

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
 Data File : BM004554.D
 Acq On : 04 Mar 2016 04:36
 Operator : SJ/UM
 Sample : H1616-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P010-SS004-01

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BM030316.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	6.852	703	708	717	rBV	68535	123818	3.88%	0.538%
2	6.934	717	722	733	rVB	65411	100723	3.16%	0.438%
3	7.157	754	760	767	rBV	93305	144414	4.53%	0.628%
4	7.593	828	834	842	rBB	128979	196663	6.17%	0.855%
5	8.369	961	966	974	rBV	75851	131935	4.14%	0.574%
6	8.757	1026	1032	1038	rBV	80899	131496	4.12%	0.572%
7	9.475	1149	1154	1162	rBB	73803	116599	3.66%	0.507%
8	10.075	1250	1256	1271	rBV	97883	189100	5.93%	0.822%
9	10.375	1299	1307	1312	rBV	197571	332734	10.44%	1.446%
10	10.428	1312	1316	1323	rVB	74204	114289	3.58%	0.497%
11	10.540	1330	1335	1357	rBB	25457	51940	1.63%	0.226%
12	13.410	1819	1823	1835	rVB5	15145	36963	1.16%	0.161%
13	13.575	1845	1851	1856	rBV2	22799	42043	1.32%	0.183%
14	13.669	1862	1867	1872	rBV	229383	328835	10.31%	1.429%
15	13.722	1872	1876	1881	rVB	110021	158581	4.97%	0.689%
16	13.951	1909	1915	1925	rBV	276813	429390	13.47%	1.867%
17	14.151	1944	1949	1953	rBV	26439	35863	1.12%	0.156%
18	14.257	1962	1967	1975	rVB2	299728	466062	14.62%	2.026%
19	14.486	2000	2006	2011	rVB6	17845	33217	1.04%	0.144%
20	14.663	2031	2036	2040	rBV4	26149	49618	1.56%	0.216%
21	14.880	2069	2073	2079	rBV3	15707	32120	1.01%	0.140%
22	15.057	2095	2103	2106	rBV2	26568	57742	1.81%	0.251%
23	15.110	2108	2112	2116	rVB	44019	55076	1.73%	0.239%
24	15.257	2132	2137	2143	rBV	351633	496181	15.56%	2.157%
25	15.416	2160	2164	2171	rBV2	58882	105975	3.32%	0.461%
26	15.974	2255	2259	2260	rBV	40979	49266	1.55%	0.214%
27	16.521	2348	2352	2357	rBV	77585	98528	3.09%	0.428%
28	16.816	2400	2402	2413	rVB6	30905	58163	1.82%	0.253%
29	17.016	2430	2436	2440	rBV	321624	456404	14.31%	1.984%
30	17.057	2440	2443	2447	rVV	39825	55238	1.73%	0.240%
31	17.116	2448	2453	2461	rVB2	319421	473098	14.84%	2.057%
32	18.821	2740	2743	2747	rBV2	45741	77973	2.45%	0.339%
33	19.427	2841	2846	2850	rBV	317953	460408	14.44%	2.001%
34	21.233	3149	3153	3158	rVB	255353	332922	10.44%	1.447%

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Integration Parameters: LSCINT.P

Integrator: RTE
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35	21.492	3191	3197	3201	rBV	574890	1178541	36.96%	5.123%
36	21.533	3201	3204	3207	rVV	621223	662767	20.79%	2.881%
37	21.598	3211	3215	3225	rVB	740478	1004218	31.50%	4.365%
38	21.703	3226	3233	3236	rBV	1480440	3188387	100.00%	13.860%
39	21.739	3236	3239	3244	rVV	1251614	1721903	54.01%	7.485%
40	21.792	3244	3248	3259	rVB	1192168	1518858	47.64%	6.602%
41	21.903	3262	3267	3269	rVV	894431	1277863	40.08%	5.555%
42	21.927	3269	3271	3277	rVB	1084684	1219261	38.24%	5.300%
43	22.074	3292	3296	3301	rVB	763316	970717	30.45%	4.220%
44	22.274	3325	3330	3336	rVB	1465001	2056721	64.51%	8.940%
45	22.380	3342	3348	3353	rVB	1125515	1579320	49.53%	6.865%
46	23.315	3503	3507	3513	rVB	169097	272204	8.54%	1.183%
47	23.450	3526	3530	3536	rVB	179326	330467	10.36%	1.437%

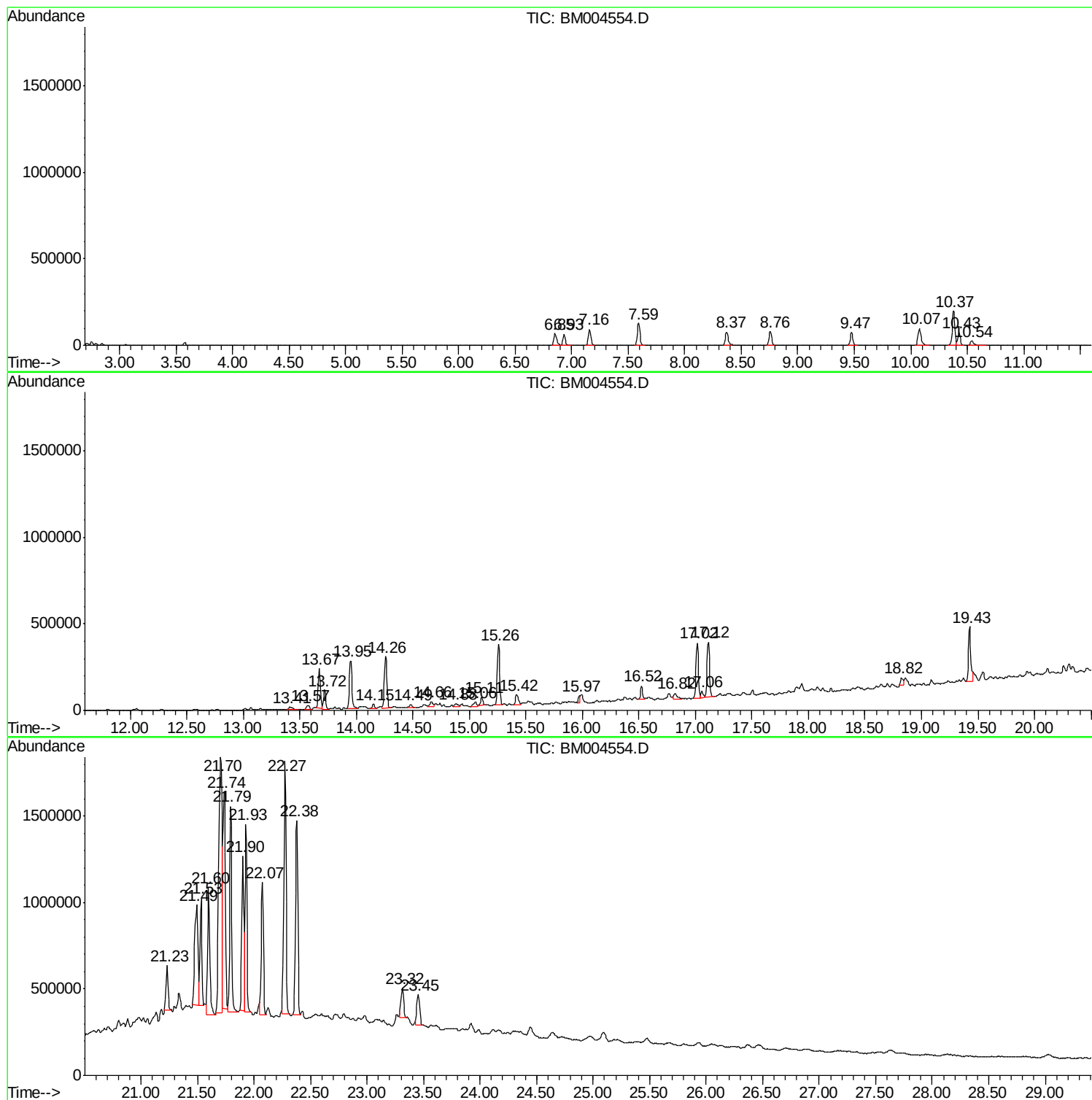
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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



