

Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
 Data File : BM004246.D
 Acq On : 13 Feb 2016 02:38
 Operator : SJ/IZ
 Sample : H1414-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9QD2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM021216.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	2.769	9	14	23	rBV	104828	169991	29.88%	2.567%
2	2.846	23	27	34	rVB	115339	135197	23.76%	2.042%
3	3.110	69	72	76	rVB	2419	2577	0.45%	0.039%
4	3.634	157	161	163	rVV	2039	1998	0.35%	0.030%
5	3.845	195	197	201	rVB	1036	983	0.17%	0.015%
6	4.751	348	351	356	rBB2	1298	1844	0.32%	0.028%
7	6.822	698	703	712	rBB	52482	80582	14.16%	1.217%
8	6.981	725	730	737	rBB	46035	67821	11.92%	1.024%
9	7.181	758	764	772	rBB	59451	92039	16.18%	1.390%
10	7.245	773	775	780	rBB	971	1246	0.22%	0.019%
11	7.645	837	843	849	rBB	103308	157427	27.67%	2.377%
12	8.357	959	964	973	rBB	70771	109646	19.27%	1.656%
13	8.804	1034	1040	1047	rBB	51988	81124	14.26%	1.225%
14	8.939	1060	1063	1068	rBB	1234	2021	0.36%	0.031%
15	9.522	1157	1162	1168	rBB	43383	65638	11.54%	0.991%
16	10.063	1249	1254	1263	rBB	67263	107577	18.91%	1.625%
17	10.433	1310	1317	1325	rBB	143335	237401	41.72%	3.585%
18	10.580	1336	1342	1354	rBB	29117	48178	8.47%	0.728%
19	13.192	1784	1786	1789	rBB	930	928	0.16%	0.014%
20	13.715	1869	1875	1879	rBV	126346	165831	29.15%	2.504%
21	13.762	1879	1883	1889	rVB	60412	80012	14.06%	1.208%
22	13.992	1917	1922	1932	rBB	142665	209838	36.88%	3.169%
23	14.204	1955	1958	1961	rBB	4212	4713	0.83%	0.071%
24	14.304	1970	1975	1982	rBB2	194000	289954	50.96%	4.379%
25	14.533	2010	2014	2025	rBB	22713	39379	6.92%	0.595%
26	15.104	2109	2111	2114	rBB	1222	1220	0.21%	0.018%
27	15.157	2117	2120	2124	rBB	6524	7645	1.34%	0.115%
28	15.304	2139	2145	2151	rBB	186057	259197	45.56%	3.914%
29	15.439	2164	2168	2174	rBB	34714	40483	7.12%	0.611%
30	15.545	2184	2186	2188	rBV	1538	1595	0.28%	0.024%
31	15.562	2188	2189	2192	rVB	1308	1082	0.19%	0.016%
32	16.021	2264	2267	2274	rBB2	8180	12121	2.13%	0.183%
33	16.209	2296	2299	2301	rBB	749	604	0.11%	0.009%
34	16.815	2399	2402	2405	rBB	4917	4916	0.86%	0.074%

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35	16.845	2405	2407	2409	rBV	512	590	0.10%	0.009%
36	16.868	2409	2411	2413	rVB	785	655	0.12%	0.010%
37	17.056	2437	2443	2448	rBV	229768	311811	54.80%	4.709%
38	17.098	2448	2450	2454	rVB	3480	3886	0.68%	0.059%
39	17.156	2454	2460	2469	rBB	177905	251949	44.28%	3.805%
40	17.556	2525	2528	2530	rBB	2185	2297	0.40%	0.035%
41	17.950	2593	2595	2599	rBB	6973	7412	1.30%	0.112%
42	18.245	2644	2645	2648	rBB	1498	1066	0.19%	0.016%
43	19.097	2787	2790	2791	rBV	1379	1384	0.24%	0.021%
44	19.127	2792	2795	2799	rVB	6569	7132	1.25%	0.108%
45	19.244	2810	2815	2819	rBV2	3285	4886	0.86%	0.074%
46	19.350	2828	2833	2836	rBV	1316	1982	0.35%	0.030%
47	19.462	2846	2852	2865	rVB2	223478	298069	52.39%	4.501%
48	19.862	2918	2920	2923	rVB	714	610	0.11%	0.009%
49	19.980	2935	2940	2943	rBV3	2338	4216	0.74%	0.064%
50	20.009	2943	2945	2948	rVB2	2553	2596	0.46%	0.039%
51	20.068	2953	2955	2959	rVB2	1334	1574	0.28%	0.024%
52	20.109	2959	2962	2963	rBV	1309	1042	0.18%	0.016%
53	20.133	2963	2966	2967	rBV	912	638	0.11%	0.010%
54	20.192	2973	2976	2978	rVB	999	1057	0.19%	0.016%
55	20.221	2978	2981	2984	rBV3	1321	2323	0.41%	0.035%
56	20.256	2984	2987	2990	rBV	2173	2735	0.48%	0.041%
57	20.297	2992	2994	3000	rBV3	3618	4713	0.83%	0.071%
58	20.391	3008	3010	3013	rVB3	1938	1815	0.32%	0.027%
59	20.444	3013	3019	3024	rBV	27303	32231	5.66%	0.487%
60	20.509	3026	3030	3034	rVB4	2594	3443	0.61%	0.052%
61	20.544	3034	3036	3039	rVB3	1014	904	0.16%	0.014%
62	20.568	3039	3040	3042	rBV	1143	748	0.13%	0.011%
63	20.586	3042	3043	3045	rBV	1506	706	0.12%	0.011%
64	20.650	3052	3054	3055	rBV	809	761	0.13%	0.011%
65	20.715	3063	3065	3068	rBV4	1322	1194	0.21%	0.018%
66	20.756	3068	3072	3075	rBV4	1246	1945	0.34%	0.029%
67	20.791	3075	3078	3079	rBV3	1566	1366	0.24%	0.021%
68	20.821	3081	3083	3085	rBV2	1038	909	0.16%	0.014%
69	20.868	3089	3091	3094	rVB4	1261	1550	0.27%	0.023%
70	20.897	3094	3096	3097	rBV2	1714	1202	0.21%	0.018%
71	20.921	3097	3100	3103	rBV3	1682	2215	0.39%	0.033%

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72	20.974	3103	3109	3118	rBV	43306	69872	12.28%	1.055%
73	21.191	3145	3146	3149	rVB3	2415	1310	0.23%	0.020%
74	21.268	3153	3159	3164	rBV	291739	378930	66.60%	5.723%
75	21.580	3209	3212	3217	rVB	13947	14526	2.55%	0.219%
76	21.862	3256	3260	3271	rBV	73385	116273	20.44%	1.756%
77	22.297	3332	3334	3338	rVB3	6495	6287	1.10%	0.095%
78	22.427	3353	3356	3360	rVB3	7433	10317	1.81%	0.156%
79	22.797	3416	3419	3423	rBV	12675	15379	2.70%	0.232%
80	22.850	3424	3428	3445	rVB	39046	84465	14.85%	1.276%
81	23.356	3507	3514	3524	rBV	173827	317310	55.77%	4.792%
82	23.497	3531	3538	3545	rVB2	195391	359204	63.13%	5.425%
83	23.991	3618	3622	3629	rVB	17957	30866	5.42%	0.466%
84	24.091	3633	3639	3649	rBV8	7439	21245	3.73%	0.321%
85	24.727	3744	3747	3756	rVB7	10396	22664	3.98%	0.342%
86	26.509	4042	4050	4062	rBV6	23752	70224	12.34%	1.061%
87	27.409	4195	4203	4219	rVB5	59082	212364	37.32%	3.207%
88	27.773	4255	4265	4278	rBV4	132921	517646	90.98%	7.817%
89	27.909	4278	4288	4298	rVV3	150149	568975	100.00%	8.593%
90	28.020	4299	4307	4327	rVB8	64575	297938	52.36%	4.499%
91	28.967	4460	4468	4479	rVB9	15349	57507	10.11%	0.868%

