

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
 Data File : BM007401.D
 Acq On : 03 Sep 2016 11:22
 Operator : UM/SJ
 Sample : H4639-15
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BD391

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-BM082316.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	6.963	735	744	759	rBV	574915	1018149	12.13%	1.124%
2	7.128	766	772	778	rBV	427387	656687	7.82%	0.725%
3	7.328	800	806	815	rVB	573667	940221	11.20%	1.038%
4	7.792	876	885	893	rBV	749485	1262138	15.03%	1.393%
5	8.492	998	1004	1012	rBV	708328	1148784	13.68%	1.268%
6	8.945	1074	1081	1088	rBV	595502	1007005	11.99%	1.112%
7	9.663	1197	1203	1214	rBV	527102	918158	10.94%	1.014%
8	10.198	1287	1294	1303	rVB	793004	1381493	16.45%	1.525%
9	10.574	1348	1358	1364	rBV	1147041	2064668	24.59%	2.279%
10	10.627	1364	1367	1375	rVB	112389	173346	2.06%	0.191%
11	10.716	1375	1382	1392	rBV	578150	1090583	12.99%	1.204%
12	11.598	1528	1532	1539	rBV	58435	88058	1.05%	0.097%
13	11.998	1593	1600	1606	rBV	58522	91204	1.09%	0.101%
14	12.239	1635	1641	1650	rVB	145925	251439	2.99%	0.278%
15	12.463	1672	1679	1686	rBV	112379	188445	2.24%	0.208%
16	13.245	1805	1812	1820	rBV2	111678	212910	2.54%	0.235%
17	13.586	1865	1870	1876	rBV4	51306	127874	1.52%	0.141%
18	13.745	1892	1897	1901	rBV2	99700	168427	2.01%	0.186%
19	13.798	1901	1906	1907	rVV	93347	124294	1.48%	0.137%
20	13.833	1907	1912	1916	rVV	1686947	2384560	28.40%	2.632%
21	13.880	1916	1920	1928	rVB	653694	985952	11.74%	1.088%
22	14.115	1954	1960	1972	rVB	1802812	2878217	34.28%	3.177%
23	14.321	1991	1995	2002	rVB	130515	168504	2.01%	0.186%
24	14.427	2002	2013	2019	rBV2	1898731	3126452	37.24%	3.451%
25	14.627	2041	2047	2057	rBV	506512	903848	10.76%	0.998%
26	14.821	2076	2080	2086	rVB	199204	318840	3.80%	0.352%
27	14.898	2087	2093	2097	rVV	55348	87687	1.04%	0.097%
28	15.045	2113	2118	2123	rVV3	52064	96016	1.14%	0.106%
29	15.215	2142	2147	2150	rVV4	95554	168220	2.00%	0.186%
30	15.274	2153	2157	2162	rVB	171005	230671	2.75%	0.255%
31	15.415	2174	2181	2187	rBV	2426891	3632549	43.26%	4.010%
32	15.462	2187	2189	2194	rVB4	79714	102076	1.22%	0.113%
33	15.539	2194	2202	2208	rBV	711193	1044444	12.44%	1.153%
34	15.680	2216	2226	2233	rBV4	87891	237445	2.83%	0.262%

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35	15.768	2236	2241	2246	rVB	114585	195945	2.33%	0.216%
36	15.915	2258	2266	2274	rVB2	104434	269157	3.21%	0.297%
37	15.998	2274	2280	2287	rVB3	53514	132531	1.58%	0.146%
38	16.139	2299	2304	2305	rBV	276709	380927	4.54%	0.421%
39	16.709	2396	2401	2403	rBV2	93315	150868	1.80%	0.167%
40	16.745	2403	2407	2410	rVB3	77818	109411	1.30%	0.121%
41	16.821	2416	2420	2427	rVB2	215814	346703	4.13%	0.383%
42	16.927	2428	2438	2444	rVV	1012982	1582698	18.85%	1.747%
43	16.980	2445	2447	2458	rVB4	136130	244409	2.91%	0.270%
44	17.168	2473	2479	2483	rBV	2812659	4062478	48.38%	4.485%
45	17.209	2483	2486	2490	rVB	874169	1138160	13.56%	1.256%
46	17.268	2490	2496	2500	rBV	2785031	3987595	47.49%	4.402%
47	17.351	2506	2510	2515	rBV4	65386	115795	1.38%	0.128%
48	17.480	2529	2532	2538	rVB2	129338	191177	2.28%	0.211%
49	17.545	2539	2543	2556	rBV3	503768	971951	11.58%	1.073%
50	17.662	2559	2563	2566	rVV	190697	222009	2.64%	0.245%
51	17.745	2574	2577	2581	rVB	109294	139435	1.66%	0.154%
52	17.845	2590	2594	2599	rBV5	50505	86994	1.04%	0.096%
53	18.056	2623	2630	2633	rBV2	436969	898304	10.70%	0.992%
54	18.086	2633	2635	2640	rVB	474924	615905	7.34%	0.680%
55	18.227	2655	2659	2664	rBV2	255626	429818	5.12%	0.475%
56	18.274	2664	2667	2673	rVB	202313	265582	3.16%	0.293%
57	18.356	2676	2681	2684	rBV2	276872	362873	4.32%	0.401%
58	18.392	2684	2687	2691	rVV4	121207	176895	2.11%	0.195%
59	18.433	2691	2694	2699	rVB4	110141	152273	1.81%	0.168%
60	18.545	2709	2713	2716	rBV	315716	420200	5.00%	0.464%
61	18.586	2716	2720	2727	rVB3	328071	497057	5.92%	0.549%
62	18.762	2747	2750	2751	rBV3	86316	93316	1.11%	0.103%
63	18.786	2752	2754	2758	rVV3	156901	208198	2.48%	0.230%
64	18.839	2760	2763	2766	rVV	151608	212646	2.53%	0.235%
65	18.880	2767	2770	2773	rVB	156420	186772	2.22%	0.206%
66	18.962	2780	2784	2786	rBV	317311	463509	5.52%	0.512%
67	19.056	2797	2800	2803	rVB	151858	172403	2.05%	0.190%
68	19.103	2804	2808	2815	rVV2	287546	437660	5.21%	0.483%
69	19.227	2824	2829	2835	rVB	2298196	3118733	37.14%	3.443%
70	19.286	2836	2839	2843	rBV	161658	226629	2.70%	0.250%
71	19.562	2880	2886	2889	rBV	4439851	6609811	78.72%	7.297%

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72	19.592	2889	2891	2899	rVB	1697168	1964810	23.40%	2.169%
73	19.662	2899	2903	2907	rBV2	271944	389103	4.63%	0.430%
74	19.909	2941	2945	2953	rBV5	234164	622197	7.41%	0.687%
75	20.386	3023	3026	3029	rBV	154082	229054	2.73%	0.253%
76	20.597	3059	3062	3065	rVB	274355	308910	3.68%	0.341%
77	20.833	3098	3102	3108	rBV	633720	892665	10.63%	0.985%
78	21.056	3137	3140	3148	rVB3	878950	1494730	17.80%	1.650%
79	21.127	3149	3152	3158	rVB	438512	607492	7.24%	0.671%
80	21.350	3183	3190	3194	rBV2	4695462	8396506	100.00%	9.269%
81	21.386	3194	3196	3203	rVB	1583446	2018747	24.04%	2.229%
82	22.938	3455	3460	3465	rBV	1838734	3290260	39.19%	3.632%
83	23.427	3538	3543	3547	rBV	1084577	1974603	23.52%	2.180%
84	23.480	3547	3552	3557	rVV2	2501298	4328380	51.55%	4.778%
85	23.621	3570	3576	3582	rBV	2537575	4909645	58.47%	5.420%