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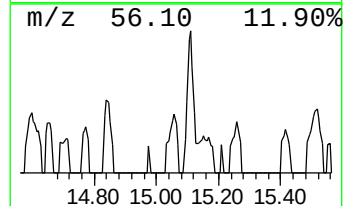
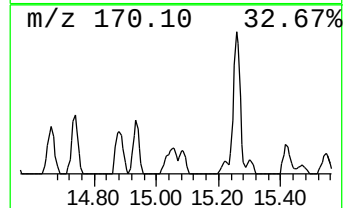
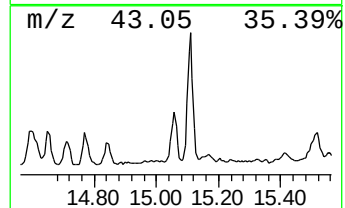
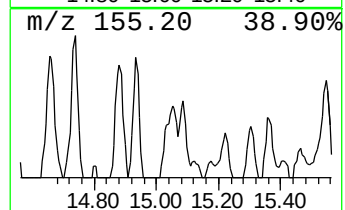
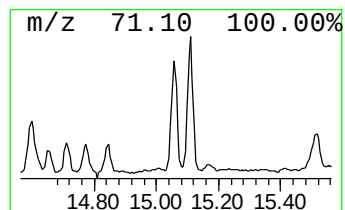
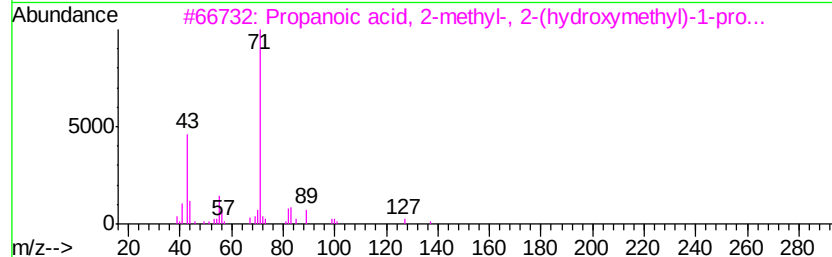
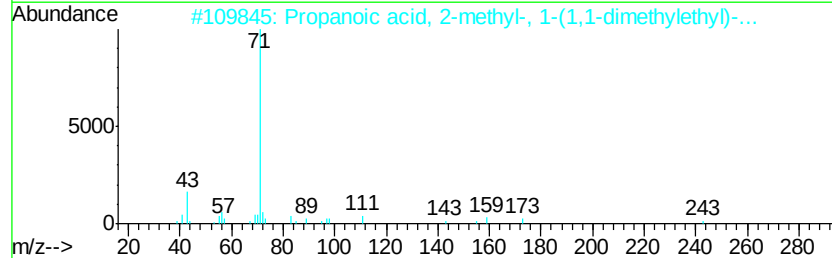
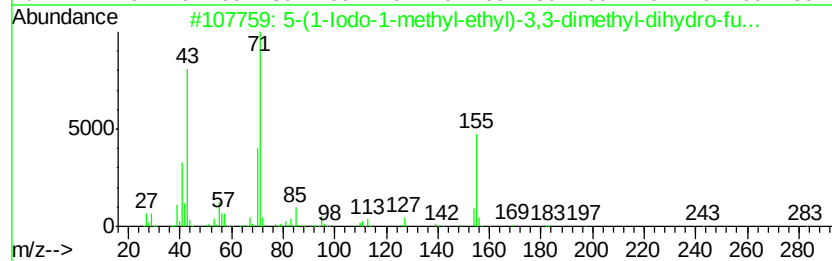
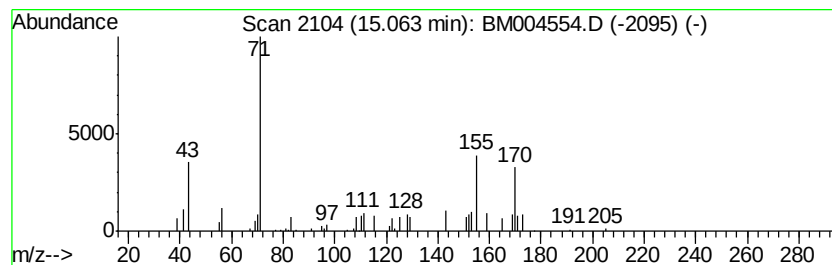
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.06	2.48 ng/ul	57742	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-(1-Iodo-1-methyl-ethyl)-3,3-di...	282	C9H15IO2	1000190-61-9	47
2		Propanoic acid, 2-methyl-, 1-(1,...	286	C16H30O4	074381-40-1	38
3		Propanoic acid, 2-methyl-, 2-(hy...	216	C12H24O3	074367-32-1	35
4		3-Buten-2-ol, 1-bromo-2-methyl-	164	C5H9BrO	036219-40-6	35
5		Pyrano[3,2-b]pyran-2(3H)-one, he...	156	C8H12O3	060378-32-7	32



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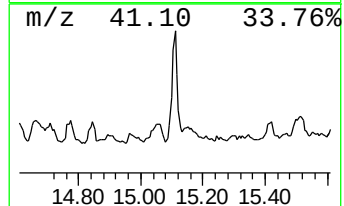
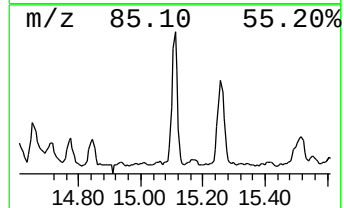
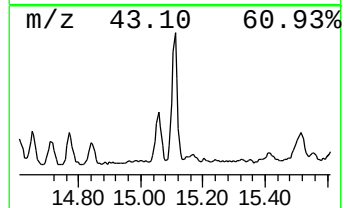
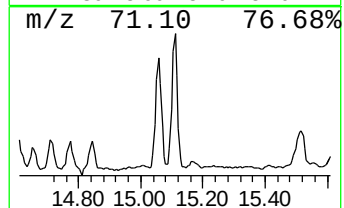
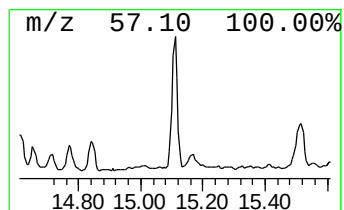
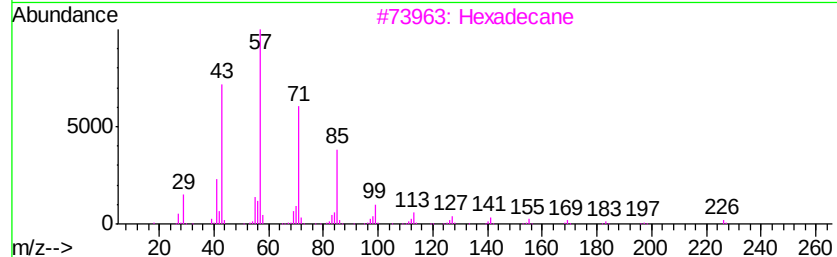
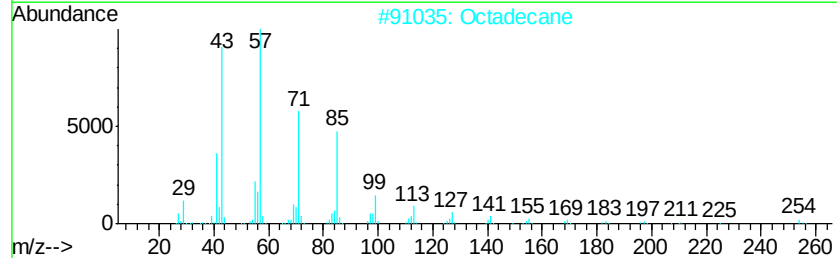
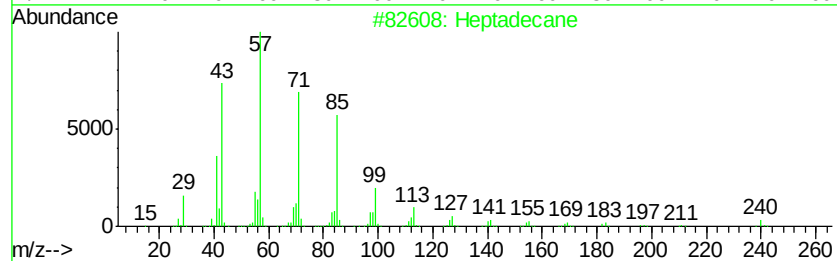
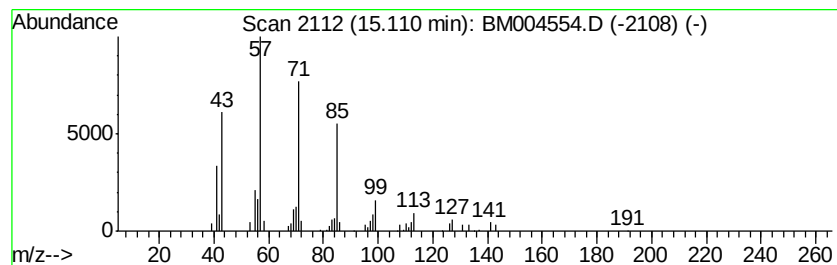
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	2.36 ng/ul	55076	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane	240	C17H36	000629-78-7	91
2		Octadecane	254	C18H38	000593-45-3	91
3		Hexadecane	226	C16H34	000544-76-3	90
4		Pentadecane	212	C15H32	000629-62-9	90
5		Octadecane	254	C18H38	000593-45-3	90



