

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
 Data File : BG025350.D
 Acq On : 20 Dec 2016 12:10
 Operator : UM/SJ
 Sample : H5905-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DA811

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	2.867	3	5	17	rVB7	26413	63450	0.24%	0.120%
2	3.249	65	70	78	rVB5	27503	47652	0.18%	0.090%
3	3.343	78	86	90	rBV5	20999	36662	0.14%	0.069%
4	3.583	121	127	140	rBV	483515	728010	2.80%	1.371%
5	4.106	211	216	224	rBV6	24421	45876	0.18%	0.086%
6	6.321	585	593	604	rBV2	136517	301050	1.16%	0.567%
7	7.379	765	773	789	rBV2	254441	558176	2.15%	1.051%
8	7.538	794	800	812	rVB	183117	313053	1.20%	0.590%
9	7.755	829	837	847	rBV	250948	464162	1.78%	0.874%
10	8.219	909	916	931	rBV3	266856	613934	2.36%	1.156%
11	8.936	1028	1038	1054	rVV2	303324	619074	2.38%	1.166%
12	9.406	1110	1118	1128	rBV2	269403	514400	1.98%	0.969%
13	10.129	1232	1241	1250	rBV2	260232	496733	1.91%	0.936%
14	10.675	1326	1334	1351	rBV2	340774	704024	2.71%	1.326%
15	11.051	1388	1398	1412	rVB	372110	712263	2.74%	1.342%
16	11.204	1415	1424	1434	rBV	88974	215584	0.83%	0.406%
17	13.225	1759	1768	1775	rBV7	30955	60939	0.23%	0.115%
18	13.337	1779	1787	1796	rVB9	11013	31660	0.12%	0.060%
19	13.625	1827	1836	1842	rBV5	14981	33232	0.13%	0.063%
20	14.036	1894	1906	1915	rBV8	27948	86454	0.33%	0.163%
21	14.165	1922	1928	1933	rBV7	19175	37330	0.14%	0.070%
22	14.242	1933	1941	1945	rVV	663796	1044102	4.01%	1.967%
23	14.283	1945	1948	1956	rVB	163836	266291	1.02%	0.502%
24	14.547	1985	1993	2007	rVB	674620	1126341	4.33%	2.121%
25	14.682	2011	2016	2027	rBV5	19707	35776	0.14%	0.067%
26	14.847	2034	2044	2056	rBV	633553	1020915	3.92%	1.923%
27	15.082	2071	2084	2097	rBV3	169716	458908	1.76%	0.864%
28	15.229	2102	2109	2118	rVV3	25800	72293	0.28%	0.136%
29	15.317	2118	2124	2129	rVB9	12787	32460	0.12%	0.061%
30	15.464	2141	2149	2152	rBV8	11829	28402	0.11%	0.053%
31	15.587	2164	2170	2174	rBV3	21617	40406	0.16%	0.076%
32	15.634	2174	2178	2185	rVB5	53137	85550	0.33%	0.161%
33	15.834	2198	2212	2223	rBV	944263	1547328	5.95%	2.914%
34	15.969	2230	2235	2242	rBV	262223	436175	1.68%	0.822%

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Integration Parameters: LSCINT.P

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35	16.034	2243	2246	2250	rVV5	21325	34223	0.13%	0.064%
36	16.081	2250	2254	2258	rVB7	16097	26770	0.10%	0.050%
37	16.186	2266	2272	2276	rVB8	17913	35725	0.14%	0.067%
38	16.492	2318	2324	2339	rBV5	58090	147393	0.57%	0.278%
39	16.874	2385	2389	2397	rVV9	14389	36112	0.14%	0.068%
40	17.279	2452	2458	2462	rBV	59550	78418	0.30%	0.148%
41	17.332	2462	2467	2473	rVB8	15534	32176	0.12%	0.061%
42	17.591	2503	2511	2517	rBV	776461	1237973	4.76%	2.332%
43	17.691	2522	2528	2536	rBV2	974326	1546705	5.94%	2.913%
44	18.020	2579	2584	2594	rVB	80780	117864	0.45%	0.222%
45	18.319	2627	2635	2640	rBV	13089	36293	0.14%	0.068%
46	18.431	2649	2654	2662	rBV4	139670	245713	0.94%	0.463%
47	18.513	2663	2668	2674	rVB9	30019	57786	0.22%	0.109%
48	18.648	2683	2691	2693	rBV8	22637	48531	0.19%	0.091%
49	18.701	2696	2700	2705	rBV2	60373	77183	0.30%	0.145%
50	18.860	2719	2727	2730	rBV9	20342	42251	0.16%	0.080%
51	19.165	2775	2779	2782	rBV6	20031	35263	0.14%	0.066%
52	19.324	2796	2806	2810	rBV5	57937	159699	0.61%	0.301%
53	19.494	2831	2835	2841	rBV	192829	300590	1.16%	0.566%
54	19.659	2859	2863	2865	rBV3	42857	49857	0.19%	0.094%
55	19.688	2865	2868	2872	rVB6	56130	61878	0.24%	0.117%
56	19.735	2873	2876	2880	rVB6	37146	41897	0.16%	0.079%
57	19.894	2900	2903	2908	rVB5	62051	79155	0.30%	0.149%
58	19.964	2908	2915	2926	rBV	1285420	1998886	7.68%	3.765%
59	20.428	2990	2994	3003	rBV9	66878	191034	0.73%	0.360%
60	21.451	3164	3168	3179	rVB7	162011	394752	1.52%	0.744%
61	21.721	3205	3214	3221	rBV	16736446	26019107	100.00%	49.007%
62	21.886	3236	3242	3251	rBV	846420	1508528	5.80%	2.841%
63	22.373	3320	3325	3329	rBV3	231791	349175	1.34%	0.658%
64	25.082	3777	3786	3795	rBV3	588006	1646662	6.33%	3.101%
65	25.311	3819	3825	3837	rVB2	481669	1331300	5.12%	2.507%
66	27.097	4121	4129	4143	rVB7	342526	1256381	4.83%	2.366%
67	28.284	4323	4331	4345	rVB7	277576	989169	3.80%	1.863%

