

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
 Data File : BM004833.D  
 Acq On : 29 Mar 2016 22:51  
 Operator : UM/SJ  
 Sample : H1939-18  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9SH0

## 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-BM031516.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 6.986       | 725           | 731         | 740          | rBV      | 222274         | 353368        | 18.80%          | 1.480%        |
| 2         | 7.157       | 754           | 760         | 768          | rBV      | 190473         | 289868        | 15.42%          | 1.214%        |
| 3         | 7.357       | 788           | 794         | 803          | rVB      | 225971         | 358357        | 19.06%          | 1.501%        |
| 4         | 7.828       | 867           | 874         | 881          | rBV      | 220786         | 354274        | 18.84%          | 1.484%        |
| 5         | 8.522       | 986           | 992         | 1001         | rBV      | 307088         | 487150        | 25.91%          | 2.040%        |
| 6         | 8.980       | 1062          | 1070        | 1082         | rBV      | 225934         | 364206        | 19.37%          | 1.525%        |
| 7         | 9.698       | 1186          | 1192        | 1201         | rVB      | 177670         | 294336        | 15.66%          | 1.233%        |
| 8         | 10.233      | 1277          | 1283        | 1297         | rBV      | 300008         | 494848        | 26.32%          | 2.073%        |
| 9         | 10.616      | 1341          | 1348        | 1361         | rVB      | 353973         | 598493        | 31.84%          | 2.507%        |
| 10        | 13.868      | 1895          | 1901        | 1905         | rBV      | 568944         | 814237        | 43.31%          | 3.410%        |
| 11        | 13.915      | 1905          | 1909        | 1917         | rVB      | 165925         | 228001        | 12.13%          | 0.955%        |
| 12        | 14.151      | 1943          | 1949        | 1961         | rVB      | 656738         | 989558        | 52.64%          | 4.145%        |
| 13        | 14.357      | 1978          | 1984        | 1989         | rBV      | 60628          | 76865         | 4.09%           | 0.322%        |
| 14        | 14.462      | 1992          | 2002        | 2013         | rVV2     | 571785         | 893737        | 47.54%          | 3.743%        |
| 15        | 14.651      | 2026          | 2034        | 2048         | rBV      | 218102         | 326329        | 17.36%          | 1.367%        |
| 16        | 14.798      | 2052          | 2059        | 2065         | rBV2     | 8945           | 20074         | 1.07%           | 0.084%        |
| 17        | 14.898      | 2072          | 2076        | 2083         | rVV      | 47327          | 73592         | 3.91%           | 0.308%        |
| 18        | 14.974      | 2085          | 2089        | 2095         | rVB4     | 13441          | 21241         | 1.13%           | 0.089%        |
| 19        | 15.309      | 2141          | 2146        | 2151         | rVB      | 76592          | 95129         | 5.06%           | 0.398%        |
| 20        | 15.451      | 2164          | 2170        | 2180         | rVB      | 847908         | 1286566       | 68.44%          | 5.389%        |
| 21        | 15.568      | 2184          | 2190        | 2203         | rVB      | 245880         | 334735        | 17.81%          | 1.402%        |
| 22        | 15.704      | 2204          | 2213        | 2219         | rBV3     | 23964          | 61235         | 3.26%           | 0.256%        |
| 23        | 15.862      | 2234          | 2240        | 2246         | rVB4     | 9723           | 20367         | 1.08%           | 0.085%        |
| 24        | 16.168      | 2287          | 2292        | 2305         | rBV      | 187346         | 272127        | 14.48%          | 1.140%        |
| 25        | 16.345      | 2316          | 2322        | 2327         | rBV      | 225845         | 302612        | 16.10%          | 1.268%        |
| 26        | 16.515      | 2344          | 2351        | 2359         | rBV2     | 140884         | 272551        | 14.50%          | 1.142%        |
| 27        | 16.809      | 2397          | 2401        | 2405         | rBV      | 24455          | 32190         | 1.71%           | 0.135%        |
| 28        | 16.962      | 2423          | 2427        | 2431         | rBV      | 51151          | 61128         | 3.25%           | 0.256%        |
| 29        | 17.027      | 2431          | 2438        | 2444         | rVB4     | 19266          | 36284         | 1.93%           | 0.152%        |
| 30        | 17.203      | 2459          | 2468        | 2474         | rBV2     | 754637         | 1123720       | 59.77%          | 4.707%        |
| 31        | 17.268      | 2474          | 2479        | 2480         | rVV2     | 40731          | 65433         | 3.48%           | 0.274%        |
| 32        | 17.303      | 2480          | 2485        | 2495         | rVB      | 991891         | 1447647       | 77.00%          | 6.064%        |
| 33        | 17.568      | 2525          | 2530        | 2536         | rBV      | 77095          | 97554         | 5.19%           | 0.409%        |
| 34        | 17.698      | 2548          | 2552        | 2558         | rBV      | 24000          | 30733         | 1.63%           | 0.129%        |

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 17.974 | 2593 | 2599 | 2606 | rBV2 | 17402   | 37709   | 2.01%   | 0.158% |
| 36 | 18.092 | 2615 | 2619 | 2625 | rBV  | 30108   | 42027   | 2.24%   | 0.176% |
| 37 | 18.392 | 2665 | 2670 | 2674 | rBV2 | 15992   | 22357   | 1.19%   | 0.094% |
| 38 | 18.621 | 2704 | 2709 | 2719 | rVB4 | 9771    | 21248   | 1.13%   | 0.089% |
| 39 | 19.021 | 2769 | 2777 | 2783 | rBV2 | 21795   | 38845   | 2.07%   | 0.163% |
| 40 | 19.233 | 2809 | 2813 | 2814 | rBV  | 19553   | 25105   | 1.34%   | 0.105% |
| 41 | 19.262 | 2814 | 2818 | 2825 | rVB  | 71525   | 103572  | 5.51%   | 0.434% |
| 42 | 19.374 | 2833 | 2837 | 2845 | rVB2 | 15675   | 22322   | 1.19%   | 0.093% |
| 43 | 19.597 | 2868 | 2875 | 2889 | rBV2 | 1271875 | 1879953 | 100.00% | 7.874% |
| 44 | 19.980 | 2923 | 2940 | 2948 | rBV2 | 30227   | 158628  | 8.44%   | 0.664% |
| 45 | 20.139 | 2963 | 2967 | 2973 | rVB  | 49416   | 57764   | 3.07%   | 0.242% |
| 46 | 20.633 | 3046 | 3051 | 3055 | rBV  | 70512   | 80359   | 4.27%   | 0.337% |
| 47 | 20.880 | 3088 | 3093 | 3097 | rBV3 | 28715   | 42362   | 2.25%   | 0.177% |
| 48 | 21.091 | 3125 | 3129 | 3139 | rBV  | 304064  | 424055  | 22.56%  | 1.776% |
| 49 | 21.386 | 3174 | 3179 | 3185 | rBV  | 1098619 | 1552243 | 82.57%  | 6.502% |
| 50 | 21.533 | 3200 | 3204 | 3209 | rBV  | 115891  | 129909  | 6.91%   | 0.544% |
| 51 | 21.715 | 3231 | 3235 | 3238 | rBV2 | 23599   | 31763   | 1.69%   | 0.133% |
| 52 | 21.956 | 3270 | 3276 | 3281 | rBV2 | 543150  | 883821  | 47.01%  | 3.702% |
| 53 | 22.450 | 3355 | 3360 | 3371 | rBV  | 90664   | 143980  | 7.66%   | 0.603% |
| 54 | 22.580 | 3379 | 3382 | 3389 | rVB  | 37672   | 58546   | 3.11%   | 0.245% |
| 55 | 22.685 | 3396 | 3400 | 3404 | rVB  | 61121   | 81380   | 4.33%   | 0.341% |
| 56 | 22.850 | 3422 | 3428 | 3438 | rVB  | 42250   | 154718  | 8.23%   | 0.648% |
| 57 | 22.968 | 3444 | 3448 | 3452 | rBV2 | 118641  | 180343  | 9.59%   | 0.755% |
| 58 | 23.015 | 3452 | 3456 | 3468 | rVB  | 204095  | 353645  | 18.81%  | 1.481% |
| 59 | 23.538 | 3537 | 3545 | 3555 | rVB  | 920977  | 1820642 | 96.85%  | 7.626% |
| 60 | 23.685 | 3562 | 3570 | 3579 | rBV  | 659701  | 1358666 | 72.27%  | 5.691% |
| 61 | 24.215 | 3654 | 3660 | 3667 | rBV  | 99394   | 189700  | 10.09%  | 0.795% |
| 62 | 24.303 | 3668 | 3675 | 3688 | rBV  | 223051  | 506398  | 26.94%  | 2.121% |
| 63 | 24.985 | 3785 | 3791 | 3799 | rVB  | 47209   | 110187  | 5.86%   | 0.462% |
| 64 | 25.350 | 3847 | 3853 | 3862 | rBV3 | 53374   | 145445  | 7.74%   | 0.609% |
| 65 | 25.874 | 3937 | 3942 | 3951 | rVB3 | 55143   | 136433  | 7.26%   | 0.571% |
| 66 | 28.197 | 4327 | 4337 | 4353 | rVB5 | 132894  | 501866  | 26.70%  | 2.102% |

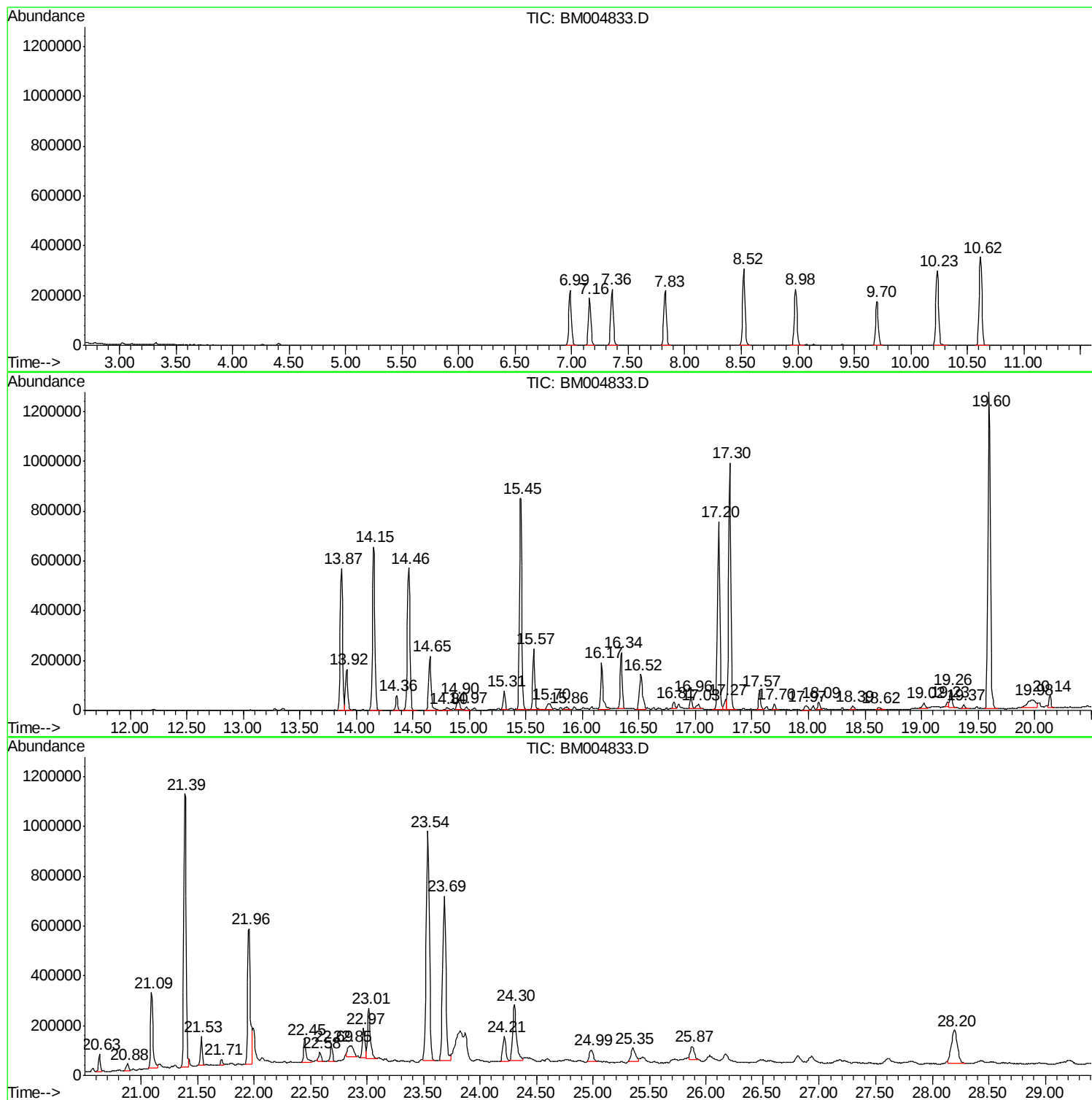
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