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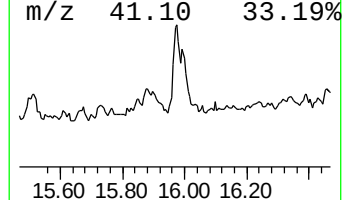
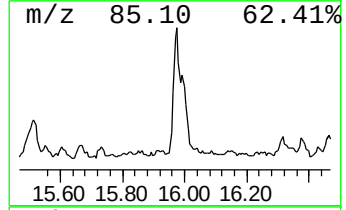
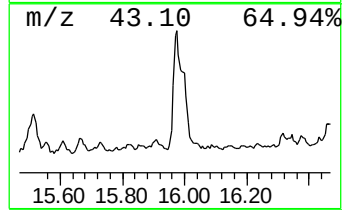
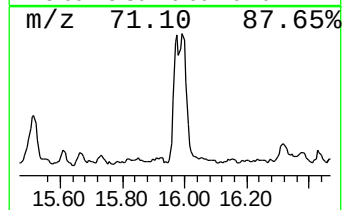
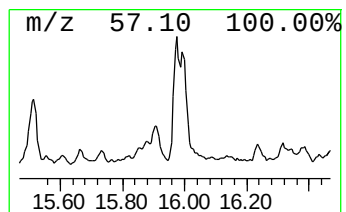
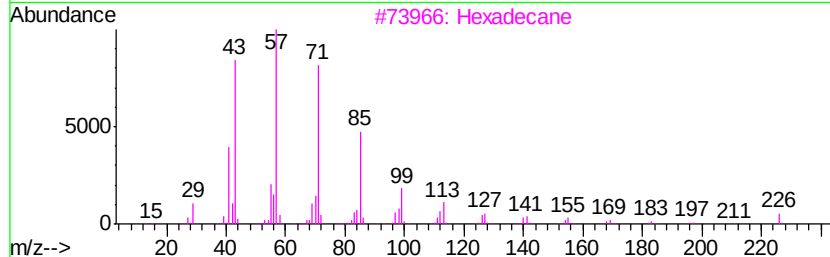
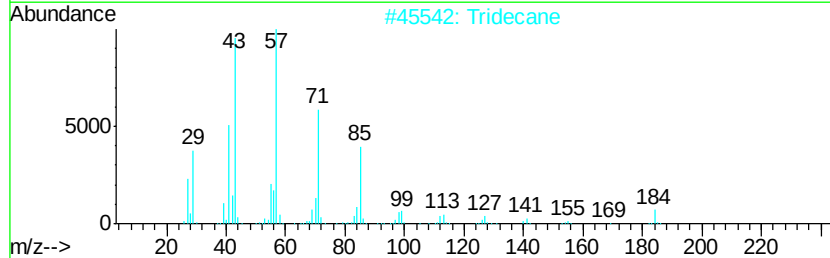
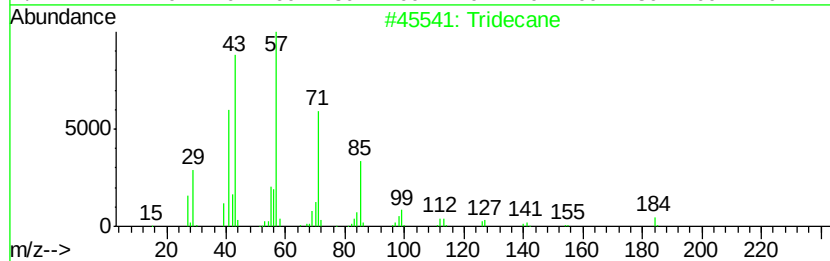
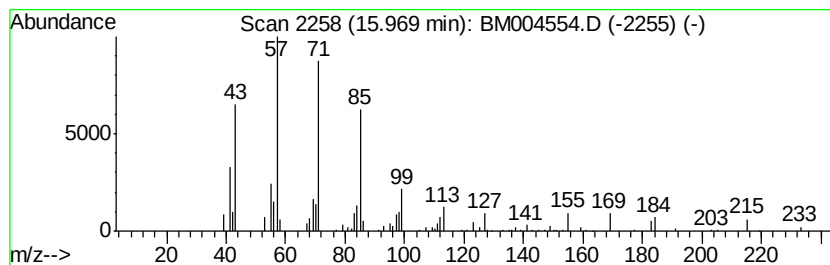
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	2.16 ng/ul	49266	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecane	184	C13H28	000629-50-5	93
2		Tridecane	184	C13H28	000629-50-5	90
3		Hexadecane	226	C16H34	000544-76-3	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heptadecane	240	C17H36	000629-78-7	90



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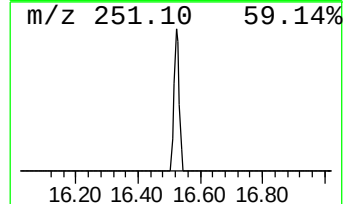
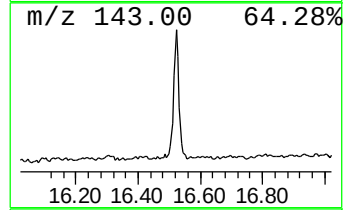
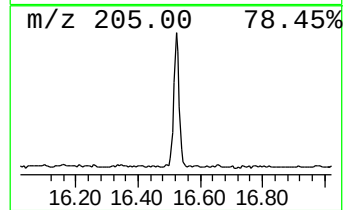
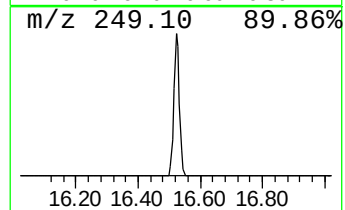
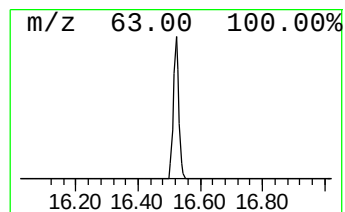
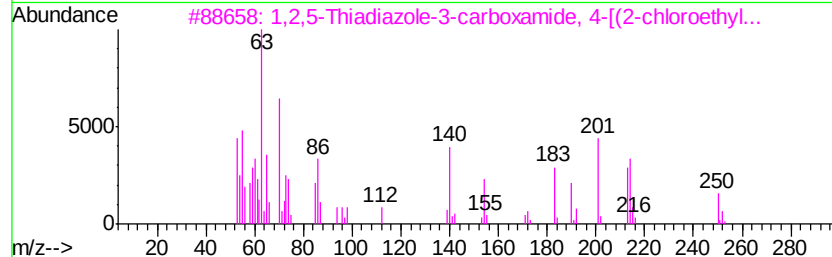
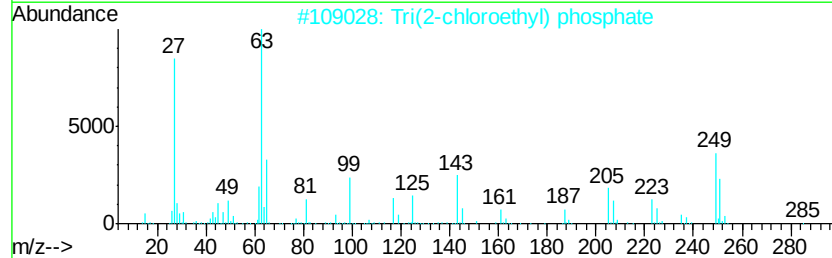
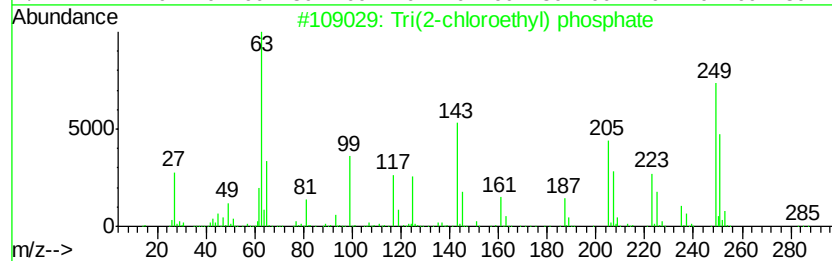
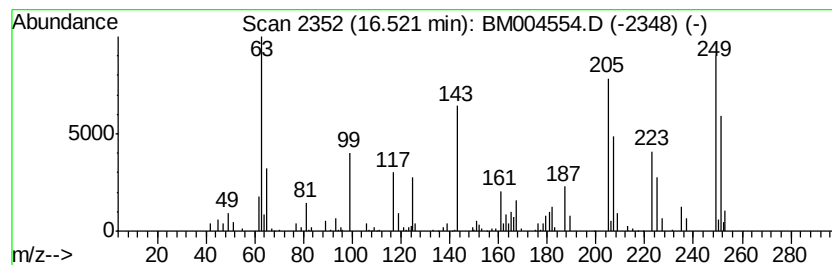
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Tri(2-chloroethyl) phosphate Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.52	4.32 ng/ul	98528	Phenanthrene-d10	17.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tri(2-chloroethyl) phosphate	284	C6H12Cl3O4P	000115-96-8	90
2			Tri(2-chloroethyl) phosphate	284	C6H12Cl3O4P	000115-96-8	90
3			1,2,5-Thiadiazole-3-carboxamide,...	250	C7H11ClN4O2S	147194-45-4	43
4			Ethene, (2-chloroethoxy)-	106	C4H7ClO	000110-75-8	38
5			1-Chloroethyl sulfone	190	C4H8Cl2O2S	010038-07-0	38



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ClientSampled :
P010-SS004-01

