

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
 Data File : BM012118.D  
 Acq On : 13 Oct 2017 02:52  
 Operator : SJ/JU  
 Sample : I5490-06  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 B-69(1-3)-092717

## 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-BM101217.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         | 4.528       | 239           | 245         | 254          | rBV      | 650030         | 1017061       | 3.20%           | 0.561%        |
| 2         | 5.457       | 396           | 403         | 414          | rBV      | 247763         | 551618        | 1.74%           | 0.304%        |
| 3         | 5.810       | 457           | 463         | 472          | rBV2     | 781175         | 1611552       | 5.07%           | 0.889%        |
| 4         | 6.310       | 536           | 548         | 554          | rBV      | 224610         | 404630        | 1.27%           | 0.223%        |
| 5         | 6.804       | 622           | 632         | 637          | rBV      | 860646         | 1287040       | 4.05%           | 0.710%        |
| 6         | 7.004       | 654           | 666         | 670          | rBV      | 1006170        | 2232489       | 7.02%           | 1.231%        |
| 7         | 7.040       | 670           | 672         | 680          | rVB3     | 387757         | 641054        | 2.02%           | 0.354%        |
| 8         | 7.163       | 683           | 693         | 705          | rBV2     | 926348         | 1873594       | 5.89%           | 1.033%        |
| 9         | 7.363       | 718           | 727         | 735          | rVB2     | 1412599        | 2742992       | 8.63%           | 1.513%        |
| 10        | 7.440       | 735           | 740         | 742          | rBV      | 377689         | 594861        | 1.87%           | 0.328%        |
| 11        | 7.651       | 772           | 776         | 781          | rVB      | 229727         | 363279        | 1.14%           | 0.200%        |
| 12        | 7.792       | 788           | 800         | 804          | rBV      | 1540509        | 2576973       | 8.11%           | 1.421%        |
| 13        | 7.834       | 804           | 807         | 814          | rVB      | 1338486        | 1968211       | 6.19%           | 1.085%        |
| 14        | 7.940       | 820           | 825         | 831          | rBV      | 545117         | 833122        | 2.62%           | 0.459%        |
| 15        | 8.045       | 837           | 843         | 850          | rBV      | 2225944        | 3765677       | 11.85%          | 2.077%        |
| 16        | 8.298       | 880           | 886         | 890          | rBV      | 514045         | 846153        | 2.66%           | 0.467%        |
| 17        | 8.357       | 890           | 896         | 910          | rVB      | 2360169        | 4975062       | 15.65%          | 2.744%        |
| 18        | 8.487       | 910           | 918         | 923          | rBV      | 2374549        | 3923367       | 12.34%          | 2.164%        |
| 19        | 8.551       | 923           | 929         | 936          | rVB3     | 1403707        | 2677850       | 8.42%           | 1.477%        |
| 20        | 8.698       | 944           | 954         | 957          | rBV2     | 1307650        | 2815658       | 8.86%           | 1.553%        |
| 21        | 8.851       | 972           | 980         | 988          | rBV4     | 540113         | 1448829       | 4.56%           | 0.799%        |
| 22        | 8.934       | 988           | 994         | 1000         | rVB      | 1266130        | 2203773       | 6.93%           | 1.215%        |
| 23        | 9.004       | 1000          | 1006        | 1010         | rBV      | 1313342        | 2468934       | 7.77%           | 1.362%        |
| 24        | 9.051       | 1010          | 1014        | 1023         | rVB2     | 1198401        | 2285526       | 7.19%           | 1.260%        |
| 25        | 9.134       | 1023          | 1028        | 1033         | rBV4     | 257136         | 579570        | 1.82%           | 0.320%        |
| 26        | 9.192       | 1033          | 1038        | 1044         | rVB4     | 266836         | 532554        | 1.68%           | 0.294%        |
| 27        | 9.263       | 1045          | 1050        | 1054         | rBV      | 186141         | 344808        | 1.08%           | 0.190%        |
| 28        | 9.334       | 1057          | 1062        | 1066         | rVV      | 1084151        | 1569941       | 4.94%           | 0.866%        |
| 29        | 9.392       | 1066          | 1072        | 1076         | rVV      | 2204858        | 3565933       | 11.22%          | 1.967%        |
| 30        | 9.434       | 1076          | 1079        | 1086         | rVB3     | 299163         | 543987        | 1.71%           | 0.300%        |
| 31        | 9.516       | 1086          | 1093        | 1096         | rBV2     | 272957         | 590182        | 1.86%           | 0.325%        |
| 32        | 9.581       | 1101          | 1104        | 1109         | rVV      | 467199         | 685737        | 2.16%           | 0.378%        |
| 33        | 9.634       | 1109          | 1113        | 1117         | rVV      | 631666         | 932566        | 2.93%           | 0.514%        |
| 34        | 9.722       | 1121          | 1128        | 1139         | rVB2     | 1000467        | 2109283       | 6.64%           | 1.163%        |