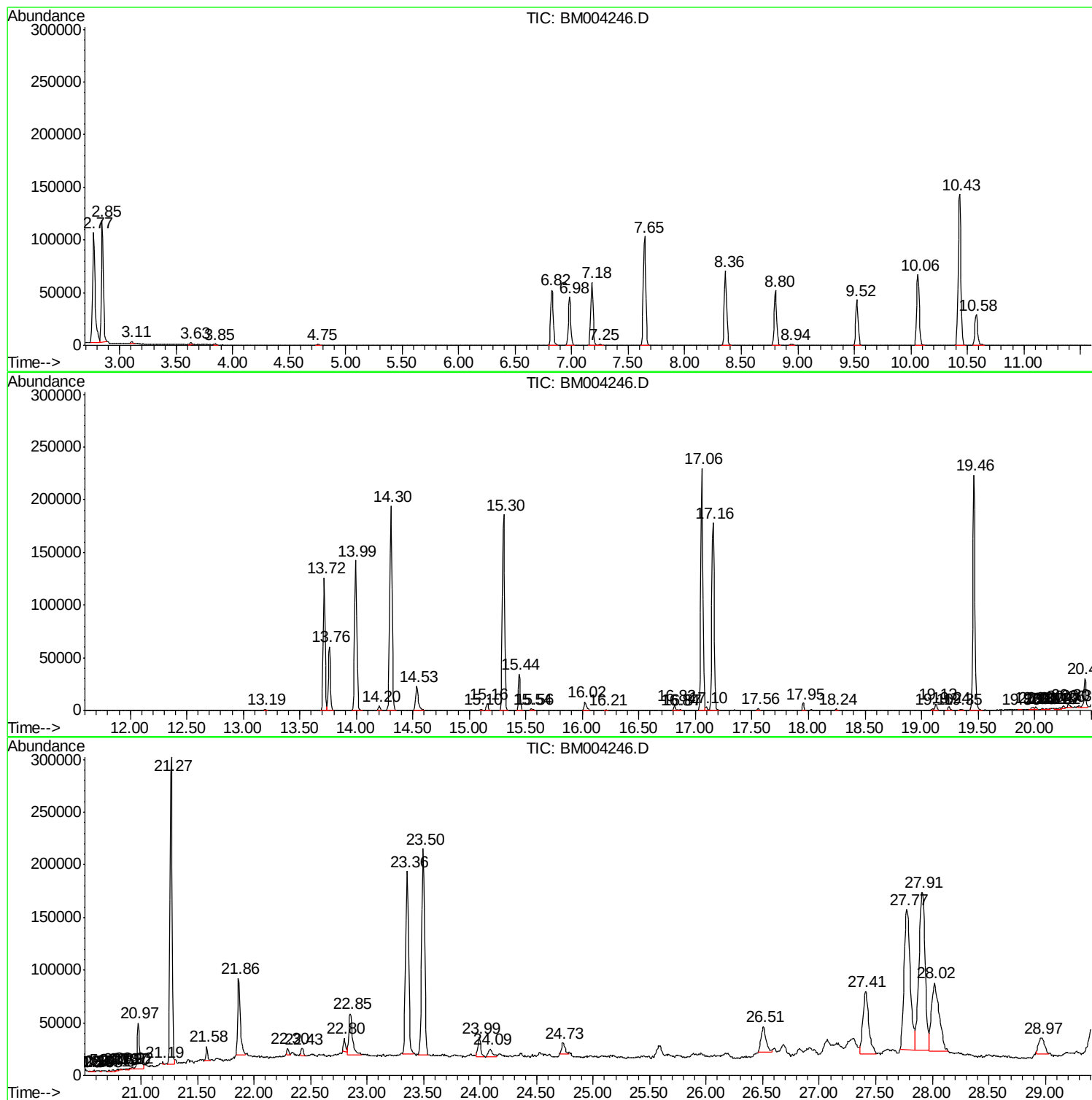
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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



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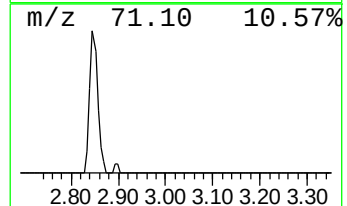
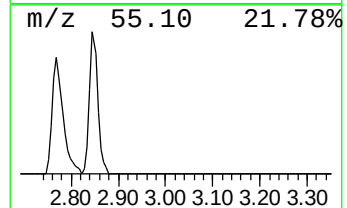
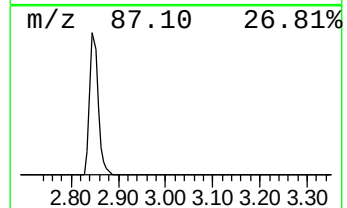
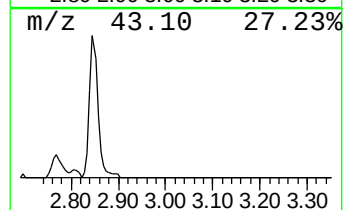
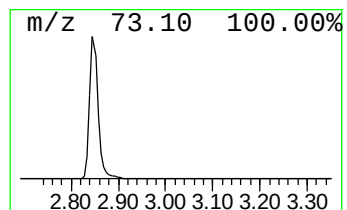
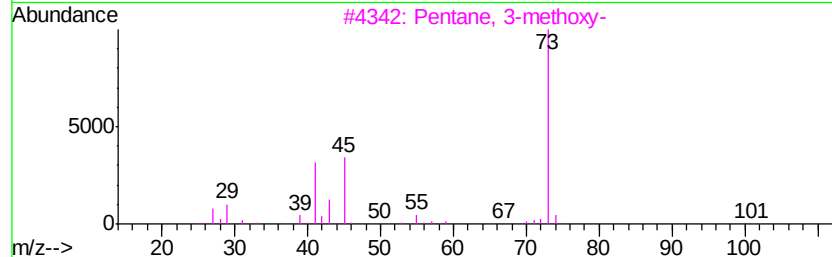
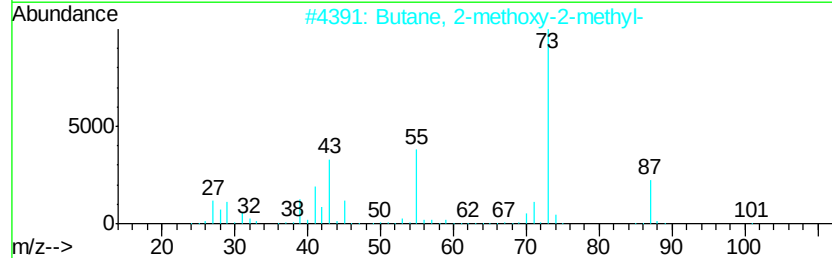
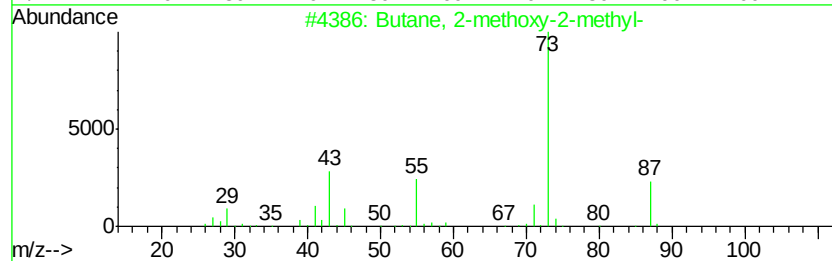
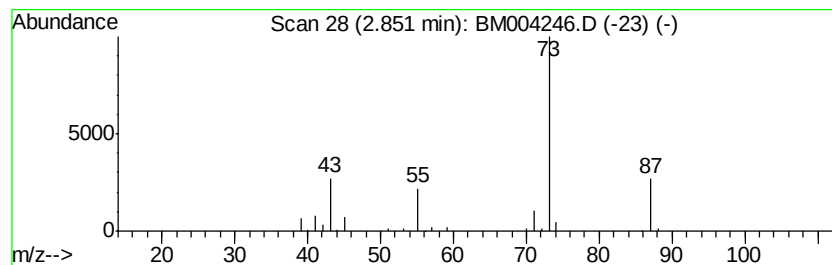
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.85	17.18 ng/ul	135197	1,4-Dichlorobenzene-d4	7.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	90
2		Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9



