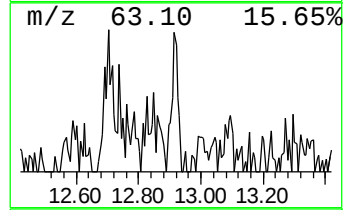
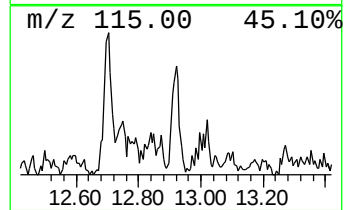
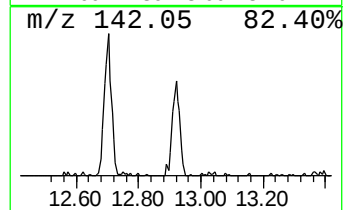
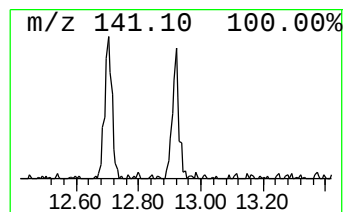
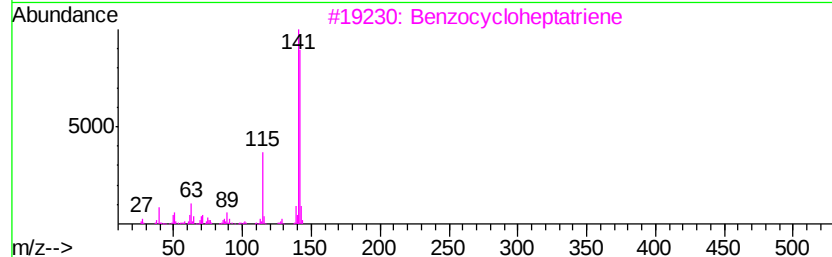
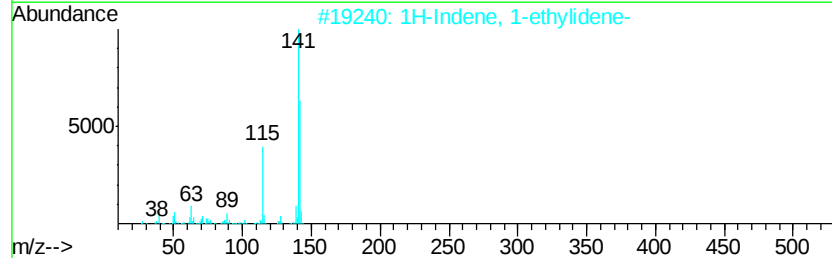
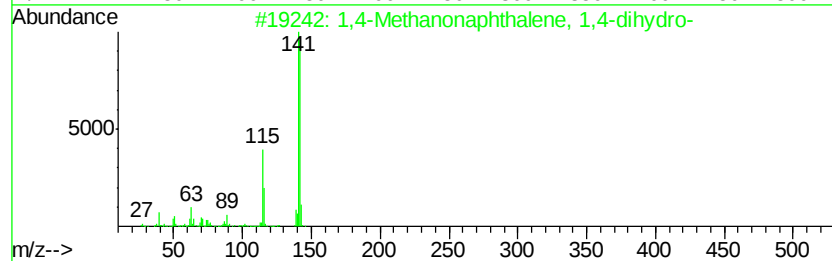
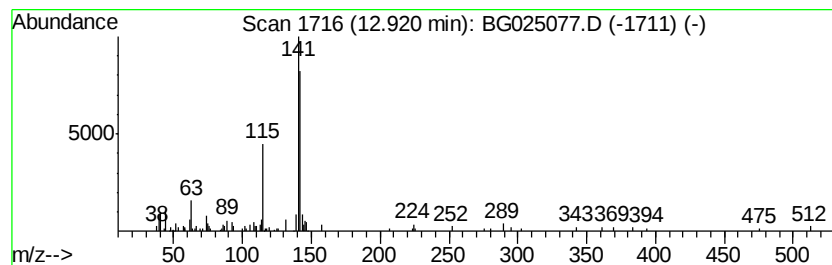
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,4-Methanonaphthalene, 1,4... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	2.51 ng/ul	83060	Naphthalene-d8	11.06

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10		004453-90-1	90
2	1H-Indene, 1-ethylidene-	142	C11H10		002471-83-2	87
3	Benzocycloheptatriene	142	C11H10		000264-09-5	87
4	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10		002443-46-1	87
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10		004453-90-1	78



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
Operator : UM/SJ