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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 9.816  | 1139 | 1144 | 1151 | rBV5 | 303173  | 613473  | 1.93%  | 0.338% |
| 36 | 9.892  | 1151 | 1157 | 1165 | rVV3 | 648208  | 1711506 | 5.38%  | 0.944% |
| 37 | 10.016 | 1169 | 1178 | 1186 | rVV  | 2568231 | 5250812 | 16.52% | 2.896% |
| 38 | 10.092 | 1186 | 1191 | 1195 | rVV  | 1905459 | 3029548 | 9.53%  | 1.671% |
| 39 | 10.134 | 1195 | 1198 | 1204 | rVB  | 1163127 | 1651437 | 5.20%  | 0.911% |
| 40 | 10.198 | 1204 | 1209 | 1215 | rBV4 | 395225  | 833522  | 2.62%  | 0.460% |
| 41 | 10.263 | 1215 | 1220 | 1228 | rVB2 | 1235422 | 2310440 | 7.27%  | 1.274% |
| 42 | 10.345 | 1228 | 1234 | 1243 | rBV3 | 601180  | 1126239 | 3.54%  | 0.621% |
| 43 | 10.586 | 1270 | 1275 | 1279 | rBV  | 378339  | 689555  | 2.17%  | 0.380% |
| 44 | 10.639 | 1279 | 1284 | 1290 | rVV  | 2094315 | 3443743 | 10.83% | 1.899% |
| 45 | 10.781 | 1303 | 1308 | 1315 | rVB4 | 209541  | 371597  | 1.17%  | 0.205% |
| 46 | 10.904 | 1325 | 1329 | 1334 | rBV3 | 165732  | 349349  | 1.10%  | 0.193% |
| 47 | 11.439 | 1414 | 1420 | 1430 | rVB5 | 201659  | 454811  | 1.43%  | 0.251% |
| 48 | 13.233 | 1721 | 1725 | 1730 | rBV  | 258012  | 362773  | 1.14%  | 0.200% |
| 49 | 13.345 | 1741 | 1744 | 1752 | rVB3 | 246185  | 359861  | 1.13%  | 0.198% |
| 50 | 13.892 | 1827 | 1837 | 1841 | rBV  | 2479357 | 3638431 | 11.45% | 2.007% |
| 51 | 13.933 | 1841 | 1844 | 1851 | rVB  | 1456342 | 2174589 | 6.84%  | 1.199% |
| 52 | 14.180 | 1880 | 1886 | 1901 | rBV  | 2558980 | 3919718 | 12.33% | 2.162% |
| 53 | 14.310 | 1903 | 1908 | 1913 | rVV4 | 204315  | 343301  | 1.08%  | 0.189% |
| 54 | 14.486 | 1932 | 1938 | 1944 | rVB2 | 2521489 | 3792690 | 11.93% | 2.092% |
| 55 | 14.704 | 1968 | 1975 | 1984 | rBV3 | 621082  | 1565877 | 4.93%  | 0.864% |
| 56 | 14.886 | 1997 | 2006 | 2014 | rVB7 | 163722  | 419194  | 1.32%  | 0.231% |
| 57 | 15.316 | 2075 | 2079 | 2084 | rVB  | 303229  | 382108  | 1.20%  | 0.211% |
| 58 | 15.480 | 2101 | 2107 | 2112 | rBV  | 2997436 | 4200933 | 13.22% | 2.317% |
| 59 | 15.616 | 2126 | 2130 | 2135 | rBV  | 293054  | 421766  | 1.33%  | 0.233% |
| 60 | 16.198 | 2221 | 2229 | 2235 | rBV4 | 579620  | 1450613 | 4.56%  | 0.800% |
| 61 | 16.304 | 2243 | 2247 | 2250 | rBV  | 253983  | 342428  | 1.08%  | 0.189% |
| 62 | 16.363 | 2254 | 2257 | 2262 | rVB2 | 184715  | 323711  | 1.02%  | 0.179% |
| 63 | 16.621 | 2297 | 2301 | 2307 | rVB4 | 293936  | 482570  | 1.52%  | 0.266% |
| 64 | 16.968 | 2356 | 2360 | 2364 | rBV  | 272581  | 367049  | 1.15%  | 0.202% |
| 65 | 17.015 | 2364 | 2368 | 2372 | rBV2 | 518666  | 775569  | 2.44%  | 0.428% |
| 66 | 17.233 | 2400 | 2405 | 2409 | rBV2 | 2575691 | 3731341 | 11.74% | 2.058% |
| 67 | 17.274 | 2409 | 2412 | 2417 | rVV  | 1002774 | 1324822 | 4.17%  | 0.731% |
| 68 | 17.333 | 2417 | 2422 | 2434 | rVV  | 2745824 | 4110349 | 12.93% | 2.267% |
| 69 | 18.121 | 2550 | 2556 | 2560 | rBV2 | 5846632 | 8628343 | 27.15% | 4.758% |
| 70 | 18.298 | 2583 | 2586 | 2590 | rBV4 | 352418  | 514865  | 1.62%  | 0.284% |
| 71 | 18.604 | 2635 | 2638 | 2643 | rBV2 | 829800  | 1004154 | 3.16%  | 0.554% |