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-BM031516.M  
Quant Title : SVOA CALIBRATION

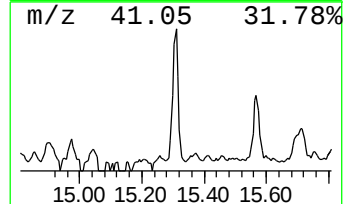
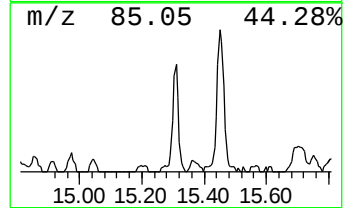
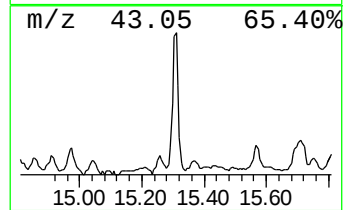
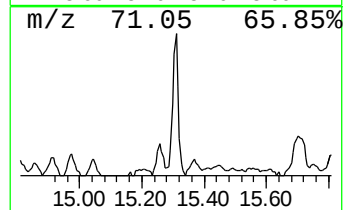
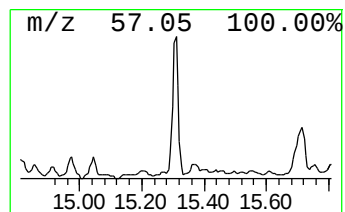
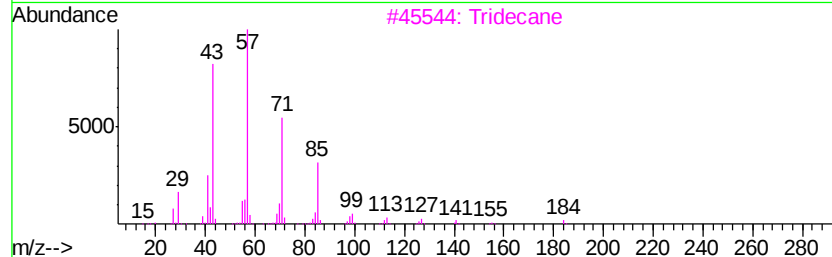
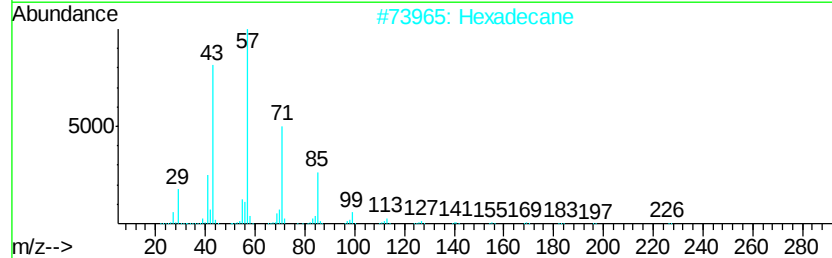
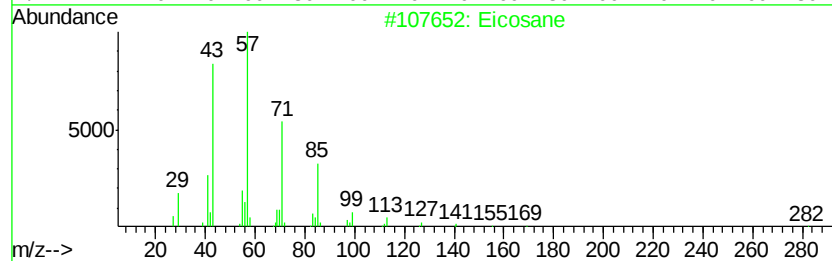
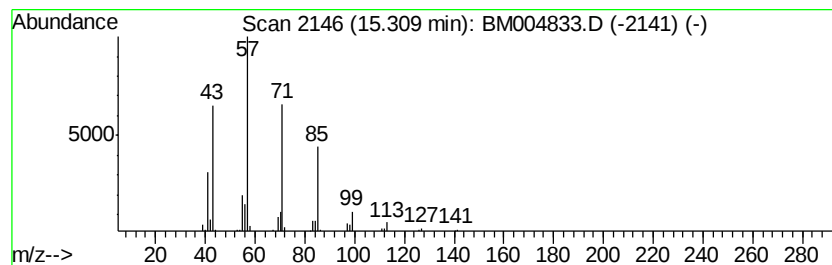
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 15.31 | 2.13 ng/ul | 95129 | Acenaphthene-d10 | 14.46 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 3    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |
| 4    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 90   |
| 5    |    |   | Heptadecane  | 240 | C17H36  | 000629-78-7 | 87   |



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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

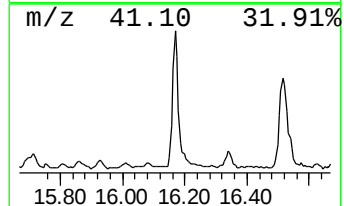
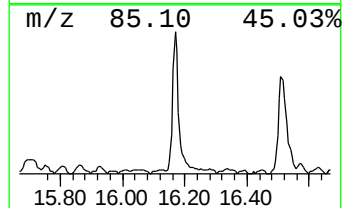
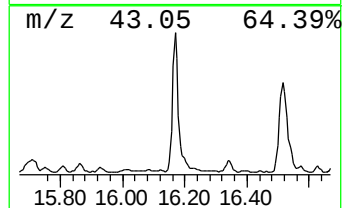
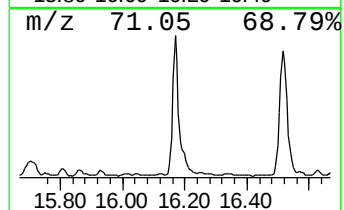
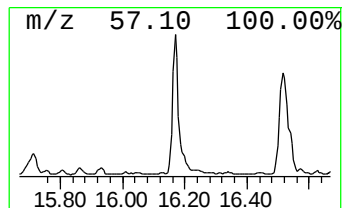
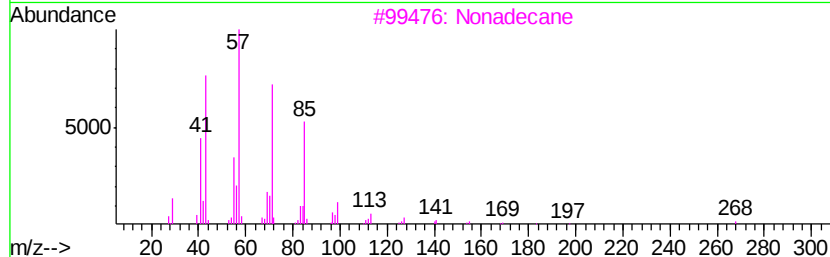
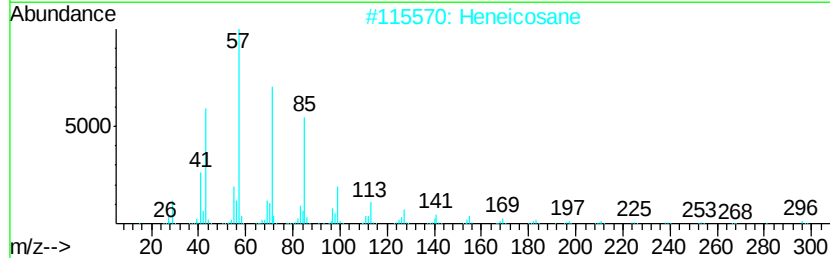
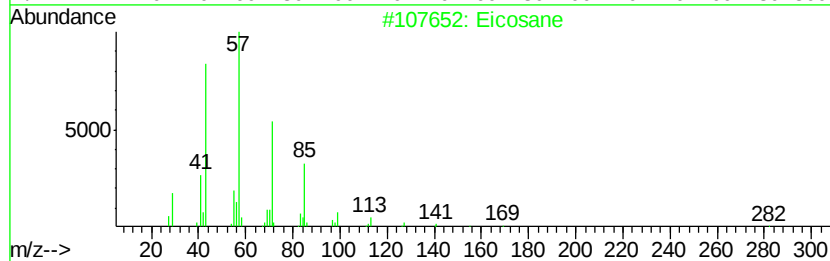
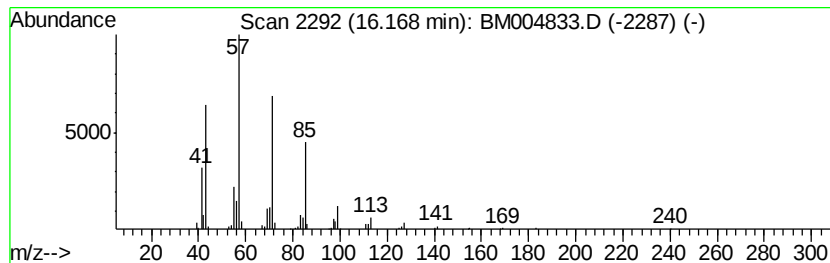
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.17 | 4.84 ng/ul | 272127 | Phenanthrene-d10 | 17.20 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 2    |    |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 90   |
| 4    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 5    |    |   | Heptacosane  | 380 | C27H56  | 000593-49-7 | 90   |



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Quant Title : SVOA CALIBRATION

