

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
 Data File : BM010464.D
 Acq On : 09 Jun 2017 21:28
 Operator : SJ/MA
 Sample : I3520-04
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 F5A36

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\SOM-EPA-BM052717.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	6.545	582	588	594	rBV4	45327	80731	1.52%	0.123%
2	7.087	672	680	691	rBV	464242	888751	16.78%	1.351%
3	7.245	702	707	715	rBV	309530	485530	9.16%	0.738%
4	7.439	734	740	750	rVB	528361	852739	16.10%	1.296%
5	7.904	811	819	827	rBV	1229092	1997995	37.71%	3.037%
6	8.181	859	866	872	rBV3	120158	193352	3.65%	0.294%
7	8.628	934	942	950	rBV	476952	816989	15.42%	1.242%
8	9.086	1010	1020	1028	rBV	456628	775746	14.64%	1.179%
9	9.798	1135	1141	1150	rBV2	416949	712537	13.45%	1.083%
10	9.963	1162	1169	1174	rBV6	24497	55840	1.05%	0.085%
11	10.333	1224	1232	1245	rBV	664136	1197237	22.60%	1.820%
12	10.469	1248	1255	1264	rBV3	37091	72155	1.36%	0.110%
13	10.704	1287	1295	1305	rBV	1554828	2642519	49.88%	4.017%
14	10.875	1318	1324	1333	rBV2	164759	320723	6.05%	0.488%
15	12.451	1586	1592	1599	rBV2	186770	352221	6.65%	0.535%
16	13.951	1841	1847	1852	rBV	1315457	1785489	33.70%	2.714%
17	13.998	1852	1855	1861	rVB	387700	519645	9.81%	0.790%
18	14.239	1887	1896	1903	rBV	1241203	1895697	35.78%	2.882%
19	14.339	1908	1913	1917	rBV4	33496	61406	1.16%	0.093%
20	14.539	1940	1947	1958	rBV2	2455915	3718867	70.19%	5.654%
21	14.780	1978	1988	2001	rBV3	301353	640277	12.09%	0.973%
22	15.527	2106	2115	2122	rBV	1804804	2588069	48.85%	3.935%
23	15.657	2131	2137	2144	rBV	473232	756918	14.29%	1.151%
24	16.233	2230	2235	2237	rVB2	49179	60488	1.14%	0.092%
25	16.968	2353	2360	2364	rBV5	121682	190146	3.59%	0.289%
26	17.286	2407	2414	2424	rVB2	3327491	4720713	89.10%	7.177%
27	17.386	2424	2431	2440	rBV	1942842	2775467	52.39%	4.219%
28	18.027	2536	2540	2544	rBV7	39524	61799	1.17%	0.094%
29	18.133	2554	2558	2567	rBV	440225	596363	11.26%	0.907%
30	19.268	2747	2751	2752	rBV4	112446	133836	2.53%	0.203%
31	19.292	2752	2755	2766	rVB10	180815	400477	7.56%	0.609%
32	19.403	2770	2774	2784	rBV	497511	756537	14.28%	1.150%
33	19.556	2796	2800	2804	rBV6	68773	132002	2.49%	0.201%
34	19.674	2814	2820	2828	rBV	2366125	3142256	59.31%	4.777%

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35	20.139	2896	2899	2903	rVB5	99797	125169	2.36%	0.190%
36	20.209	2905	2911	2914	rBV6	269590	465487	8.79%	0.708%
37	20.580	2968	2974	2979	rBV5	353197	611045	11.53%	0.929%
38	21.109	3059	3064	3068	rBV5	410477	861971	16.27%	1.310%
39	21.450	3117	3122	3129	rVB	4471032	5298014	100.00%	8.054%
40	21.968	3207	3210	3213	rBV3	370209	470681	8.88%	0.716%
41	22.009	3213	3217	3226	rVB2	1146064	2091272	39.47%	3.179%
42	22.956	3373	3378	3384	rVB	2042490	2865748	54.09%	4.357%
43	23.033	3387	3391	3401	rVB5	570917	1286198	24.28%	1.955%
44	23.533	3472	3476	3482	rVB	812441	1368219	25.83%	2.080%
45	23.638	3489	3494	3501	rVB2	1406005	2558751	48.30%	3.890%
46	23.791	3514	3520	3528	rVB2	2446056	4583773	86.52%	6.969%
47	24.197	3583	3589	3598	rVB2	2764231	5123005	96.70%	7.788%
48	24.956	3714	3718	3728	rVB2	781908	1687420	31.85%	2.565%

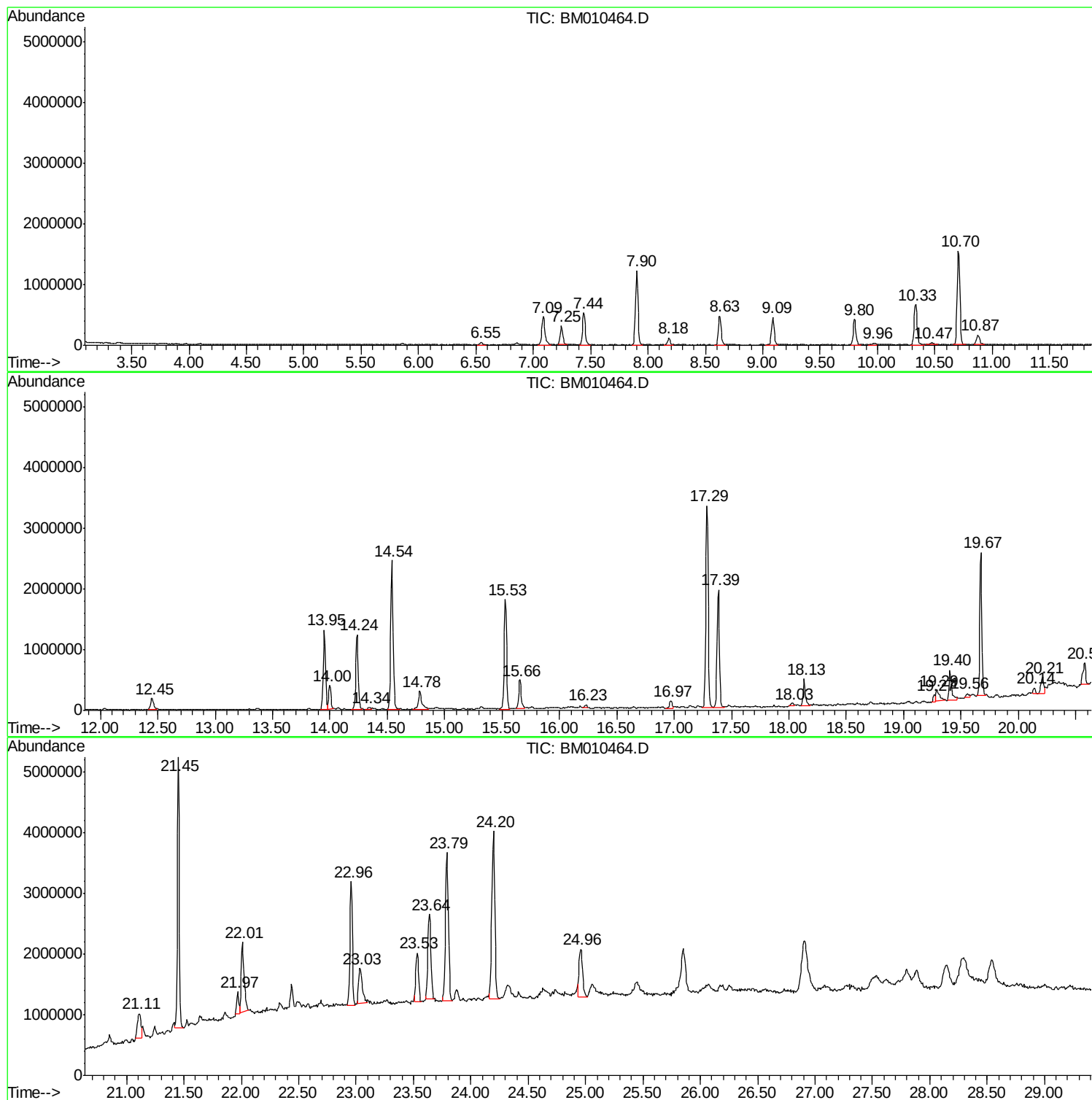
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Quant Method : Z:\HPCHEM1\BNA_M\Methods\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION