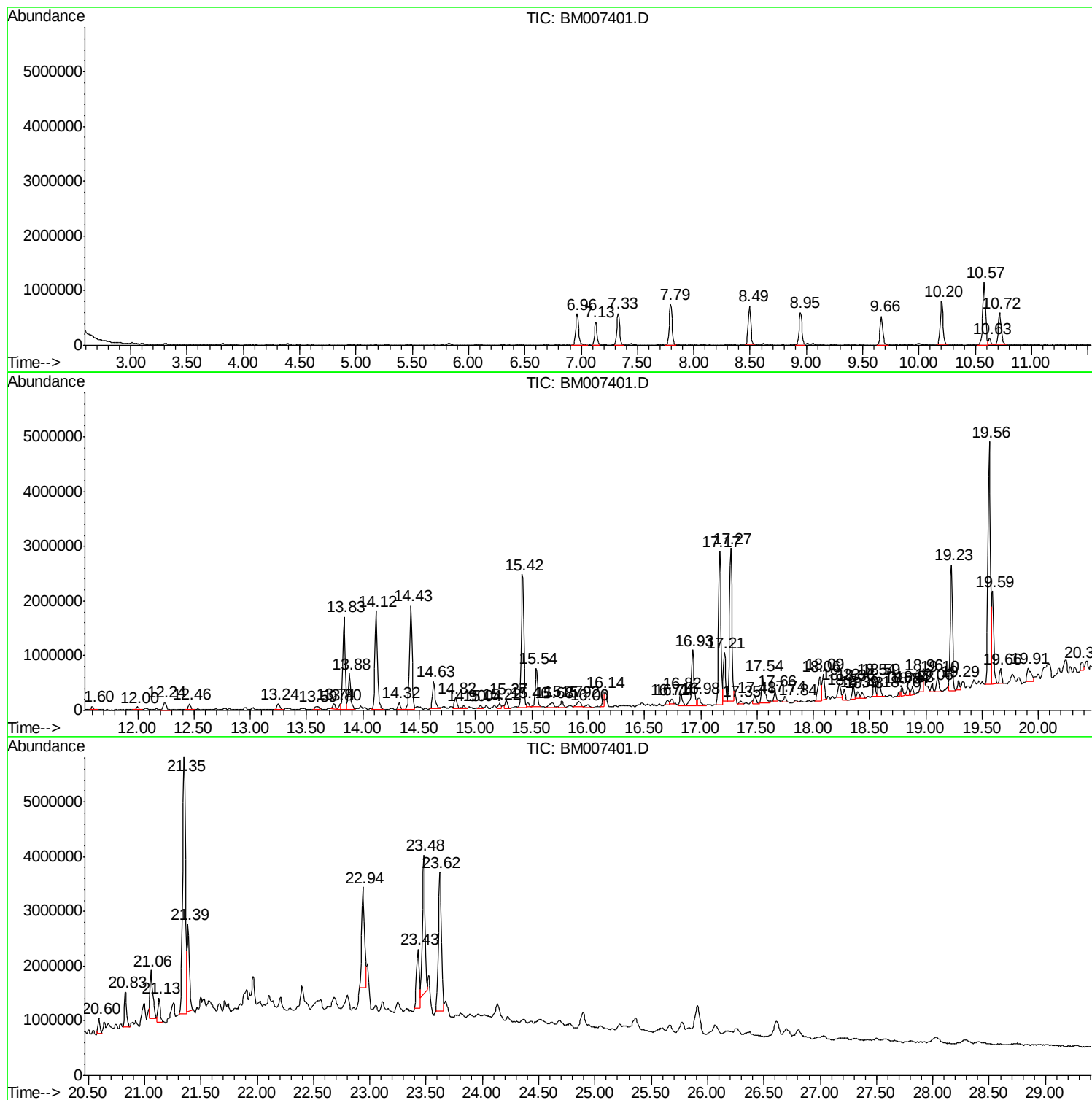
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION

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



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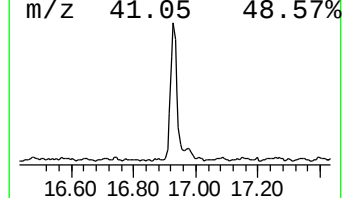
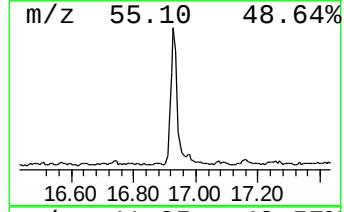
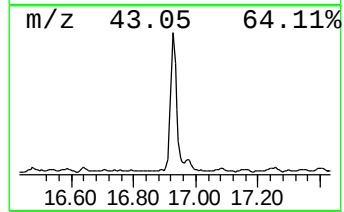
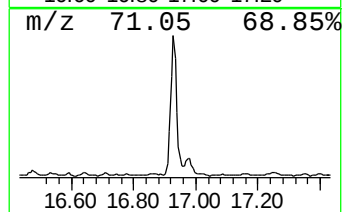
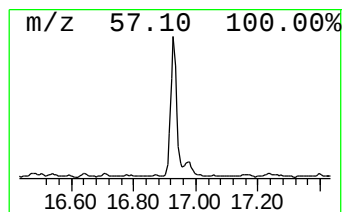
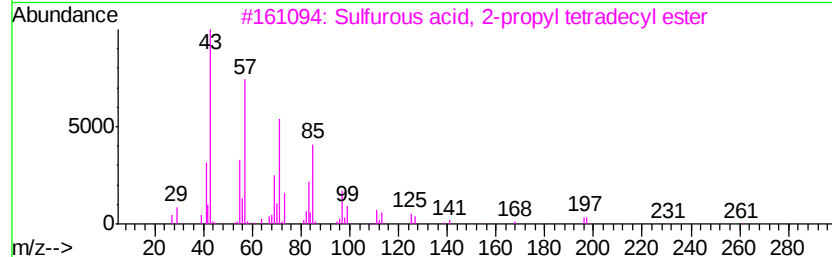
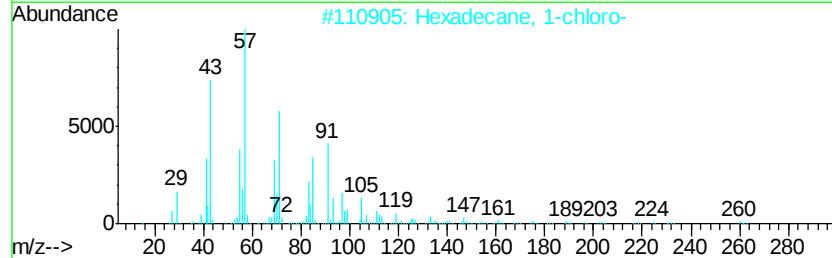
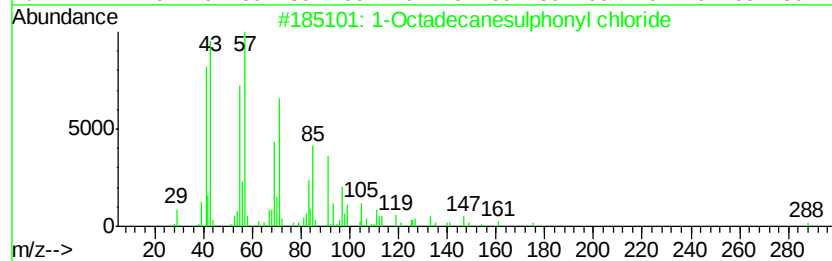
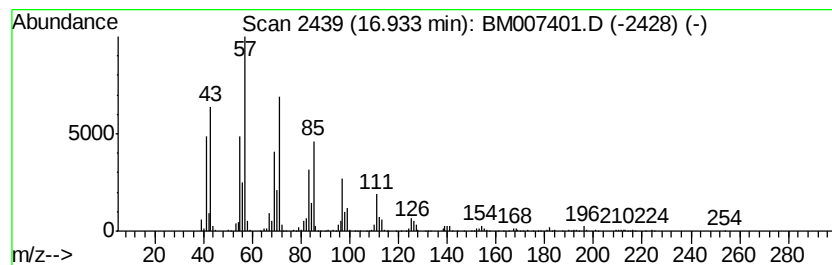
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-Octadecanesulphonyl chloride Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.93	7.79 ng/ul	1582700	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	91
2		Hexadecane, 1-chloro-	260	C16H33Cl	004860-03-1	90
3		Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1000309-12-5	86
4		1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	83
5		1-Decanol, 2-octyl-	270	C18H38O	045235-48-1	83