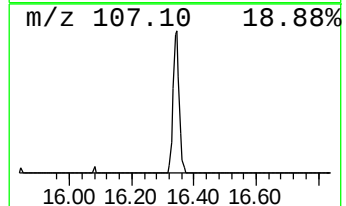
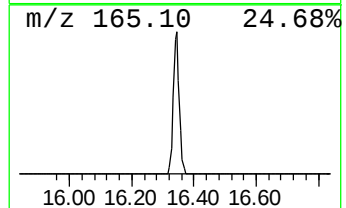
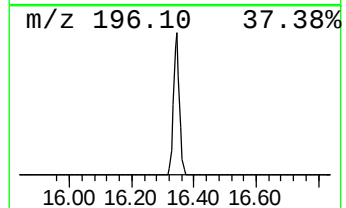
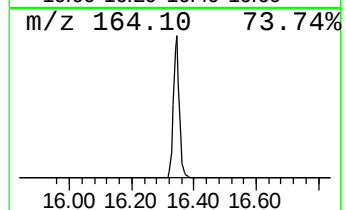
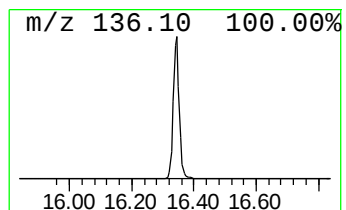
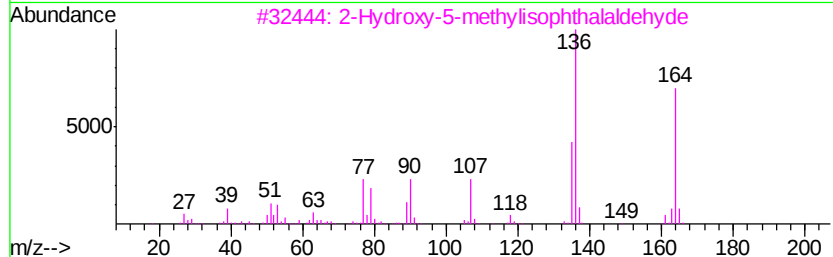
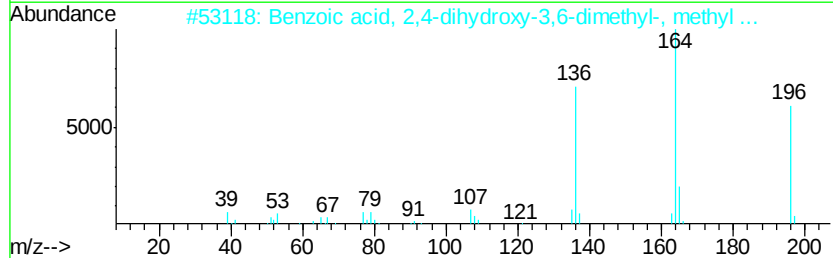
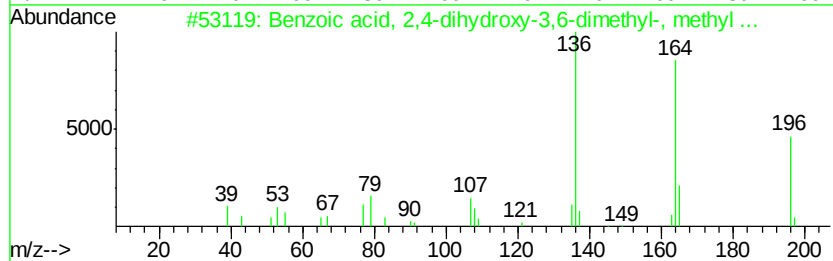
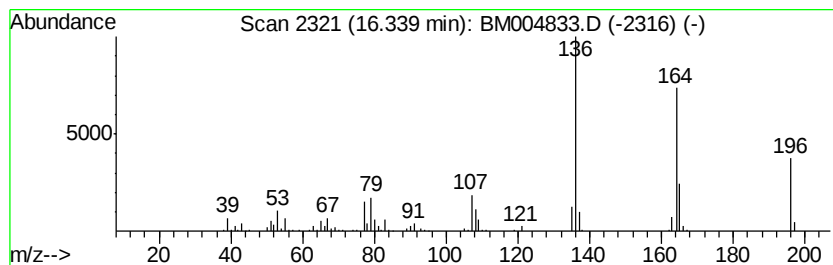
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Benzoic acid, 2,4-dihydroxy... Concentration Rank 5

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|-------|------------|--------|------------------|-------|
| 16.34 | 5.39 ng/ul | 302612 | Phenanthrene-d10 | 17.20 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | Benzoic acid, 2,4-dihydroxy-3,6-... | 196 | C10H12O4 | 004707-47-5  | 98   |
| 2       |   | Benzoic acid, 2,4-dihydroxy-3,6-... | 196 | C10H12O4 | 004707-47-5  | 94   |
| 3       |   | 2-Hydroxy-5-methylisophthalaldehyde | 164 | C9H8O3   | 007310-95-4  | 58   |
| 4       |   | Benzoic acid, 2-hydroxy-4-methox... | 196 | C10H12O4 | 000520-43-4  | 40   |
| 5       |   | Benzo-1,3,2-oxathiaborole, 2-ethyl- | 164 | C8H9BO3  | 1000161-38-9 | 40   |



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Quant Title : SVOA CALIBRATION

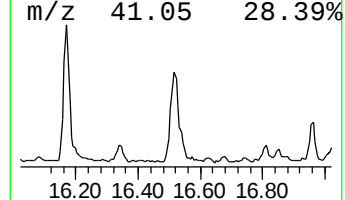
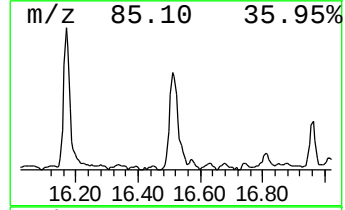
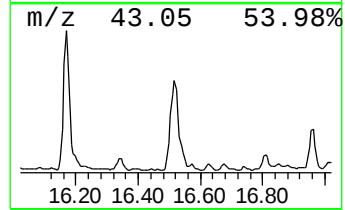
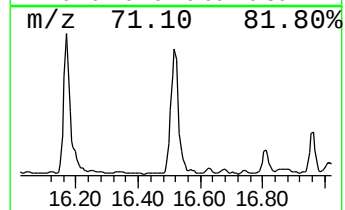
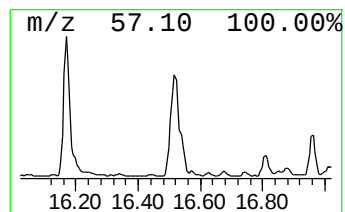
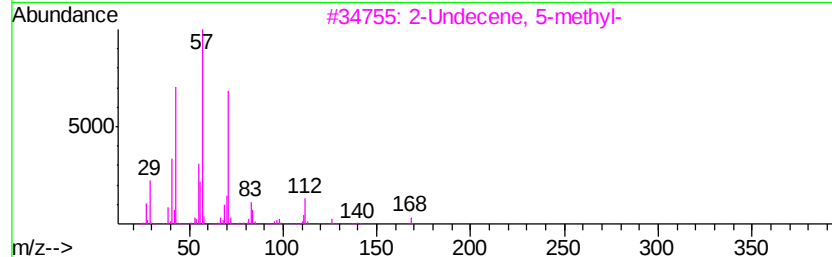
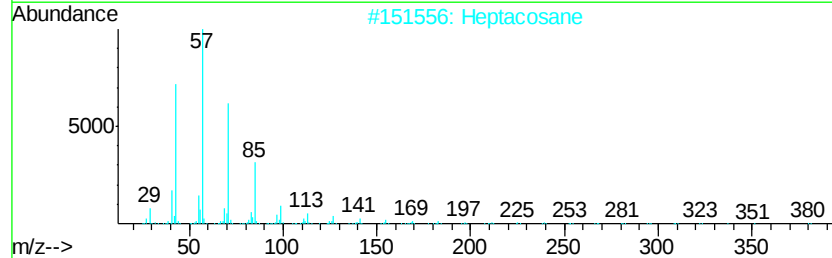
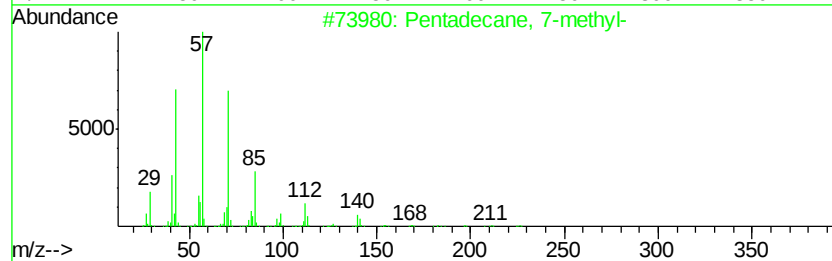
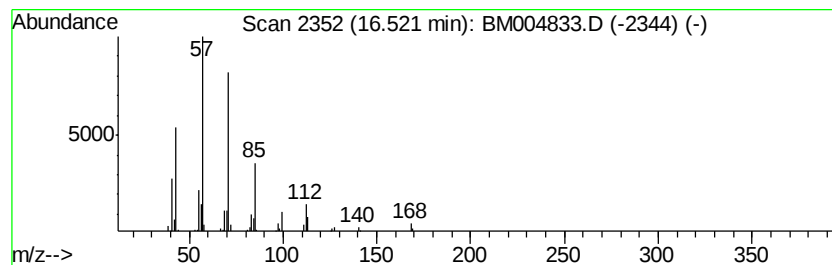
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.52 | 4.85 ng/ul | 272551 | Phenanthrene-d10 | 17.20 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 7-methyl-  | 226 | C16H34  | 006165-40-8 | 90   |
| 2    |    | Heptacosane             | 380 | C27H56  | 000593-49-7 | 64   |
| 3    |    | 2-Undecene, 5-methyl-   | 168 | C12H24  | 056851-34-4 | 64   |
| 4    |    | Heneicosane             | 296 | C21H44  | 000629-94-7 | 52   |
| 5    |    | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