Sample : H5905-03
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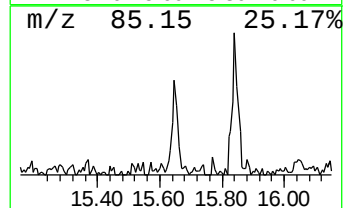
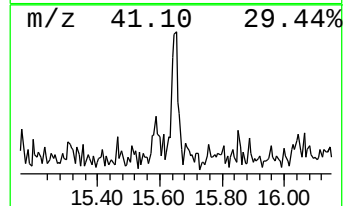
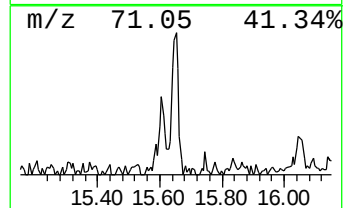
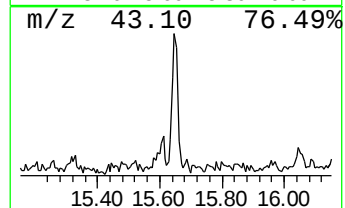
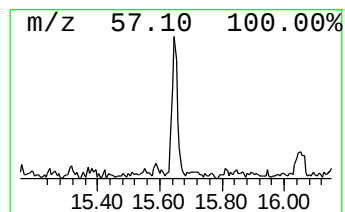
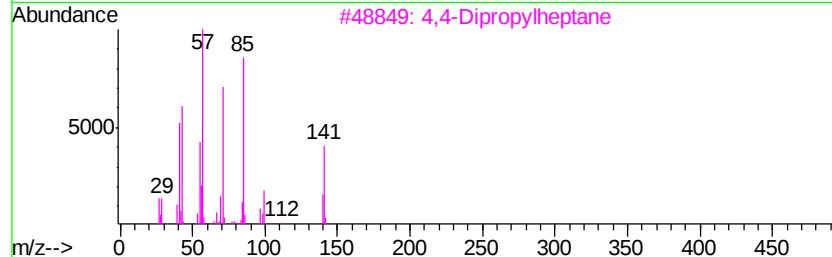
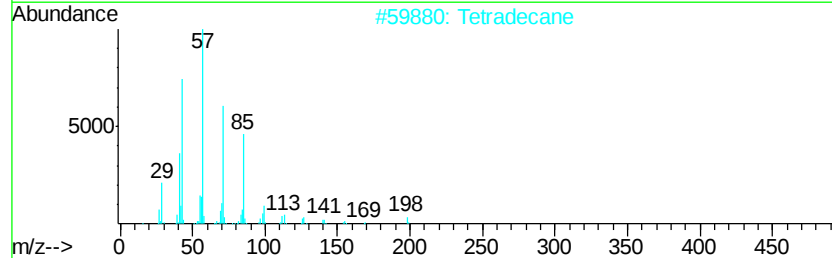
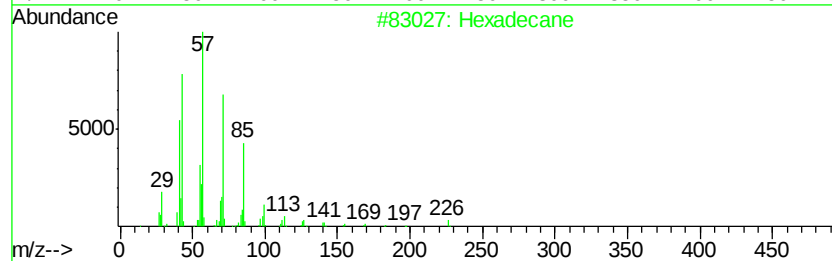
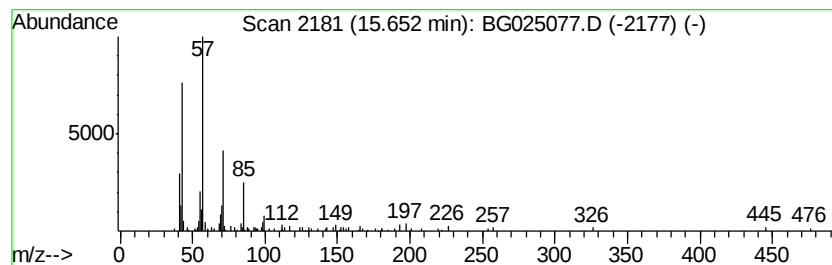
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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.65	2.77 ng/ul	144992	Acenaphthene-d10		14.86	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	52
2	Tetradecane		198	C14H30	000629-59-4	50
3	4,4-Dipropylheptane		184	C13H28	017312-72-0	50
4	Decane, 2,3,6-trimethyl-		184	C13H28	062238-12-4	50
5	Pentadecane		212	C15H32	000629-62-9	50



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