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



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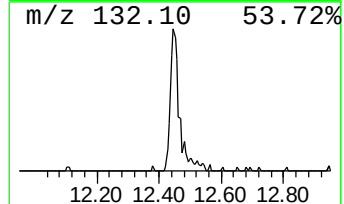
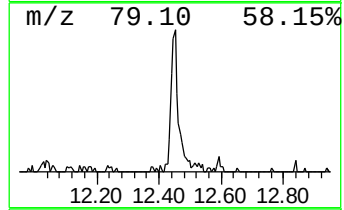
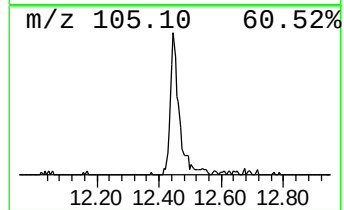
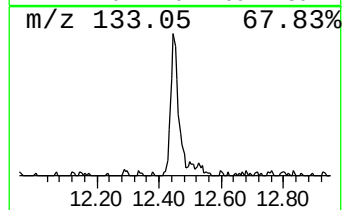
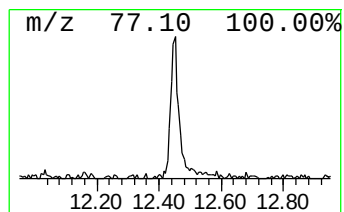
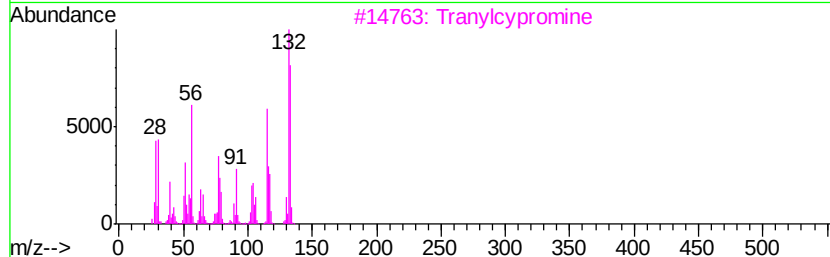
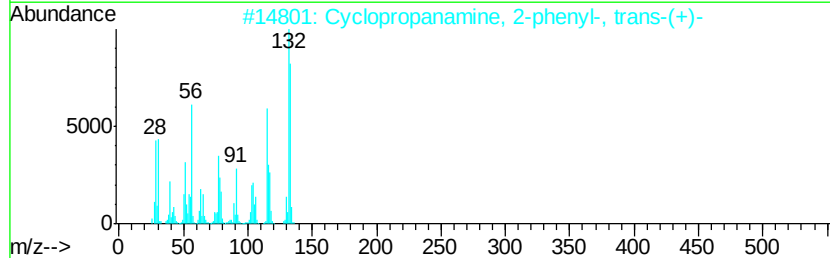
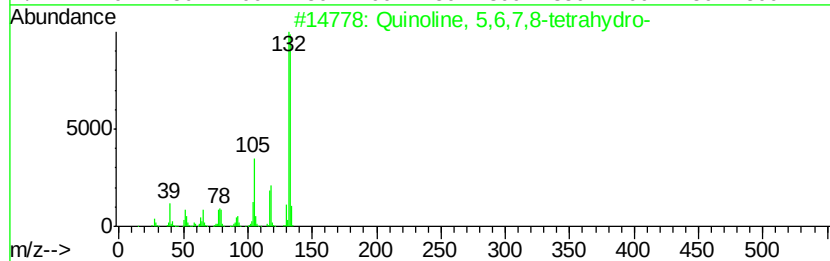
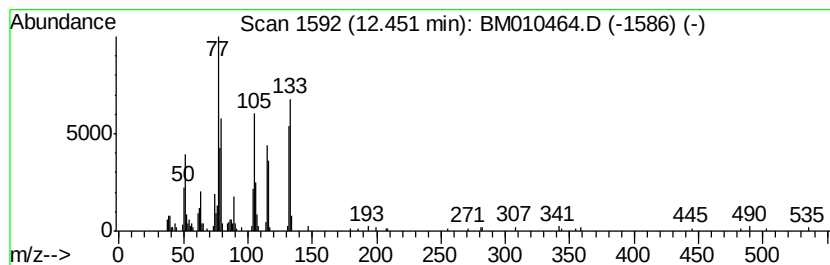
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.45	2.67 ng/ul	352221	Napthalene-d8	10.70
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Quinoline, 5,6,7,8-tetrahydro-	133 C9H11N	010500-57-9	25
2	Cyclopropanamine, 2-phenyl-, tra...	133 C9H11N	003721-28-6	22
3	Tranylcypromine	133 C9H11N	000155-09-9	22
4	Benzeneacetonitrile, 4-hydroxy-	133 C8H7NO	014191-95-8	22
5	5,6,7,8-Tetrahydroisoquinoline	133 C9H11N	1000379-32-5	22