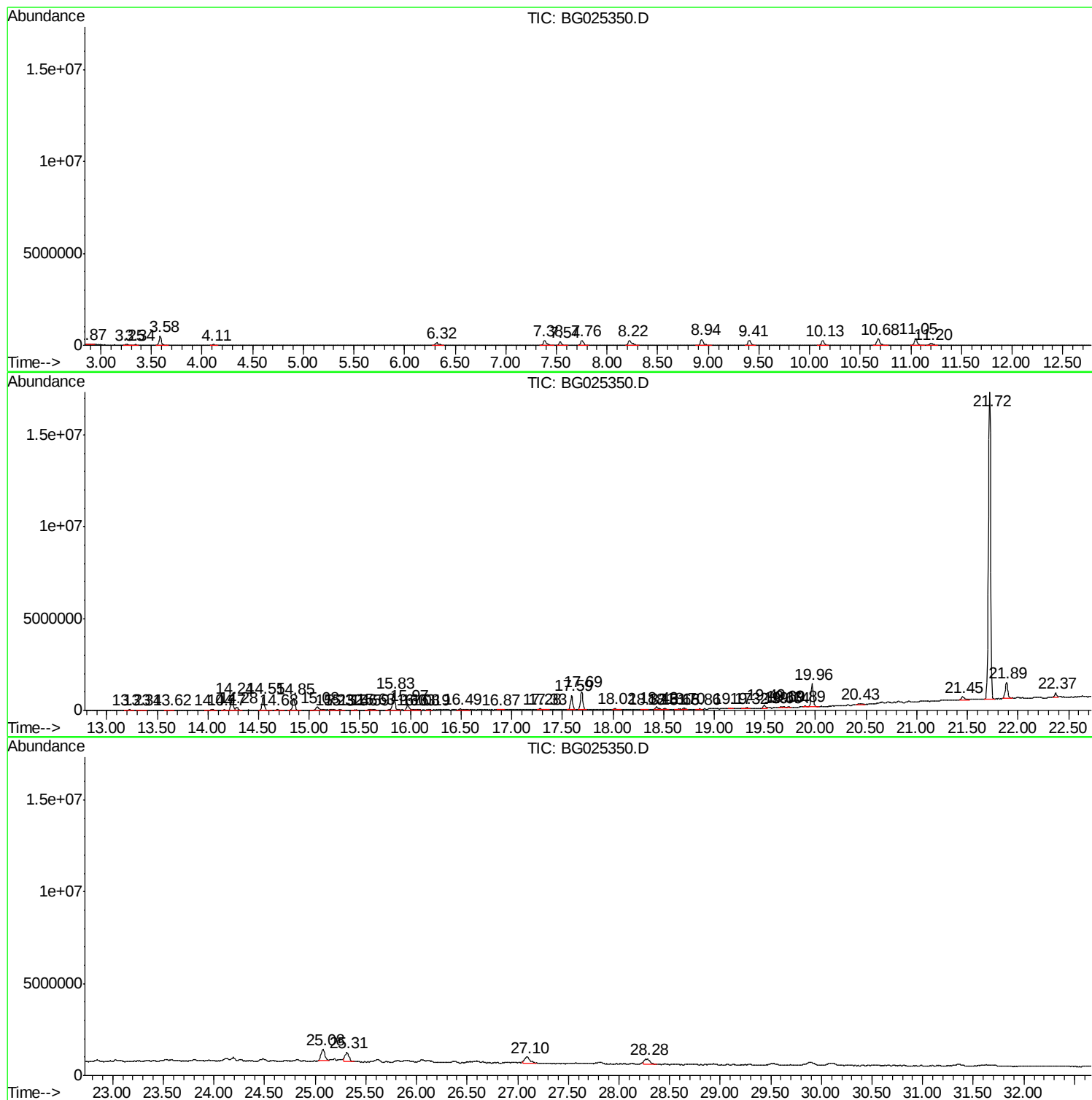
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Quant Title : SVOA CALIBRATION

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



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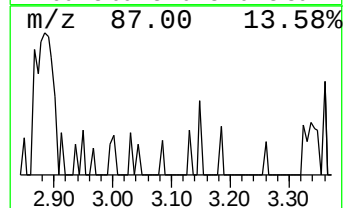
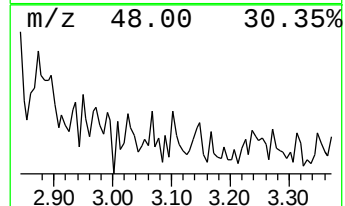
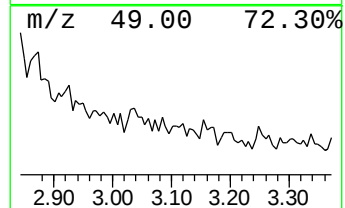
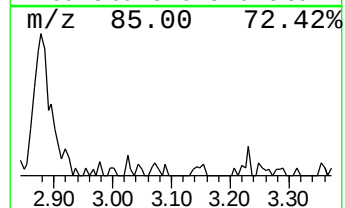
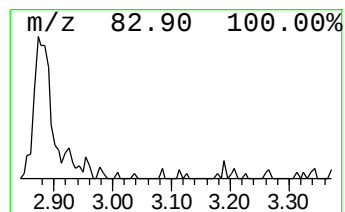
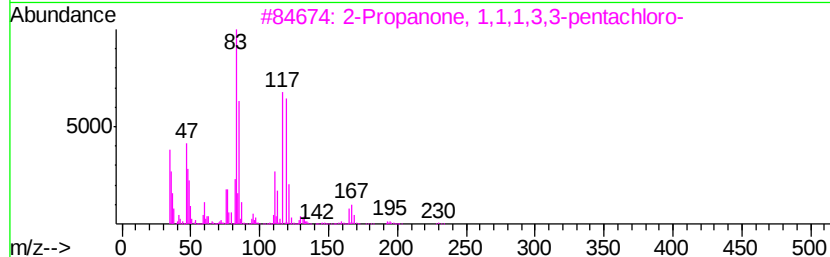
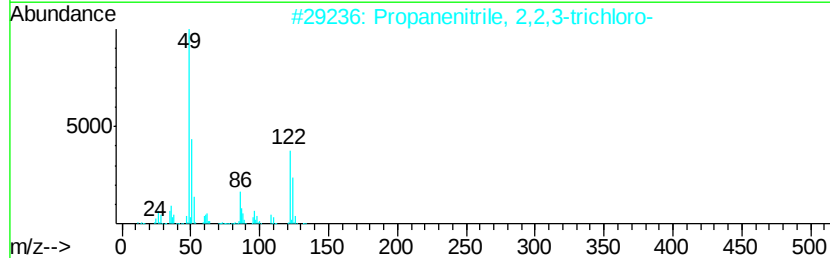
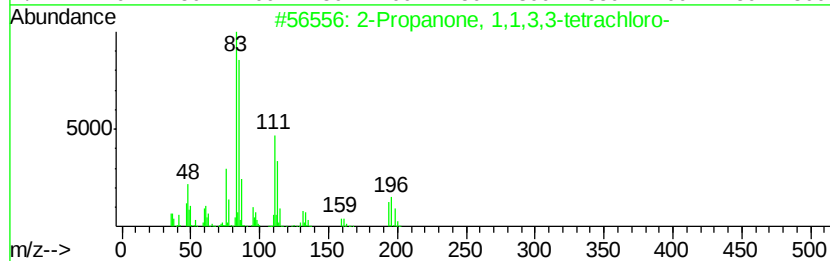
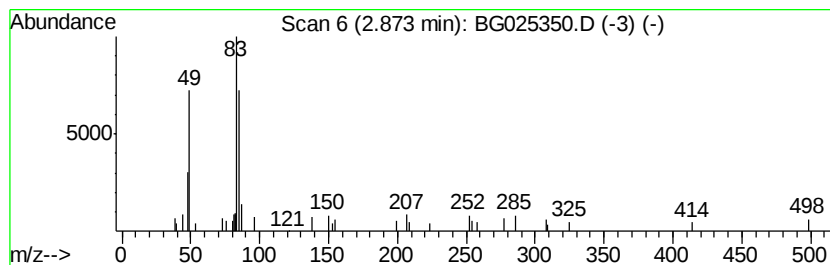
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.87	2.07 ng/ul	63450	1,4-Dichlorobenzene-d4	8.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanone, 1,1,3,3-tetrachloro-	194	C3H2Cl4O	000632-21-3	12		
2	Propanenitrile, 2,2,3-trichloro-	157	C3H2Cl3N	000813-74-1	9		
3	2-Propanone, 1,1,1,3,3-pentachloro-	228	C3HCl5O	001768-31-6	9		
4	Ethane, 2,2-dichloro-1,1,1-trifl...	152	C2HCl2F3	000306-83-2	9		
5	Bromodichloroacetaldehyde	190	C2HBrCl2O	034619-29-9	9		