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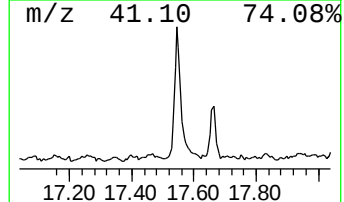
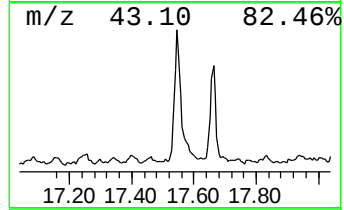
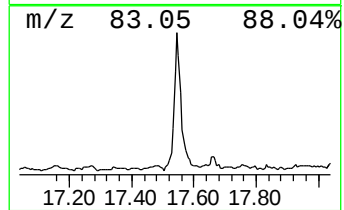
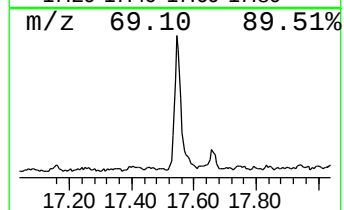
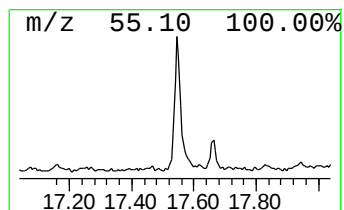
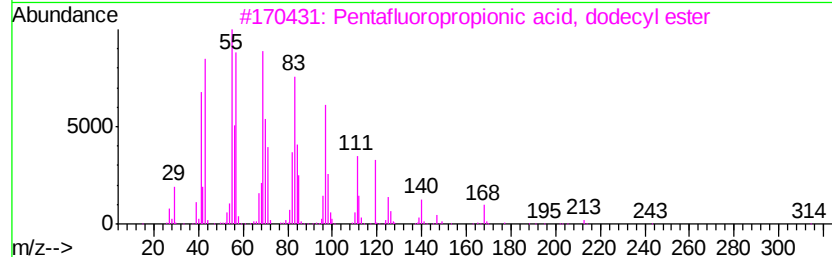
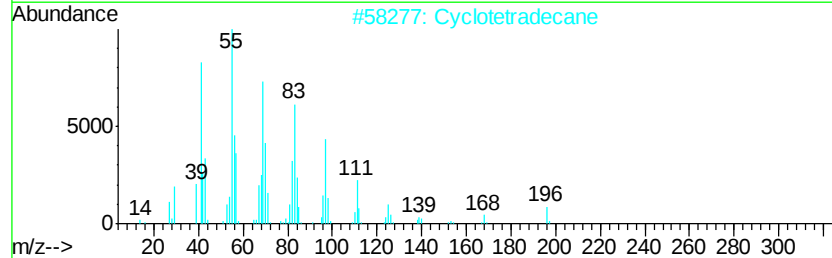
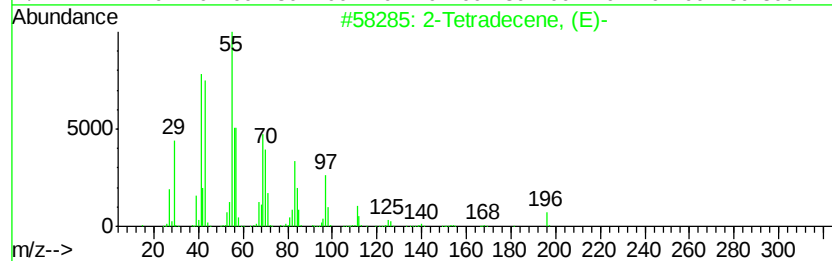
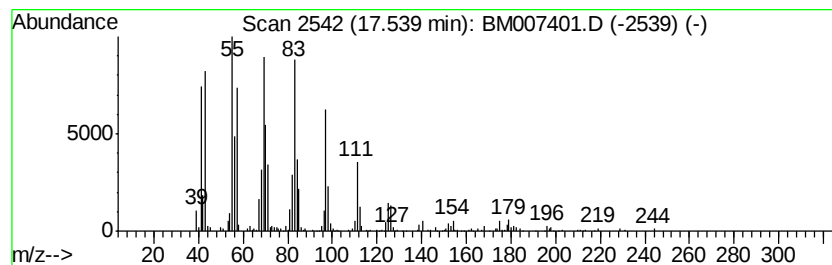
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 (DEL) Alkane: Cyclic17.54 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.54	4.79 ng/ul	971951	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Tetradecene, (E)-	196	C14H28	035953-54-9	97
2		Cyclotetradecane	196	C14H28	000295-17-0	94
3		Pentafluoropropionic acid, dodecyl...	332	C15H25F5O2	006222-04-4	94
4		Acetic acid, trifluoro-, dodecyl...	282	C14H25F3O2	006222-00-0	94
5		5-Octadecene, (E)-	252	C18H36	007206-21-5	91



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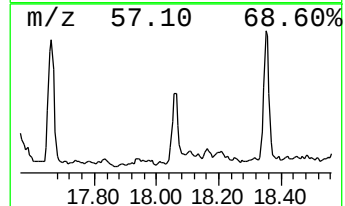
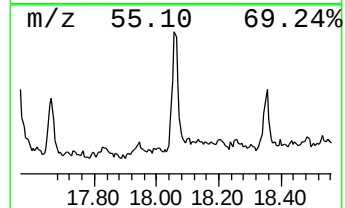
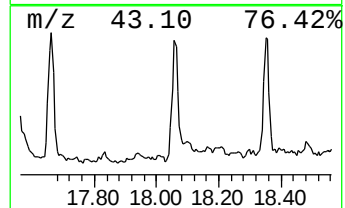
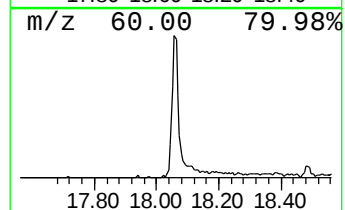
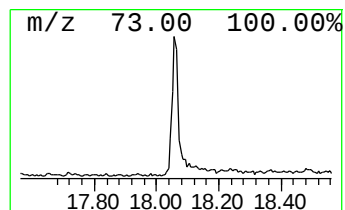
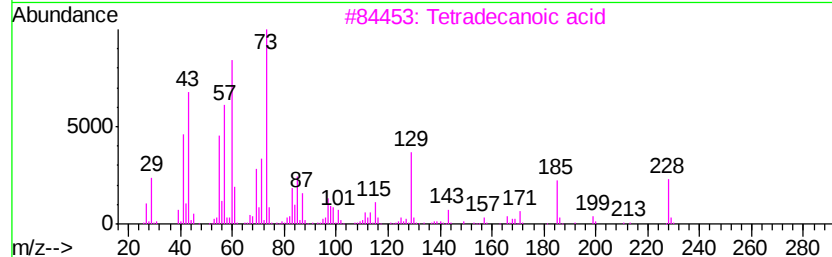
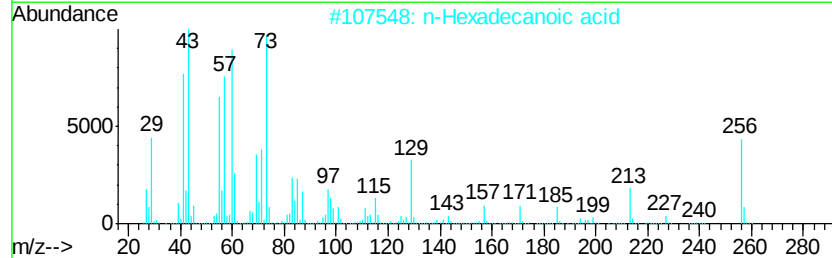
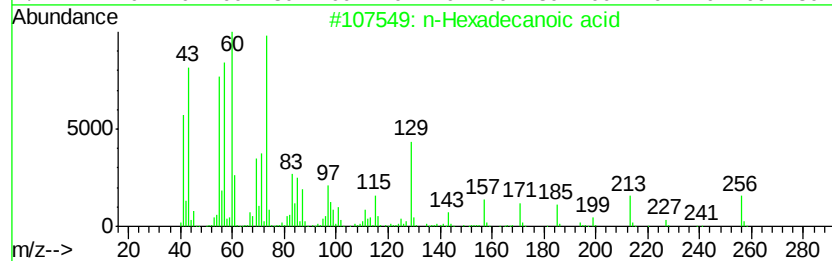
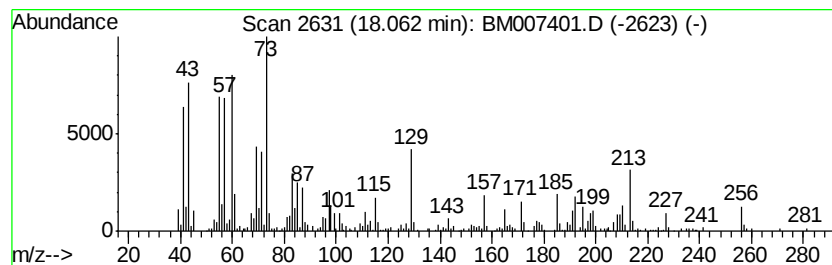
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.06	4.42 ng/ul	898304	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		Tridecanoic acid	214	C13H26O2	000638-53-9	64
5		Octadecanoic acid	284	C18H36O2	000057-11-4	64