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Sample : H5905-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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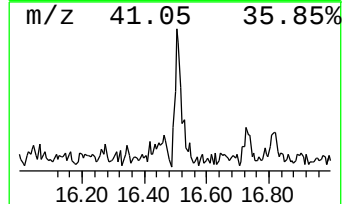
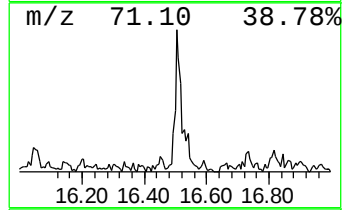
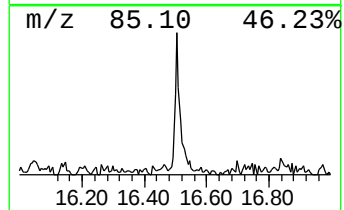
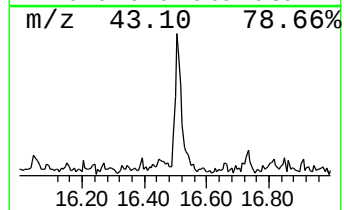
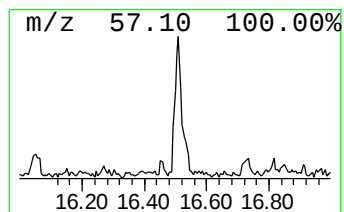
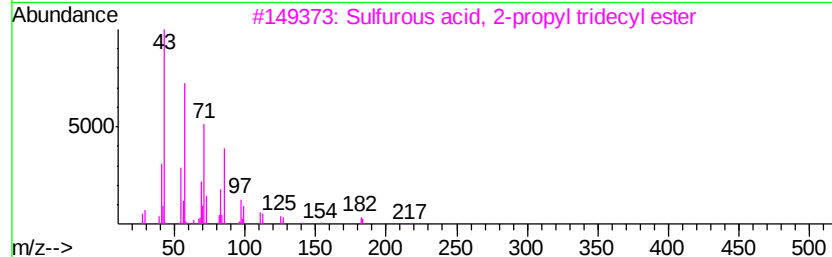
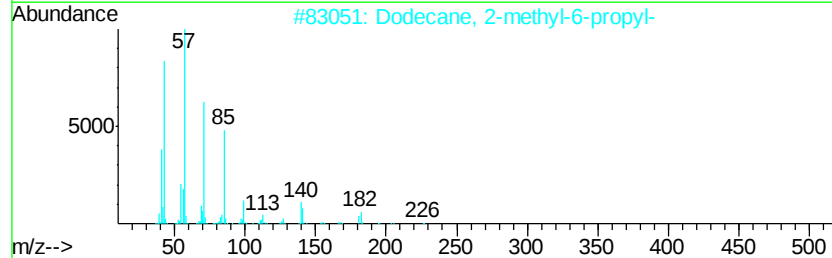
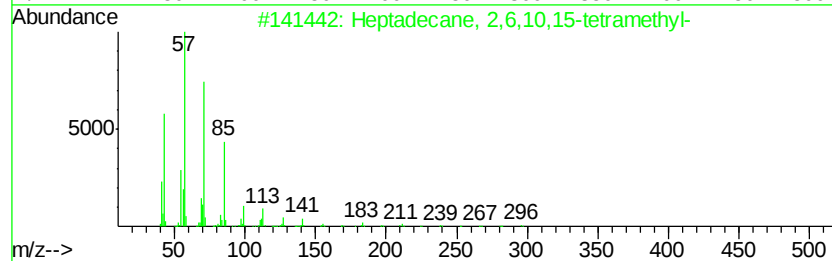
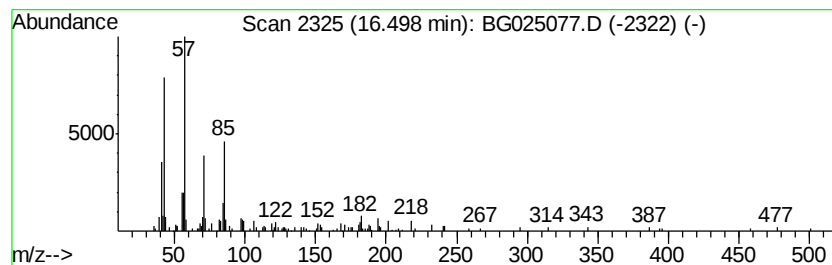
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	3.12 ng/ul	229318	Phenanthrene-d10	17.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
3			Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	80
4			1-Iodo-2-methylnonane	268	C10H21I	1000101-47-9	80