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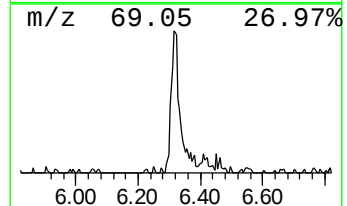
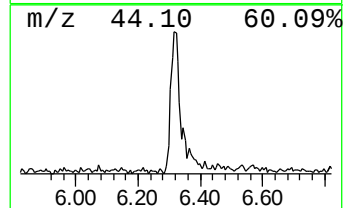
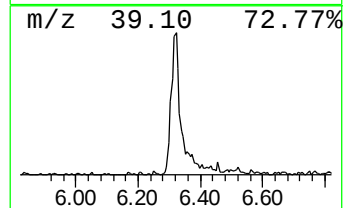
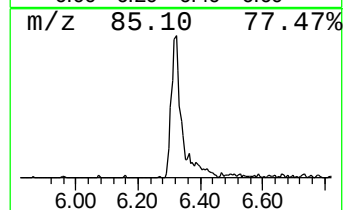
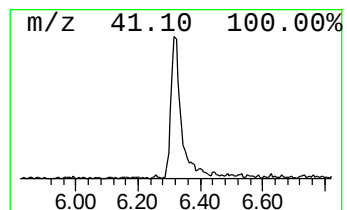
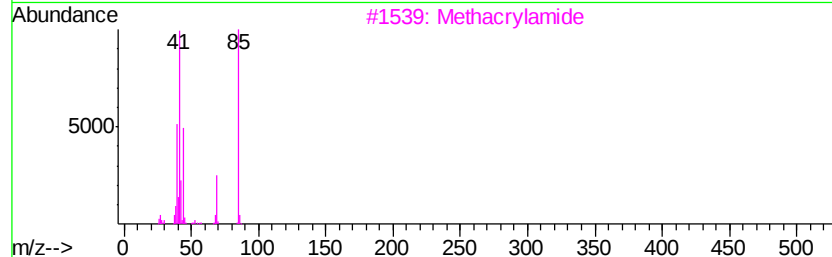
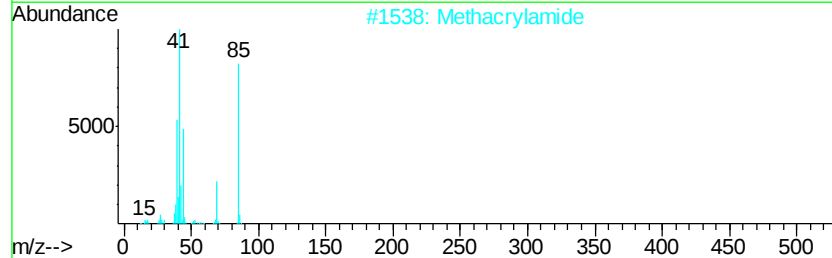
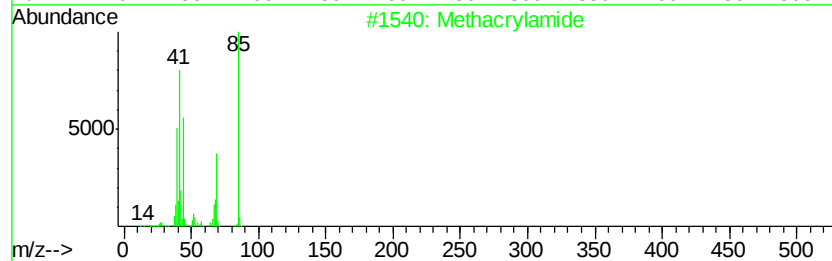
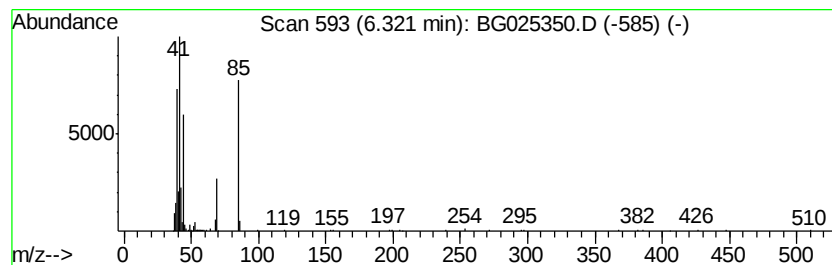
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Methacrylamide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	9.81 ng/ul	301050	1,4-Dichlorobenzene-d4	8.23

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	90
2			Methacrylamide	85	C4H7NO	000079-39-0	90
3			Methacrylamide	85	C4H7NO	000079-39-0	64
4			trans-Crotonamide	85	C4H7NO	000625-37-6	59
5			E-2-Butenoic acid, 3-iodo-	212	C4H5IO2	1000308-87-1	42



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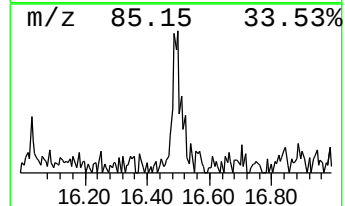
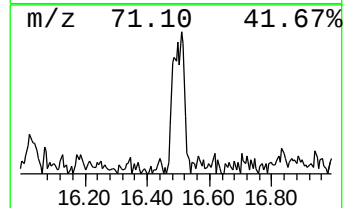
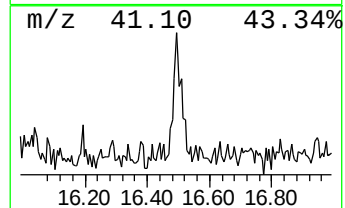
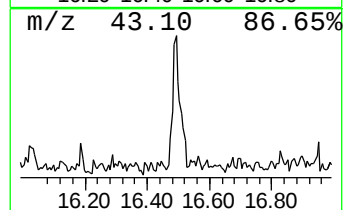
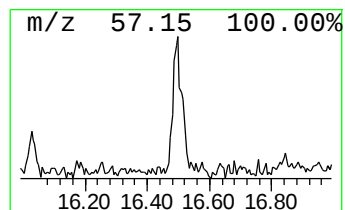
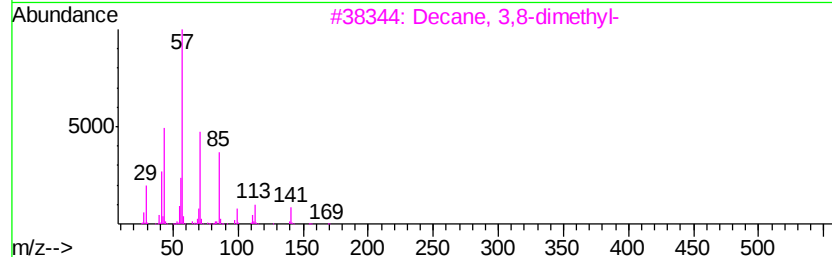
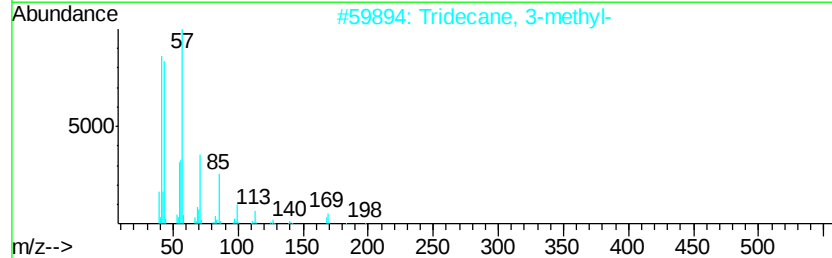
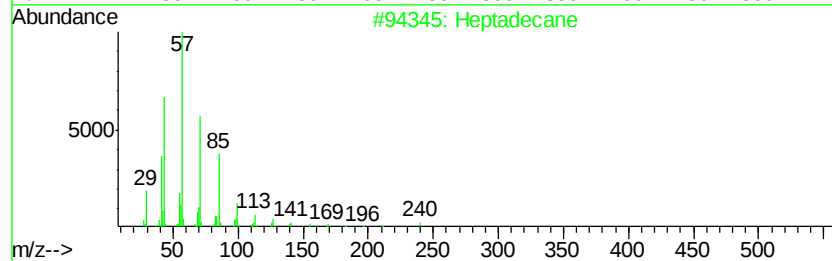
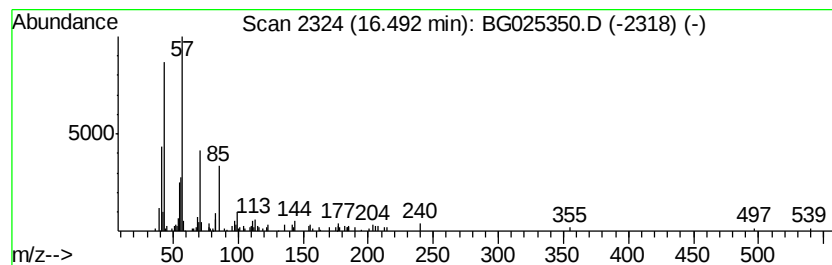
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.49	2.38 ng/ul	147393	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	74
2		Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
3		Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
4		Heptadecane	240	C17H36	000629-78-7	70
5		Octadecane	254	C18H38	000593-45-3	64