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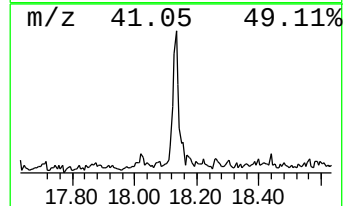
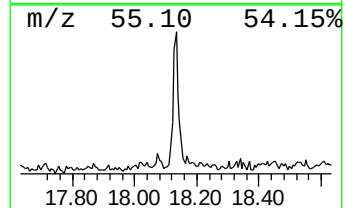
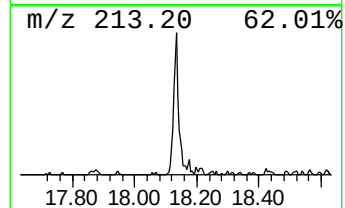
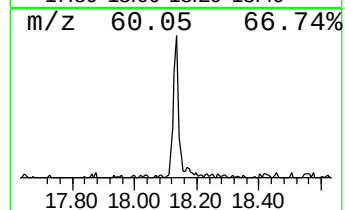
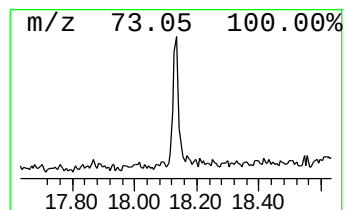
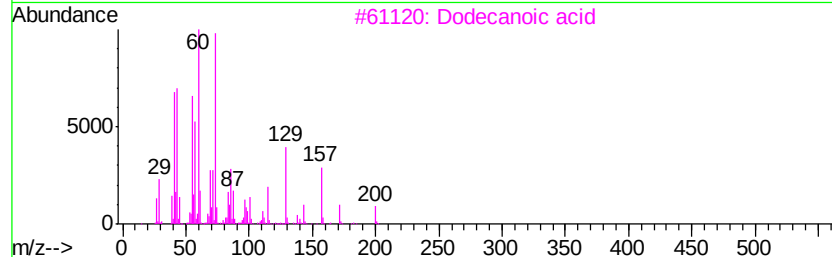
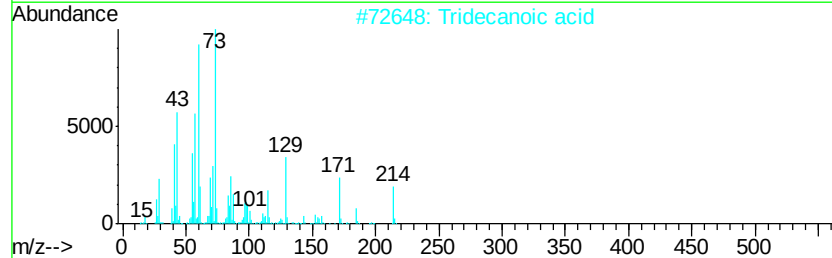
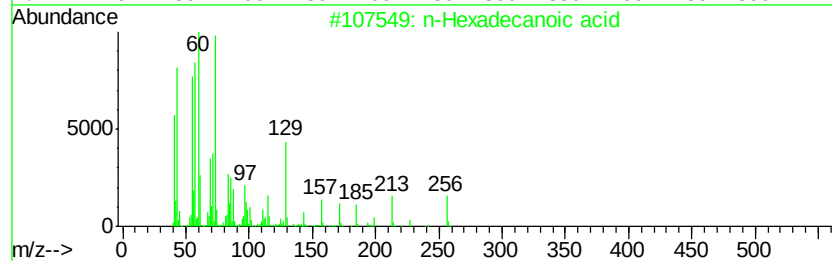
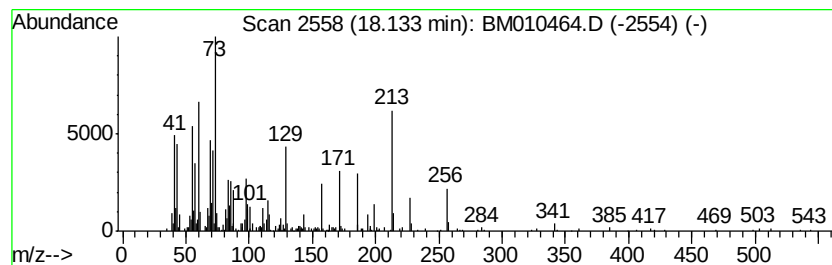
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	2.53 ng/ul	596363	Phenanthrene-d10	17.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
2		Tridecanoic acid	214	C13H26O2	000638-53-9	64
3		Dodecanoic acid	200	C12H24O2	000143-07-7	50
4		Dodecanoic acid	200	C12H24O2	000143-07-7	38
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	35



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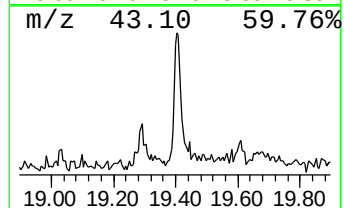
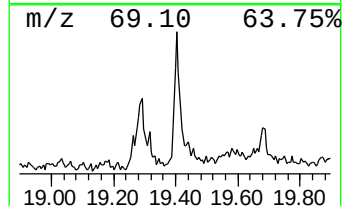
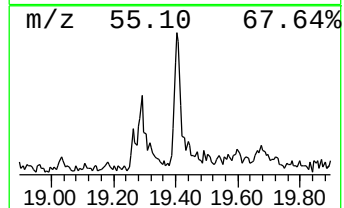
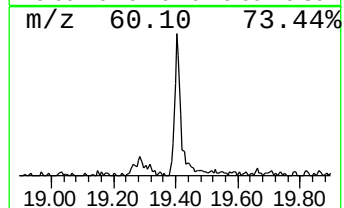
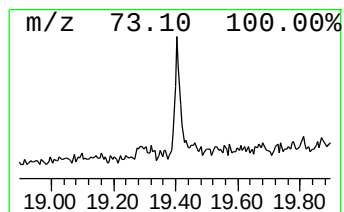
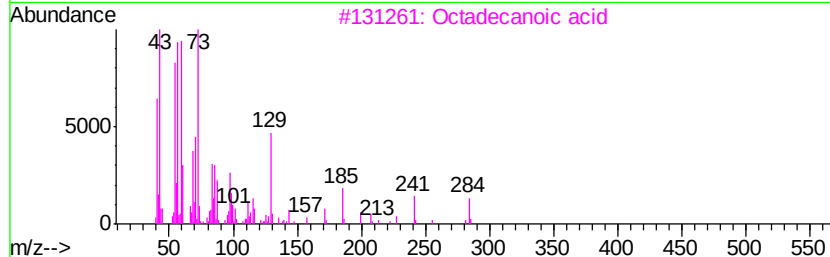
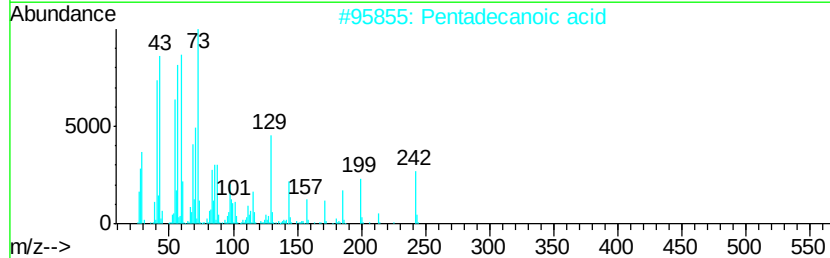
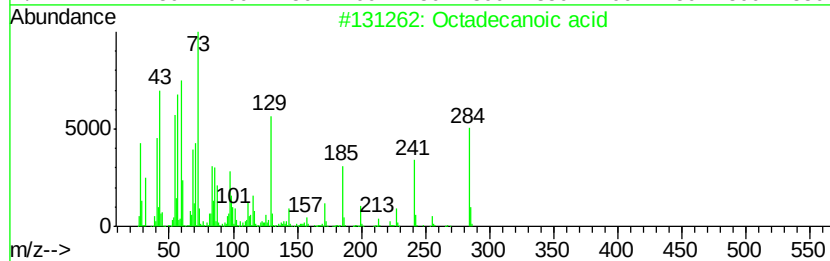
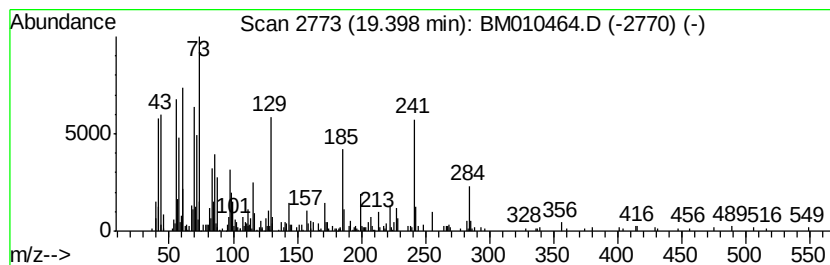
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
19.40	2.86 ng/ul	756537	Chrysene-d12	21.45		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecanoic acid	284	C18H36O2	000057-11-4	89
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3		Octadecanoic acid	284	C18H36O2	000057-11-4	64
4		n-Decanoic acid	172	C10H20O2	000334-48-5	55
5		Undecanoic acid	186	C11H22O2	000112-37-8	45