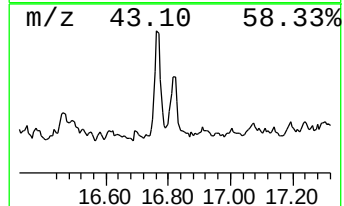
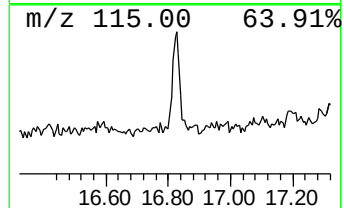
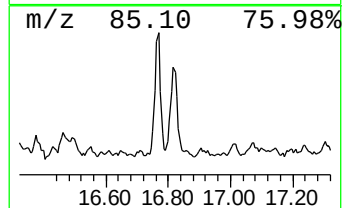
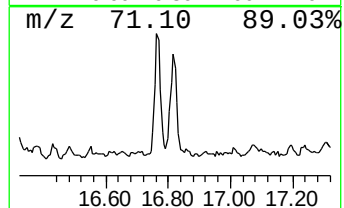
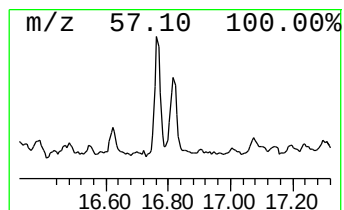
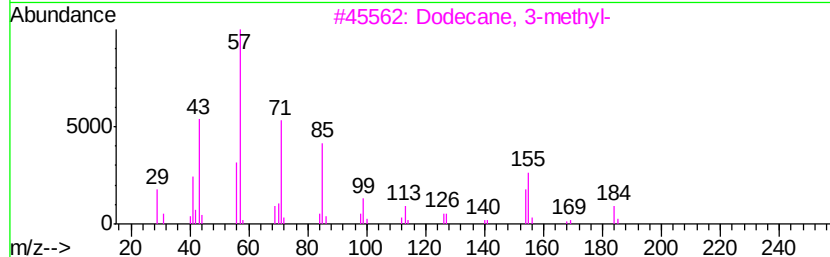
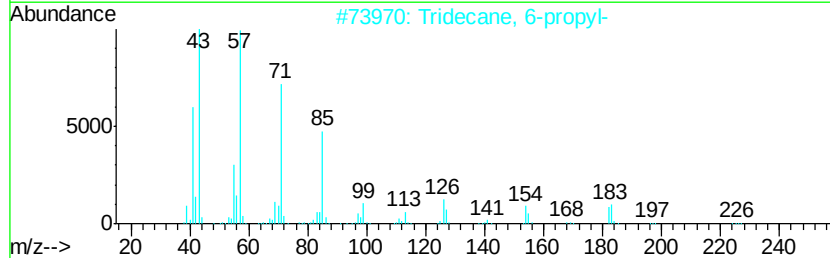
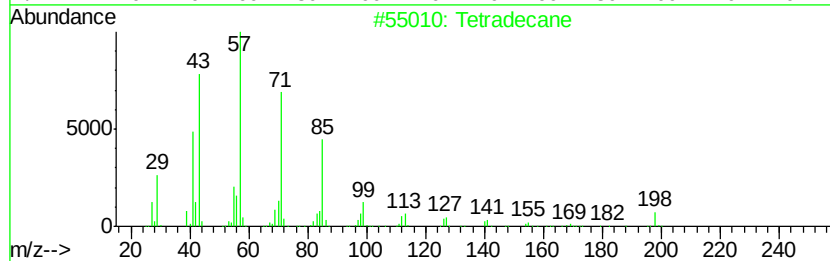
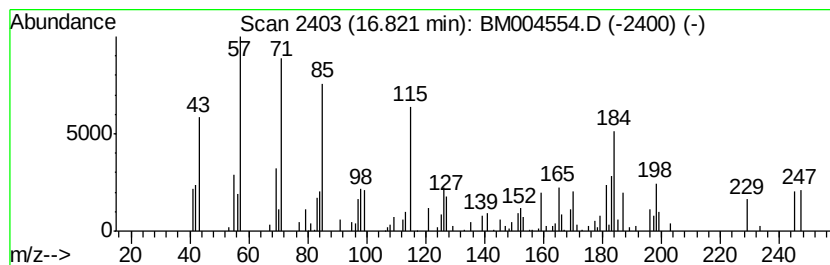
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.82	2.55 ng/ul	58163	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	45
2		Tridecane, 6-propyl-	226	C16H34	055045-10-8	43
3		Dodecane, 3-methyl-	184	C13H28	017312-57-1	38
4		Pentadecane, 3-methyl-	226	C16H34	002882-96-4	38
5		Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	35



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Data File : BM004554.D
Acq On : 04 Mar 2016 04:36
Operator : SJ/UM
Sample : H1616-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P010-SS004-01

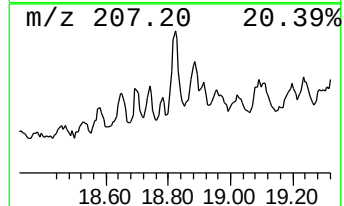
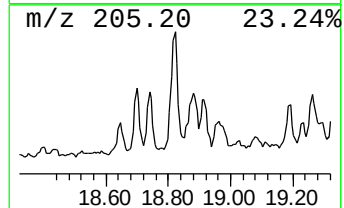
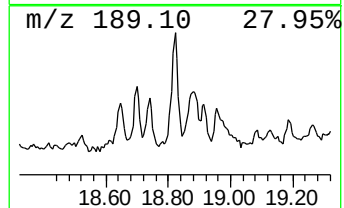
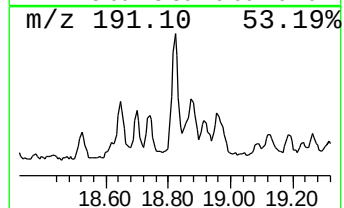
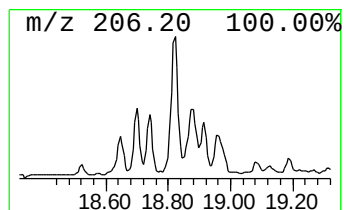
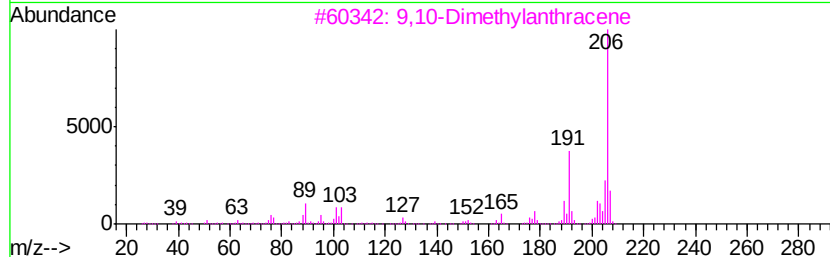
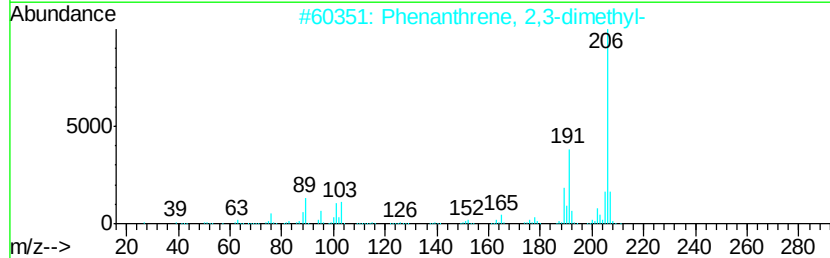
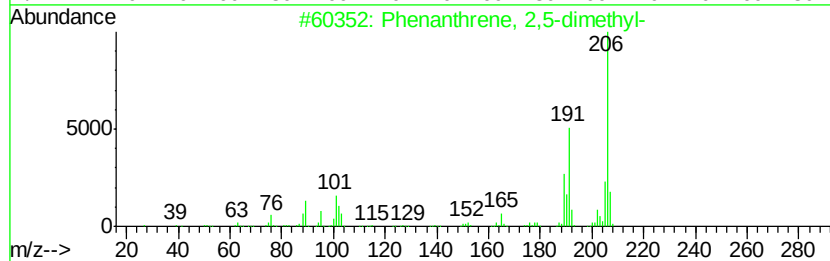
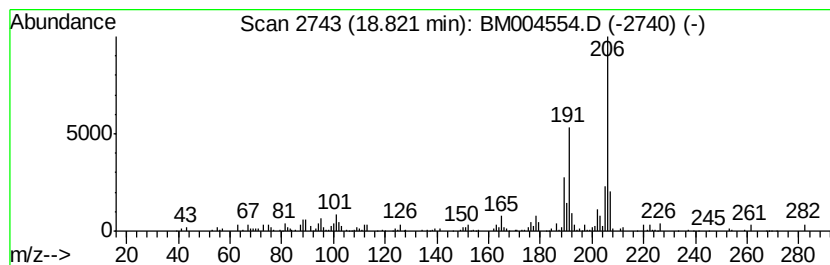
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenanthrene, 2,5-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	3.42 ng/ul	77973	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
4		Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	90
5		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	89