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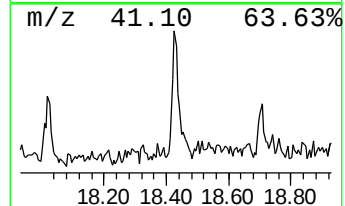
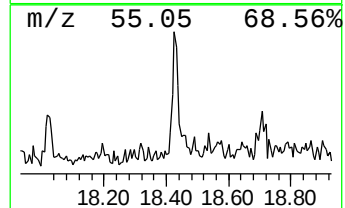
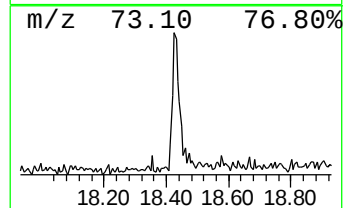
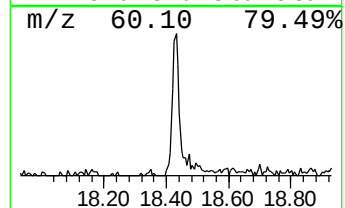
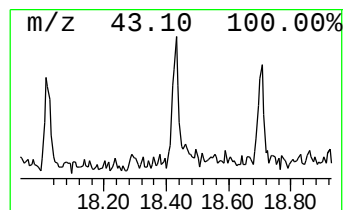
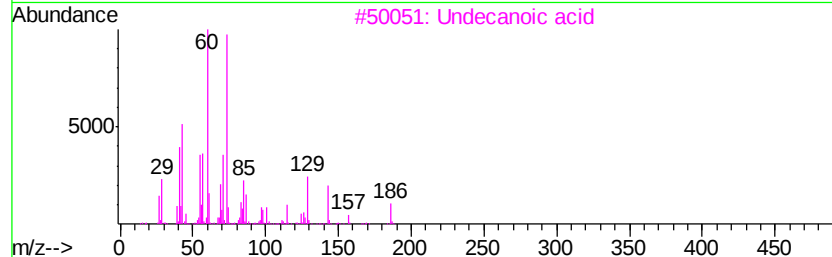
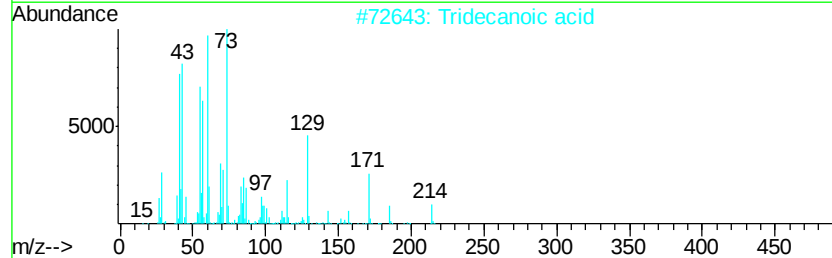
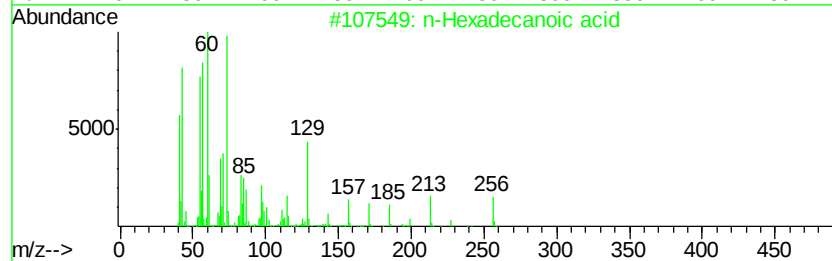
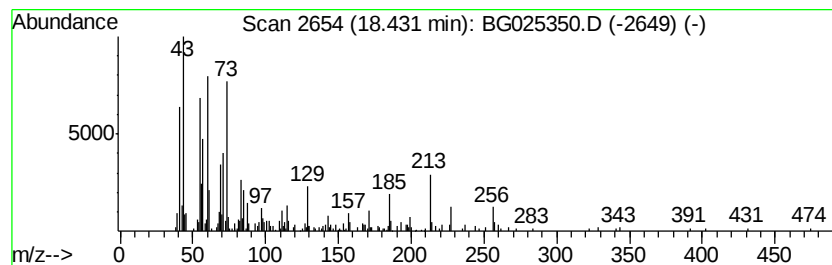
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.43	3.97 ng/ul	245713	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2		Tridecanoic acid	214	C13H26O2	000638-53-9	83
3		Undecanoic acid	186	C11H22O2	000112-37-8	81
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	74



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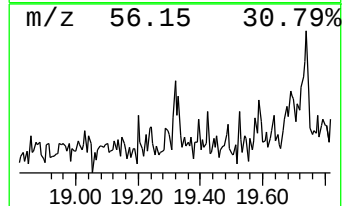
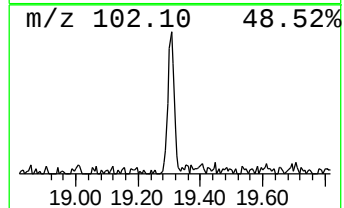
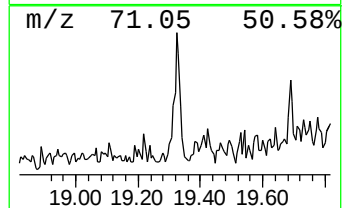
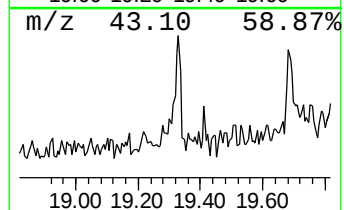
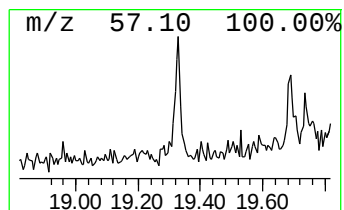
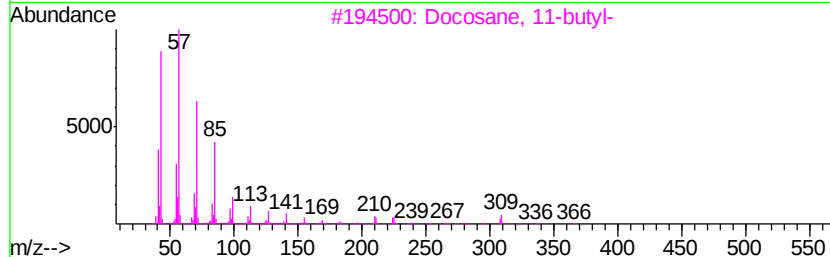
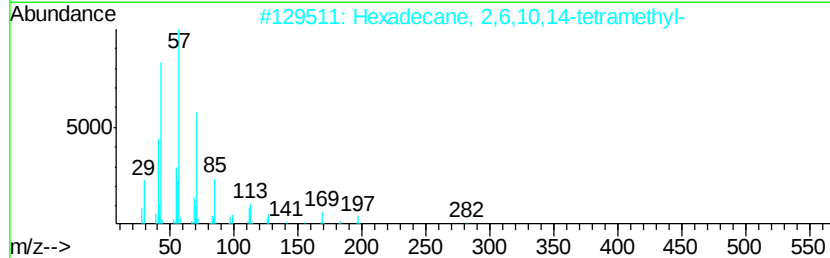
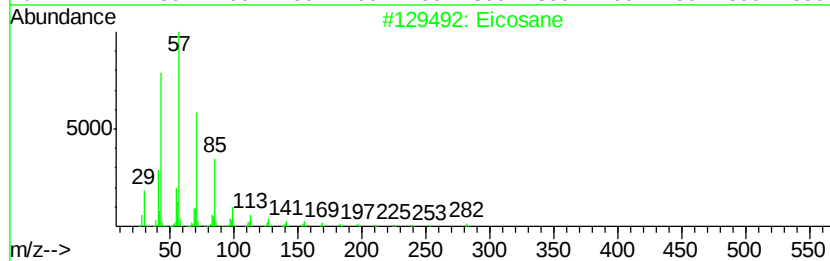
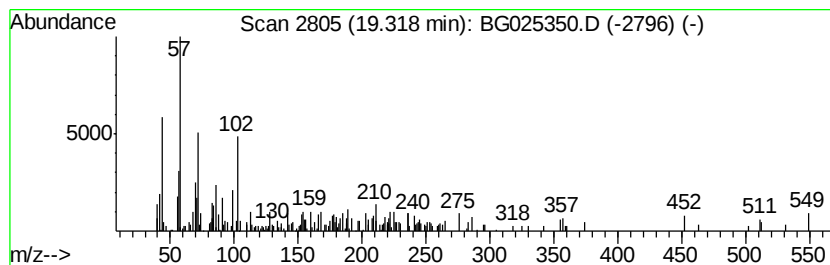
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.32	2.58 ng/ul	159699	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	55
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	55
3		Docosane, 11-butyl-	366	C26H54	013475-76-8	46
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	46
5		Heptacosane	380	C27H56	000593-49-7	46