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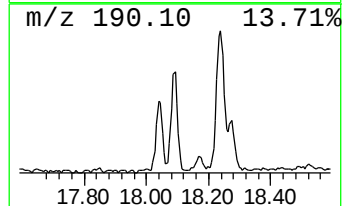
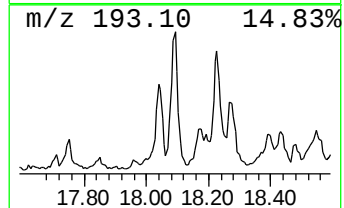
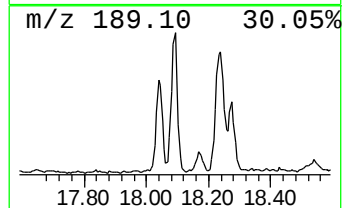
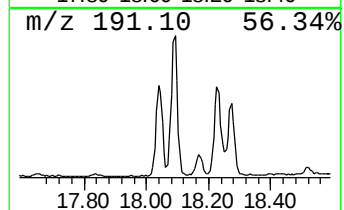
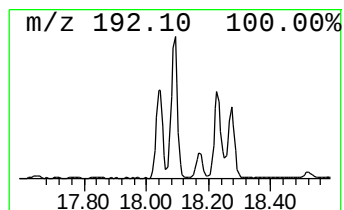
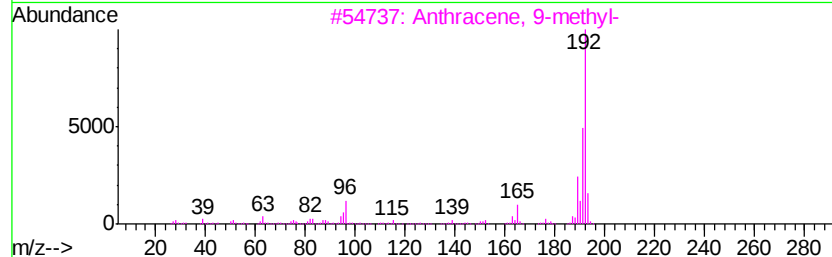
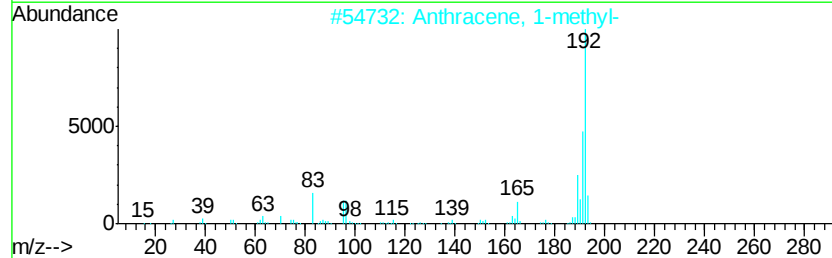
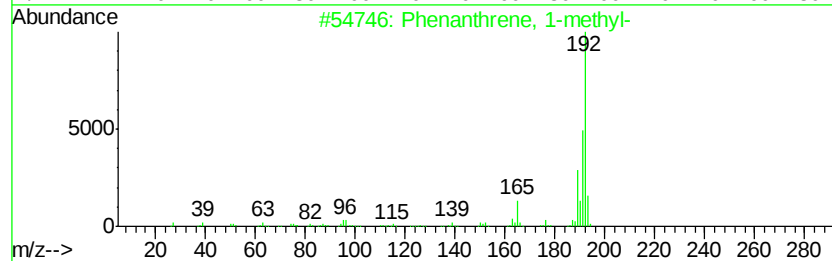
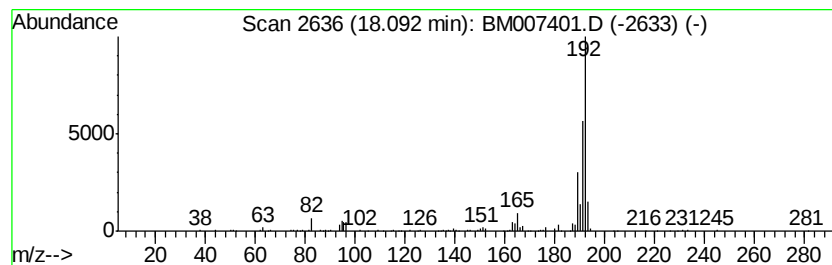
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.09	3.03 ng/ul	615905	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	90
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	90