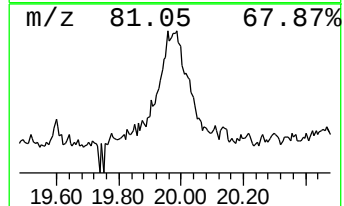
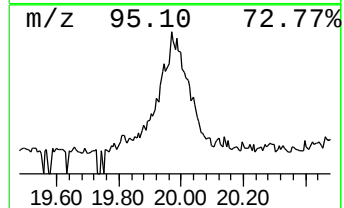
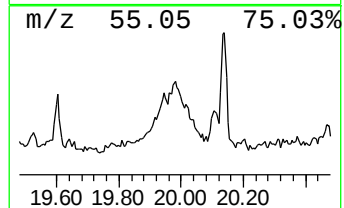
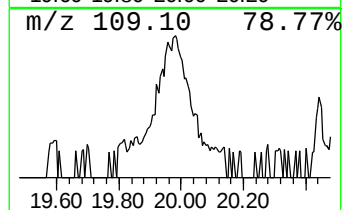
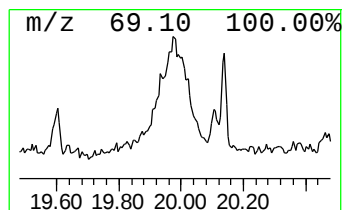
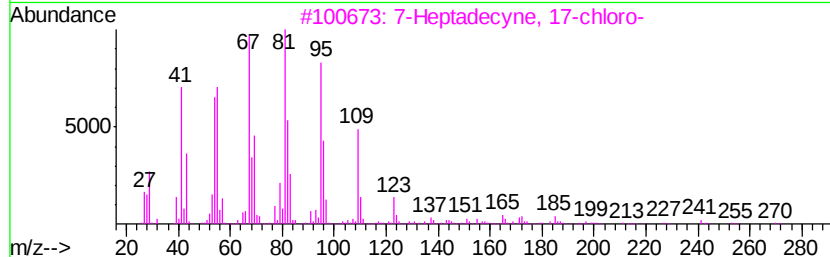
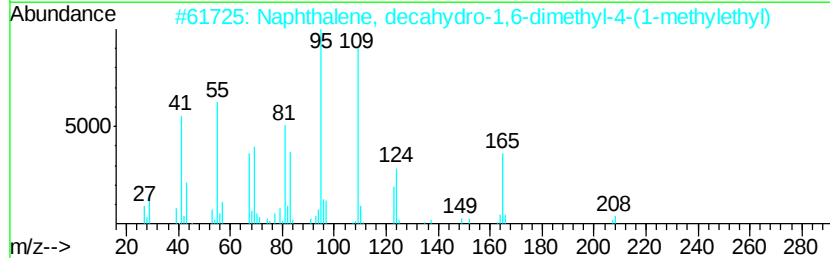
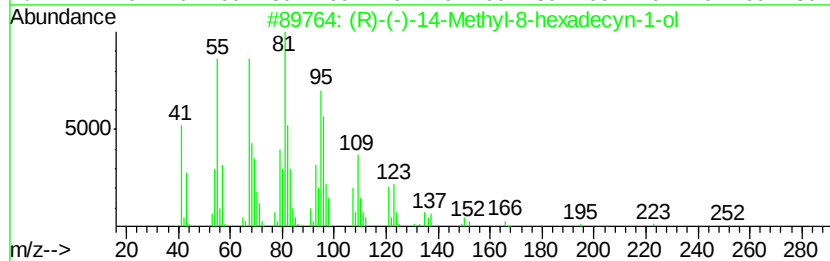
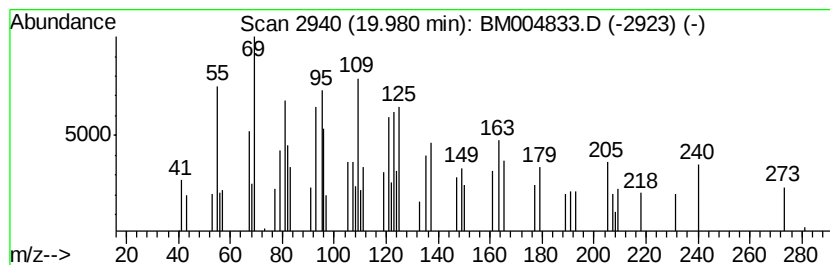
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 unknown-01 Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.98 | 2.04 ng/ul | 158628 | Chrysene-d12     | 21.39 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | (R)-(-)-14-Methyl-8-hexadecyn-1-ol  | 252 | C17H32O      | 064566-18-3  | 30      |      |      |
| 2    | Naphthalene, decahydro-1,6-dimet... | 208 | C15H28       | 034315-85-0  | 30      |      |      |
| 3    | 7-Heptadecyne, 17-chloro-           | 270 | C17H31Cl     | 056554-75-7  | 27      |      |      |
| 4    | 5-Acetoxymethyl-2,6,10-trimethyl... | 282 | C17H30O3     | 1000144-12-3 | 27      |      |      |
| 5    | Longifolenaldehyde                  | 220 | C15H24O      | 019890-84-7  | 25      |      |      |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

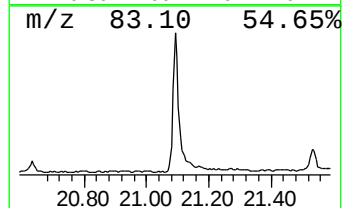
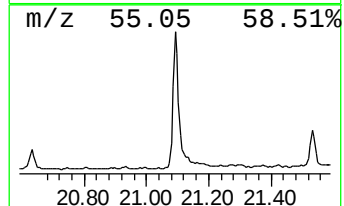
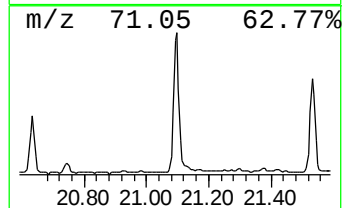
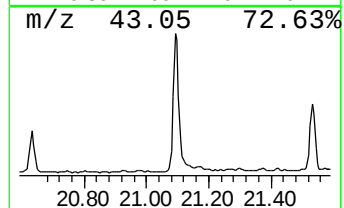
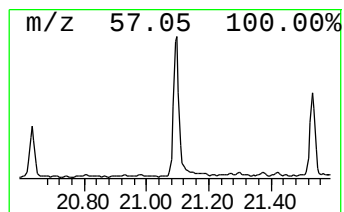
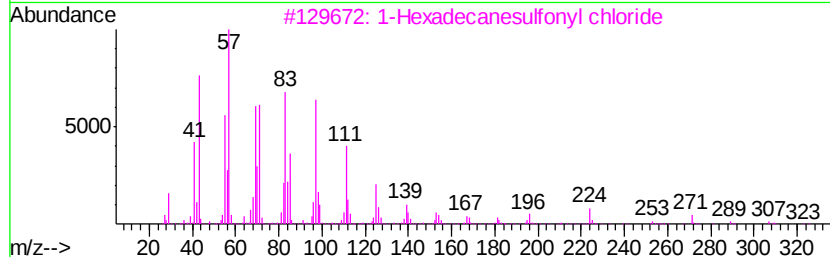
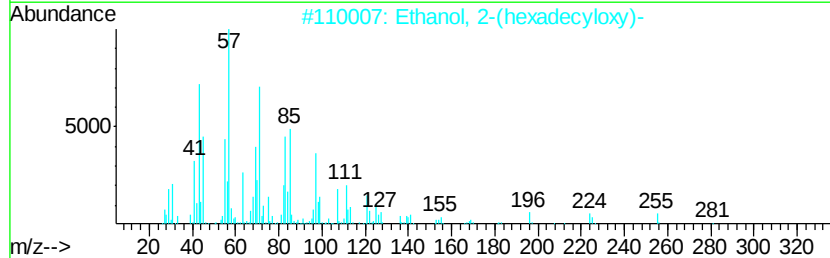
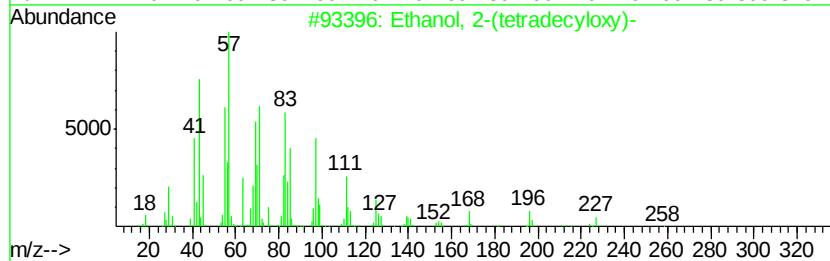
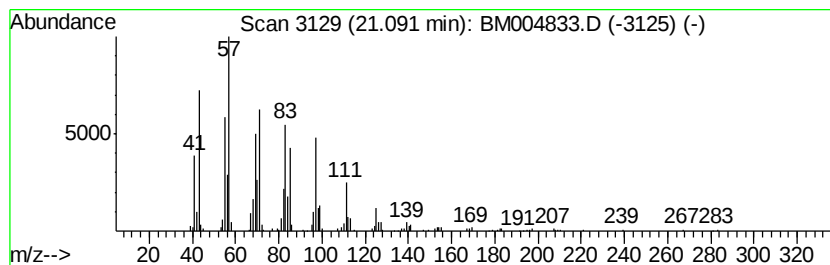
Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 6 Ethanol, 2-(tetradecyloxy)- Concentration Rank 4

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.        |              |      |
|---------|------------|-------------------------------------|------------------|-------------|--------------|------|
| 21.09   | 5.46 ng/ul | 424055                              | Chrysene-d12     | 21.39       |              |      |
| Hit# of | 5          | Tentative ID                        | MW               | MolForm     | CAS#         | Qual |
| 1       |            | Ethanol, 2-(tetradecyloxy)-         | 258              | C16H34O2    | 002136-70-1  | 91   |
| 2       |            | Ethanol, 2-(hexadecyloxy)-          | 286              | C18H38O2    | 002136-71-2  | 91   |
| 3       |            | 1-Hexadecanesulfonyl chloride       | 324              | C16H33ClO2S | 038775-38-1  | 87   |
| 4       |            | Pentafluoropropionic acid, hepta... | 402              | C20H35F5O2  | 1000283-04-2 | 87   |
| 5       |            | Heptafluorobutanoic acid, heptad... | 452              | C21H35F7O2  | 1000282-97-3 | 87   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