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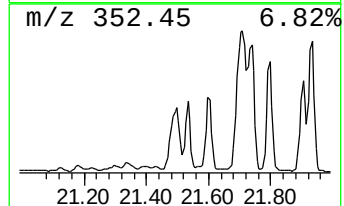
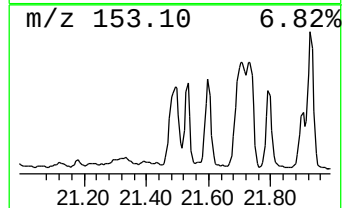
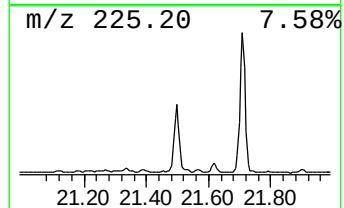
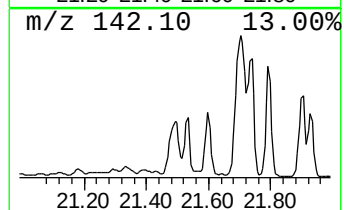
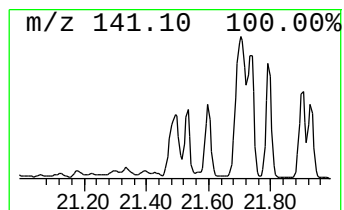
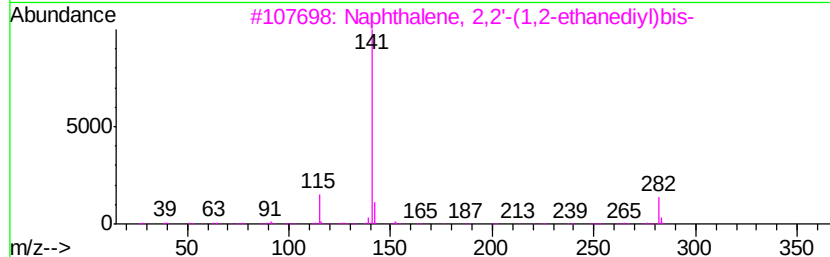
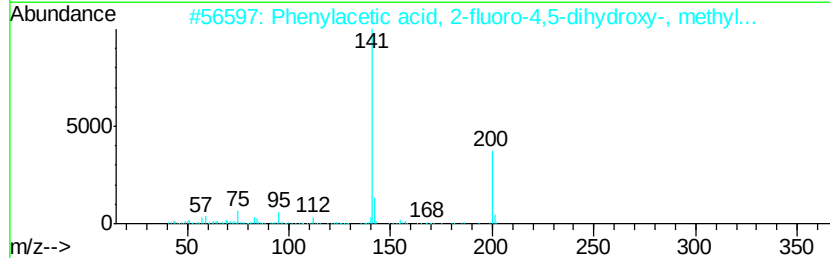
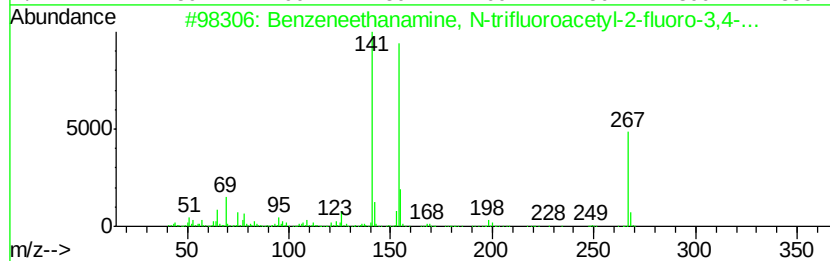
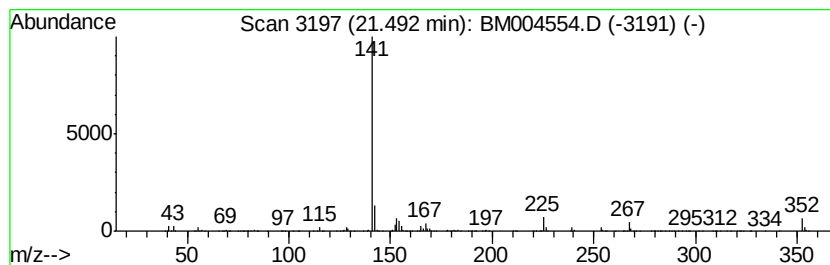
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.49	70.80 ng/ul	1178540	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneethanamine, N-trifluoroac...	267	C10H9F4N03	059043-77-5	38
2		Phenylacetic acid, 2-fluoro-4,5-...	200	C9H9F04	141523-25-3	9
3		Naphthalene, 2,2'-(1,2-ethanediyl...	282	C22H18	021969-45-9	9
4		1-Naphthalenepropionic acid	200	C13H12O2	003243-42-3	9
5		1-But-3-enyl naphthalene	182	C14H14	002489-88-5	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
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Sample : H1616-15
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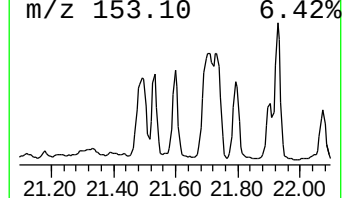
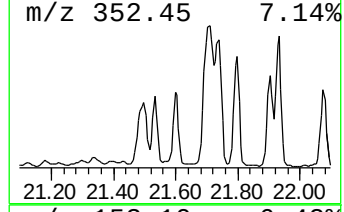
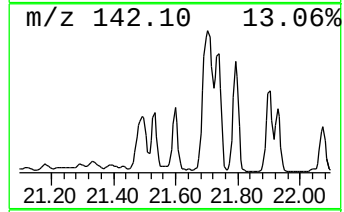
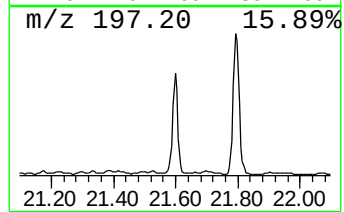
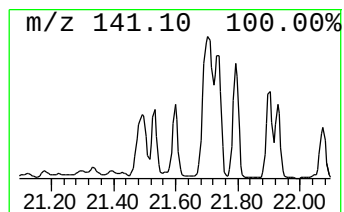
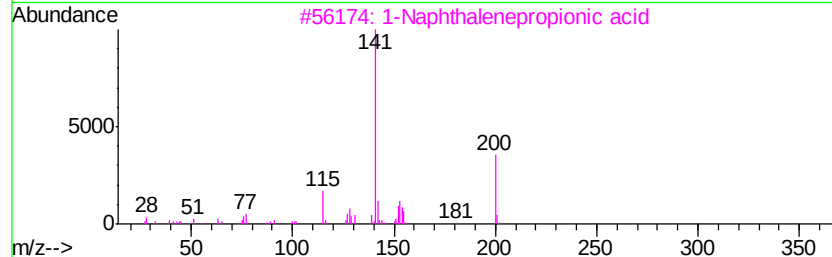
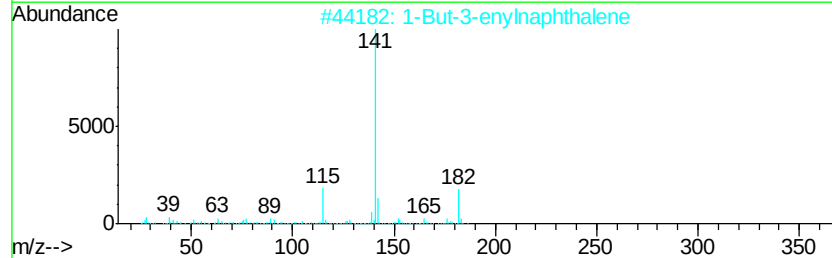
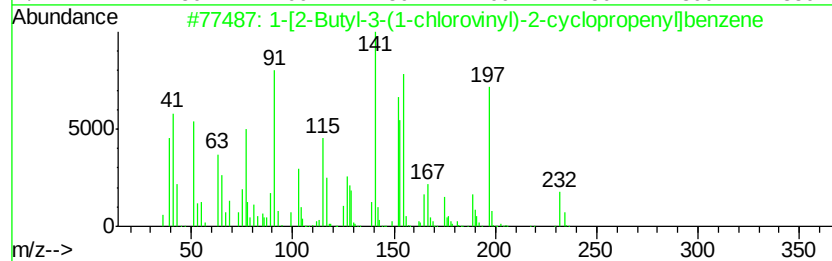
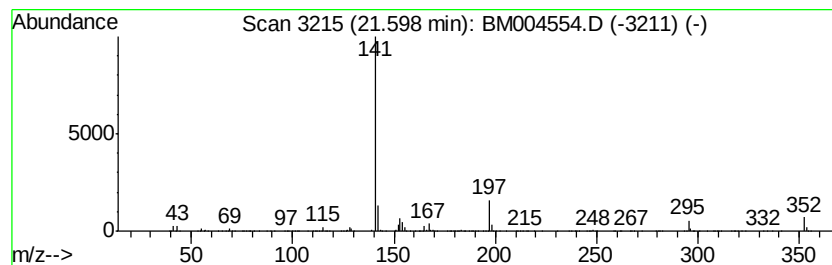
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-[2-Butyl-3-(1-chlorovinyl)... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.60	60.33 ng/ul	1004220	Chrysene-d12	21.23		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-[2-Butyl-3-(1-chlorovinyl)-2-c...	232	C15H17Cl	1000261-30-1	59	
2	1-But-3-enylnaphthalene	182	C14H14	002489-88-5	9	
3	1-Naphthalenepropionic acid	200	C13H12O2	003243-42-3	9	
4	Pentanamide, N-(3-chloro-4-methy...	239	C13H18ClNO	002307-68-8	9	
5	3-(1-Ethoxy-ethoxy)-4,4,4-triflu...	258	C10H17F3O4	095605-52-0	9	