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
Operator : UM/SJ
Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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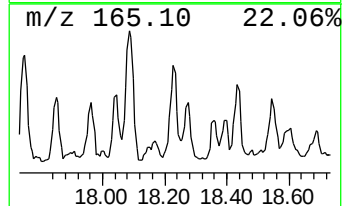
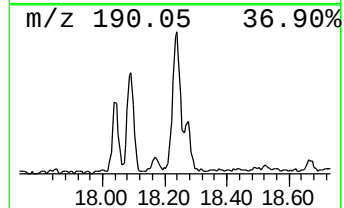
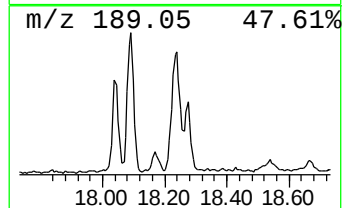
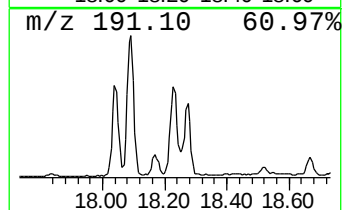
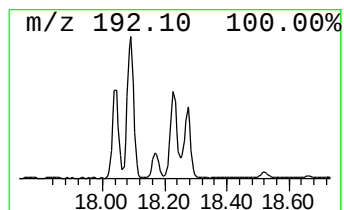
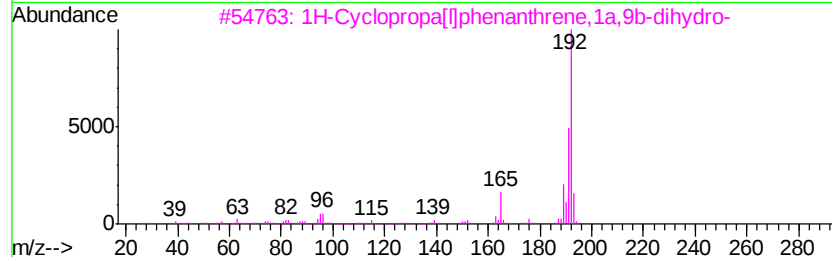
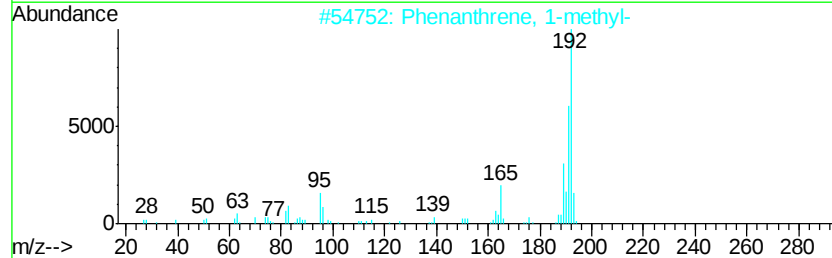
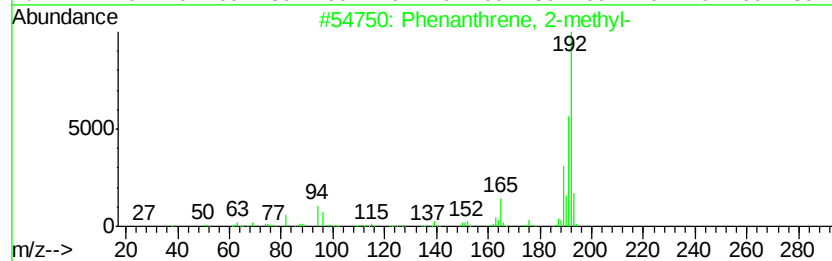
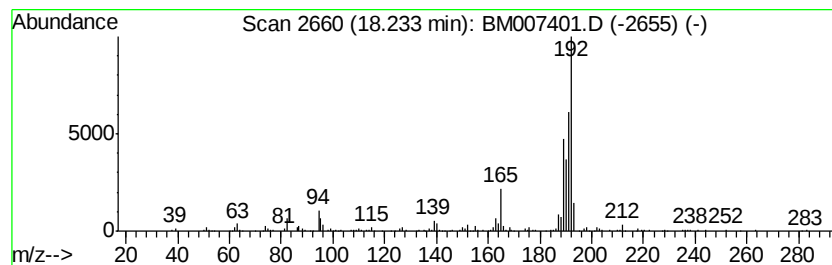
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.23	2.12 ng/ul	429818	Phenanthrene-d10	17.17

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	89
3		1H-Cyclopropa[1]phenanthrene, 1a, ...	192	C15H12	000949-41-7	81
4		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	81
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	81



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
Operator : UM/SJ
Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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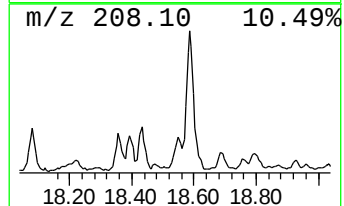
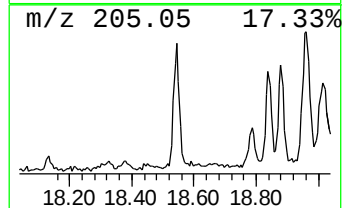
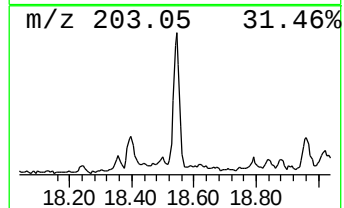
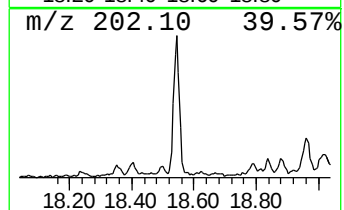
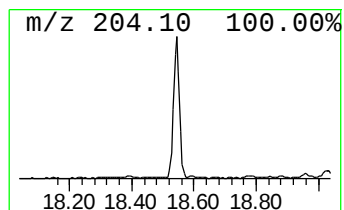
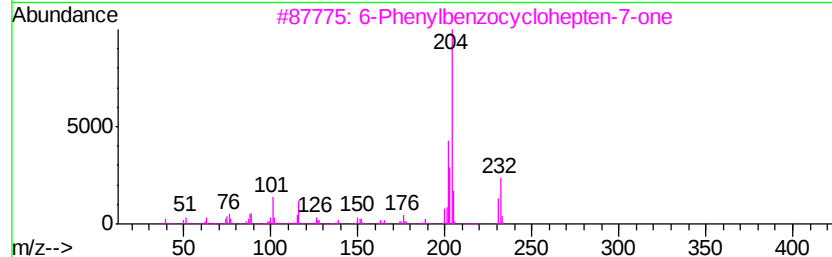
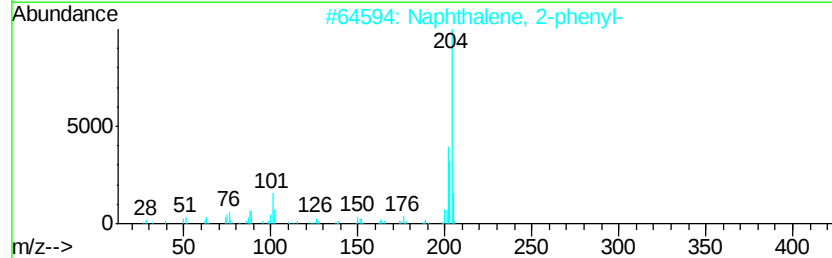
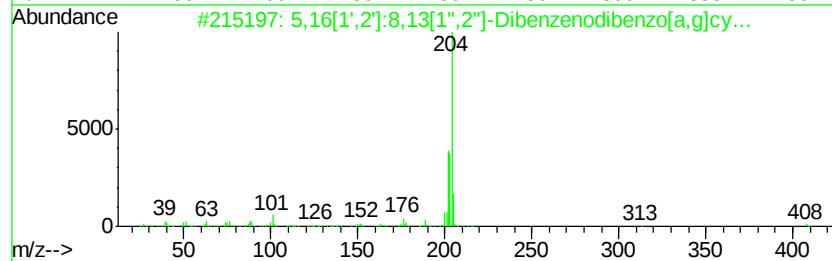
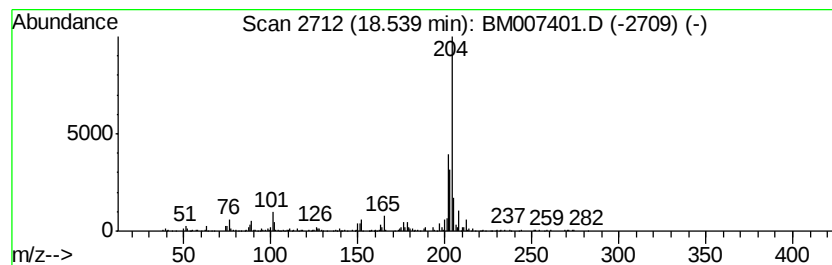
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.54	2.07 ng/ul	420200	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	83
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
3		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	64
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	58
5		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	53



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
Operator : UM/SJ
Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