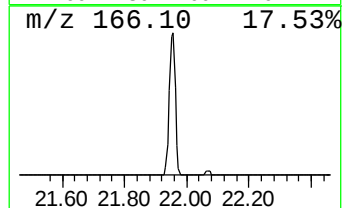
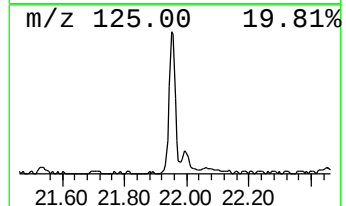
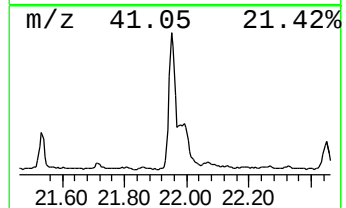
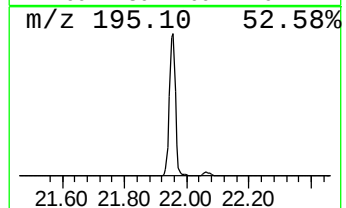
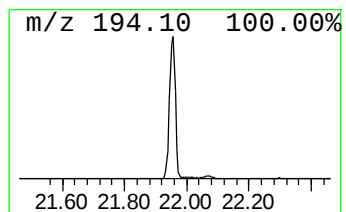
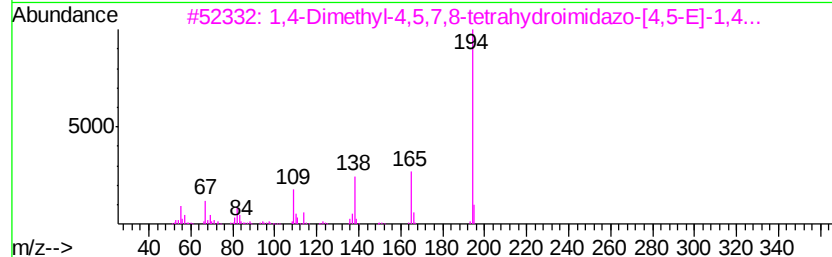
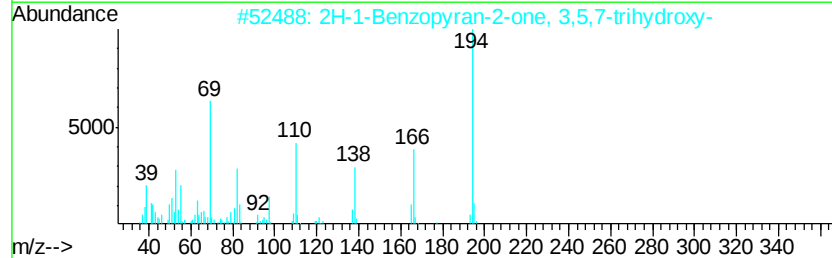
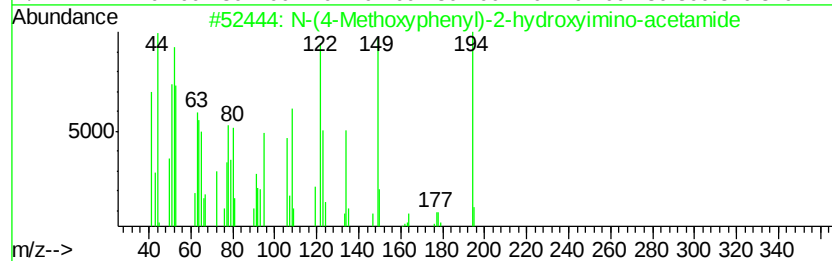
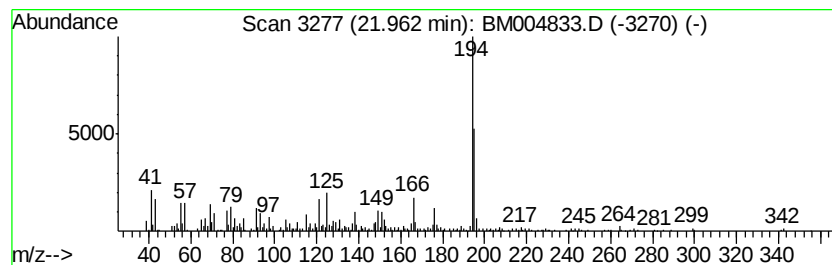
Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 7 N-(4-Methoxyphenyl)-2-hydro... Concentration Rank 1

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 21.96     | 11.39 ng/ul                         | 883821 | Chrysene-d12     | 21.39        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | N-(4-Methoxyphenyl)-2-hydroxyimi... | 194    | C9H10N2O3        | 1000143-61-3 | 95   |
| 2         | 2H-1-Benzopyran-2-one, 3,5,7-tri... | 194    | C9H6O5           | 022065-07-2  | 43   |
| 3         | 1,4-Dimethyl-4,5,7,8-tetrahydroi... | 194    | C8H10N4O2        | 130063-15-9  | 43   |
| 4         | Isophthaldiamidoxime                | 194    | C8H10N4O2        | 015325-51-6  | 42   |
| 5         | 4-Amino-6-methyl-2-methylthiopyr... | 194    | C8H10N4S         | 097337-20-7  | 38   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleID :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

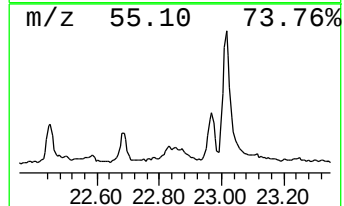
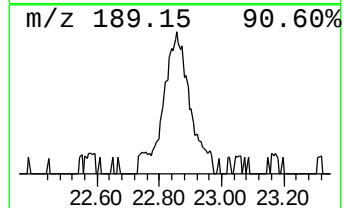
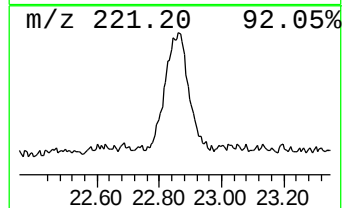
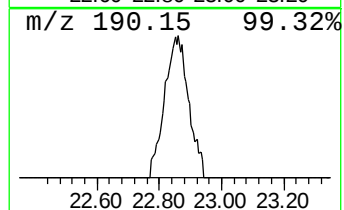
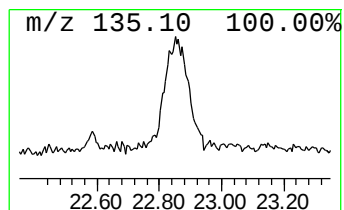
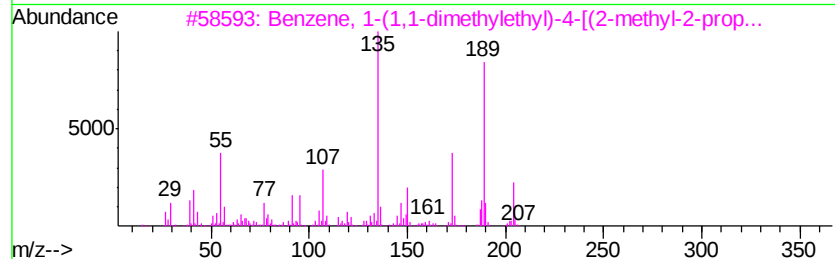
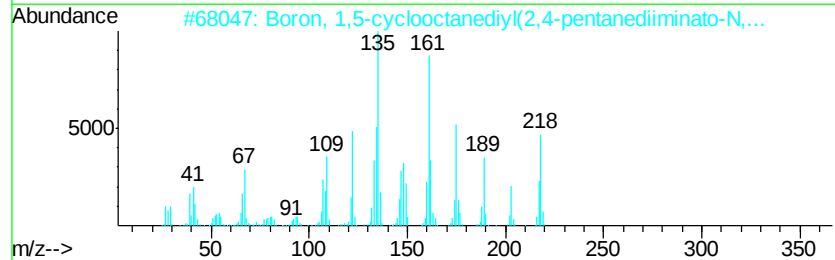
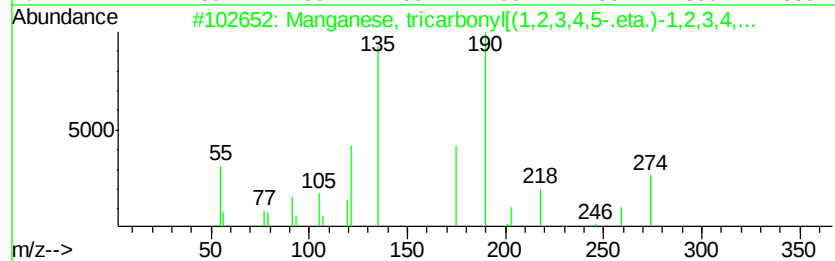
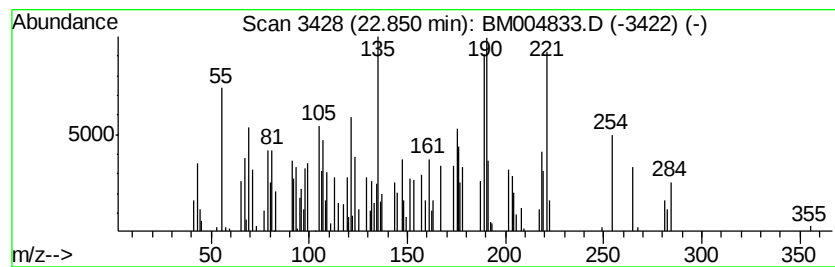
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 unknown-02 Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.85 | 2.28 ng/ul | 154718 | Perylene-d12     | 23.69 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Manganese, tricarbonyl[(1,2,3,4,... | 274 | C13H15MnO3 | 034807-89-1  | 27   |
| 2    |    | Boron, 1,5-cyclooctanediyl(2,4-p... | 218 | C13H23BN2  | 136704-99-9  | 22   |
| 3    |    | Benzene, 1-(1,1-dimethylethyl)-4... | 204 | C14H20O    | 054932-87-5  | 18   |
| 4    |    | .gamma.-Neoclovene                  | 204 | C15H24     | 1000156-11-7 | 15   |
| 5    |    | 6S-2,3,8,8-Tetramethyltricyclo[5... | 204 | C15H24     | 137235-48-4  | 15   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

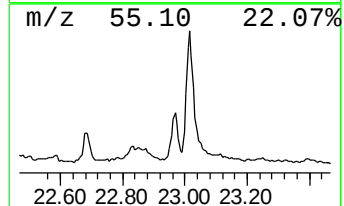
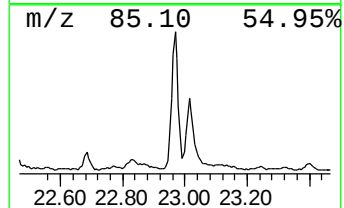
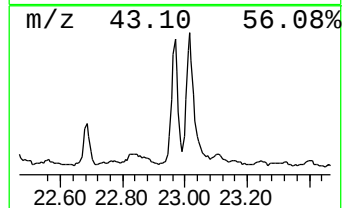
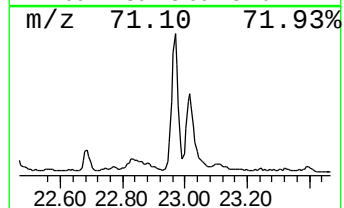
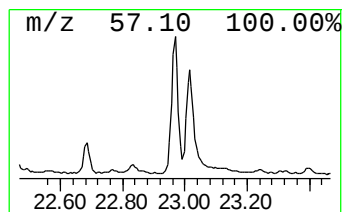
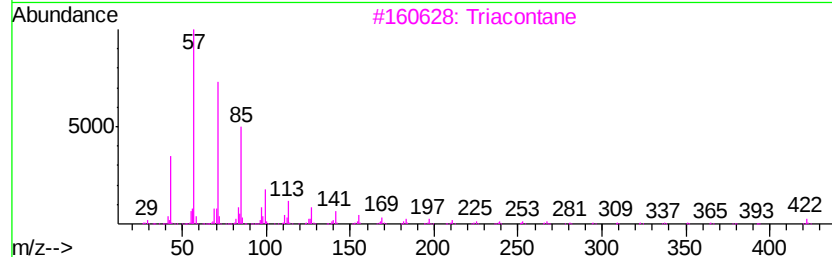
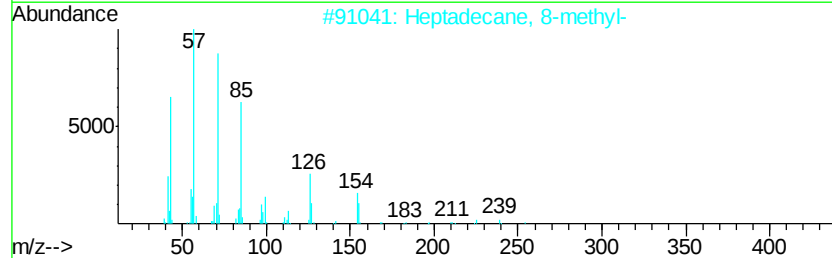
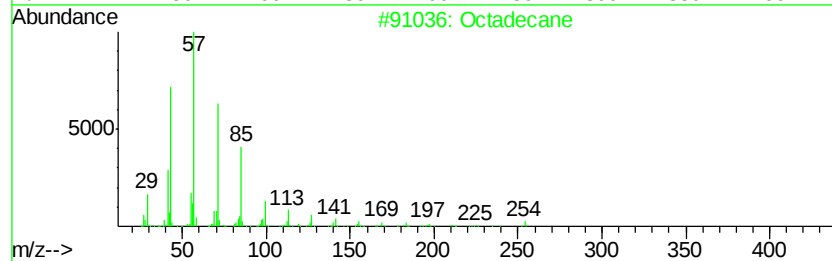
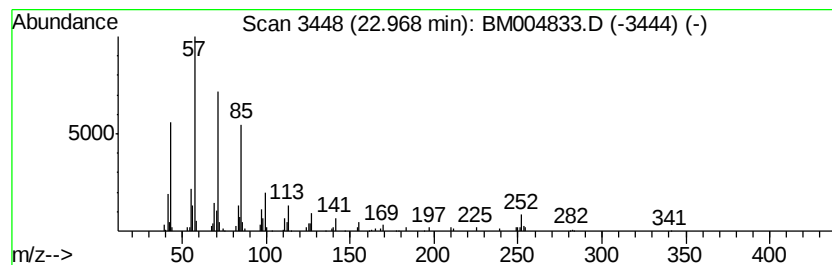
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.97 | 2.65 ng/ul | 180343 | Perylene-d12     | 23.69 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane             | 254 | C18H38  | 000593-45-3 | 93   |
| 2    |    | Heptadecane, 8-methyl- | 254 | C18H38  | 013287-23-5 | 91   |
| 3    |    | Triacontane            | 422 | C30H62  | 000638-68-6 | 90   |
| 4    |    | Heptadecane            | 240 | C17H36  | 000629-78-7 | 90   |
| 5    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