5			Hexadecane, 7-methyl-	240	C17H36	026730-20-1	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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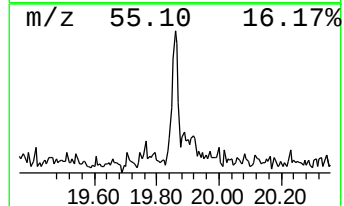
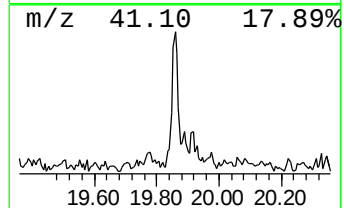
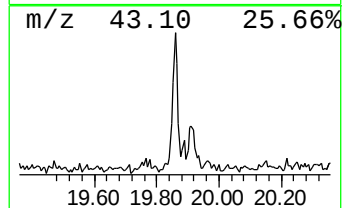
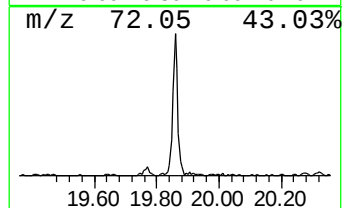
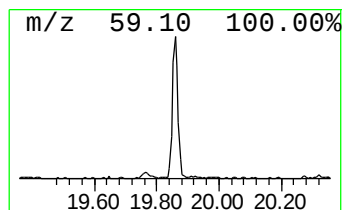
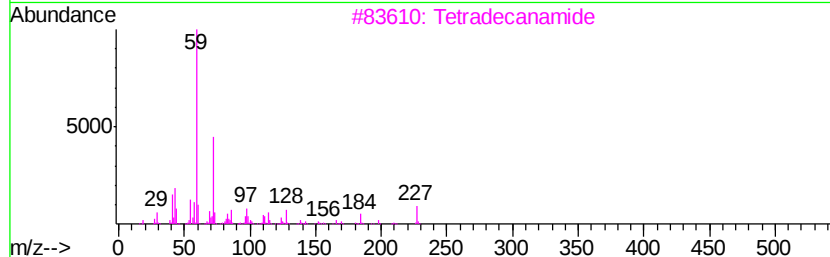
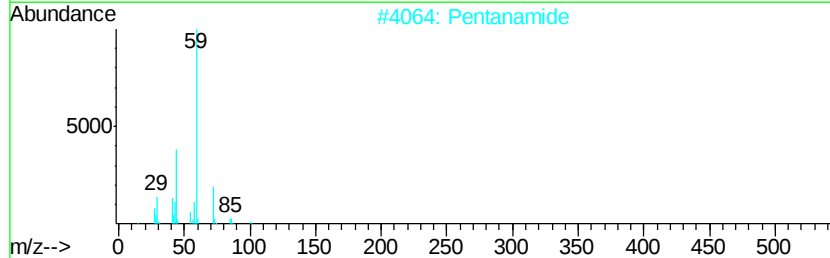
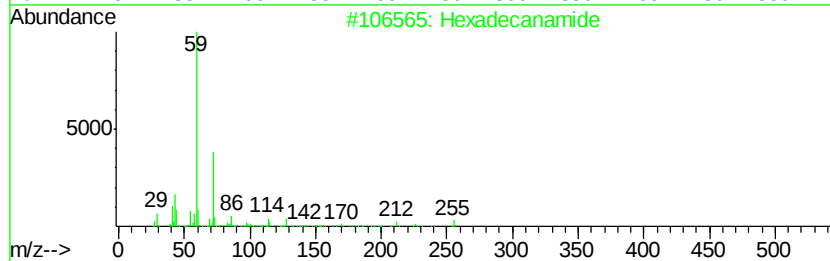
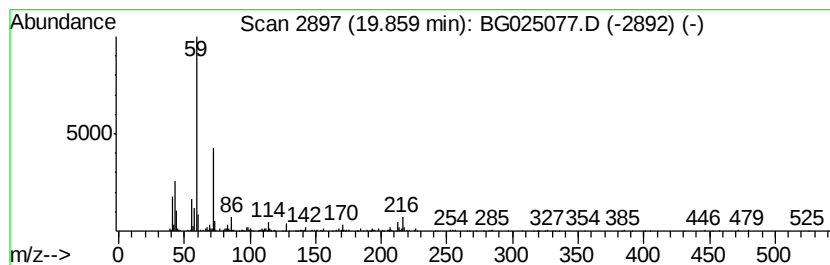
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexadecanamide Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.86	3.09 ng/ul	303899	Chrysene-d12	21.89

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanamide		255	C16H33NO	000629-54-9	80
2	Pentanamide		101	C5H11NO	000626-97-1	64
3	Tetradecanamide		227	C14H29NO	000638-58-4	56
4	9-Octadecenamide, (Z)-		281	C18H35NO	000301-02-0	56
5	Octanamide		143	C8H17NO	000629-01-6	56