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
Data File : BG025350.D
Acq On : 20 Dec 2016 12:10
Operator : UM/SJ
Sample : H5905-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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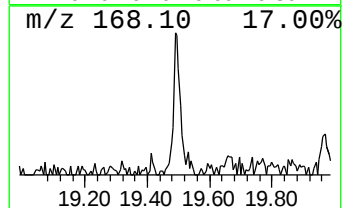
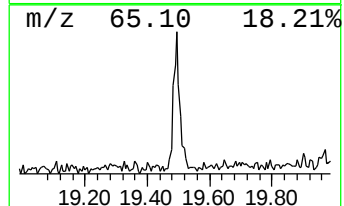
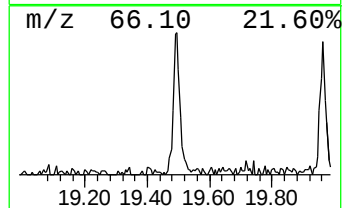
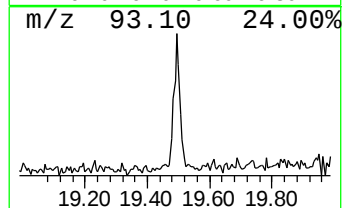
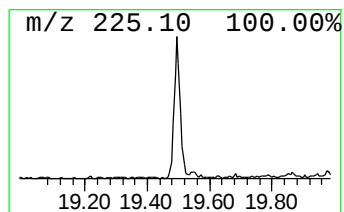
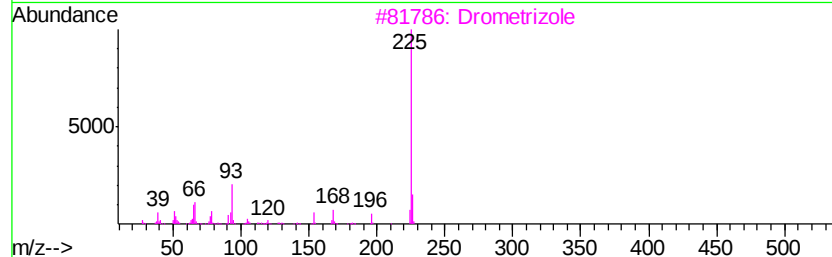
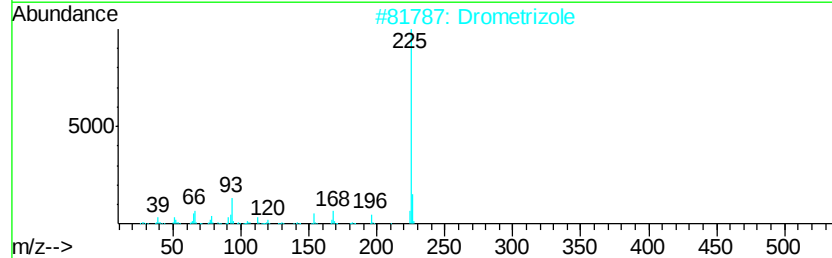
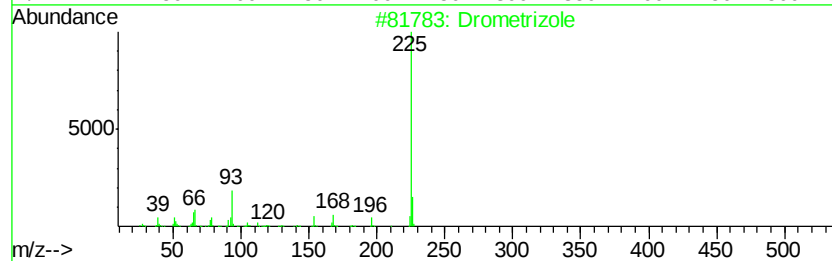
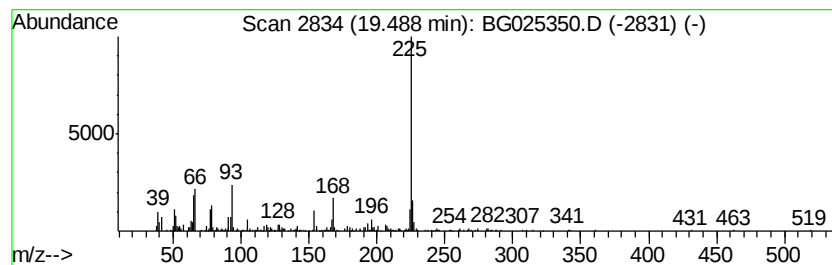
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Drometrizole Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.49	4.86 ng/ul	300590	Phenanthrene-d10	17.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Drometrizole	225	C13H11N3O	002440-22-4	95
2		Drometrizole	225	C13H11N3O	002440-22-4	94
3		Drometrizole	225	C13H11N3O	002440-22-4	94
4		2-(2H-Benzotriazol-2-yl)-5-methy...	225	C13H11N3O	004998-48-5	94
5		Drometrizole	225	C13H11N3O	002440-22-4	93