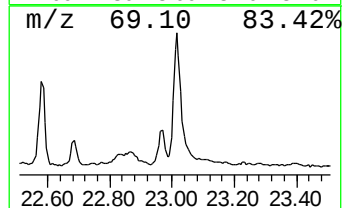
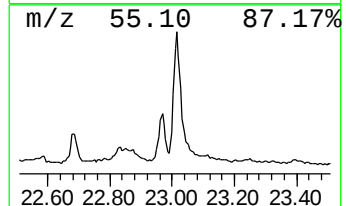
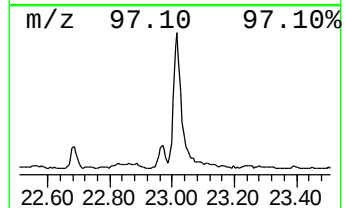
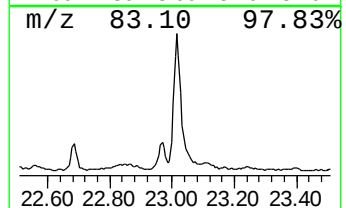
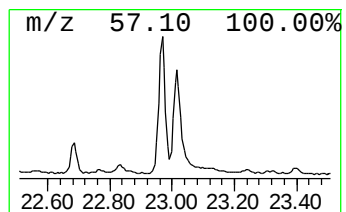
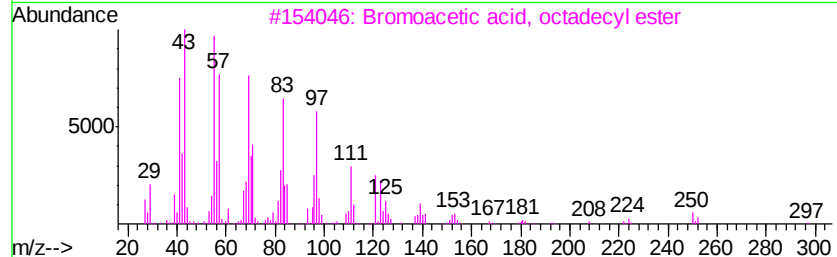
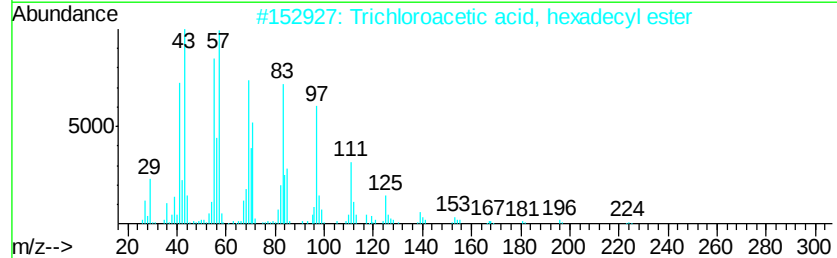
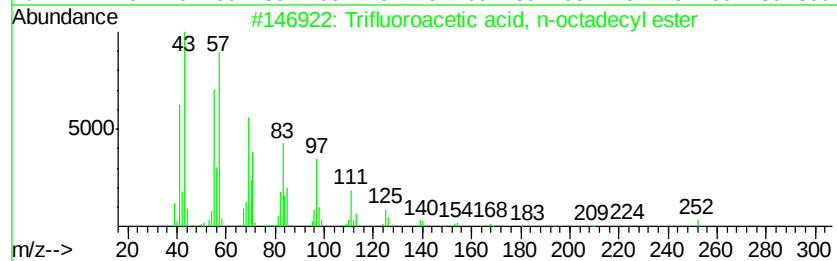
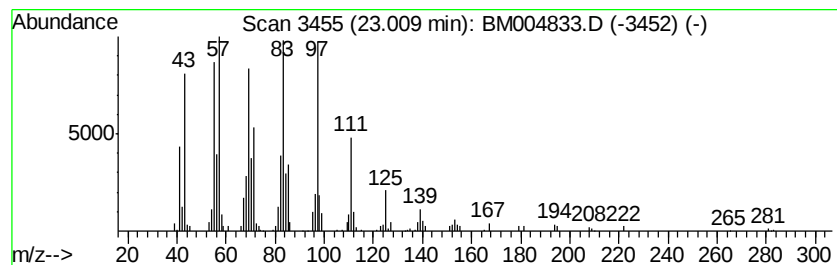
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Trifluoroacetic acid, n-oct... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.01 | 5.21 ng/ul | 353645 | Perylene-d12     | 23.69 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1  | 91   |
| 2    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 91   |
| 3    |    |   | Bromoacetic acid, octadecyl ester   | 390 | C20H39BrO2  | 018992-03-5  | 90   |
| 4    |    |   | Heptafluorobutyric acid, n-octad... | 466 | C22H37F7O2  | 000400-57-7  | 90   |
| 5    |    |   | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 81   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

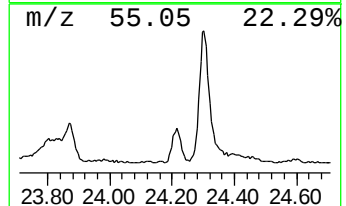
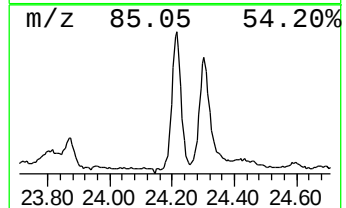
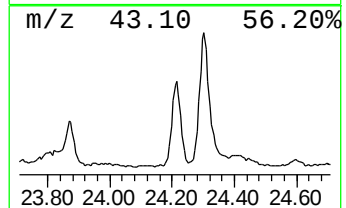
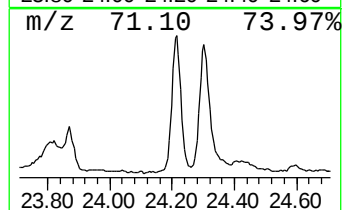
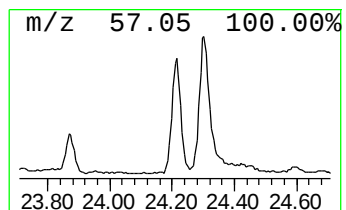
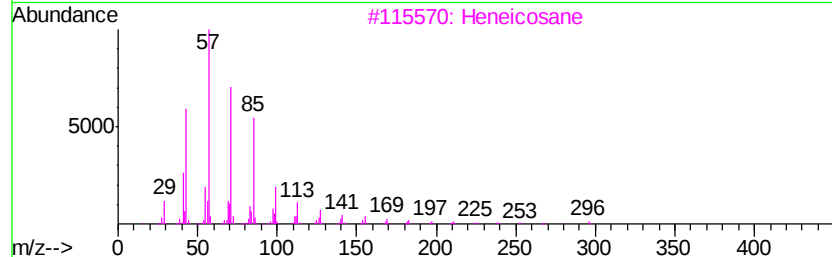
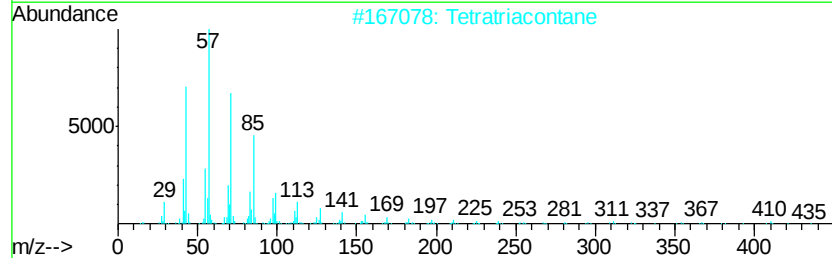
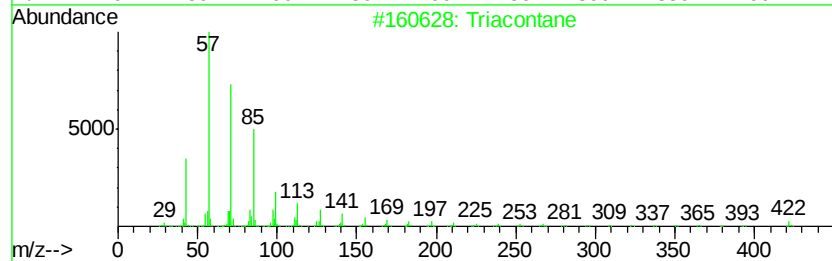
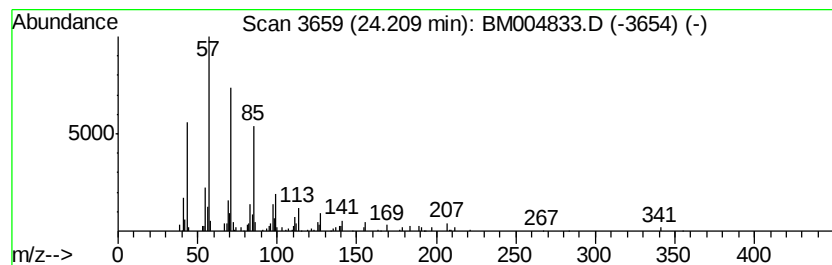
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.21 | 2.79 ng/ul | 189700 | Perylene-d12     | 23.69 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Triacontane            | 422 | C30H62  | 000638-68-6 | 91   |
| 2    |    | Tetratriacontane       | 479 | C34H70  | 014167-59-0 | 91   |
| 3    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    | Octacosane             | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    | Tetracosane, 11-decyl- | 479 | C34H70  | 055429-84-0 | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