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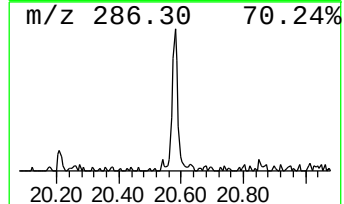
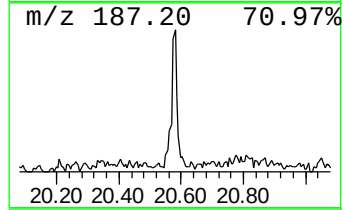
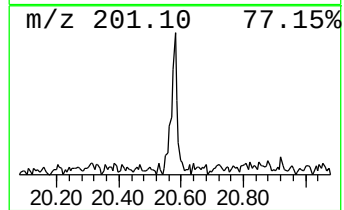
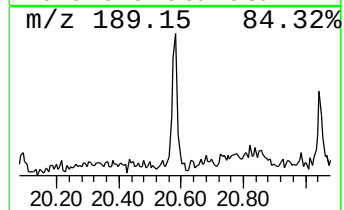
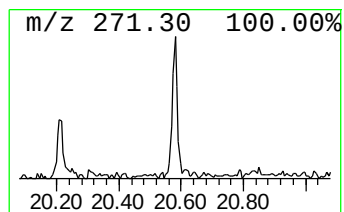
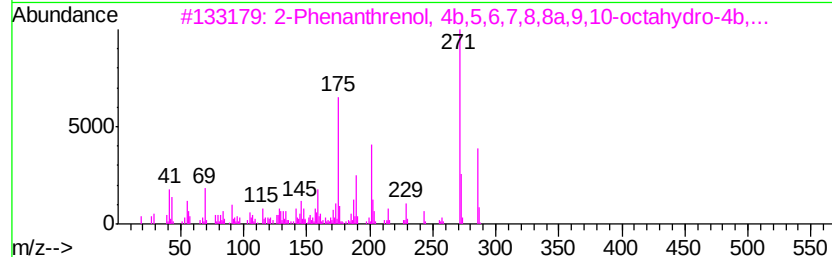
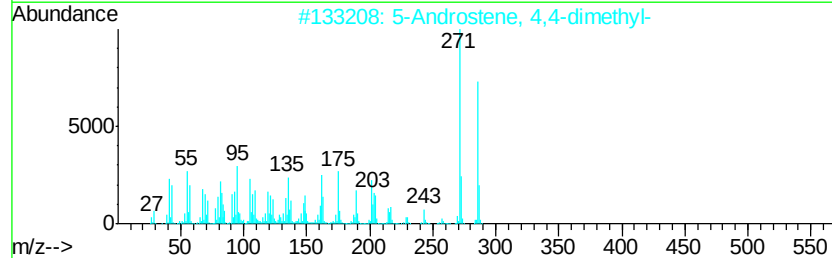
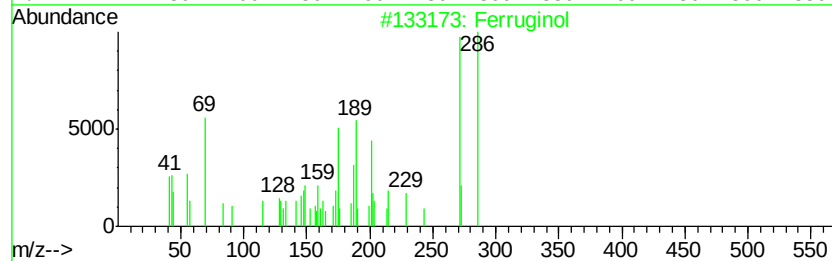
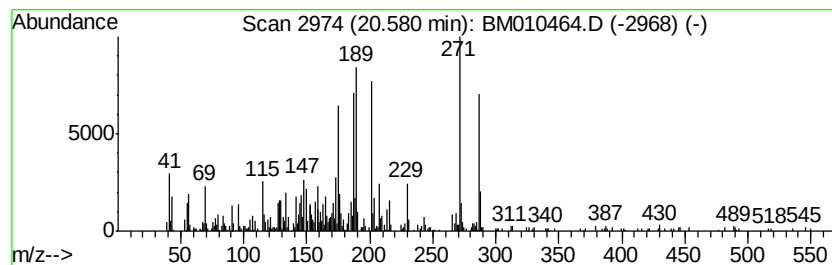
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Peak Number 4 Ferruginol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.58	2.31 ng/ul	611045	Chrysene-d12	21.45
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Ferruginol	286 C20H30O	000514-62-5 87
2		5-Androstene, 4,4-dimethyl-	286 C21H34	1000194-15-4 43
3		2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286 C20H30O	000511-15-9 30
4		Pyrrolidine, 1-[5-(1,3-benzodiox...	271 C16H17NO3	025924-78-1 30
5		5.beta.-Pregn-11-ene	286 C21H34	006673-73-0 25



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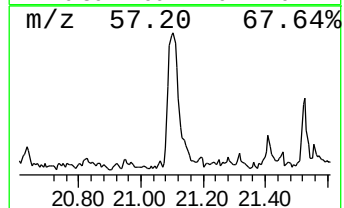
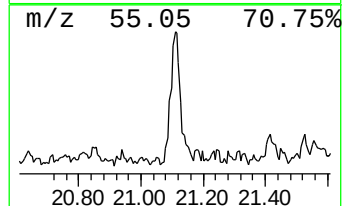
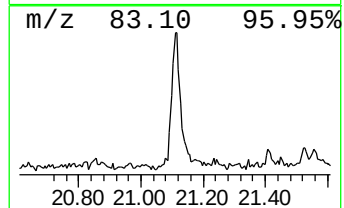
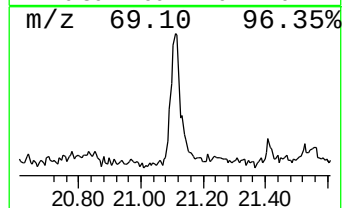
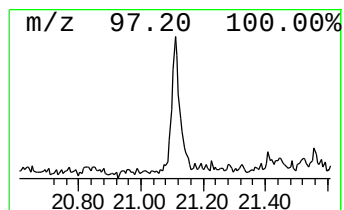
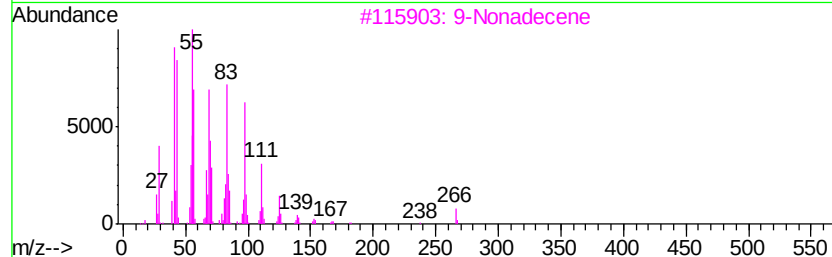
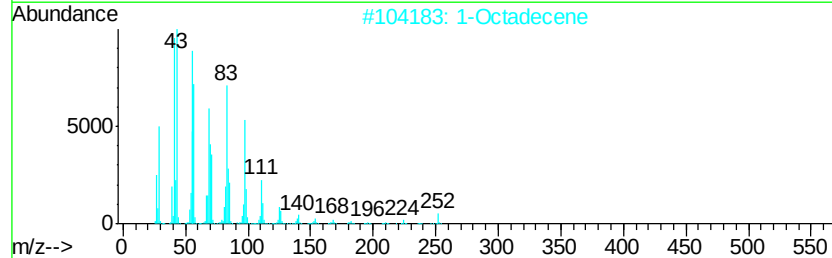
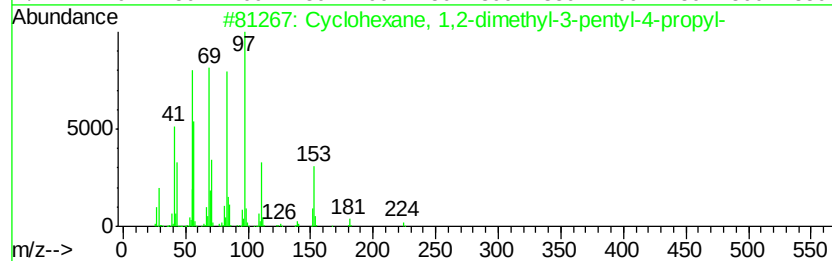
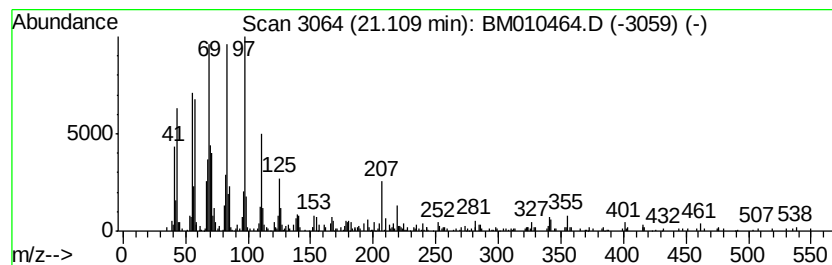
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic21.11 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.11	3.25 ng/ul	861971	Chrysene-d12	21.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,2-dimethyl-3-pent...	224	C16H32	062376-17-4	83
2		1-Octadecene	252	C18H36	000112-88-9	83
3		9-Nonadecene	266	C19H38	031035-07-1	83
4		1-Docosene	308	C22H44	001599-67-3	83
5		Cyclododecane	168	C12H24	000294-62-2	68