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
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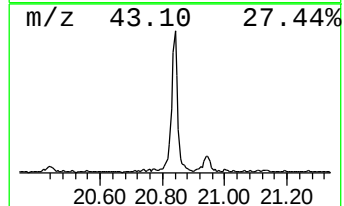
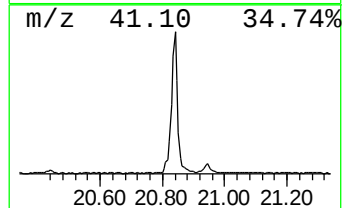
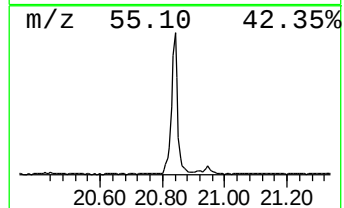
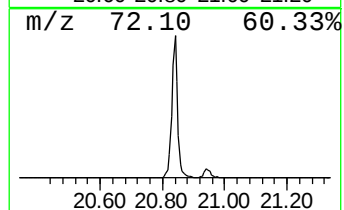
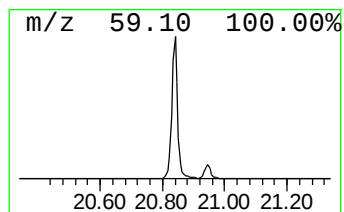
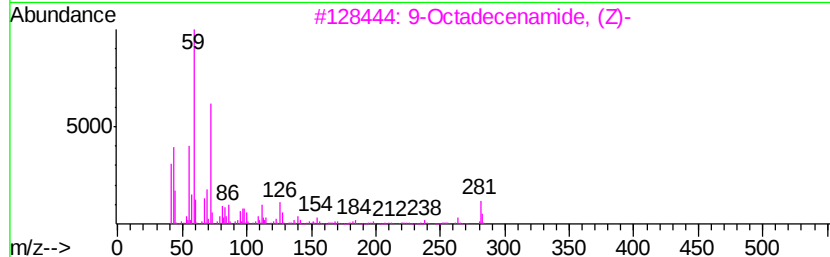
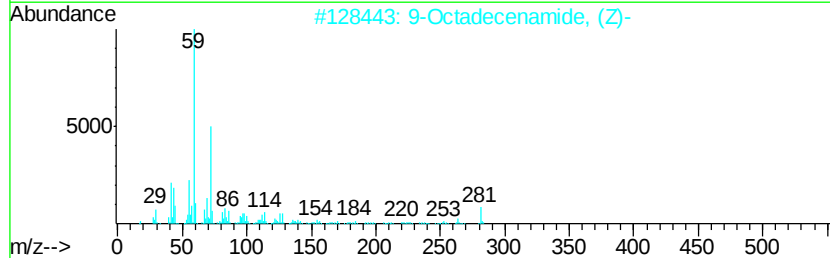
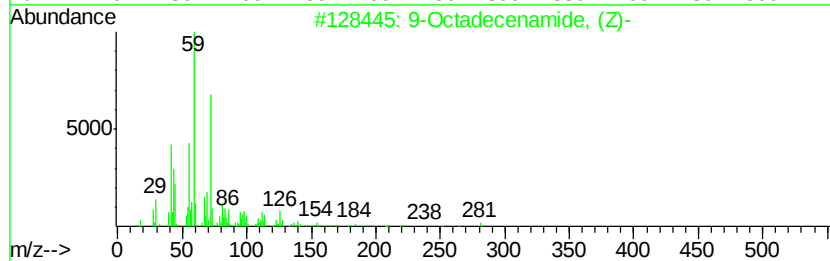
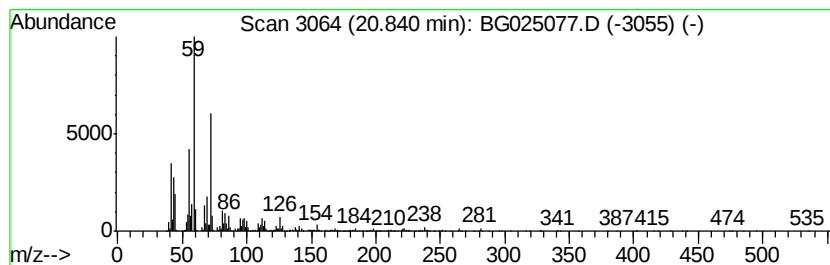
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.84	38.91 ng/ul	3831010	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	99
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
3		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	78
5		Tetradecanamide	227	C14H29NO	000638-58-4	72



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
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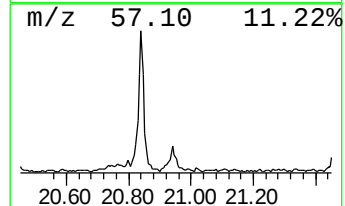