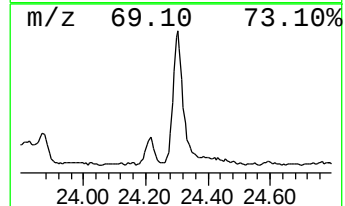
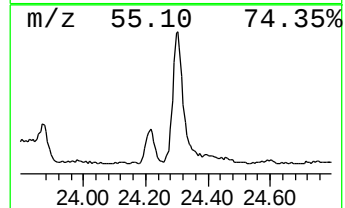
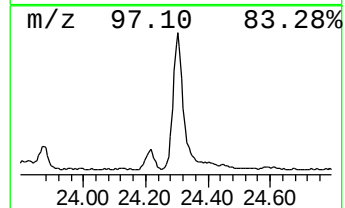
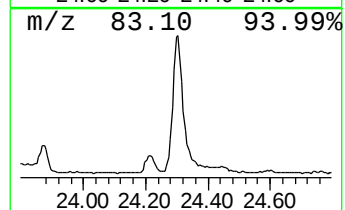
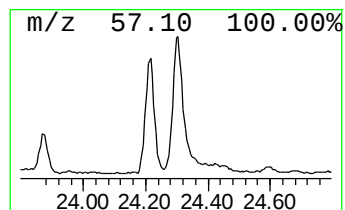
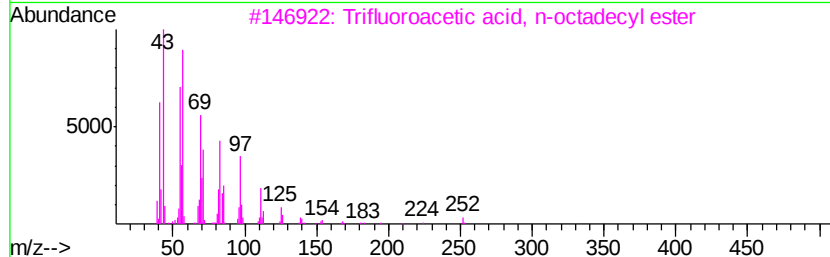
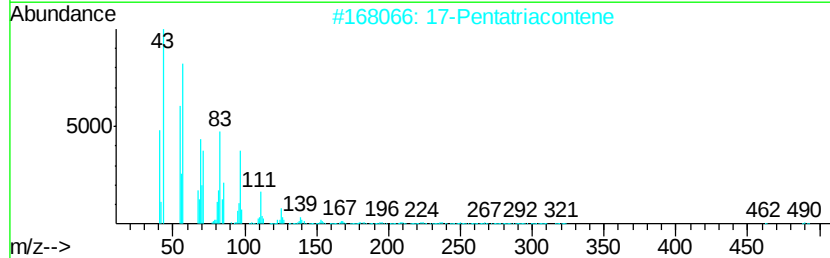
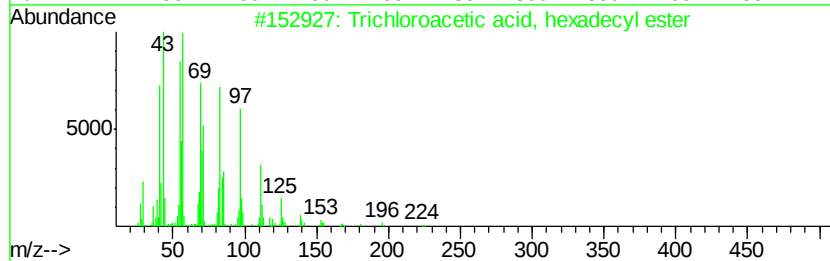
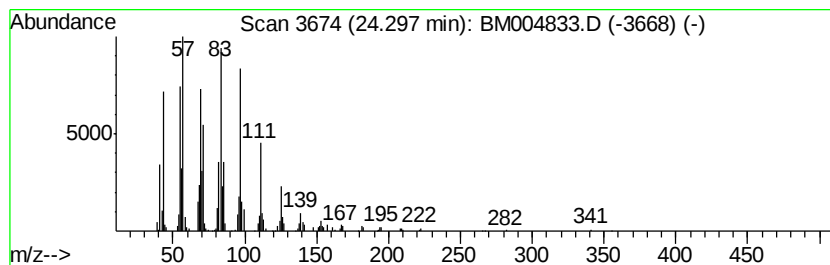
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Trichloroacetic acid, hexad... Concentration Rank 2

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.30 | 7.45 ng/ul | 506398 | Perylene-d12     | 23.69 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1 | 91   |
| 2    |    |   | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0 | 90   |
| 3    |    |   | Trifluoroacetic acid, n-octadecy... | 366 | C20H37F3O2  | 079392-43-1 | 83   |
| 4    |    |   | 1-Docosene                          | 308 | C22H44      | 001599-67-3 | 76   |
| 5    |    |   | 2-Chloropropionic acid, octadec...  | 360 | C21H41ClO2  | 088104-31-8 | 68   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

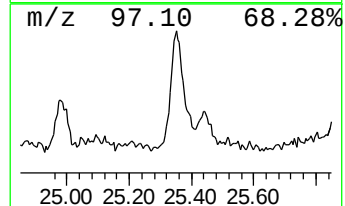
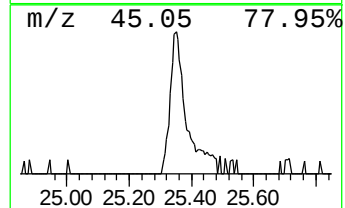
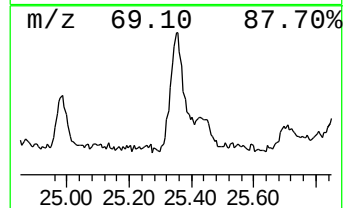
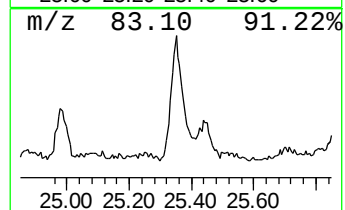
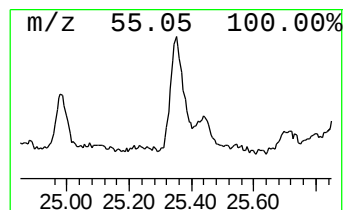
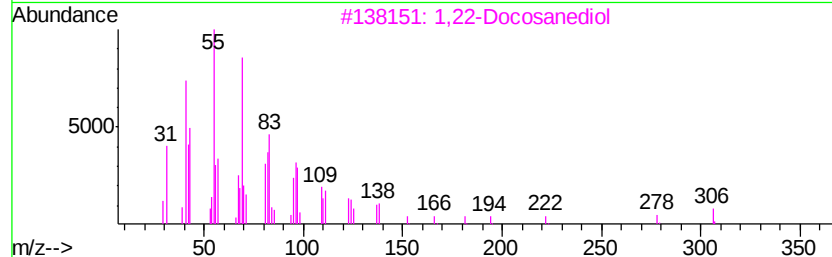
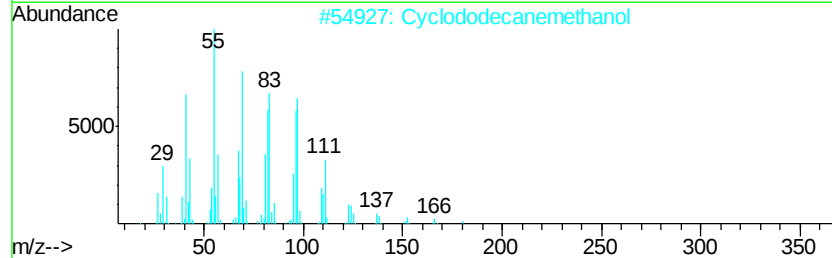
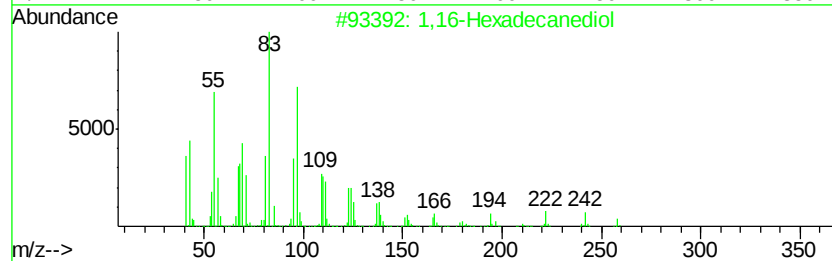
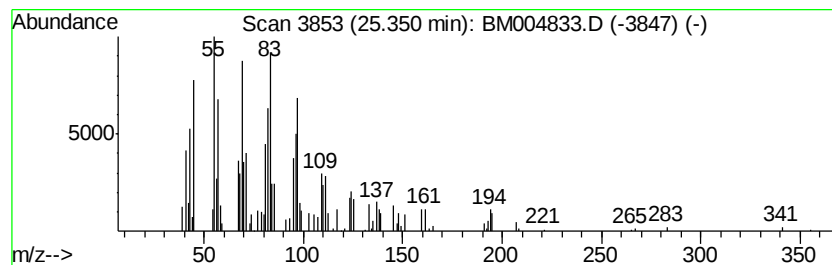
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 1,16-Hexadecanediol Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 25.35 | 2.14 ng/ul | 145445 | Perylene-d12     | 23.69 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm    | CAS#        | Qual |
|------|----|---|------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 1,16-Hexadecanediol                | 258 | C16H34O2   | 007735-42-4 | 50   |
| 2    |    |   | Cyclododecanemethanol              | 198 | C13H26O    | 001892-12-2 | 49   |
| 3    |    |   | 1,22-Docosanediol                  | 342 | C22H46O2   | 022513-81-1 | 49   |
| 4    |    |   | Bromoacetic acid, pentadecyl ester | 348 | C17H33BrO2 | 131143-01-6 | 49   |
| 5    |    |   | 1,13-Tetradecadiene                | 194 | C14H26     | 021964-49-8 | 47   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