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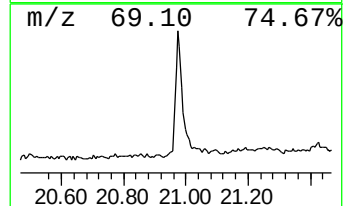
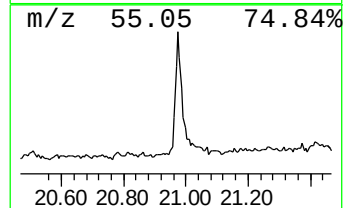
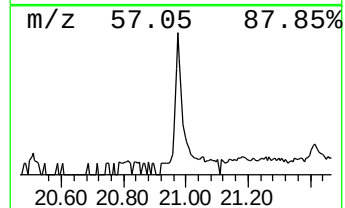
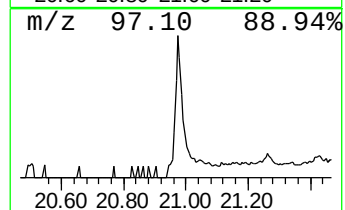
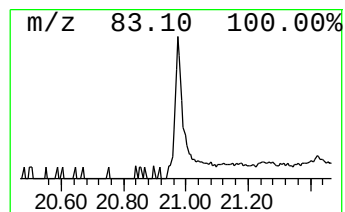
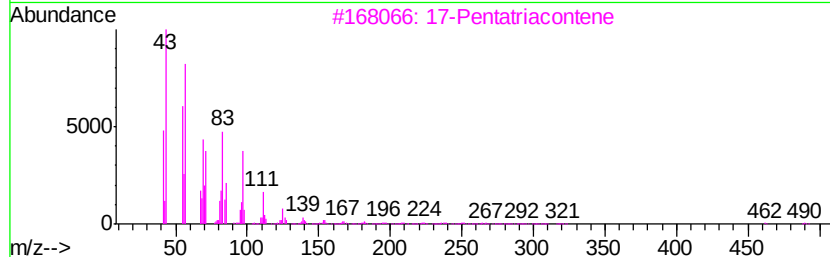
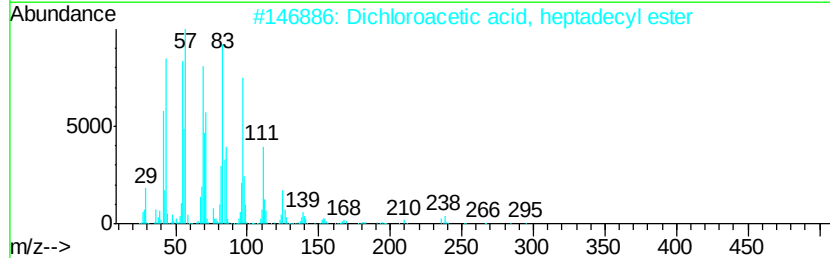
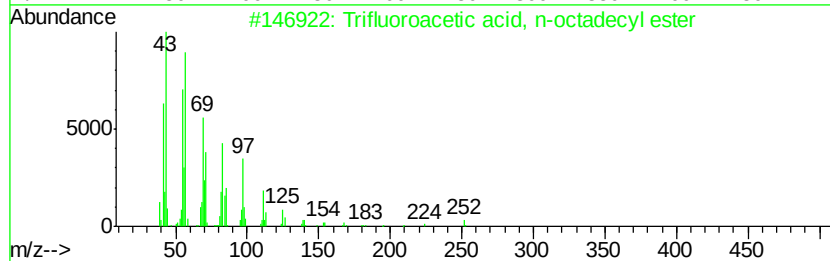
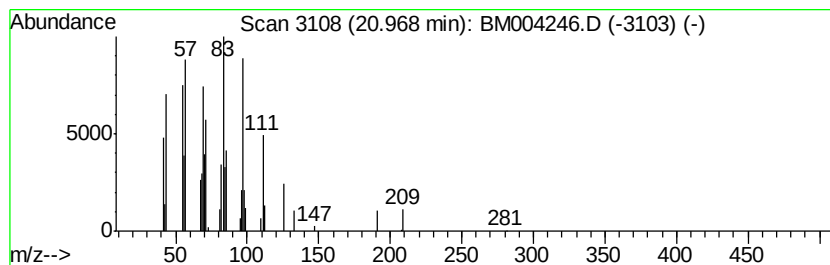
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Trifluoroacetic acid, n-oct... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.97	3.69 ng/ul	69872	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90
5		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	90