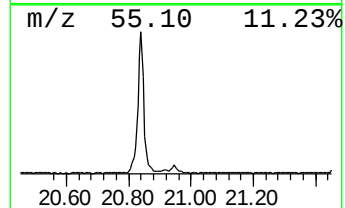
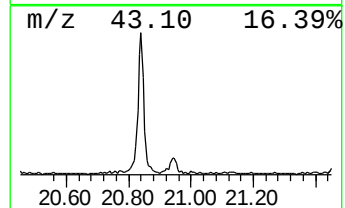
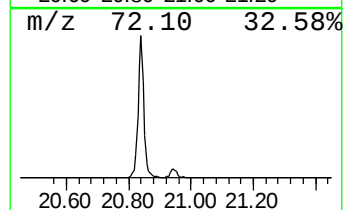
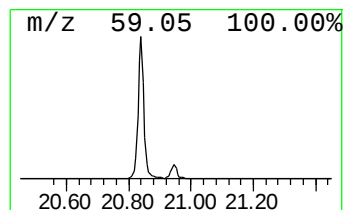
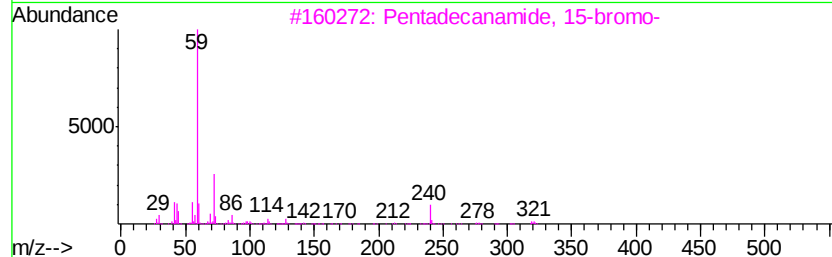
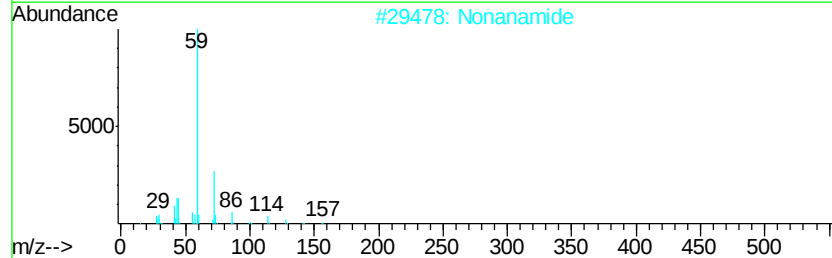
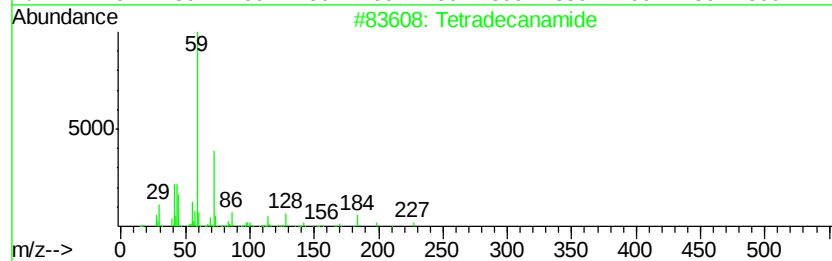
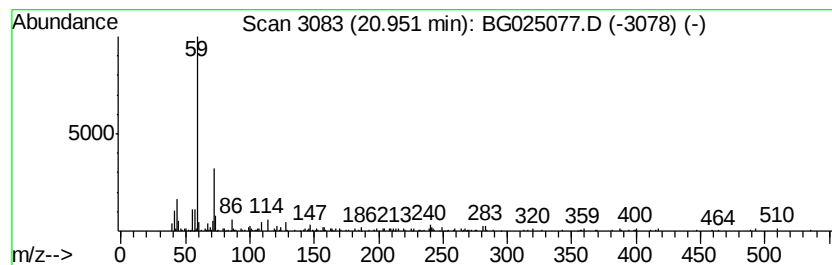
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Tetradecanamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	2.87 ng/ul	282087	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecanamide	227	C14H29NO	000638-58-4	81
2		Nonanamide	157	C9H19NO	001120-07-6	74
3		Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	72
4		Tetradecanamide	227	C14H29NO	000638-58-4	64
5		Hexadecanamide	255	C16H33NO	000629-54-9	64



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
Operator : UM/SJ
Sample : H5905-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DA805