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Data File : BM010464.D
Acq On : 09 Jun 2017 21:28
Operator : SJ/MA
Sample : I3520-04
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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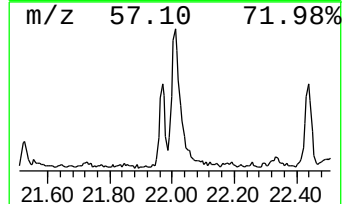
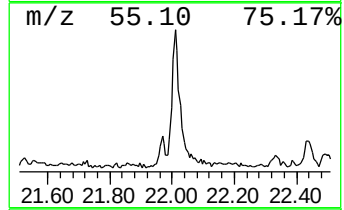
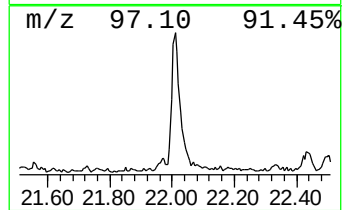
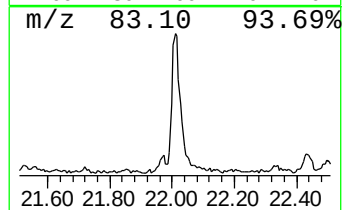
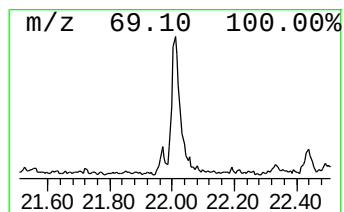
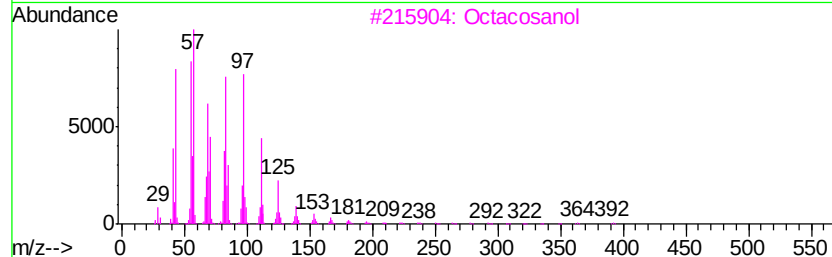
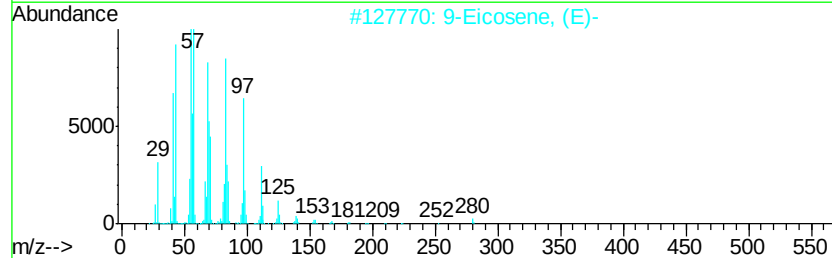
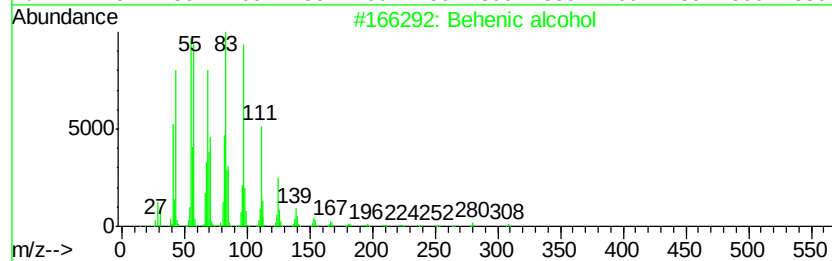
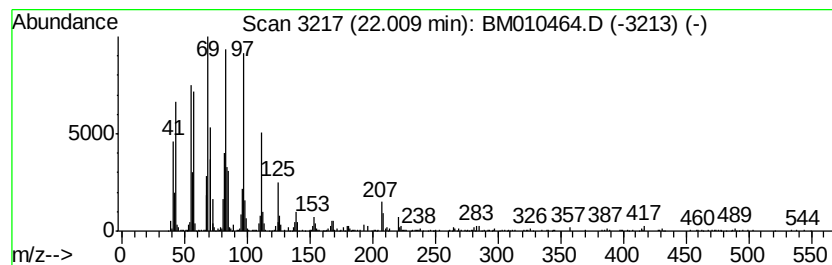
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Behenic alcohol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.01	7.89 ng/ul	2091270	Chrysene-d12	21.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Behenic alcohol	326	C22H46O	000661-19-8	90
2		9-Eicosene, (E)-	280	C20H40	074685-29-3	90
3		Octacosanol	410	C28H58O	000557-61-9	89
4		Heptacosyl acetate	438	C29H58O2	1000351-78-2	76
5		1-Docosene	308	C22H44	001599-67-3	70