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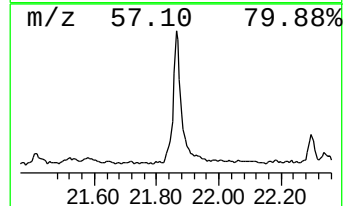
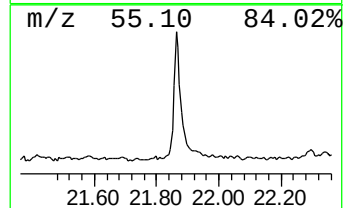
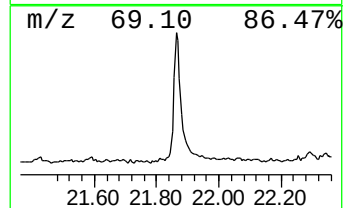
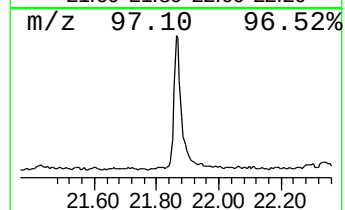
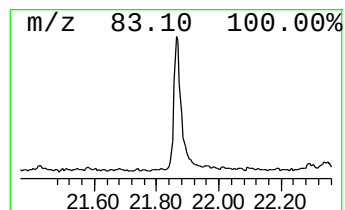
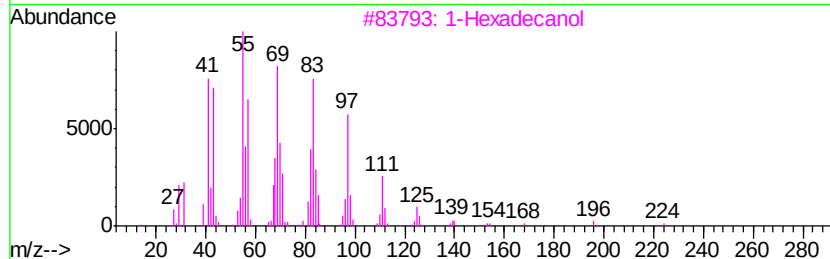
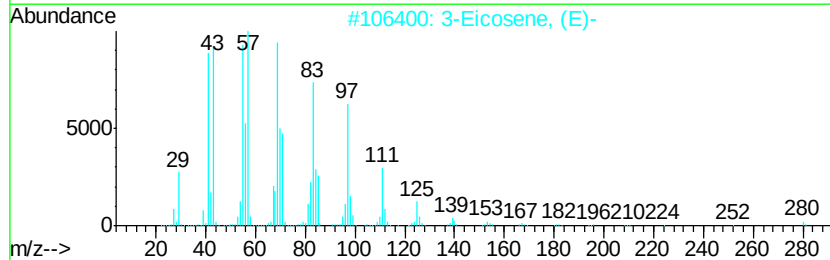
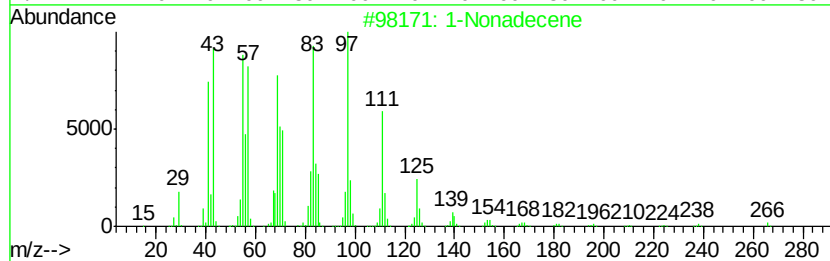
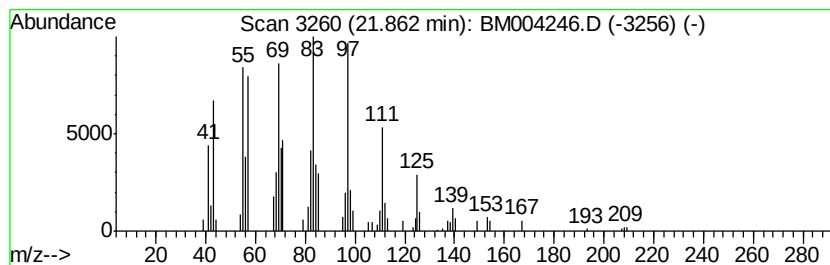
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic21.86 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.86	6.14 ng/ul	116273	Chrysene-d12	21.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Nonadecene	266	C19H38	018435-45-5	95
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
3		1-Hexadecanol	242	C16H34O	036653-82-4	91
4		Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	90
5		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



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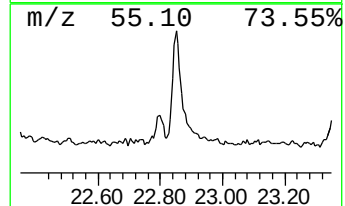
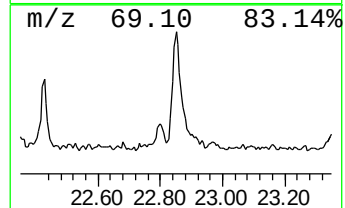
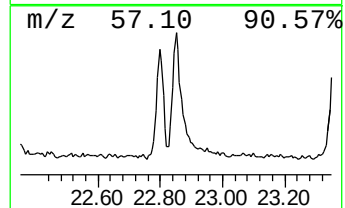
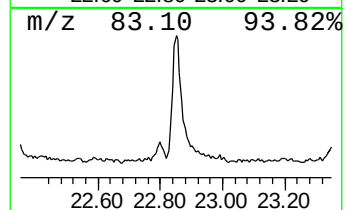
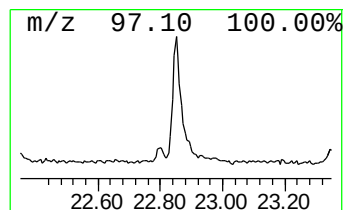
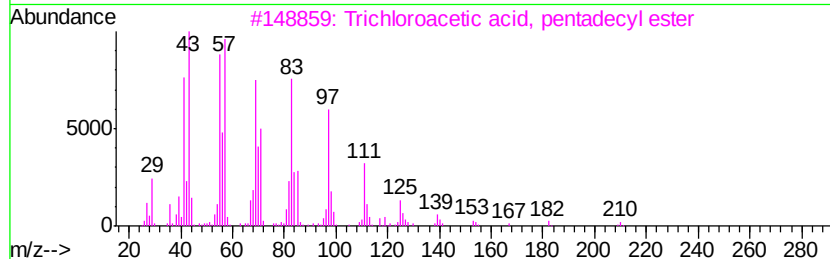
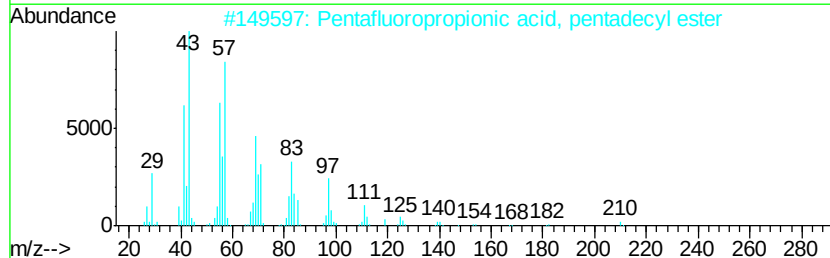
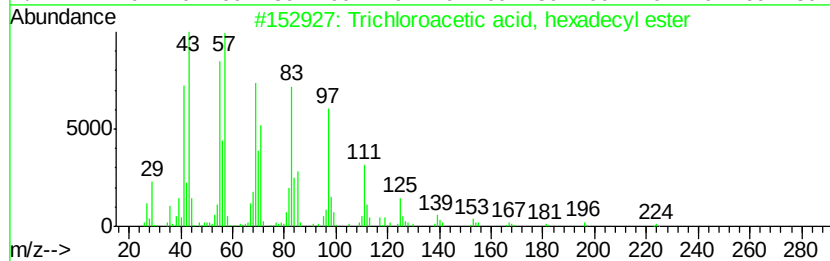
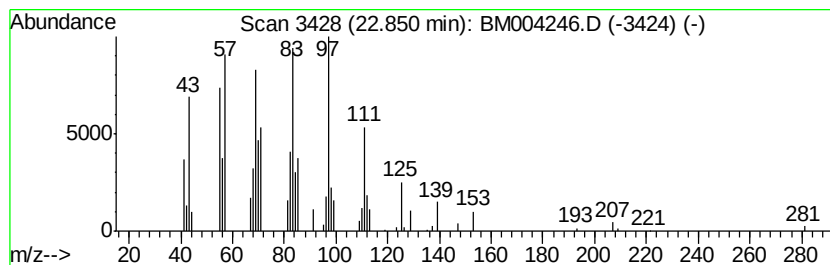
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Trichloroacetic acid, hexad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.85	4.70 ng/ul	84465	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2		Pentafluoropropionic acid, penta...	374	C18H31F5O2	1000280-07-5	91
3		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
4		1-Nonadecanol	284	C19H40O	001454-84-8	91
5		1-Octadecanol	270	C18H38O	000112-92-5	91