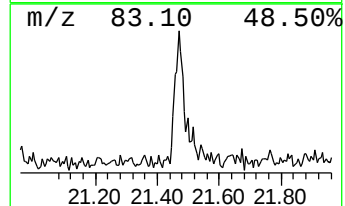
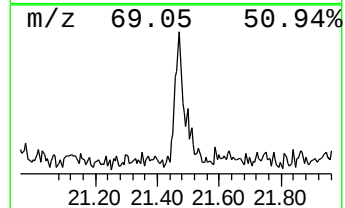
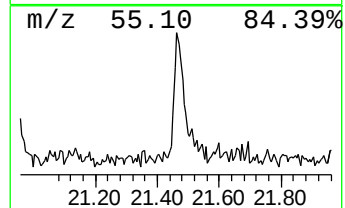
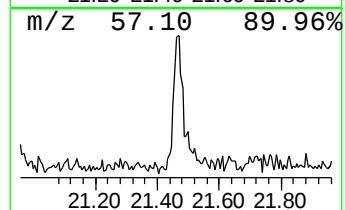
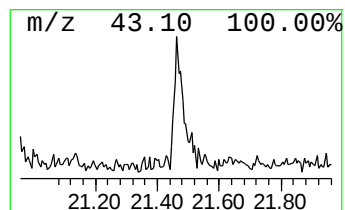
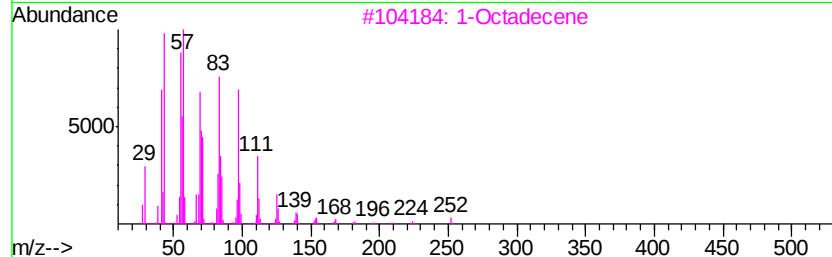
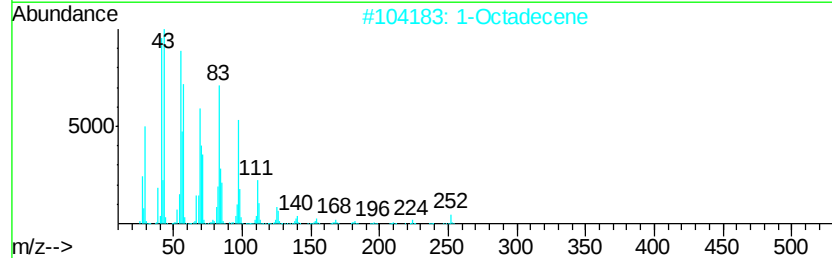
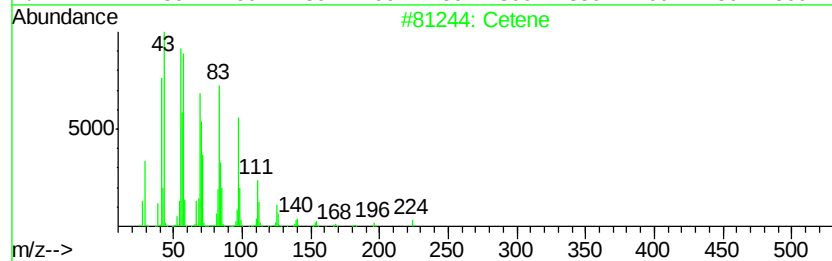
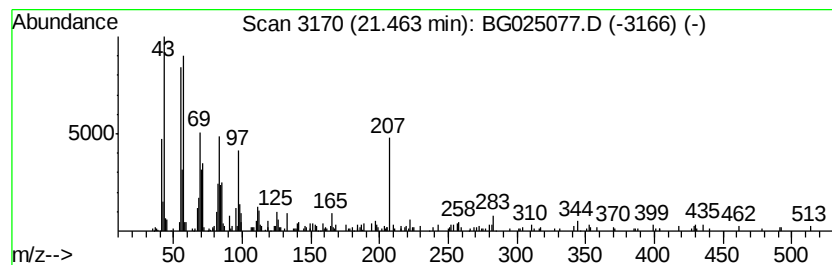
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic21.46 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.46	2.43 ng/ul	239316	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cetene	224	C16H32	000629-73-2	95
2		1-Octadecene	252	C18H36	000112-88-9	90
3		1-Octadecene	252	C18H36	000112-88-9	90
4		1-Heptadecene	238	C17H34	006765-39-5	78
5		E-15-Heptadecenal	252	C17H32O	1000130-97-9	78



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121016\
Data File : BG025077.D
Acq On : 10 Dec 2016 20:34
Operator : UM/SJ
Sample : H5905-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DA805

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.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
Phenol, 3-ethyl-	10.49	3.0	ng/ul	98140	2	11.06	662377	20.0
unknown-01	10.52	2.1	ng/ul	70968	2	11.06	662377	20.0
Benzofuran, 4,7-d...	11.46	2.9	ng/ul	97240	2	11.06	662377	20.0
3-Hydroxy-2-methy...	11.86	5.2	ng/ul	173253	2	11.06	662377	20.0
1,4-Methanonaphth...	12.92	2.5	ng/ul	83060	2	11.06	662377	20.0
(DEL) Alkane: Str...	15.65	2.8	ng/ul	144992	3	14.86	1047040	20.0
(DEL) Alkane: Str...	16.50	3.1	ng/ul	229318	4	17.60	1470580	20.0
Hexadecanamide	19.86	3.1	ng/ul	303899	5	21.89	1969170	20.0
9-Octadecenamide, ...	20.84	38.9	ng/ul	3831010	5	21.89	1969170	20.0
Tetradecanamide	20.95	2.9	ng/ul	282087	5	21.89	1969170	20.0
(DEL) Alkane: Cyc...	21.46	2.4	ng/ul	239316	5	21.89	1969170	20.0