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
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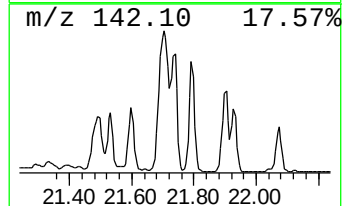
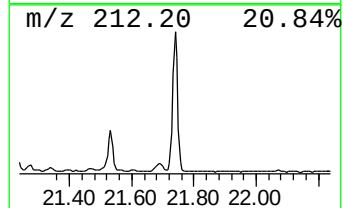
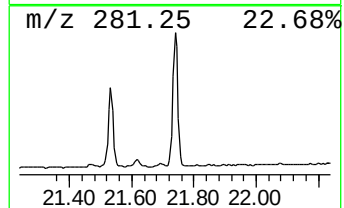
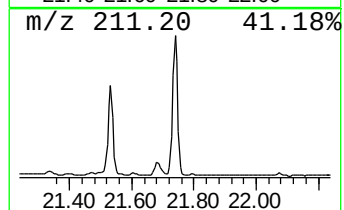
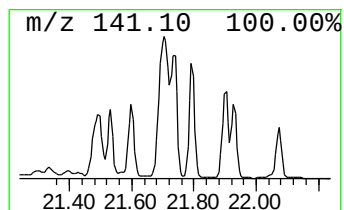
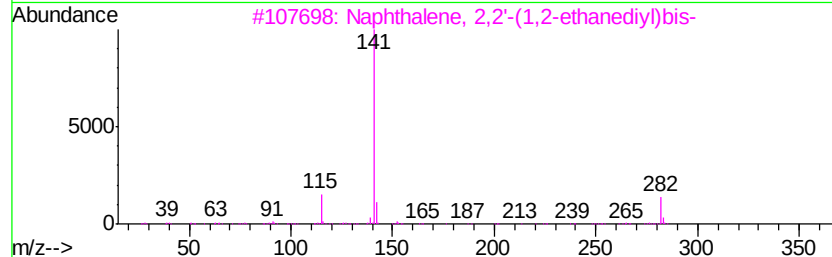
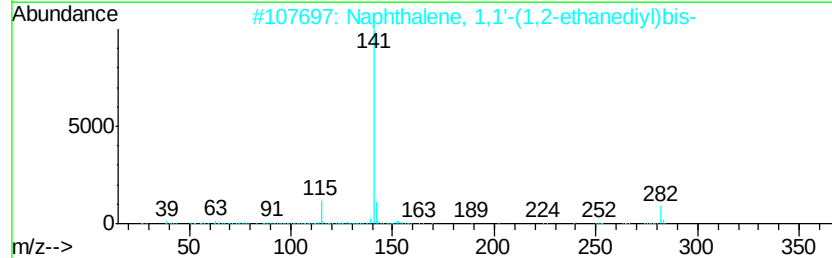
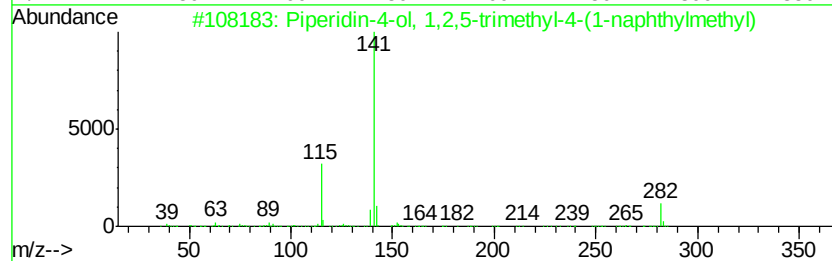
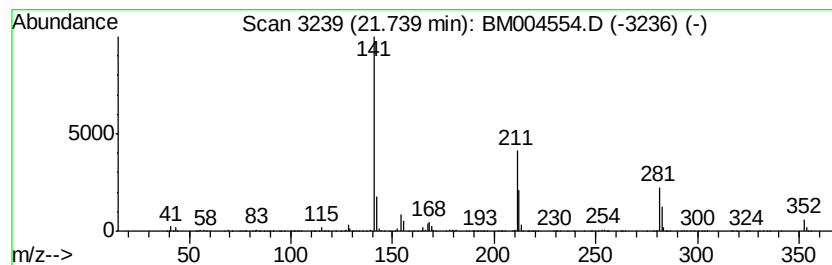
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-03 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.74	103.44 ng/ul	1721900	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Piperidin-4-ol, 1,2,5-trimethyl-...	283	C19H25NO	329789-67-5	38
2		Naphthalene, 1,1'-(1,2-ethanediyl)...	282	C22H18	015374-45-5	35
3		Naphthalene, 2,2'-(1,2-ethanediyl)...	282	C22H18	021969-45-9	35
4		Undecane, 1-(1-naphthyl)-	282	C21H30	007225-71-0	9
5		Naphthalene, 2-hexyl-	212	C16H20	002876-46-2	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
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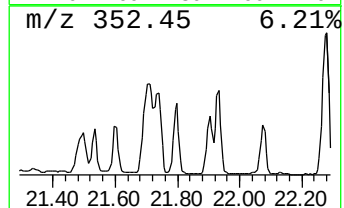
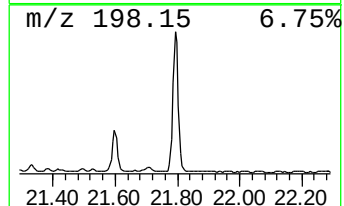
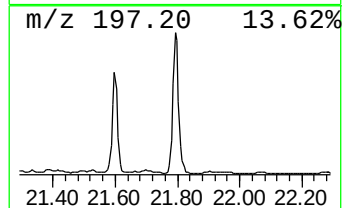
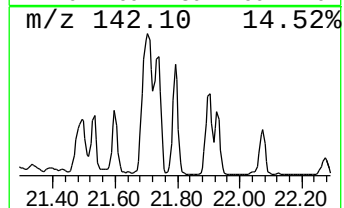
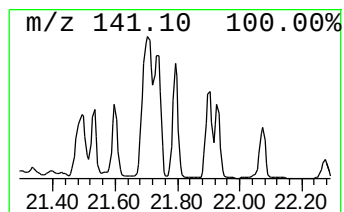
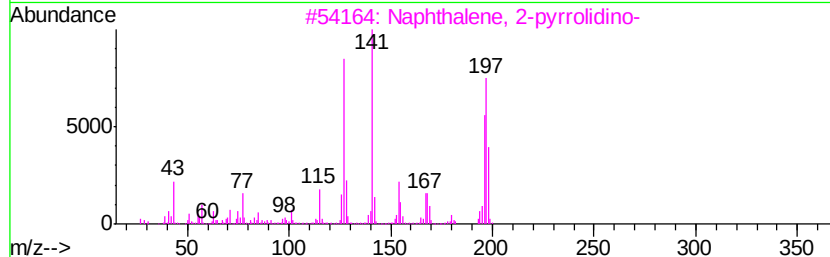
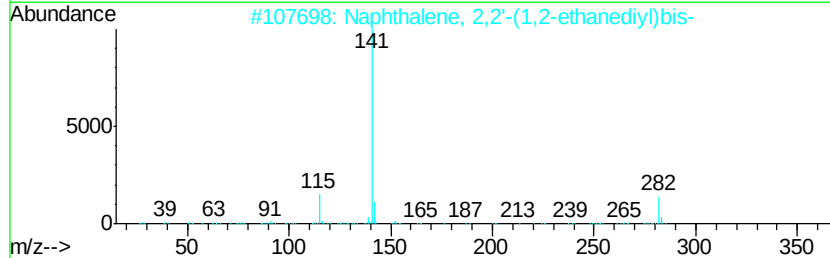
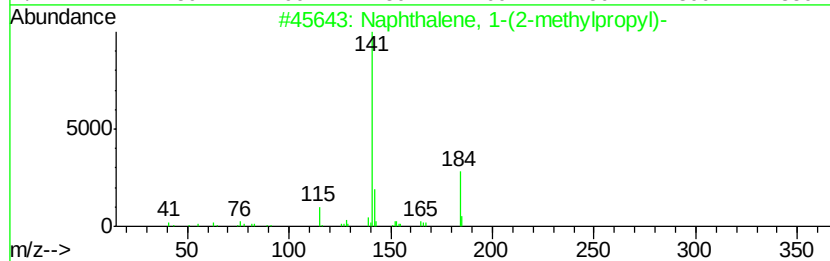
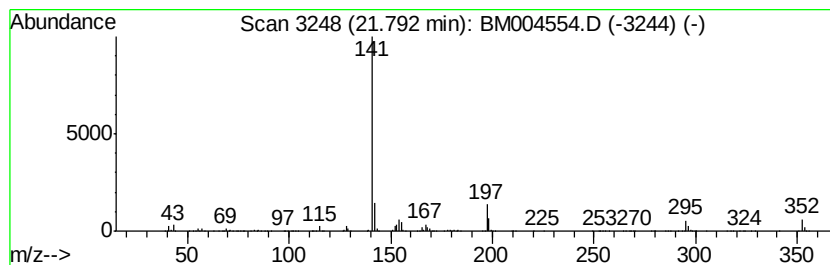
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Naphthalene, 1-(2-methylpro... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.79	91.24 ng/ul	1518860	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-(2-methylpropyl)-	184	C14H16	016727-91-6	53
2		Naphthalene, 2,2'-(1,2-ethanediyl)bis-	282	C22H18	021969-45-9	36
3		Naphthalene, 2-pyrrolidino-	197	C14H15N	013672-14-5	10
4		1-Pent-4-enynaphthalene	196	C15H16	093359-74-1	9
5		2-Naphthaleneethanol	172	C12H12O	001485-07-0	9