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
D9QD2

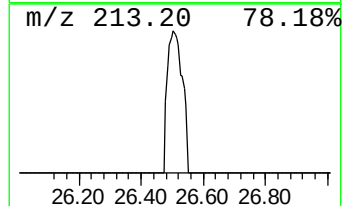
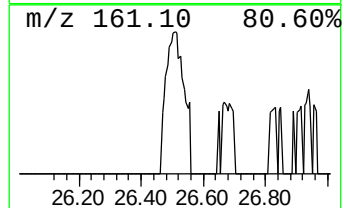
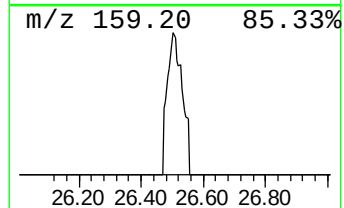
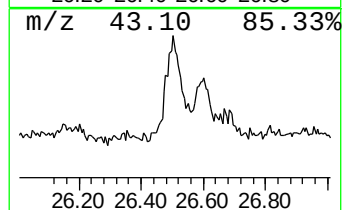
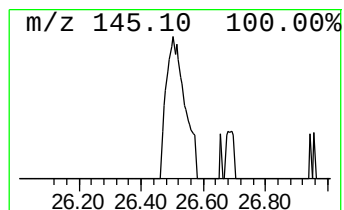
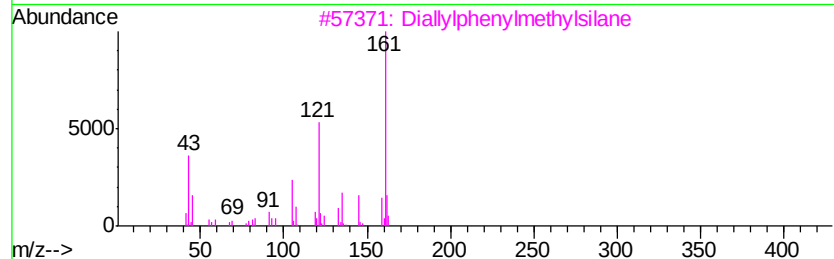
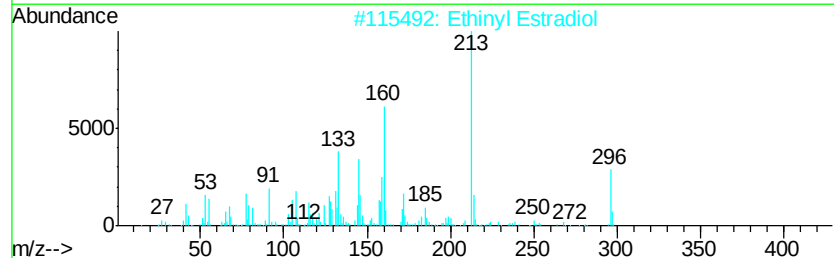
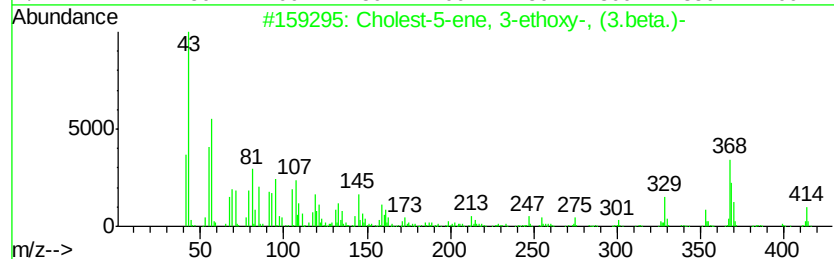
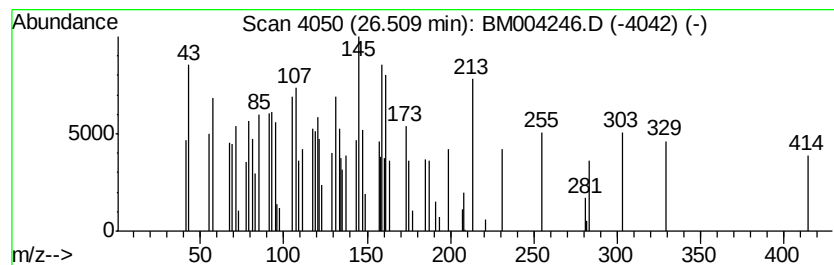
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.51	3.91 ng/ul	70224	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	30
2		Ethinyl Estradiol	296	C20H24O2	000057-63-6	16
3		Diallylphenylmethylsilane	202	C13H18Si	002633-60-5	12
4		Cholan-24-oic acid, 3,12-bis(ace...	490	C29H46O6	001181-44-8	10
5		Ledene oxide-(II)	220	C15H24O	1000159-36-7	10



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampled :
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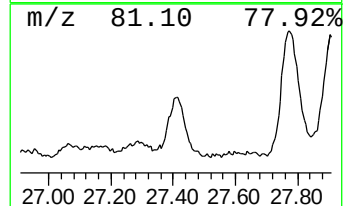
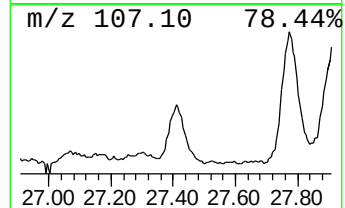
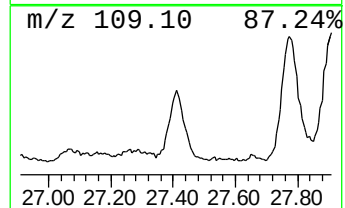
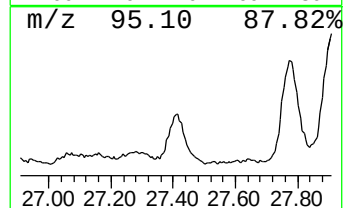
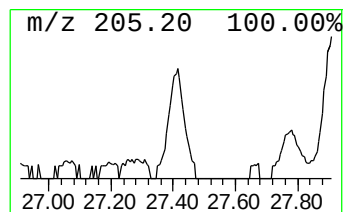
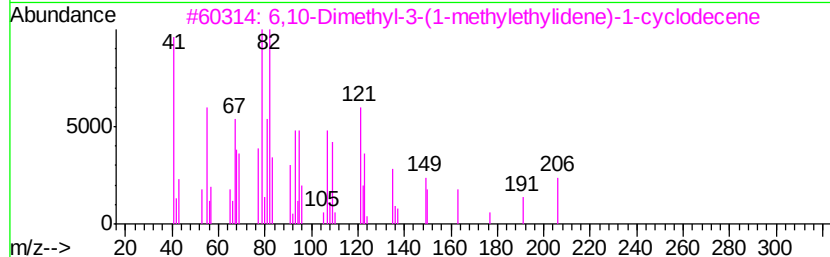
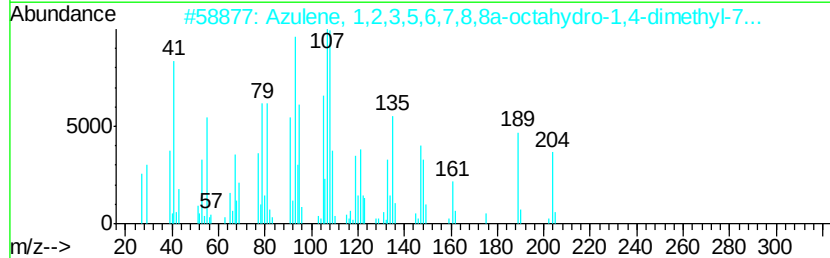
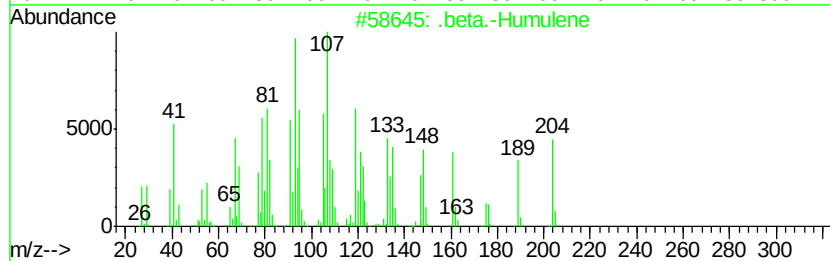
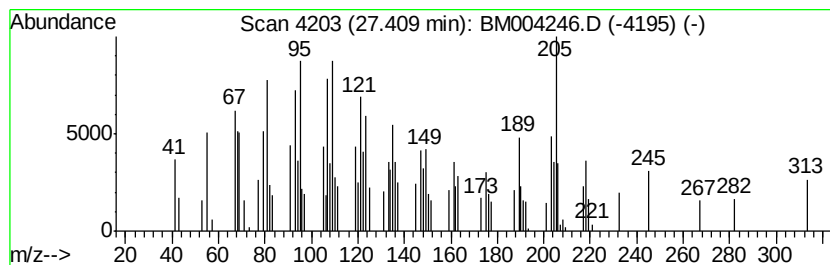
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 .beta.-Humulene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.41	11.82 ng/ul	212364	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Humulene	204	C15H24	000116-04-1	91
2		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	70
3		6,10-Dimethyl-3-(1-methylethylid...	206	C15H26	069239-71-0	53
4		Azulene, 1,2,3,5,6,7,8,8a-octahy...	204	C15H24	003691-11-0	49
5		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	46