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Sample : I3520-04
Misc :
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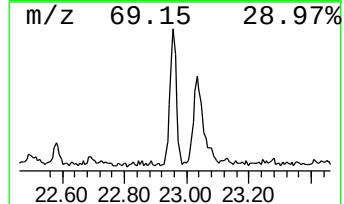
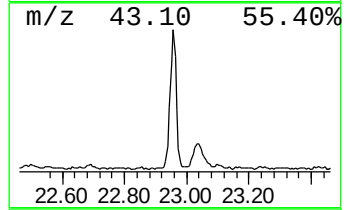
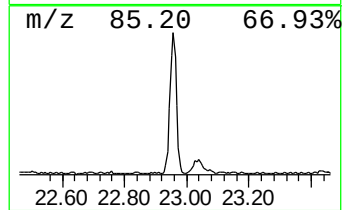
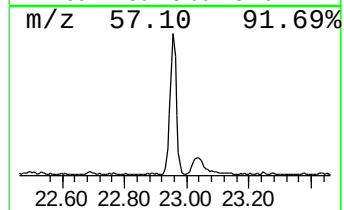
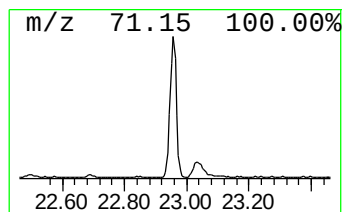
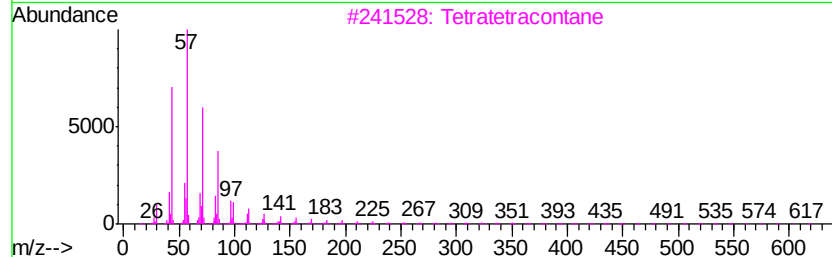
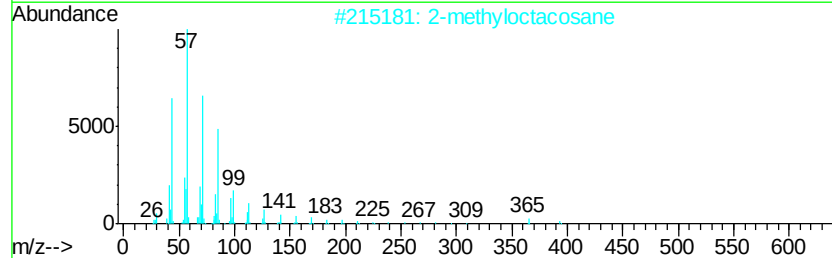
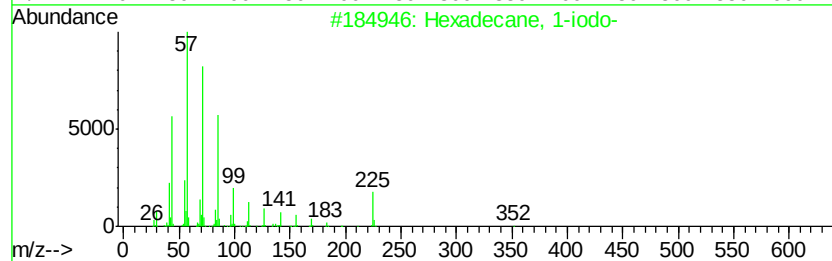
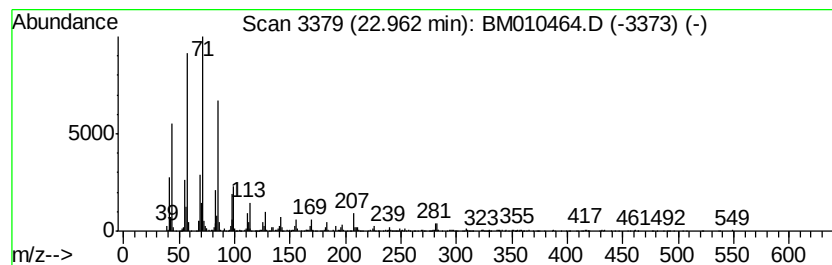
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Hexadecane, 1-iodo- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.96	12.50 ng/ul	2865750	Perylene-d12	23.79
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane, 1-iodo-	352 C16H33I	000544-77-4	96
2	2-methyloctacosane	408 C29H60	1000376-72-8	90
3	Tetratetracontane	619 C44H90	007098-22-8	90
4	Tetracosane, 11-decyl-	479 C34H70	055429-84-0	87
5	Hexacosane	366 C26H54	000630-01-3	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
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Sample : I3520-04
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Instrument :
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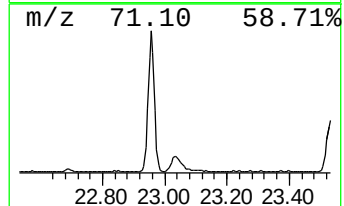
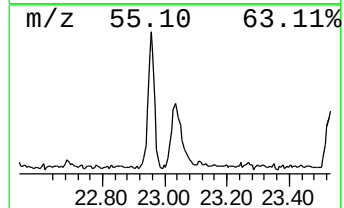
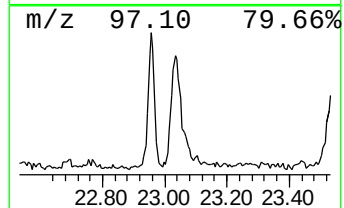
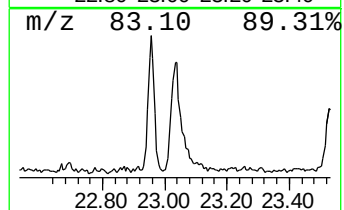
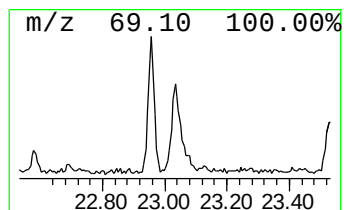
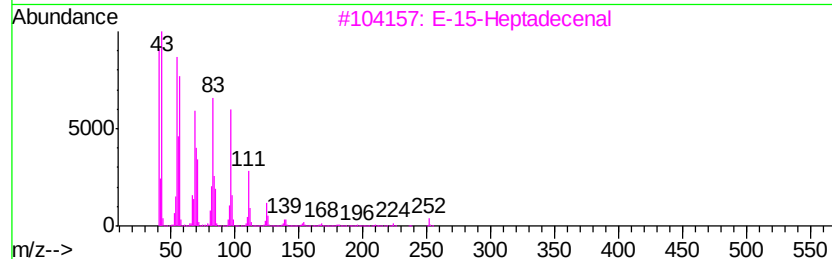
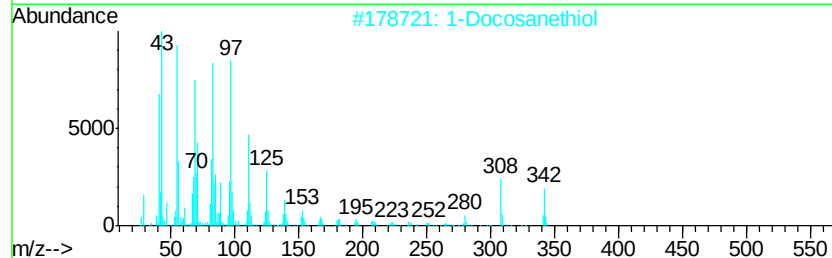
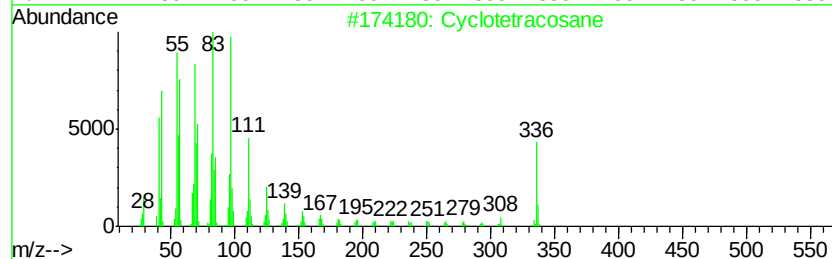
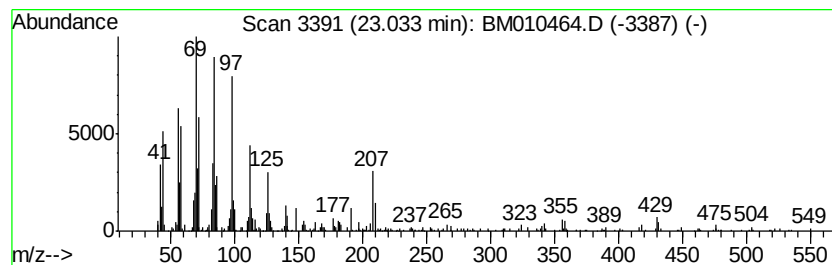
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic23.03 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.03	5.61 ng/ul	1286200	Perylene-d12	23.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetracosane	336	C24H48	000297-03-0	92
2		1-Docosanethiol	342	C22H46S	007773-83-3	87
3		E-15-Heptadecenal	252	C17H32O	1000130-97-9	86
4		4-n-Hexylthiane, S,S-dioxide	218	C11H22O2S	070928-52-8	74
5		Tetracosyl trifluoroacetate	450	C26H49F3O2	1000351-76-4	64