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
Data File : BG025350.D
Acq On : 20 Dec 2016 12:10
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Sample : H5905-11
Misc :
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Instrument :
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ClientSampleId :
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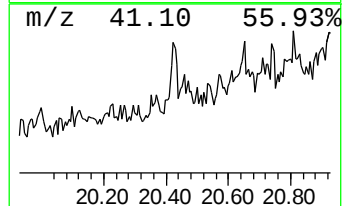
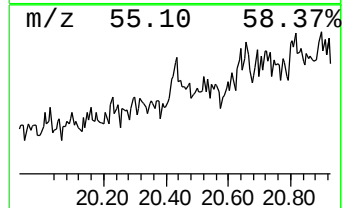
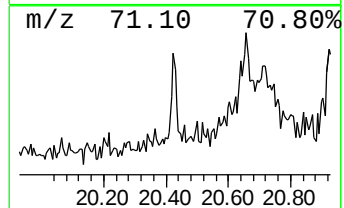
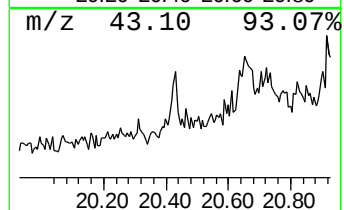
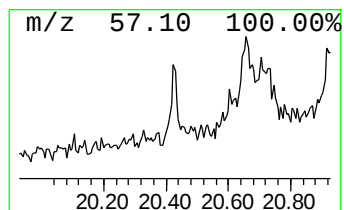
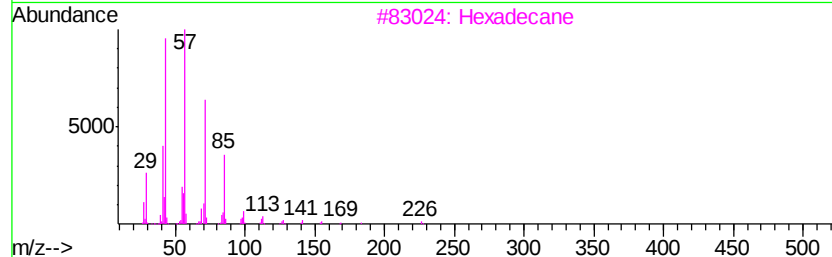
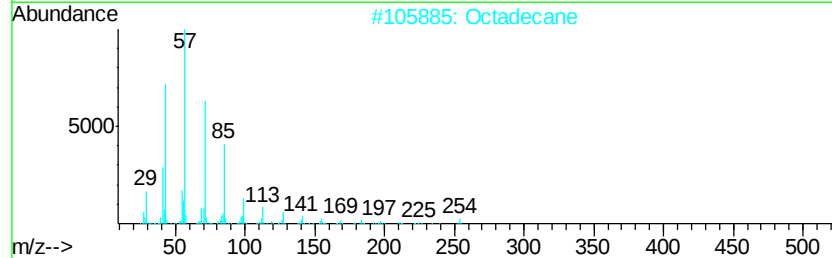
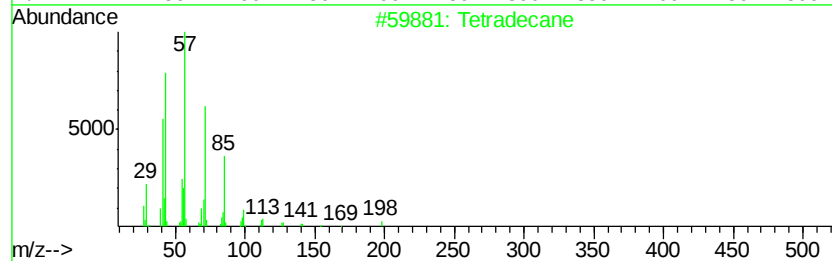
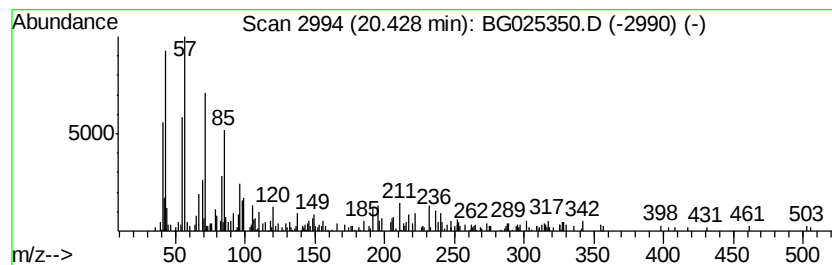
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.43	2.53 ng/ul	191034	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	45
2		Octadecane	254	C18H38	000593-45-3	43
3		Hexadecane	226	C16H34	000544-76-3	43
4		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	43
5		Nonacosane	408	C29H60	000630-03-5	43



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
Data File : BG025350.D
Acq On : 20 Dec 2016 12:10
Operator : UM/SJ
Sample : H5905-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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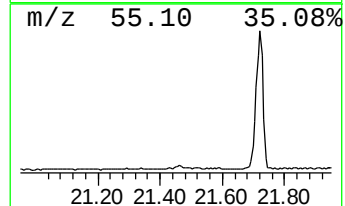
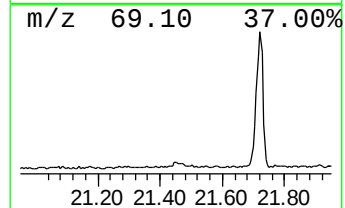
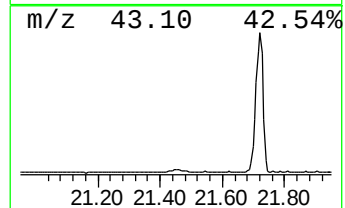
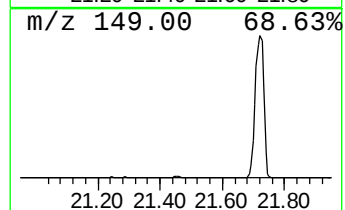
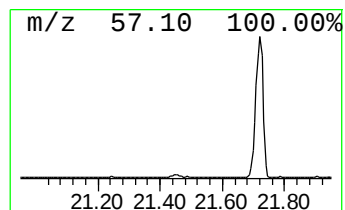
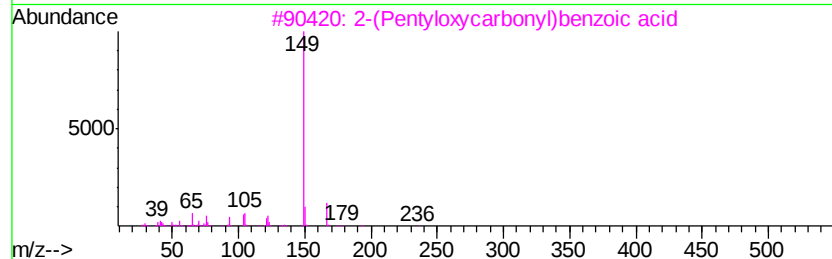
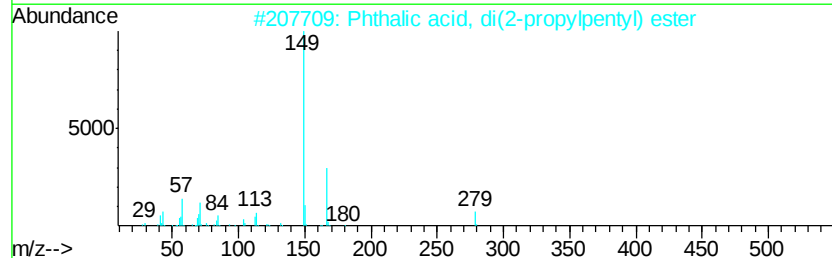
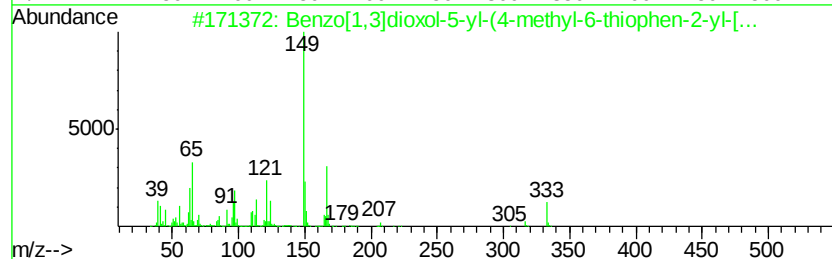
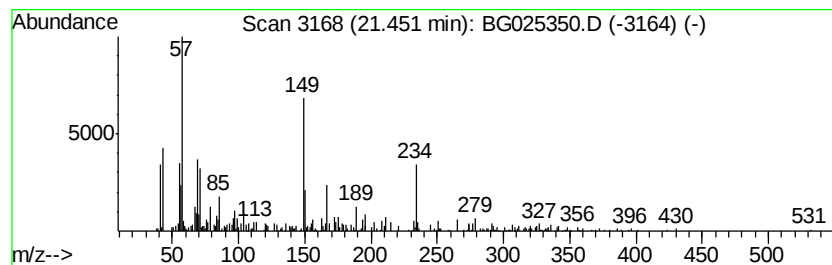
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.45	5.23 ng/ul	394752	Chrysene-d12	21.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[1,3]dioxol-5-yl-(4-methyl-...	333	C16H15N05S	1000300-13-0	43
2		Phthalic acid, di(2-propylpentyl)...	390	C24H38O4	1000377-93-5	43
3		2-(Pentyloxycarbonyl)benzoic acid	236	C13H16O4	1000373-49-7	38
4		2-((3,5,5-Trimethylhexyloxy)carb...	292	C17H24O4	1000373-83-4	35
5		Phthalic acid, heptyl oct-3-yl e...	376	C23H36O4	1000377-71-6	27