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
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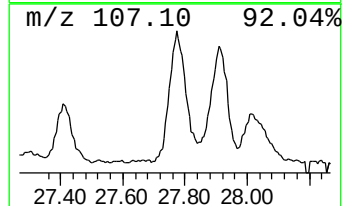
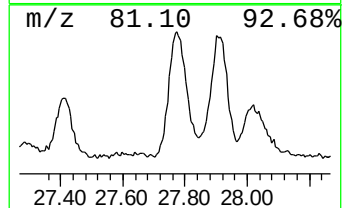
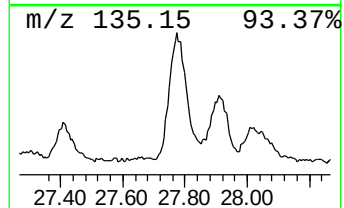
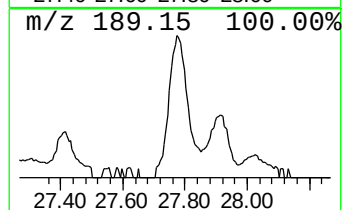
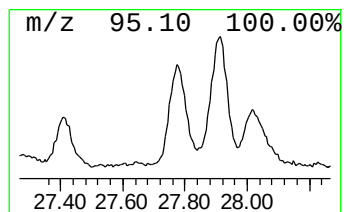
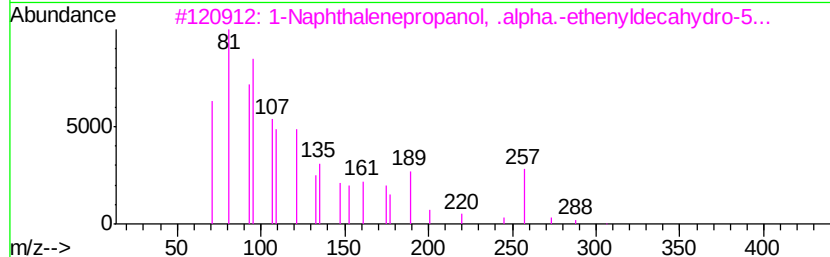
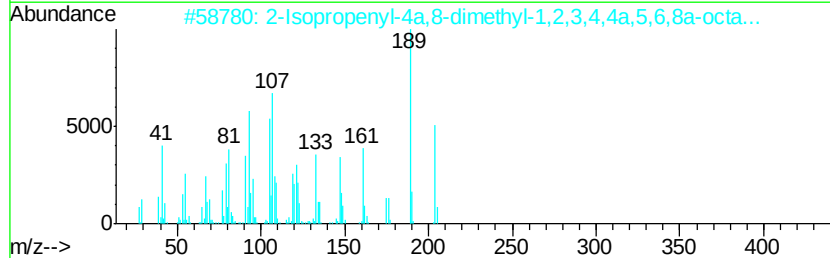
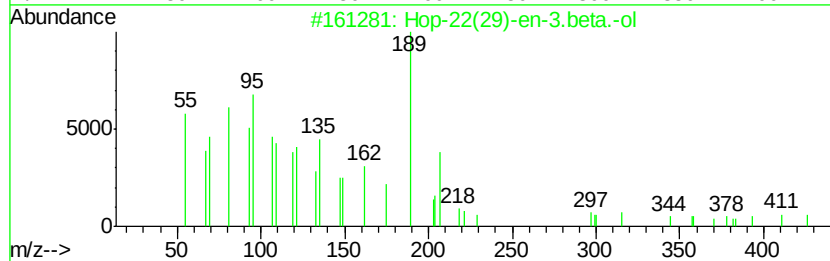
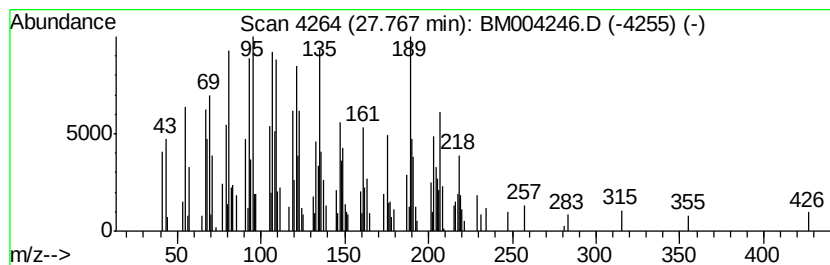
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hop-22(29)-en-3.beta.-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.		
27.77	28.82 ng/ul	517646	Perylene-d12	23.50		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hop-22(29)-en-3.beta.-ol	426	C30H50O	058801-23-3	92
2		2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	1000193-57-0	72
3		1-Naphthalenepropanol, .alpha.-e...	306	C20H34O2	003650-30-4	62
4		Cyclohexane, 1-ethenyl-1-methyl-...	204	C15H24	003242-08-8	46
5		1,5-Cyclodecadiene, 1,5-dimethyl...	204	C15H24	015423-57-1	46