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ClientSampled :
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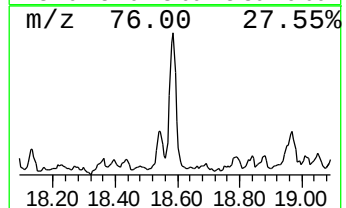
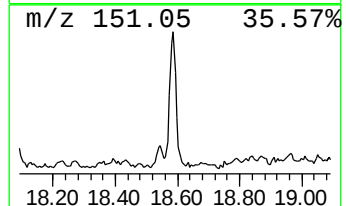
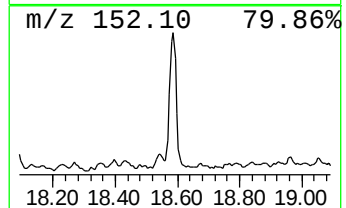
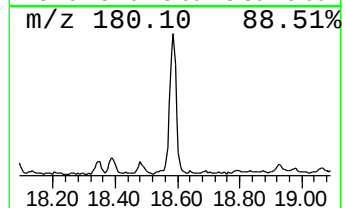
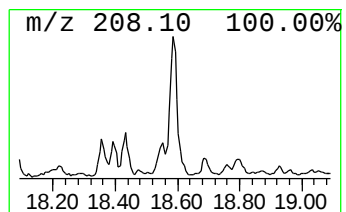
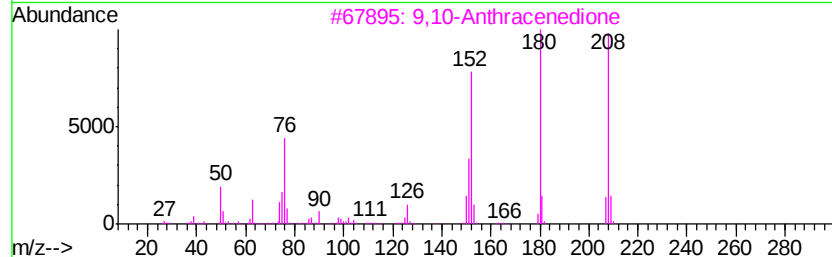
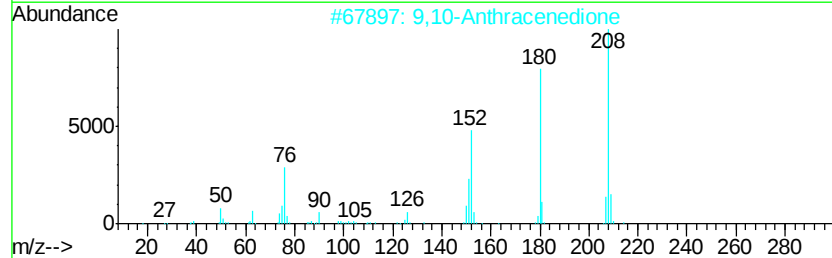
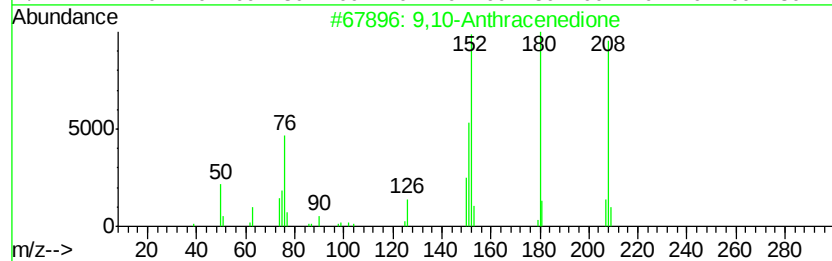
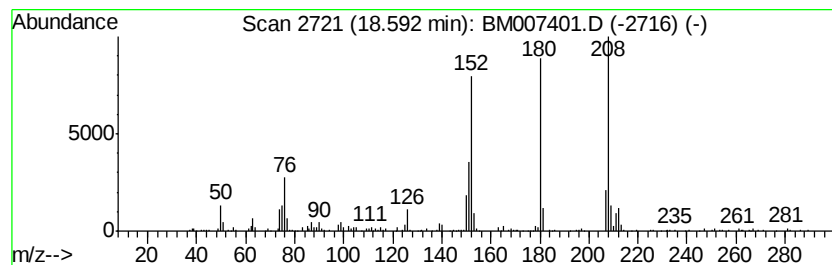
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.59	2.45 ng/ul	497057	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9,10-Anthracenedione	208	C14H8O2	000084-65-1	99
2		9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
3		9,10-Anthracenedione	208	C14H8O2	000084-65-1	91
4		Anthraquinone-1-carboxylic acid	252	C15H8O4	000602-69-7	86
5		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
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Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

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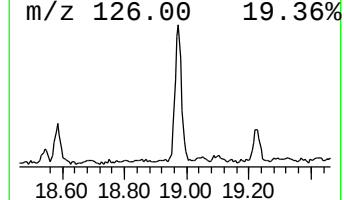
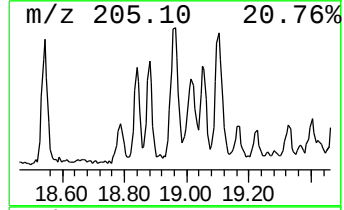
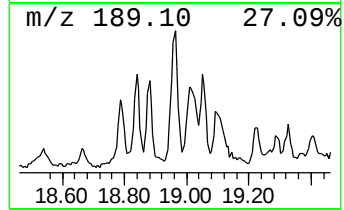
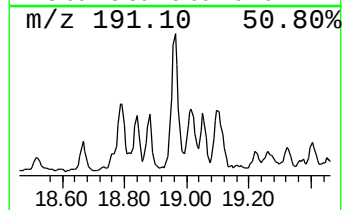
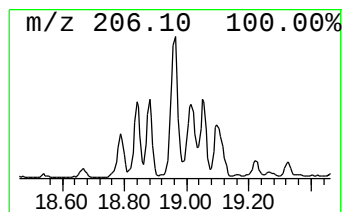
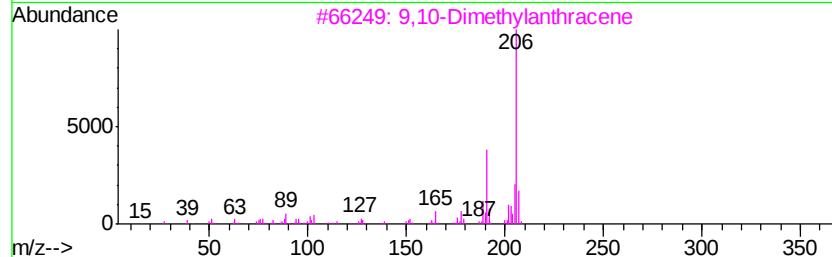
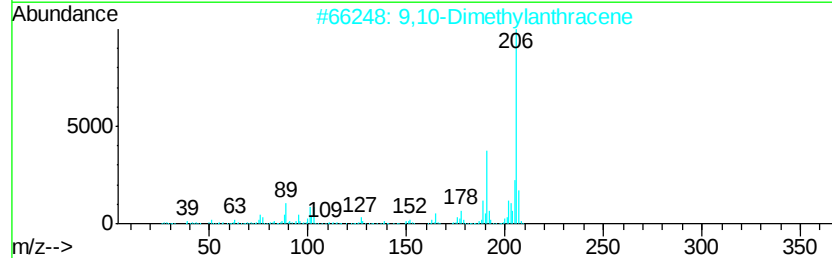
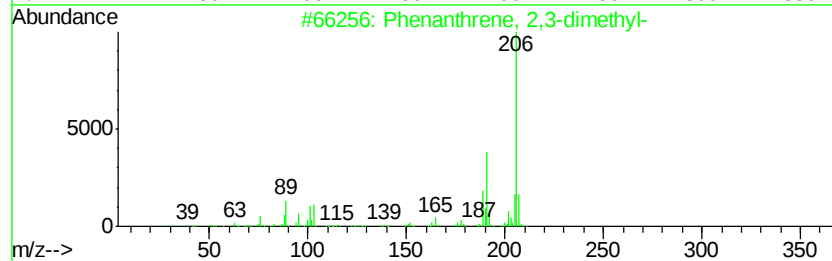
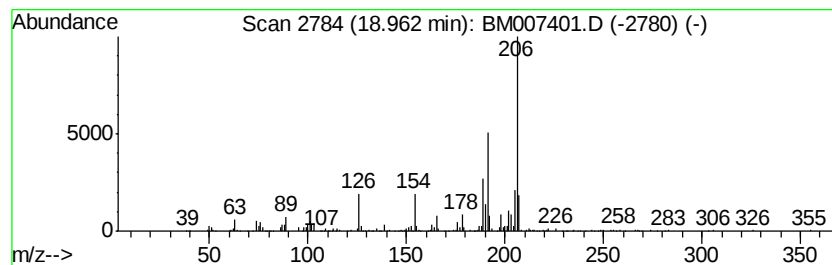
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenanthrene, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.96	2.28 ng/ul	463509	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	89
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	86
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	78
4		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	76
5		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	70