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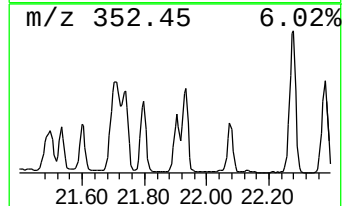
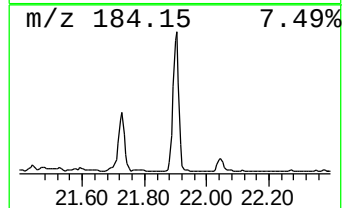
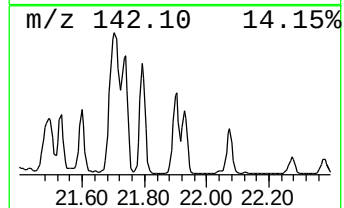
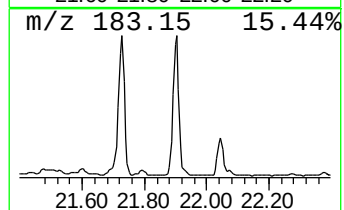
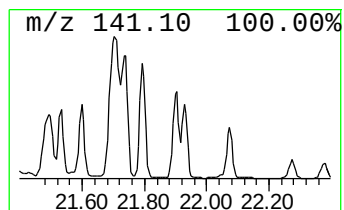
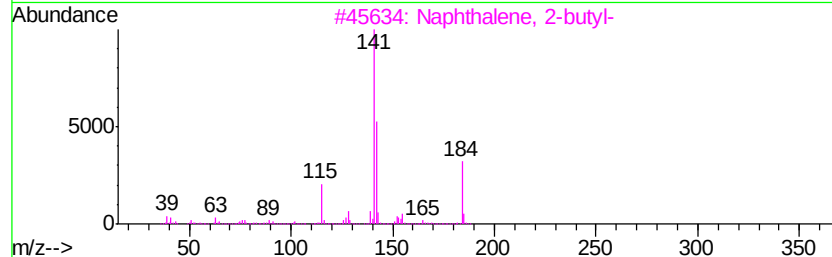
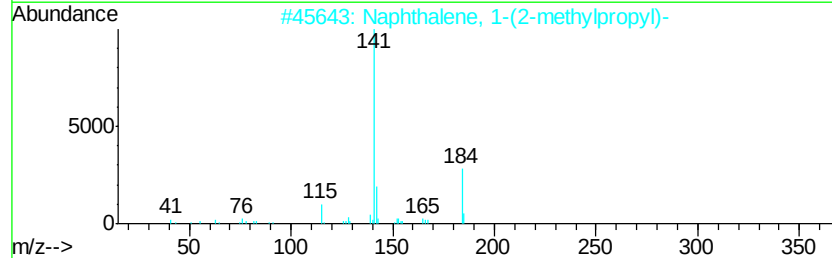
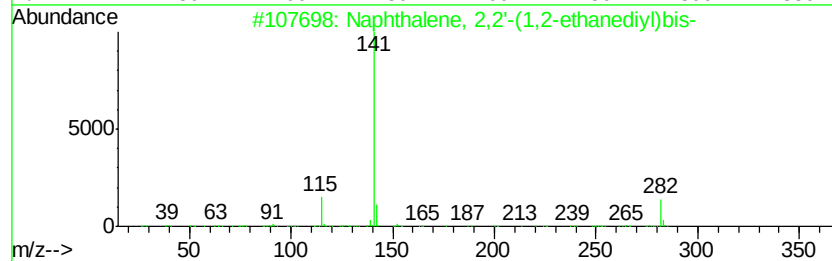
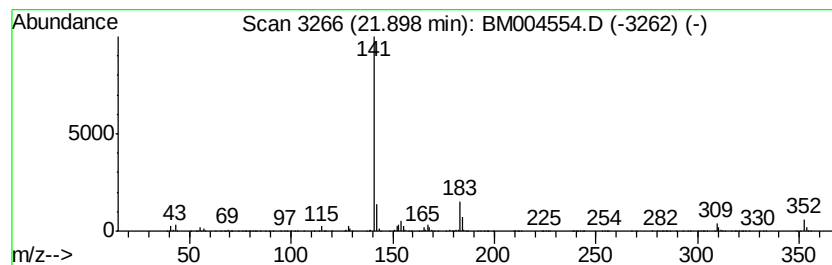
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 2,2'-(1,2-etha... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.90	76.77 ng/ul	1277860	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,2'-(1,2-ethanediy...	282	C22H18	021969-45-9	58
2		Naphthalene, 1-(2-methylpropyl)-	184	C14H16	016727-91-6	27
3		Naphthalene, 2-butyl-	184	C14H16	001134-62-9	9
4		Methyl 8-oxo-cis-2-nonenoate	184	C10H16O3	028297-04-3	9
5		2-Butyl-4-ethyl-5-methylthiazole	183	C10H17NS	052414-88-7	9



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ClientSampleId :
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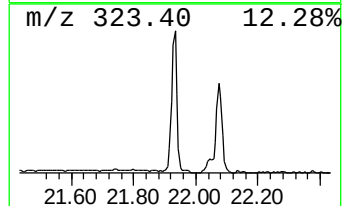
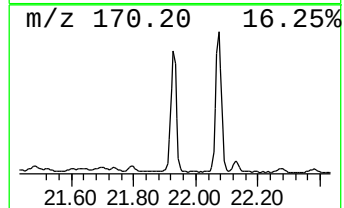
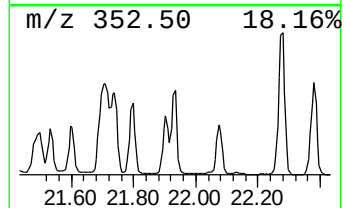
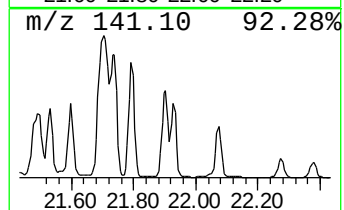
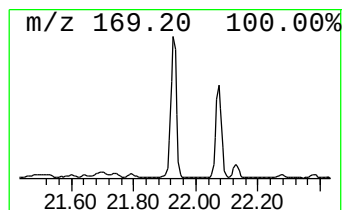
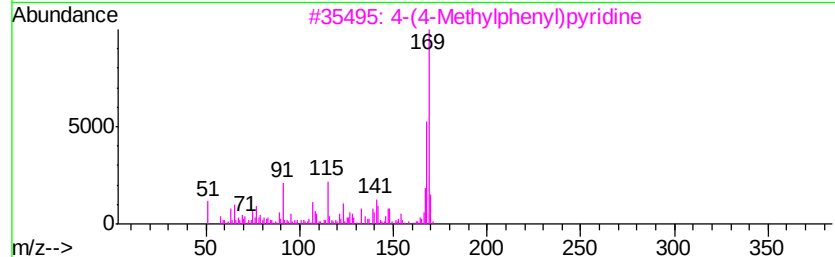
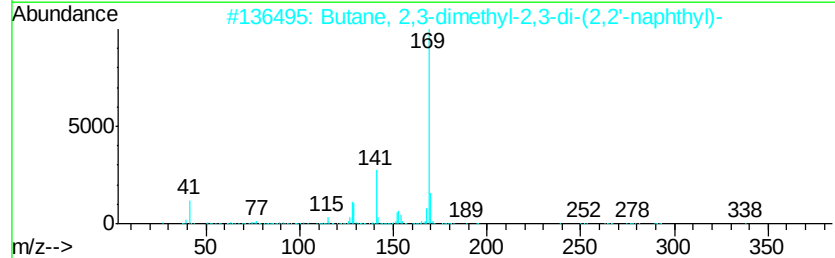
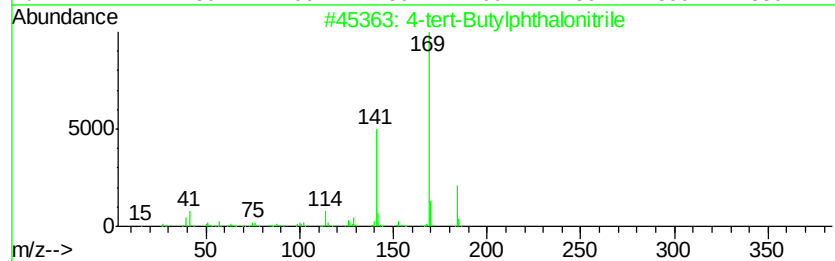
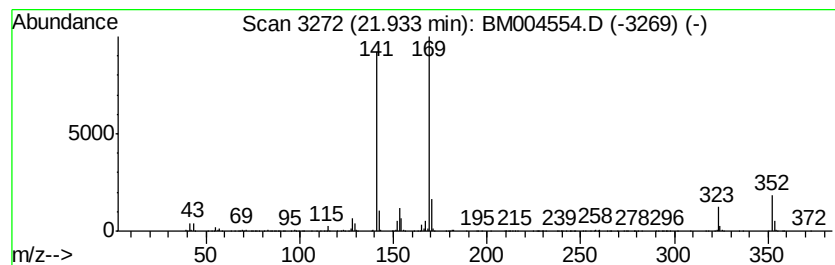
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 4-tert-Butylphthalonitrile Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.93	73.25 ng/ul	1219260	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-tert-Butylphthalonitrile	184	C12H12N2	032703-80-3	50
2		Butane, 2,3-dimethyl-2,3-di-(2,2...	338	C26H26	1000150-93-8	39
3		4-(4-Methylphenyl)pyridine	169	C12H11N	004423-10-3	38
4		Bicyclo[2.2.1]hept-2-ene, 1,7,7-...	212	C16H20	007070-09-9	38
5		Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	37



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004554.D
Acq On : 04 Mar 2016 04:36
Operator : SJ/UM
Sample : H1616-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
P010-SS004-01