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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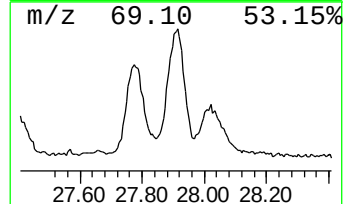
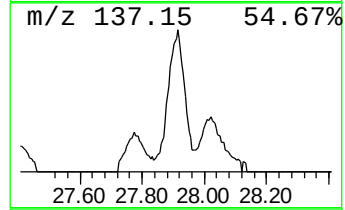
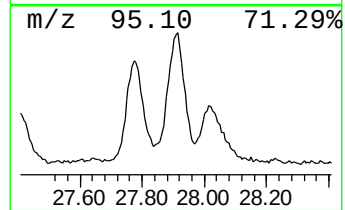
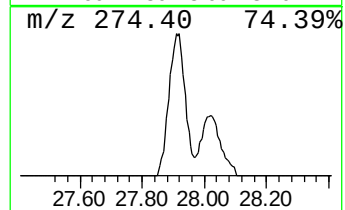
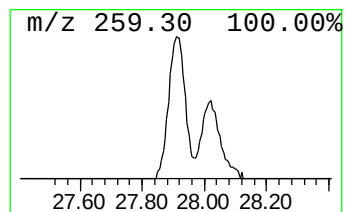
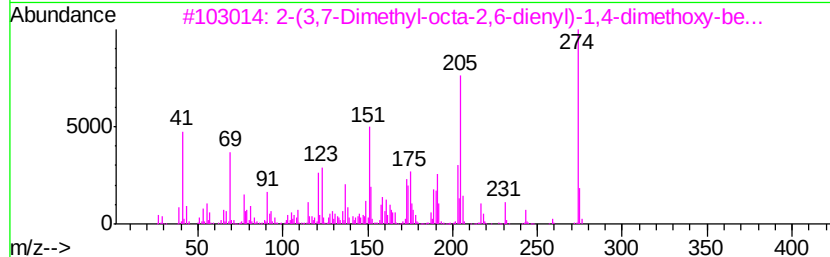
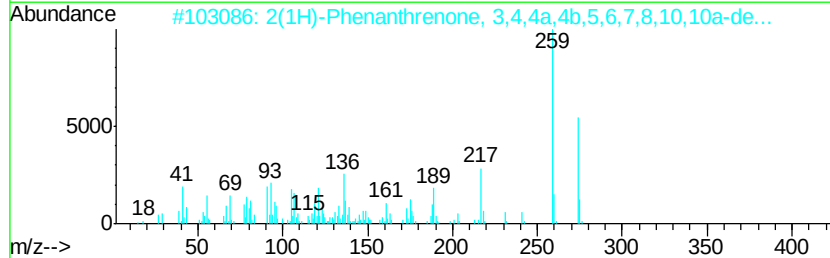
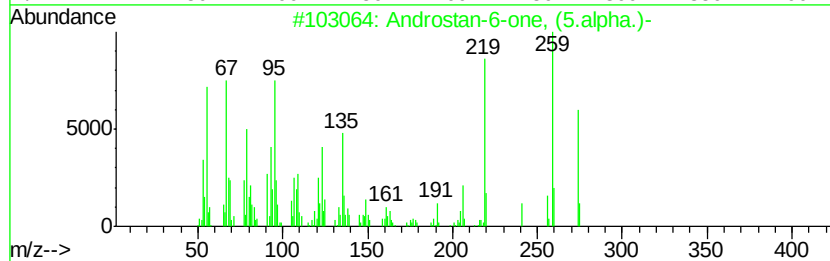
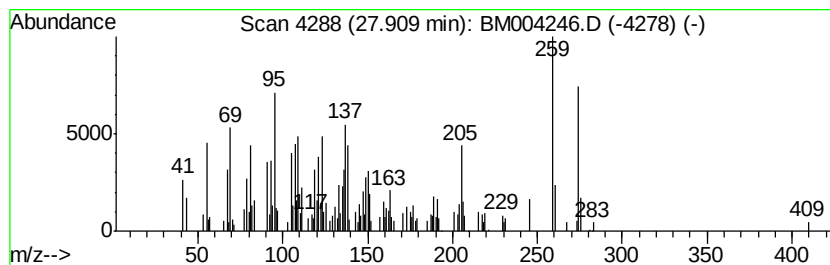
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Androstan-6-one, (5.alpha.)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.91	31.68 ng/ul	568975	Perylene-d12	23.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Androstan-6-one, (5.alpha.)-	274	C19H30O	003676-06-0	58
2		2(1H)-Phenanthrenone, 3,4,4a,4b,...	274	C19H30O	007715-44-8	45
3		2-(3,7-Dimethyl-octa-2,6-dienvl)...	274	C18H26O2	107697-34-7	40
4		Androst-16-en-3-ol, (3.alpha.,5....	274	C19H30O	001153-51-1	38
5		Androst-7-en-3-ol #	274	C19H30O	1000148-48-7	30



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

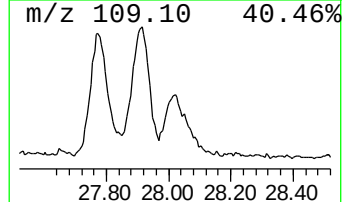
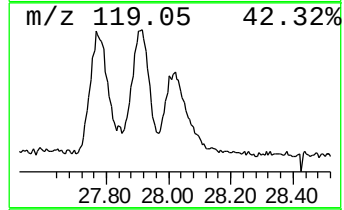
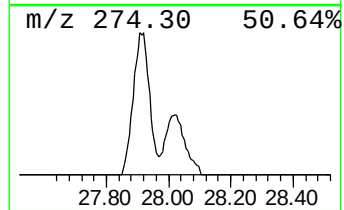
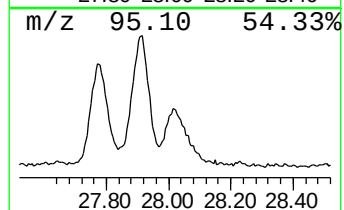
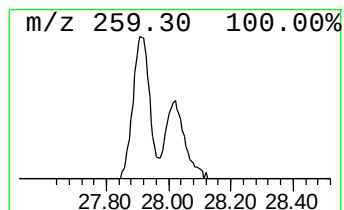
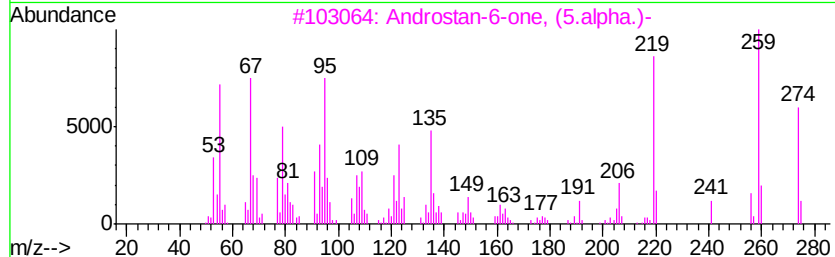
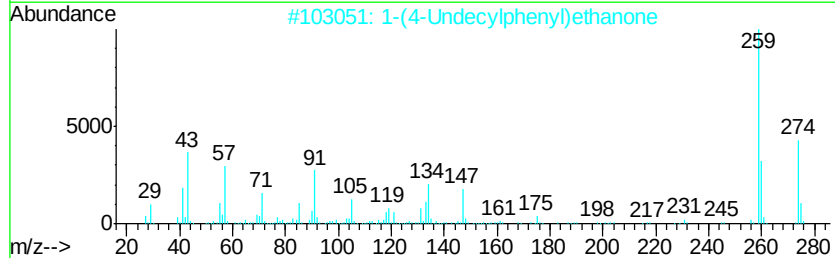
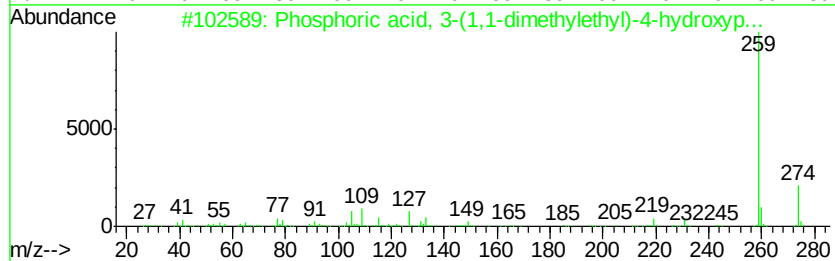
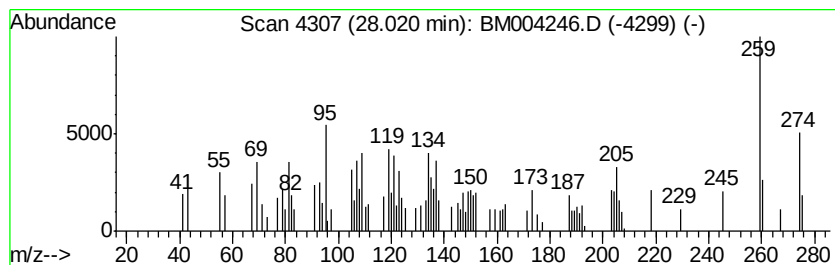
Instrument :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.02	16.59 ng/ul	297938	Perylene-d12	23.50
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Phosphoric acid, 3-(1,1-dimethyl...	274 C12H19O5P	094420-61-8 43
2		1-(4-Undecylphenyl)ethanone	274 C19H30O	1000185-92-4 38
3		Androstan-6-one, (5.alpha.)-	274 C19H30O	003676-06-0 38
4		Androstan-6-one, (5.alpha.)-	274 C19H30O	003676-06-0 38
5		2(1H)-Phenanthrenone, 3,4,4a,4b,...	274 C19H30O	007715-44-8 35