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Peak Location: TOP

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Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Title : SVOA CALIBRATION

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|----|--------|------|------|------|-----|----------|----------|---------|---------|
| 72 | 19.292 | 2748 | 2755 | 2768 | rBV | 18565728 | 31784899 | 100.00% | 17.529% |
| 73 | 19.403 | 2769 | 2774 | 2785 | rVV | 8463450  | 11788109 | 37.09%  | 6.501%  |
| 74 | 19.627 | 2807 | 2812 | 2815 | rBV | 3583039  | 5425061  | 17.07%  | 2.992%  |
| 75 | 19.656 | 2815 | 2817 | 2825 | rVB | 2035525  | 2647446  | 8.33%   | 1.460%  |
| 76 | 20.392 | 2939 | 2942 | 2947 | rVB | 2064827  | 2609833  | 8.21%   | 1.439%  |
| 77 | 21.309 | 3095 | 3098 | 3102 | rVB | 2523211  | 3038280  | 9.56%   | 1.676%  |

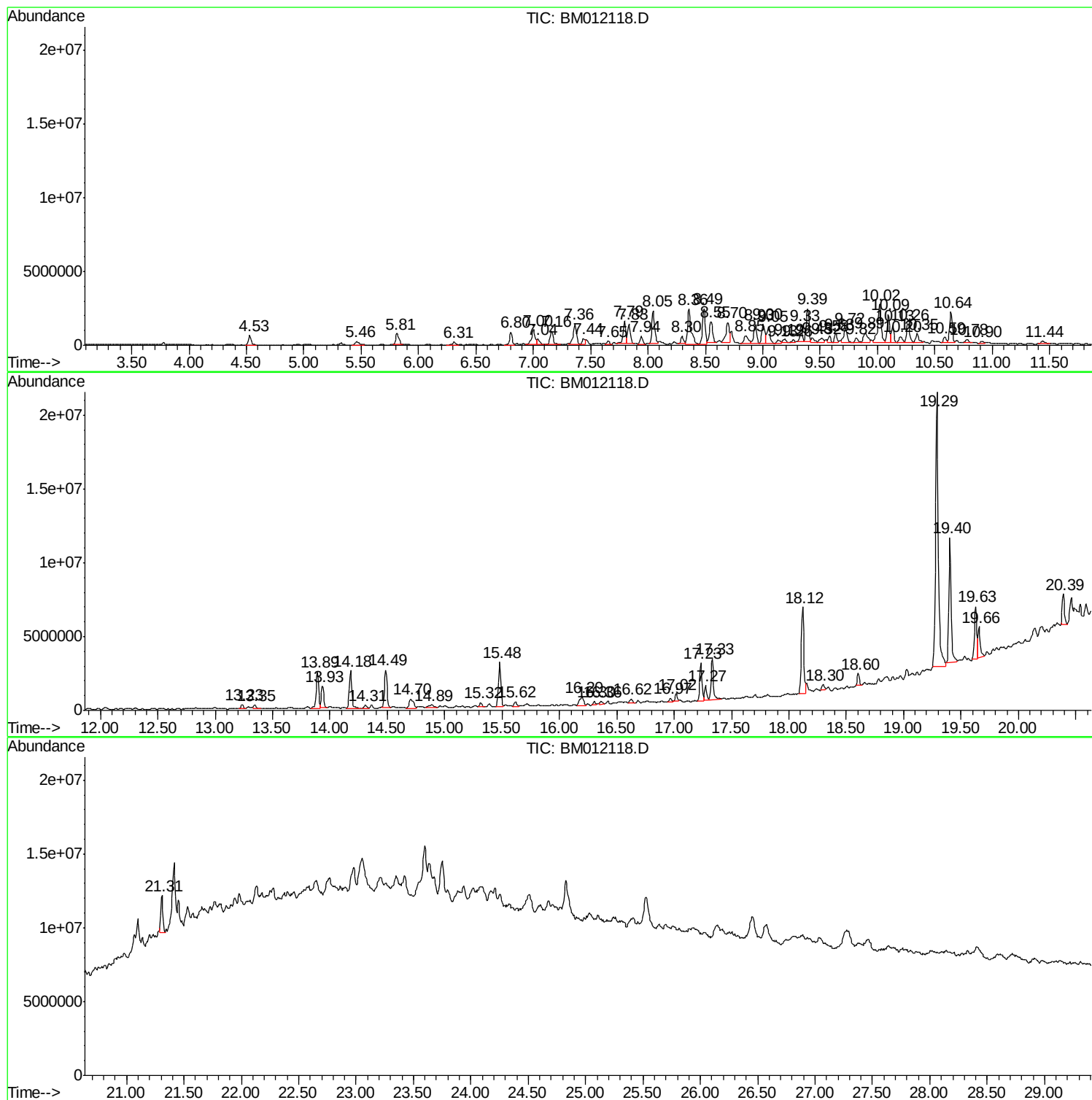
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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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