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
Operator : UM/SJ
Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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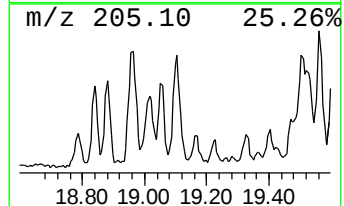
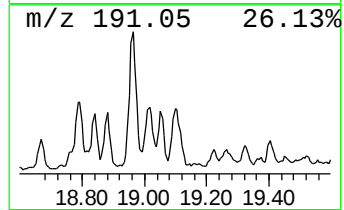
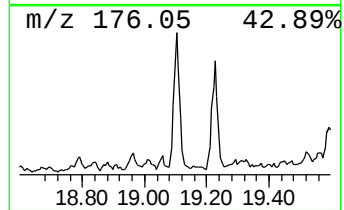
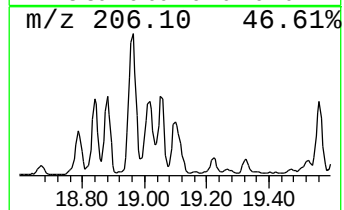
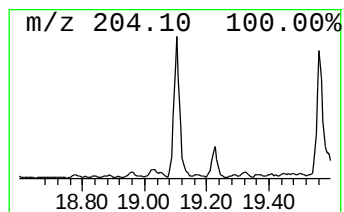
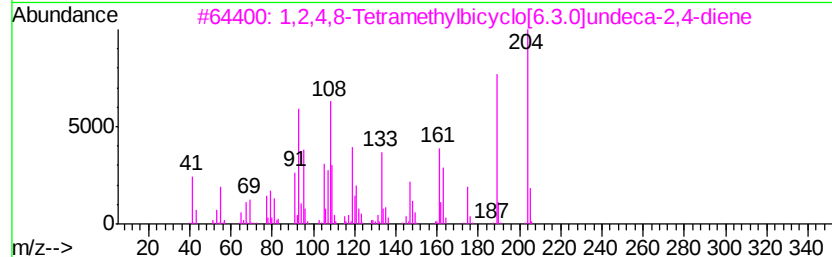
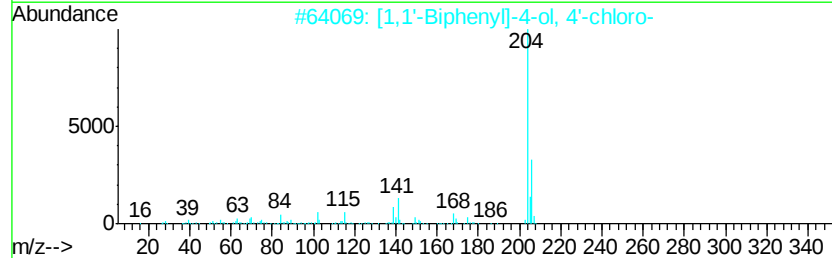
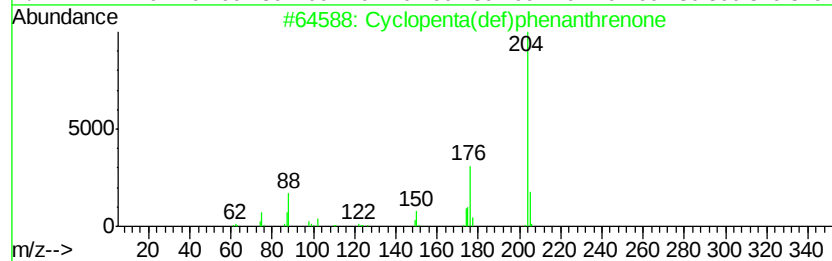
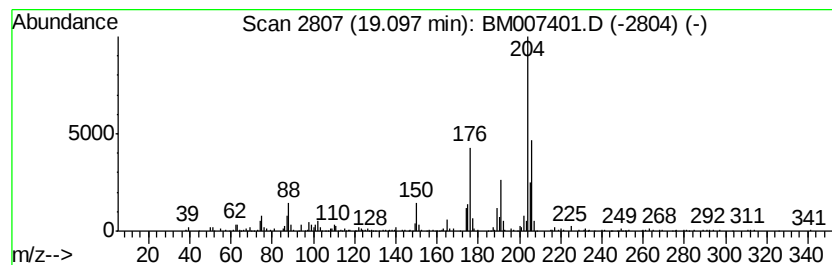
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cyclopenta(def)phenanthrenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	2.15 ng/ul	437660	Phenanthrene-d10	17.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2		[1,1'-Biphenyl]-4-ol, 4'-chloro-	204	C12H9ClO	028034-99-3	43
3		1,2,4,8-Tetramethylbicyclo[6.3.0...]	204	C15H24	137235-51-9	41
4		2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	38
5		2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	38



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
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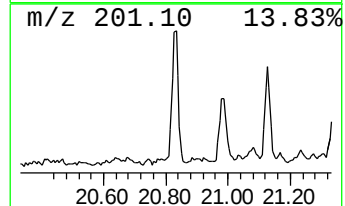
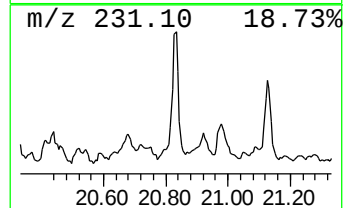
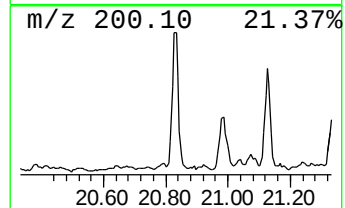
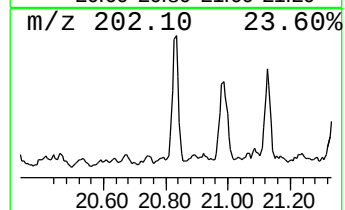
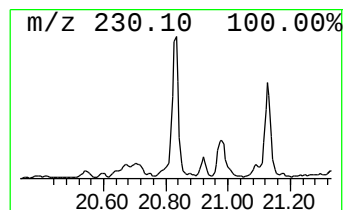
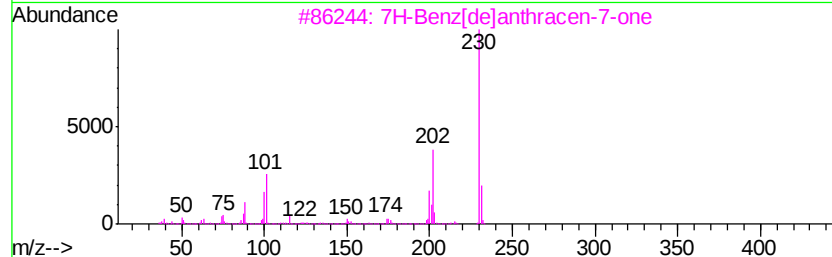
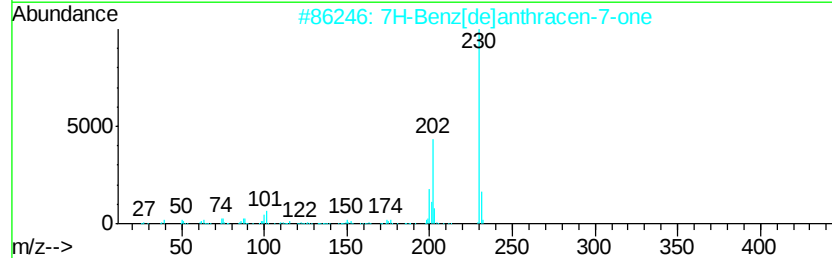
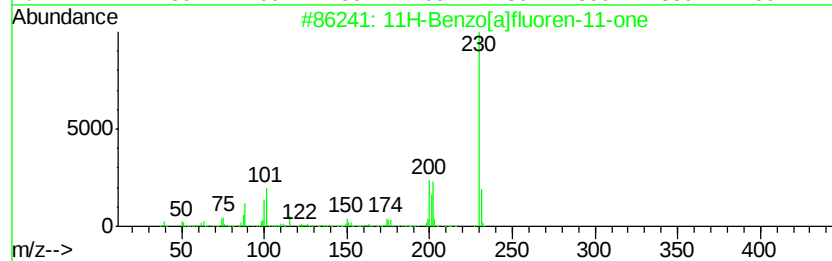
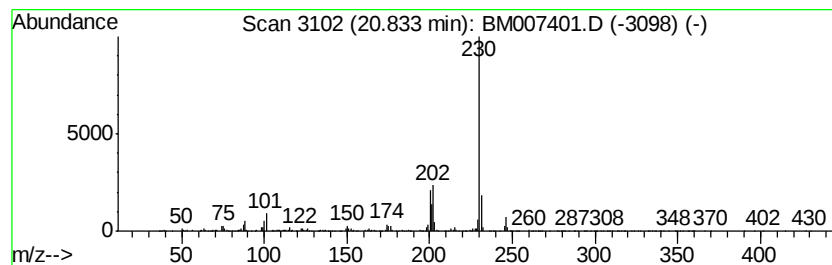
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluoren-11-one Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.83	2.13 ng/ul	892665	Chrysene-d12	21.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
4		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
5		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	81