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
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Sample : I3520-04
Misc :
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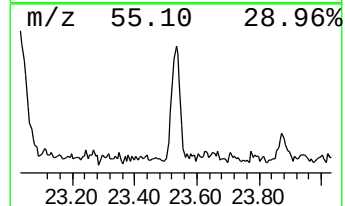
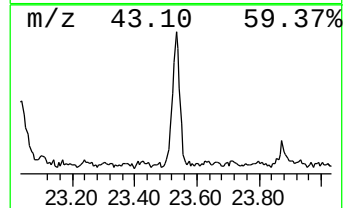
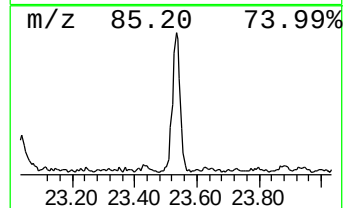
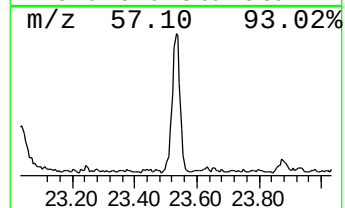
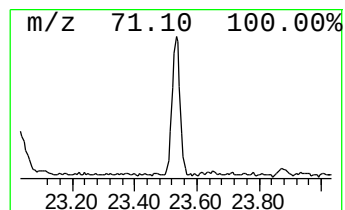
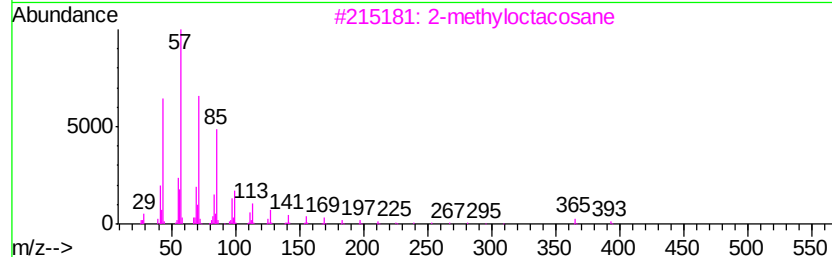
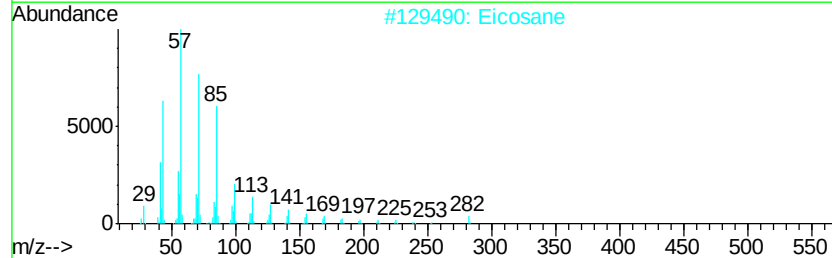
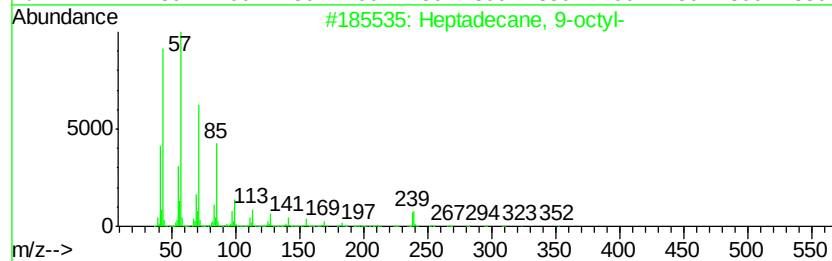
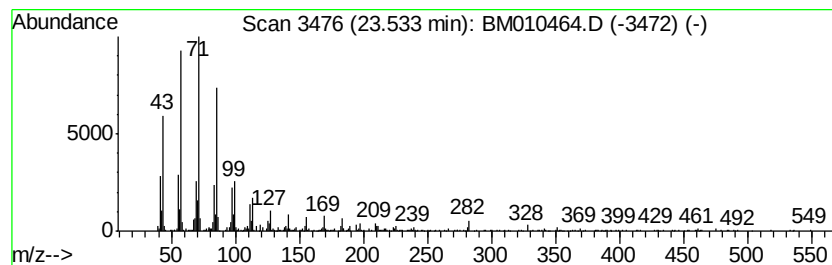
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.53	5.97 ng/ul	1368220	Perylene-d12	23.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2		Eicosane	282	C20H42	000112-95-8	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	90
4		Tetratetracontane	619	C44H90	007098-22-8	90
5		2-methylhexacosane	380	C27H56	1000376-72-7	87



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060917\
Data File : BM010464.D
Acq On : 09 Jun 2017 21:28
Operator : SJ/MA
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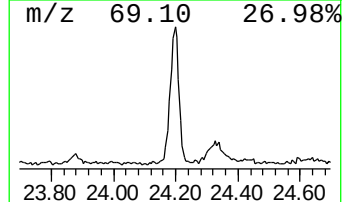
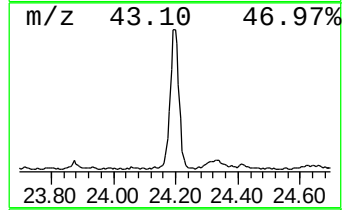
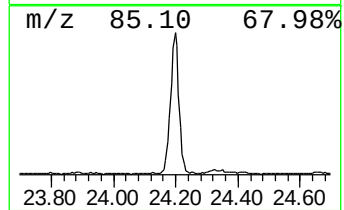
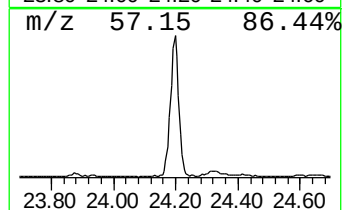
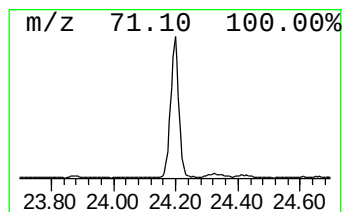
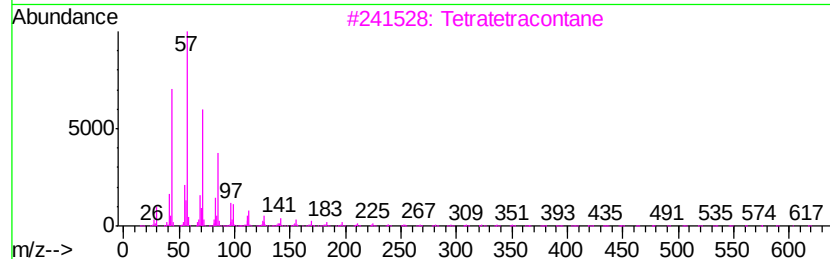
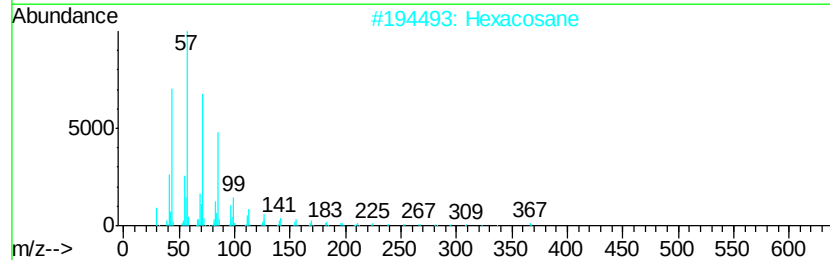
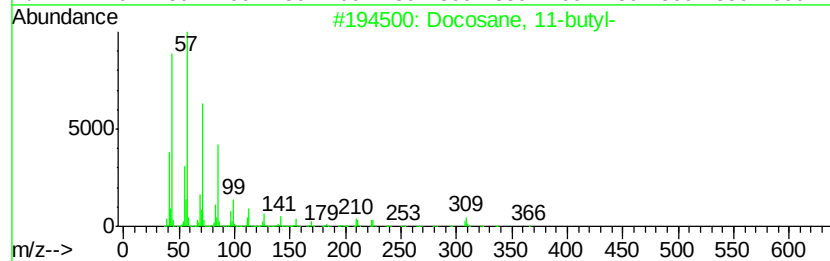
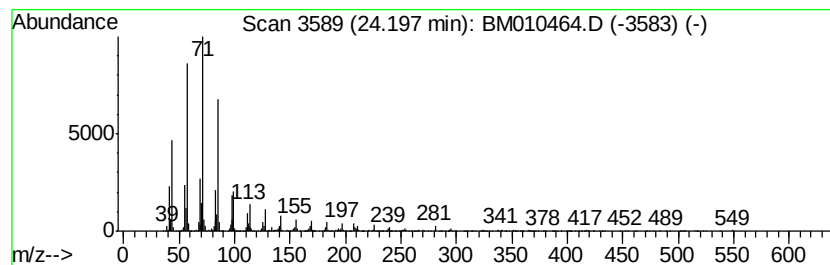
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.20	22.35 ng/ul	5123010	Perylene-d12	23.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane, 11-butyl-	366	C26H54	013475-76-8	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		Tetratetracontane	619	C44H90	007098-22-8	90
4		2-methyloctacosane	408	C29H60	1000376-72-8	90
5		Hentriacontane	437	C31H64	000630-04-6	90