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
Data File : BG025350.D
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Sample : H5905-11
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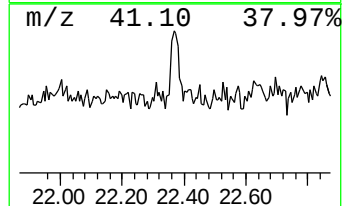
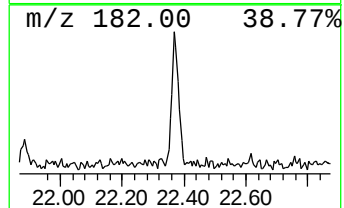
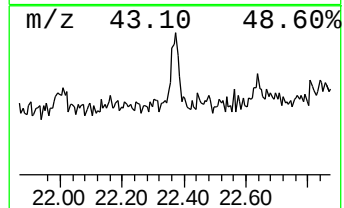
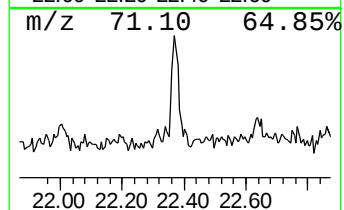
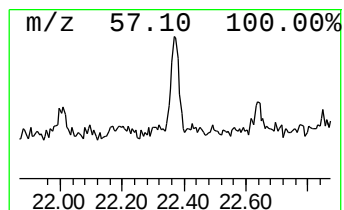
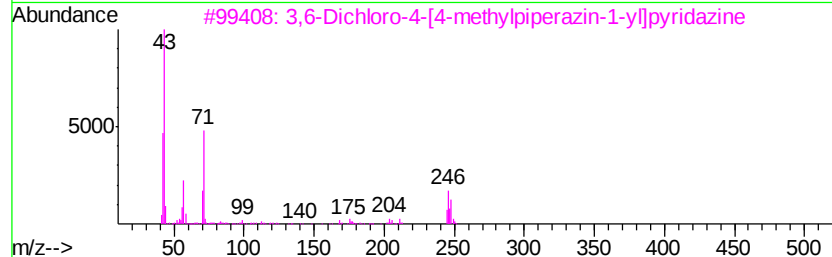
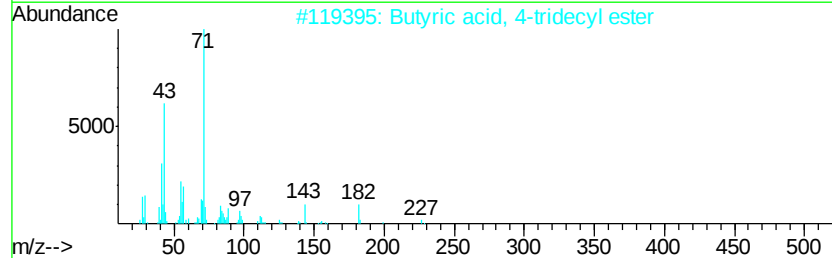
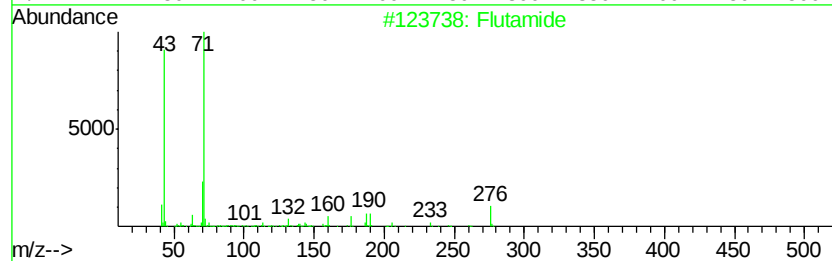
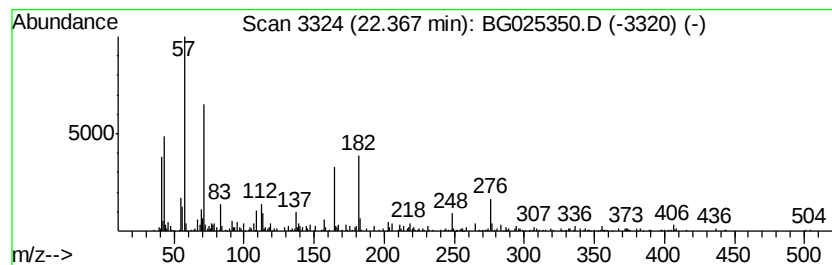
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	4.63 ng/ul	349175	Chrysene-d12	21.89
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Flutamide	276 C11H11F3N2O3	013311-84-7	27
2	Butyric acid, 4-tridecyl ester	270 C17H34O2	1000280-56-9	10
3	3,6-Dichloro-4-[4-methylpiperazi...	246 C9H12Cl2N4	1000255-94-7	10
4	n-Propyl sec-butyl disulfide	164 C7H16S2	059849-54-6	9
5	Disulfide, ethyl isopentyl	164 C7H16S2	072437-60-6	9



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
Data File : BG025350.D
Acq On : 20 Dec 2016 12:10
Operator : UM/SJ
Sample : H5905-11
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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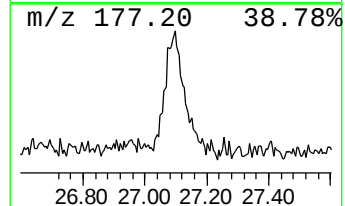
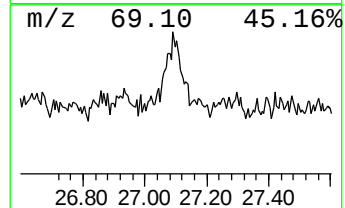
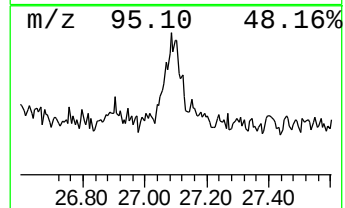
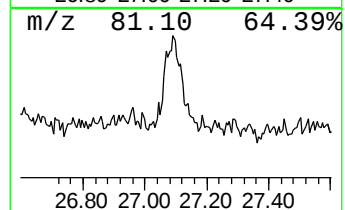
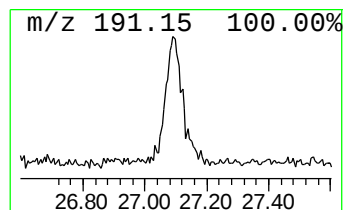
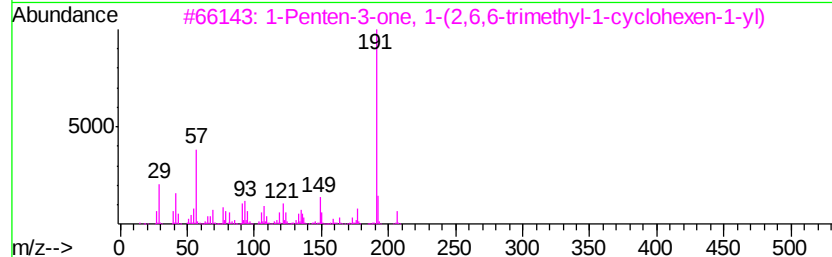
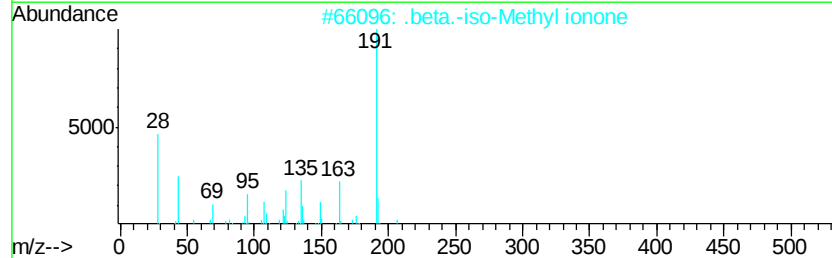
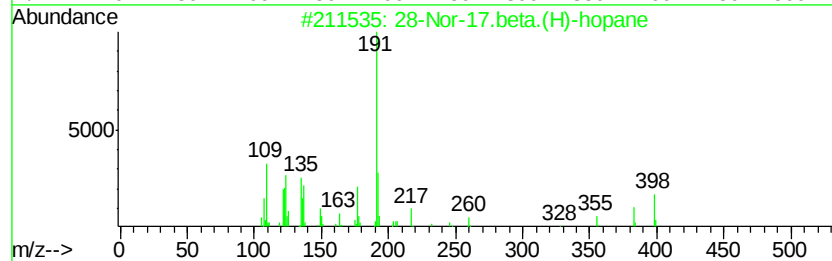
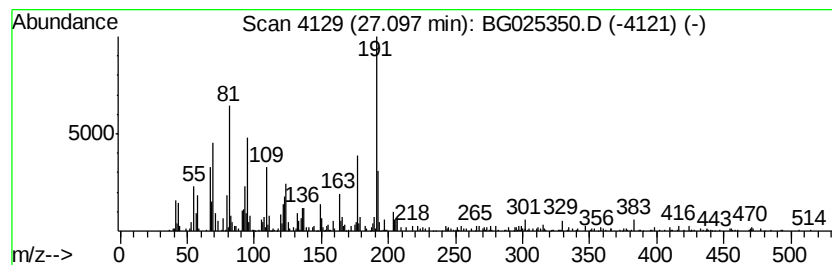
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 28-Nor-17.beta.(H)-hopane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.10	18.87 ng/ul	1256380	Perylene-d12	25.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	68
2		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	62
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	42
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
5		5-Fluoro-2-trifluoromethylbenzoi...	376	C20H28F4O2	1000338-54-5	35