Data Path : Z:\HPCHEM1\BNA M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IZ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

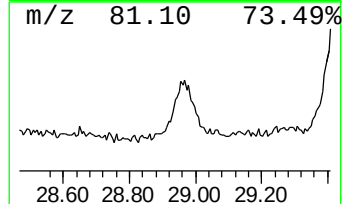
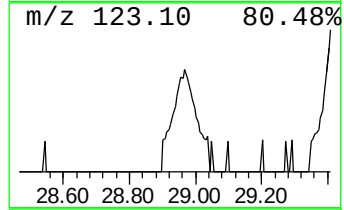
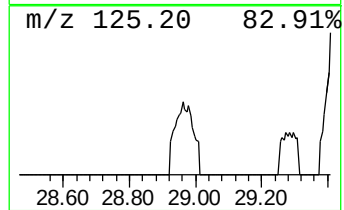
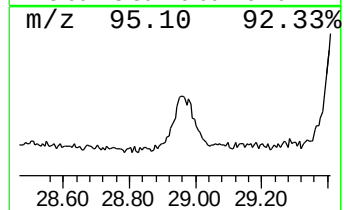
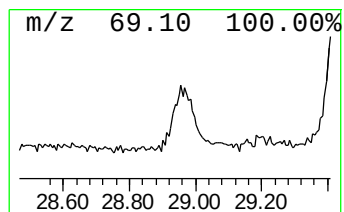
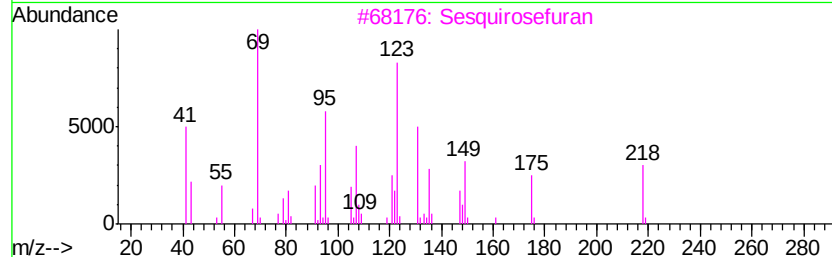
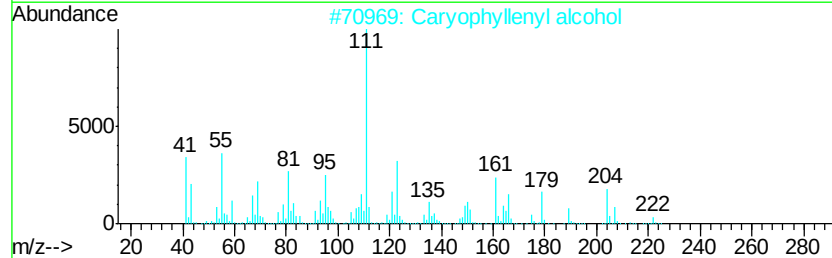
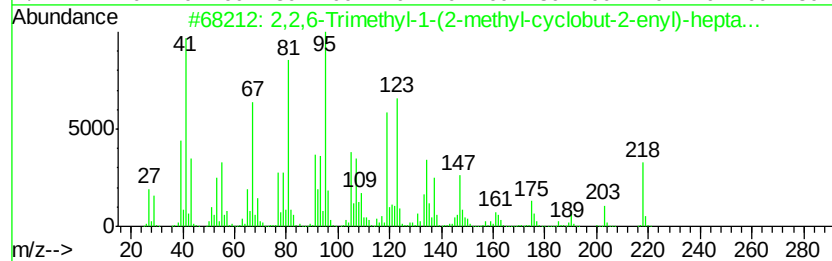
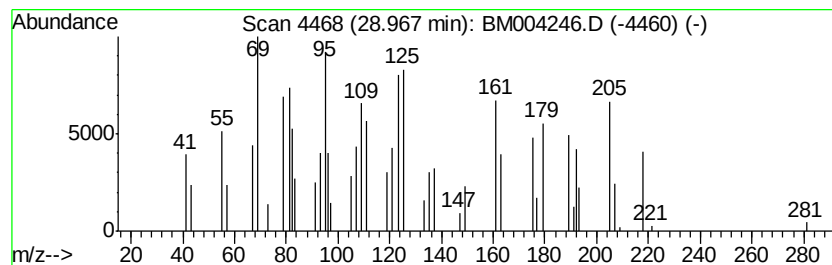
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-03 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
28.97	3.20 ng/ul	57507	Perylene-d12	23.50		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	30	
2	Carvophyllenyl alcohol	222	C15H26O	1000131-90-4	25	
3	Sesquirosefuran	218	C15H22O	039007-93-7	22	
4	3-Cyclopentylpropionic acid, 3-c...	216	C11H17ClO2	1000292-47-2	22	
5	1,5-Dimethyl-bicyclo[3.2.0]hepta...	168	C10H16O2	1000190-69-8	17	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021216\
Data File : BM004246.D
Acq On : 13 Feb 2016 02:38
Operator : SJ/IJ
Sample : H1414-07
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
D9QD2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM021216.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.85	17.2	ng/ul	135197	1	7.65	157427	20.0
Trifluoroacetic a...	20.97	3.7	ng/ul	69872	5	21.27	378930	20.0
(DEL) Alkane: Cyc...	21.86	6.1	ng/ul	116273	5	21.27	378930	20.0
Trichloroacetic a...	22.85	4.7	ng/ul	84465	6	23.50	359204	20.0
unknown-01	26.51	3.9	ng/ul	70224	6	23.50	359204	20.0
.beta.-Humulene	27.41	11.8	ng/ul	212364	6	23.50	359204	20.0
Hop-22(29)-en-3.b...	27.77	28.8	ng/ul	517646	6	23.50	359204	20.0
Androstan-6-one, ...	27.91	31.7	ng/ul	568975	6	23.50	359204	20.0
unknown-02	28.02	16.6	ng/ul	297938	6	23.50	359204	20.0
unknown-03	28.97	3.2	ng/ul	57507	6	23.50	359204	20.0