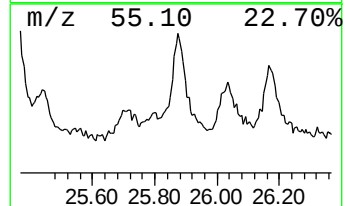
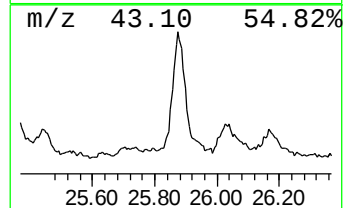
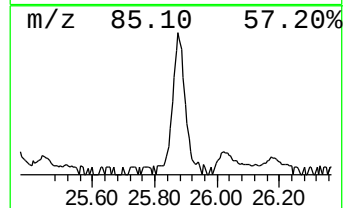
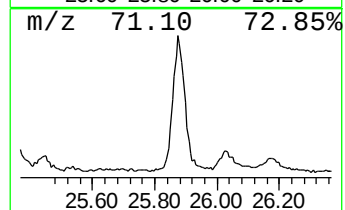
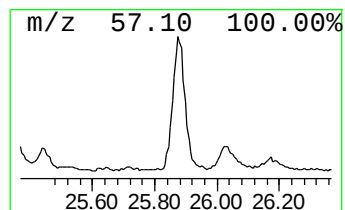
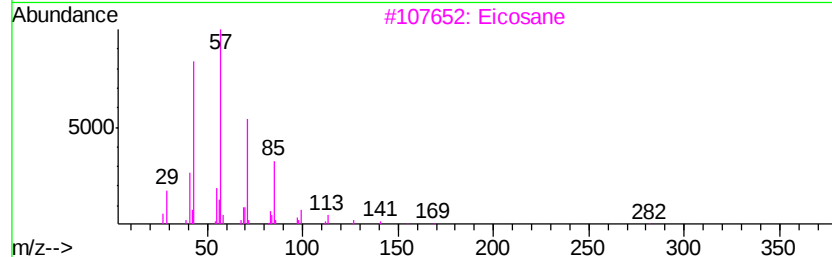
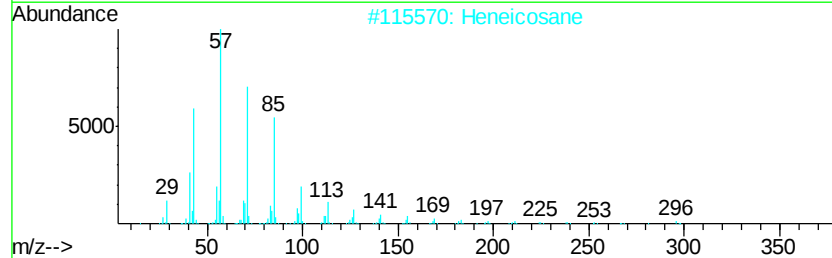
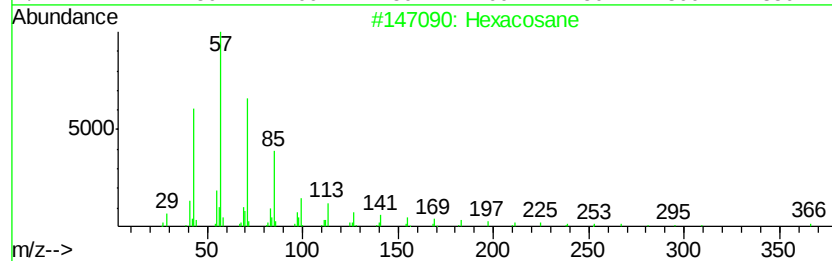
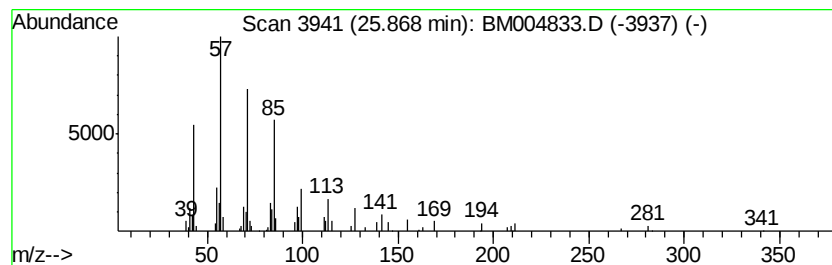
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 25.87 | 2.01 ng/ul | 136433 | Perylene-d12     | 23.69 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Hexacosane   | 366 | C26H54  | 000630-01-3 | 91   |
| 2    |    |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 4    |    |   | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 5    |    |   | Pentacosane  | 352 | C25H52  | 000629-99-2 | 91   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
Data File : BM004833.D  
Acq On : 29 Mar 2016 22:51  
Operator : UM/SJ  
Sample : H1939-18  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

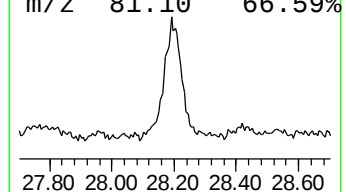
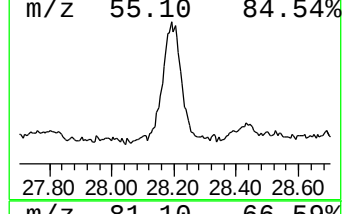
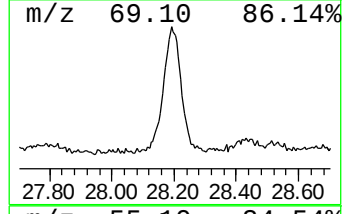
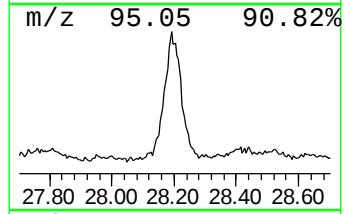
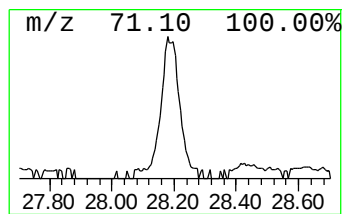
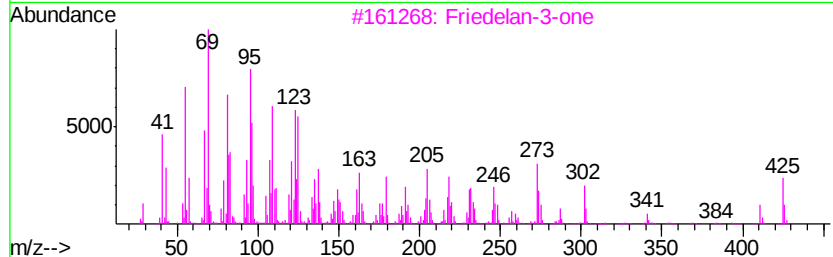
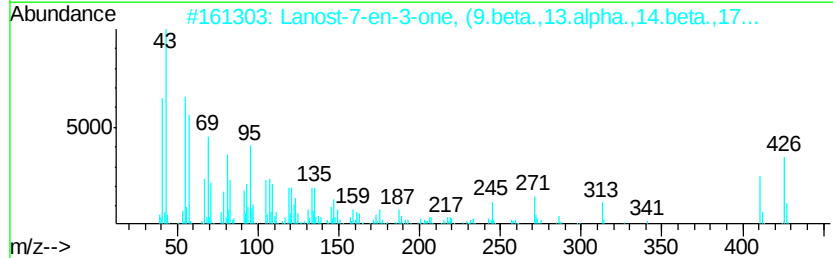
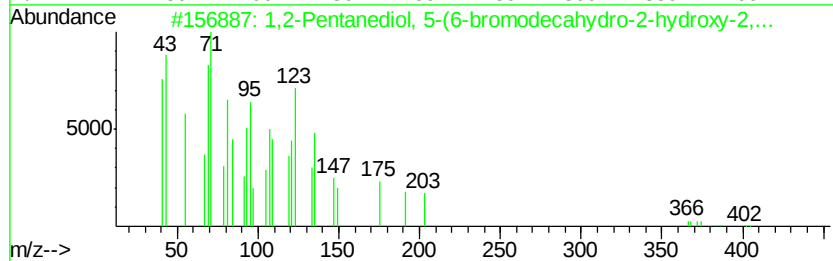
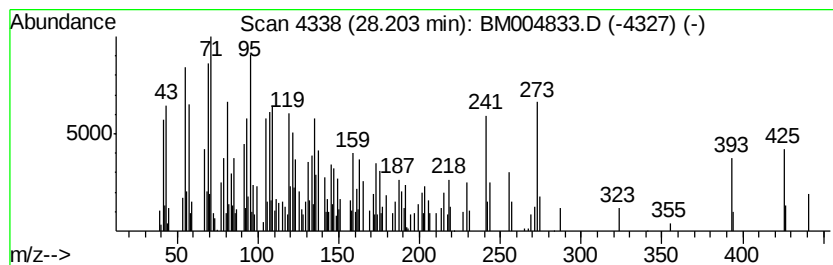
Instrument :  
BNA\_M  
ClientSampleId :  
D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 15 1,2-Pentanediol, 5-(6-bromo... Concentration Rank 3

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.         |      |      |
|---------|-------------------------------------|--------------|------------------|--------------|------|------|
| 28.20   | 7.39 ng/ul                          | 501866       | Perylene-d12     | 23.69        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | 1,2-Pentanediol, 5-(6-bromodecah... | 402          | C20H35BrO3       | 115346-29-7  | 64   |      |
| 2       | Lanost-7-en-3-one, (9.beta.,13.a... | 426          | C30H50O          | 055515-18-9  | 43   |      |
| 3       | Friedelan-3-one                     | 426          | C30H50O          | 000559-74-0  | 38   |      |
| 4       | 9-Thiabicyclo[3.3.1]nonane, 2,6-... | 274          | C14H18N4S        | 1000141-52-1 | 38   |      |
| 5       | Lanosterol                          | 426          | C30H50O          | 000079-63-0  | 38   |      |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM032916\  
 Data File : BM004833.D  
 Acq On : 29 Mar 2016 22:51  
 Operator : UM/SJ  
 Sample : H1939-18  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9SH0

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM031516.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 |
| (DEL) Alkane: Str... | 15.31 | 2.1     | ng/ul | 95129    | 3                     | 14.46 | 893737  | 20.0 |
| (DEL) Alkane: Str... | 16.17 | 4.8     | ng/ul | 272127   | 4                     | 17.20 | 1123720 | 20.0 |
| Benzoic acid, 2,4... | 16.34 | 5.4     | ng/ul | 302612   | 4                     | 17.20 | 1123720 | 20.0 |
| (DEL) Alkane: Str... | 16.52 | 4.8     | ng/ul | 272551   | 4                     | 17.20 | 1123720 | 20.0 |
| unknown-01           | 19.98 | 2.0     | ng/ul | 158628   | 5                     | 21.39 | 1552240 | 20.0 |
| Ethanol, 2-(tetra... | 21.09 | 5.5     | ng/ul | 424055   | 5                     | 21.39 | 1552240 | 20.0 |
| N-(4-Methoxypheny... | 21.96 | 11.4    | ng/ul | 883821   | 5                     | 21.39 | 1552240 | 20.0 |
| unknown-02           | 22.85 | 2.3     | ng/ul | 154718   | 6                     | 23.69 | 1358670 | 20.0 |
| (DEL) Alkane: Str... | 22.97 | 2.6     | ng/ul | 180343   | 6                     | 23.69 | 1358670 | 20.0 |
| Trifluoroacetic a... | 23.01 | 5.2     | ng/ul | 353645   | 6                     | 23.69 | 1358670 | 20.0 |
| (DEL) Alkane: Str... | 24.21 | 2.8     | ng/ul | 189700   | 6                     | 23.69 | 1358670 | 20.0 |
| Trichloroacetic a... | 24.30 | 7.5     | ng/ul | 506398   | 6                     | 23.69 | 1358670 | 20.0 |
| 1,16-Hexadecanediol  | 25.35 | 2.1     | ng/ul | 145445   | 6                     | 23.69 | 1358670 | 20.0 |
| (DEL) Alkane: Str... | 25.87 | 2.0     | ng/ul | 136433   | 6                     | 23.69 | 1358670 | 20.0 |
| 1,2-Pentanediol, ... | 28.20 | 7.4     | ng/ul | 501866   | 6                     | 23.69 | 1358670 | 20.0 |