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
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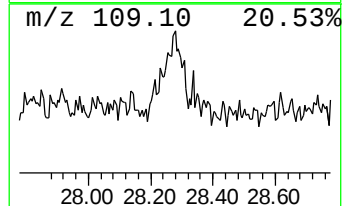
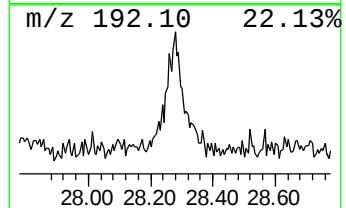
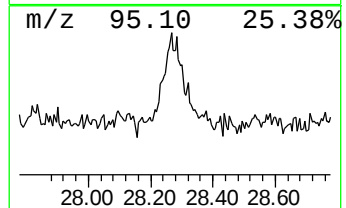
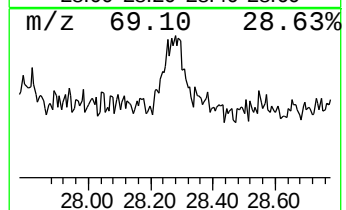
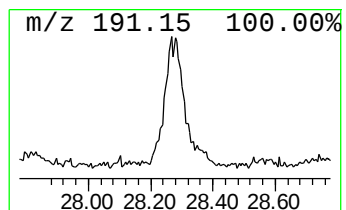
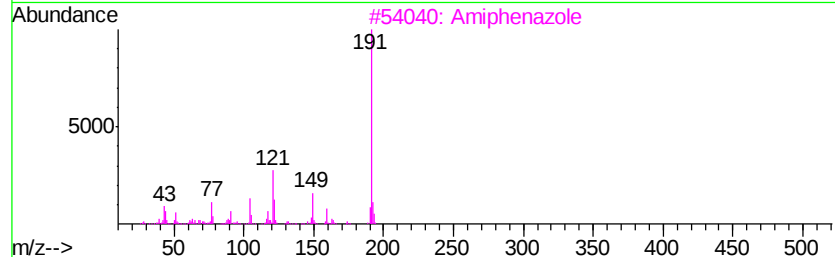
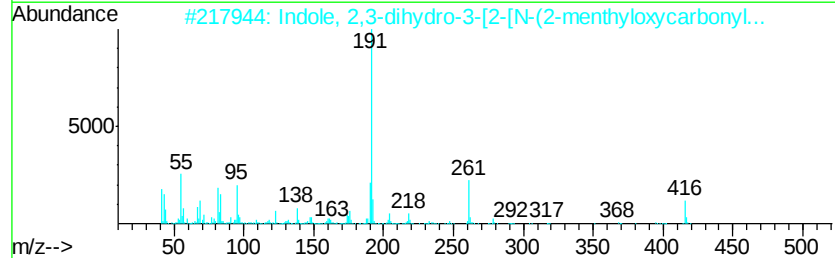
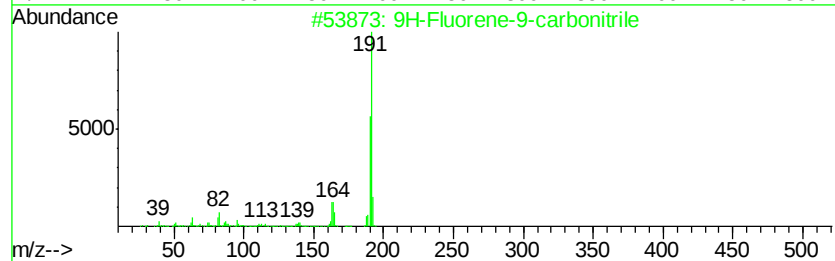
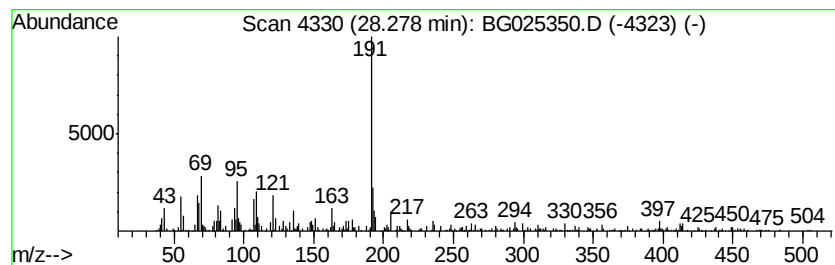
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 9H-Fluorene-9-carbonitrile Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
28.28	14.86 ng/ul	989169	Perylene-d12	25.31		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	59	
2	Indole, 2,3-dihydro-3-[2-[N-(2-m...	416	C25H40N2O3	1000124-26-9	50	
3	Amiphenazole	191	C9H9N3S	000490-55-1	50	
4	4H-1,2,4-Triazole, 3-ethyl-5-(3,...	191	C9H13N5	328532-29-2	49	
5	Isophthalic acid, 2-methyloct-5-...	330	C20H26O4	1000343-92-3	49	



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG121916\
 Data File : BG025350.D
 Acq On : 20 Dec 2016 12:10
 Operator : UM/SJ
 Sample : H5905-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DA811

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
unknown-01	2.87	2.1	ng/ul	63450	1	8.23	613934	20.0
Methacrylamide	6.32	9.8	ng/ul	301050	1	8.23	613934	20.0
(DEL) Alkane: Str...	16.49	2.4	ng/ul	147393	4	17.59	1237970	20.0
n-Hexadecanoic acid	18.43	4.0	ng/ul	245713	4	17.59	1237970	20.0
(DEL) Alkane: Str...	19.32	2.6	ng/ul	159699	4	17.59	1237970	20.0
Drometrizole	19.49	4.9	ng/ul	300590	4	17.59	1237970	20.0
(DEL) Alkane: Str...	20.43	2.5	ng/ul	191034	5	21.89	1508530	20.0
unknown-02	21.45	5.2	ng/ul	394752	5	21.89	1508530	20.0
unknown-03	22.37	4.6	ng/ul	349175	5	21.89	1508530	20.0
28-Nor-17.beta.(H...	27.10	18.9	ng/ul	1256380	6	25.31	1331300	20.0
9H-Fluorene-9-car...	28.28	14.9	ng/ul	989169	6	25.31	1331300	20.0