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
Operator : UM/SJ
Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BD391

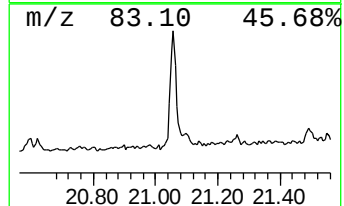
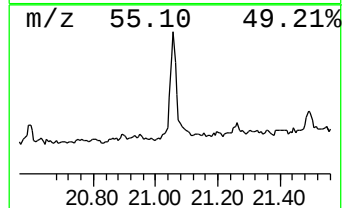
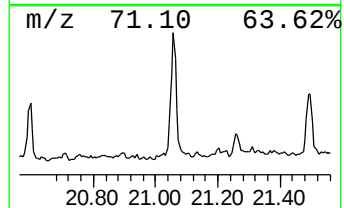
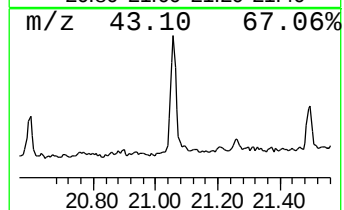
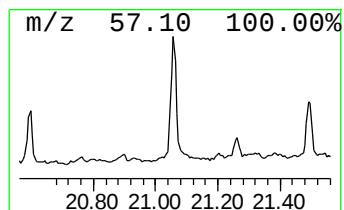
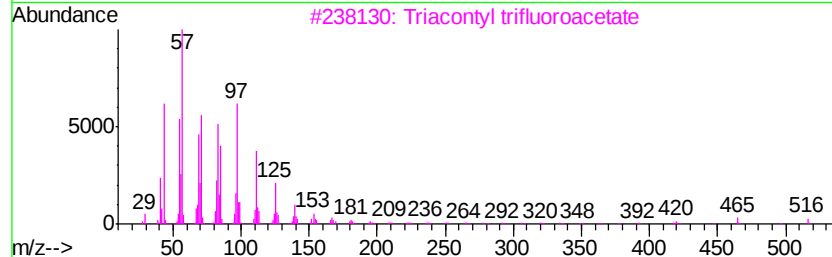
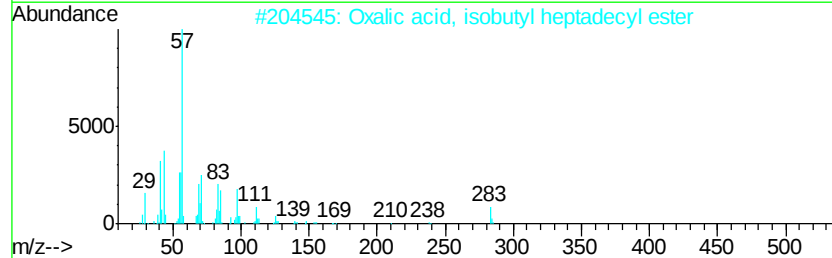
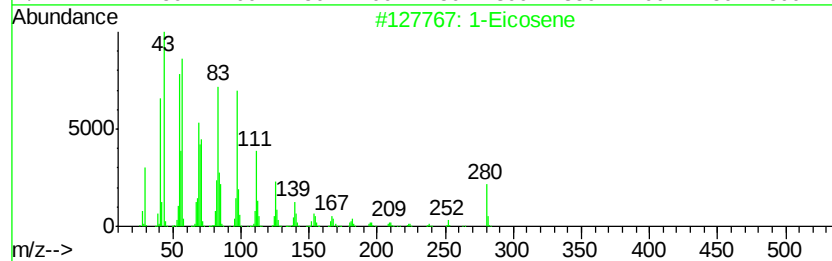
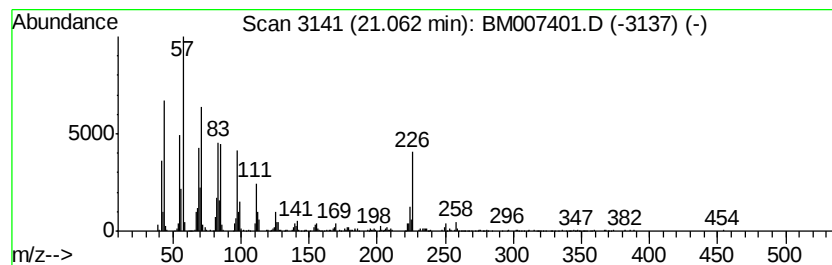
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic21.06 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.06	3.56 ng/ul	1494730	Chrysene-d12	21.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Eicosene	280	C20H40	003452-07-1	92
2		Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	91
3		Triacetyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	91
4		Oxalic acid, isobutyl pentadecyl...	356	C21H40O4	1000309-38-0	91
5		Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM090316\
Data File : BM007401.D
Acq On : 03 Sep 2016 11:22
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Sample : H4639-15
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BD391

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Octadecanesulph...	16.93	7.8	ng/ul	1582700	4	17.17	4062480	20.0
(DEL) Alkane: Cyc...	17.54	4.8	ng/ul	971951	4	17.17	4062480	20.0
n-Hexadecanoic acid	18.06	4.4	ng/ul	898304	4	17.17	4062480	20.0
Phenanthrene, 1-m...	18.09	3.0	ng/ul	615905	4	17.17	4062480	20.0
Phenanthrene, 2-m...	18.23	2.1	ng/ul	429818	4	17.17	4062480	20.0
5,16[1',2']:8,13[...	18.54	2.1	ng/ul	420200	4	17.17	4062480	20.0
9,10-Anthracenedione	18.59	2.5	ng/ul	497057	4	17.17	4062480	20.0
Phenanthrene, 2,3...	18.96	2.3	ng/ul	463509	4	17.17	4062480	20.0
Cyclopenta(def)ph...	19.10	2.1	ng/ul	437660	4	17.17	4062480	20.0
11H-Benzo[a]fluor...	20.83	2.1	ng/ul	892665	5	21.35	8396510	20.0
(DEL) Alkane: Cyc...	21.06	3.6	ng/ul	1494730	5	21.35	8396510	20.0