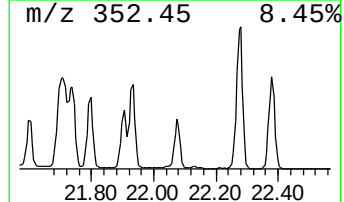
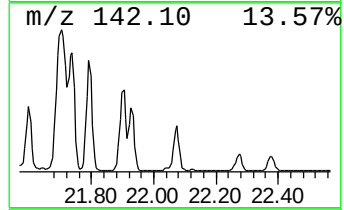
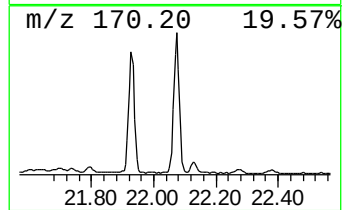
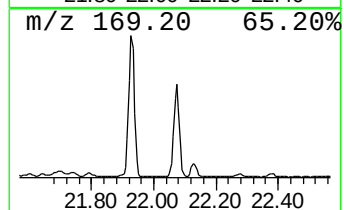
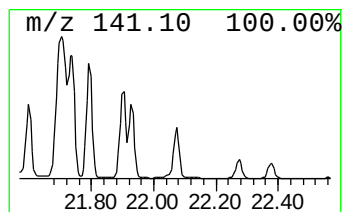
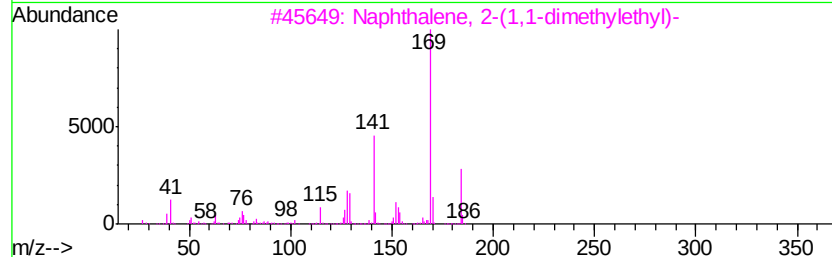
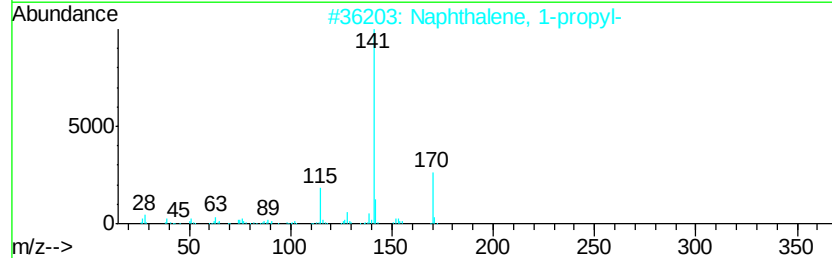
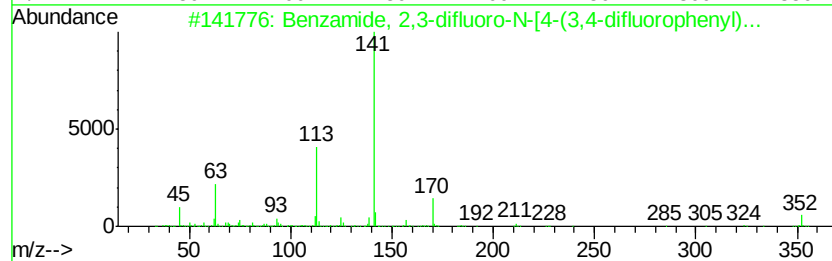
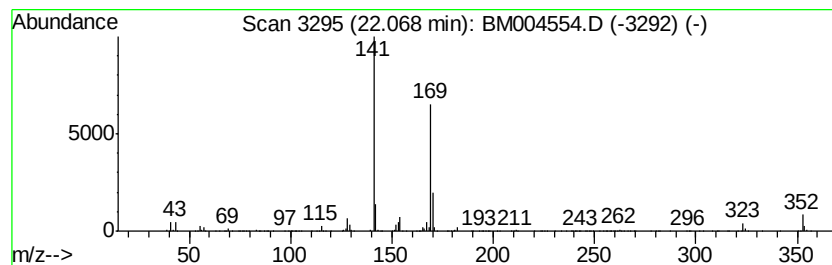
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzamide, 2,3-difluoro-N-[... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.07	58.32 ng/ul	970717	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, 2,3-difluoro-N-[4-(3,...	352	C16H8F4N2OS	1000277-49-1	50
2		Naphthalene, 1-propyl-	170	C13H14	002765-18-6	47
3		Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	37
4		4-Dodecene-6,8,10-triyn-3-one, (E)-	170	C12H10O	069698-50-6	35
5		4-tert-Butylphthalonitrile	184	C12H12N2	032703-80-3	28



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004554.D
Acq On : 04 Mar 2016 04:36
Operator : SJ/UM
Sample : H1616-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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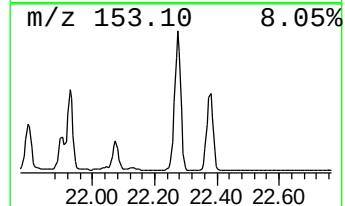
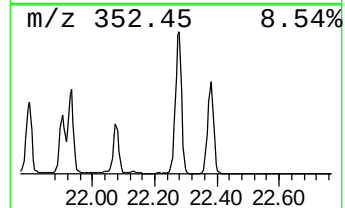
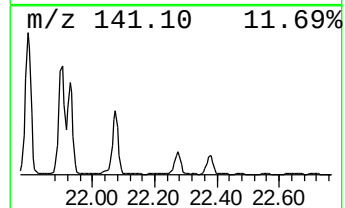
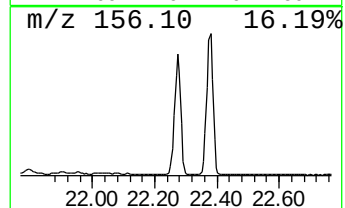
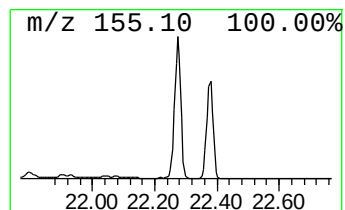
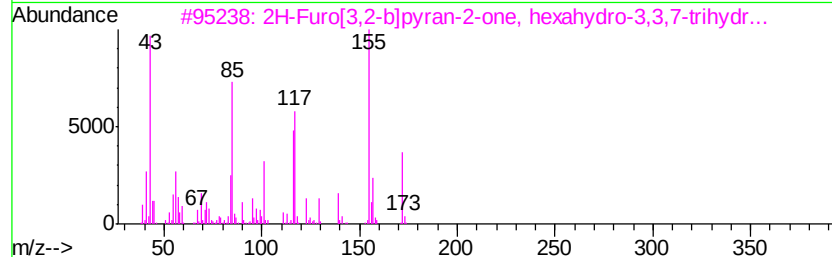
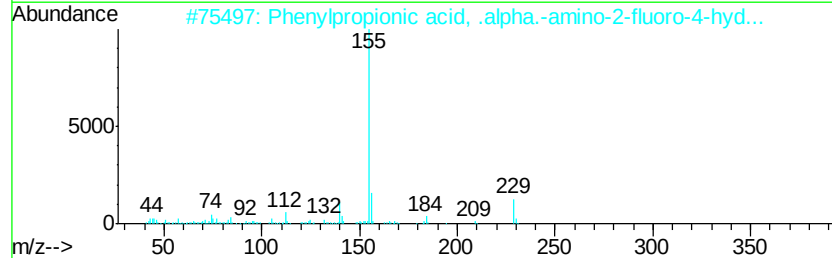
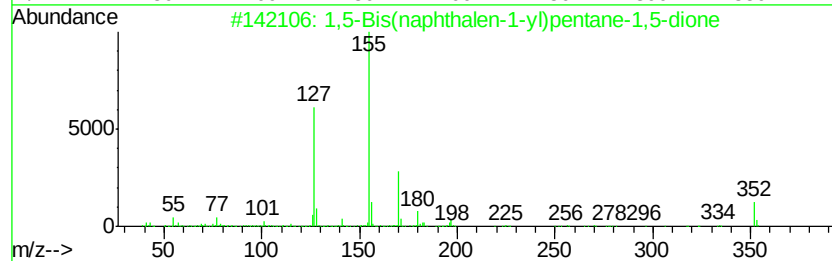
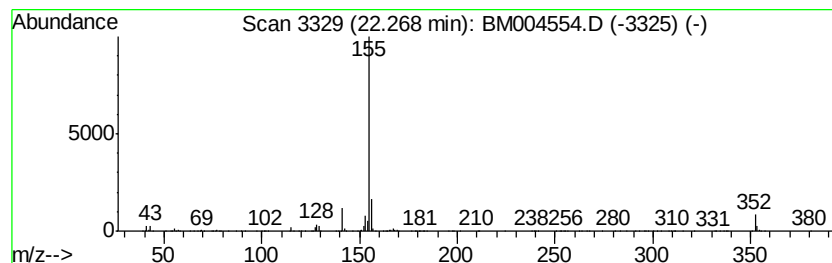
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 1,5-Bis(naphthalen-1-yl)pen... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.27	123.56 ng/ul	2056720	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,5-Bis(naphthalen-1-yl)pentane-...	352	C25H20O2	1000210-52-9	72
2		Phenylpropionic acid, .alpha.-am...	229	C10H12FN04	1000126-07-3	36
3		2H-Furo[3,2-b]pyran-2-one, hexah...	262	C11H18O7	050392-34-2	33
4		Dodecyl cis-9,10-epoxyoctadecanoate	466	C30H58O3	092332-53-1	12
5		3,7-Dimethyl-oct-6-enoic acid, (...)	323	C22H29NO	1000194-84-6	10



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
 Data File : BM004554.D
 Acq On : 04 Mar 2016 04:36
 Operator : SJ/UM
 Sample : H1616-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P010-SS004-01

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.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
unknown-01	15.06	2.5	ng/ul	57742	3	14.26	466062	20.0
(DEL) Alkane: Str...	15.11	2.4	ng/ul	55076	3	14.26	466062	20.0
(DEL) Alkane: Str...	15.97	2.2	ng/ul	49266	4	17.02	456404	20.0
Tri(2-chloroethyl...	16.52	4.3	ng/ul	98528	4	17.02	456404	20.0
(DEL) Alkane: Str...	16.82	2.5	ng/ul	58163	4	17.02	456404	20.0
Phenanthrene, 2,5...	18.82	3.4	ng/ul	77973	4	17.02	456404	20.0
unknown-02	21.49	70.8	ng/ul	1178540	5	21.23	332922	20.0
1-[2-Butyl-3-(1-c...	21.60	60.3	ng/ul	1004220	5	21.23	332922	20.0
unknown-03	21.74	103.4	ng/ul	1721900	5	21.23	332922	20.0
Naphthalene, 1-(2...	21.79	91.2	ng/ul	1518860	5	21.23	332922	20.0
Naphthalene, 2,2'...	21.90	76.8	ng/ul	1277860	5	21.23	332922	20.0
4-tert-Butylphtha...	21.93	73.3	ng/ul	1219260	5	21.23	332922	20.0
Benzamide, 2,3-di...	22.07	58.3	ng/ul	970717	5	21.23	332922	20.0
1,5-Bis(naphthale...	22.27	123.6	ng/ul	2056720	5	21.23	332922	20.0