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Sample : I3520-04
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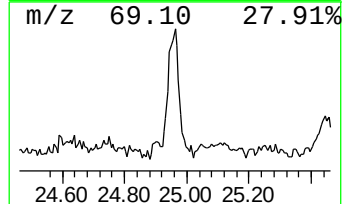
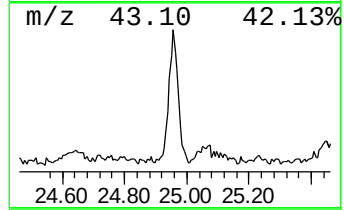
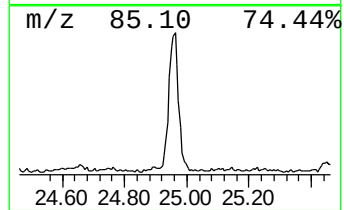
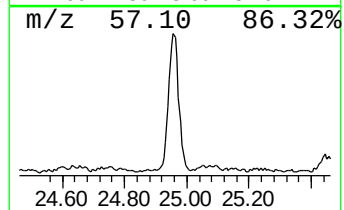
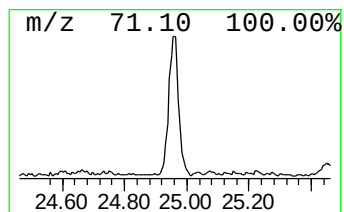
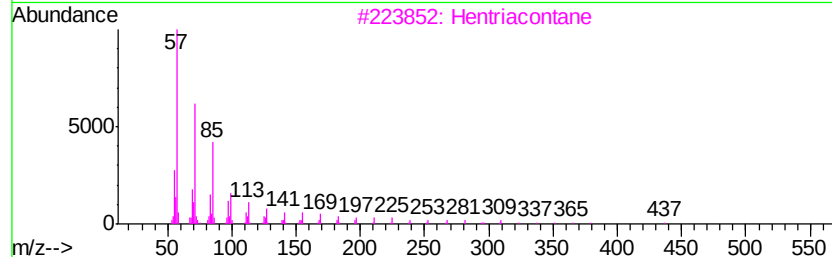
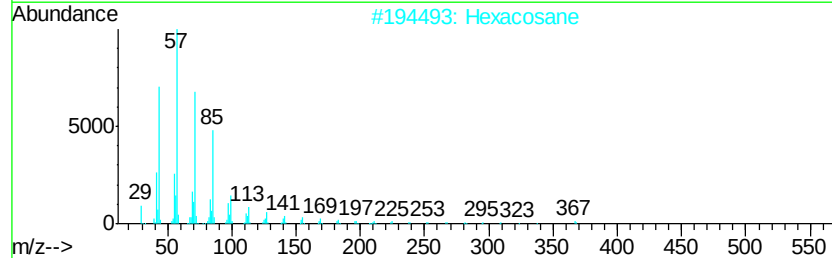
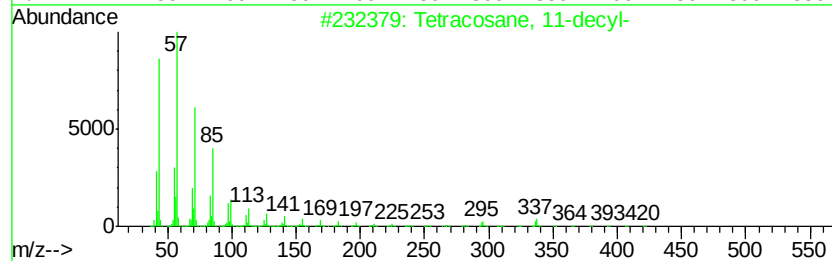
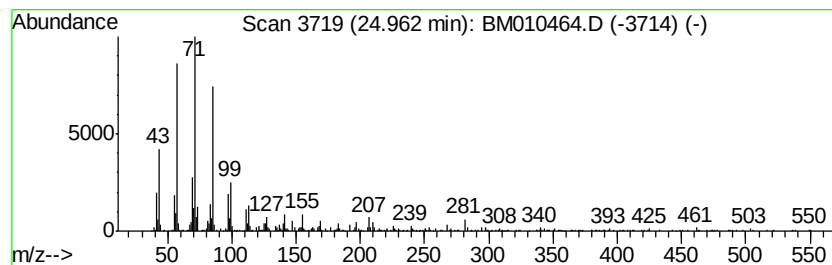
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.96	7.36 ng/ul	1687420	Perylene-d12	23.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetracosane, 11-decyl-	479	C34H70	055429-84-0	86
2		Hexacosane	366	C26H54	000630-01-3	83
3		Hentriacontane	437	C31H64	000630-04-6	83
4		Hentriacontane	437	C31H64	000630-04-6	83
5		Hexatriacontane	507	C36H74	000630-06-8	81



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Misc :
ALS Vial : 19 Sample Multiplier: 1

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

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					#	RT	Resp	Conc
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n-Hexadecanoic acid	18.13	2.5	ng/ul	596363	4	17.29	4720710	20.0
Octadecanoic acid	19.40	2.9	ng/ul	756537	5	21.45	5298010	20.0
Ferruginol	20.58	2.3	ng/ul	611045	5	21.45	5298010	20.0
(DEL) Alkane: Cyc...	21.11	3.3	ng/ul	861971	5	21.45	5298010	20.0
Behenic alcohol	22.01	7.9	ng/ul	2091270	5	21.45	5298010	20.0
Hexadecane, 1-iodo-	22.96	12.5	ng/ul	2865750	6	23.79	4583770	20.0
(DEL) Alkane: Cyc...	23.03	5.6	ng/ul	1286200	6	23.79	4583770	20.0
(DEL) Alkane: Str...	23.53	6.0	ng/ul	1368220	6	23.79	4583770	20.0
(DEL) Alkane: Str...	24.20	22.4	ng/ul	5123010	6	23.79	4583770	20.0
(DEL) Alkane: Str...	24.96	7.4	ng/ul	1687420	6	23.79	4583770	20.0