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Quant Title : SVOA CALIBRATION

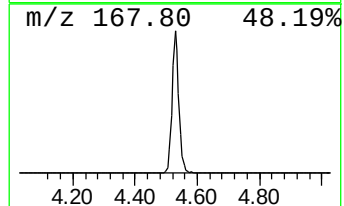
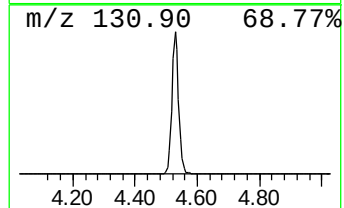
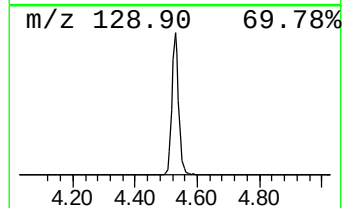
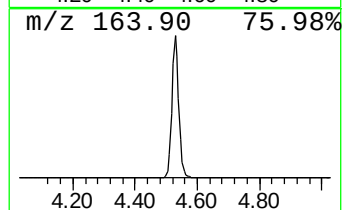
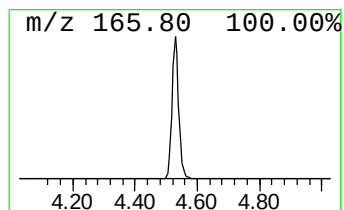
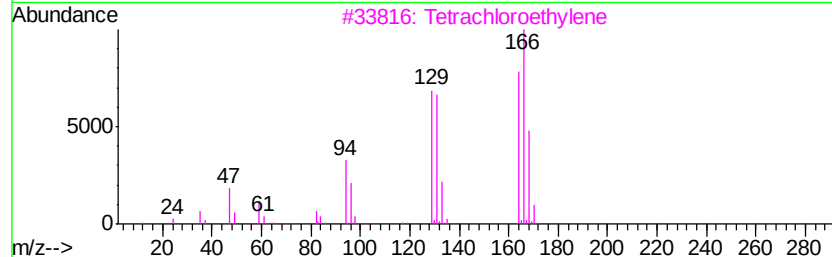
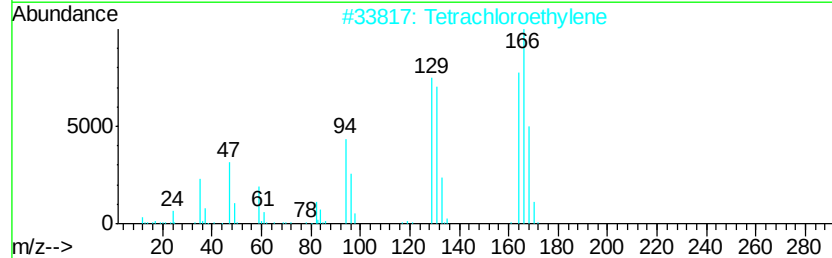
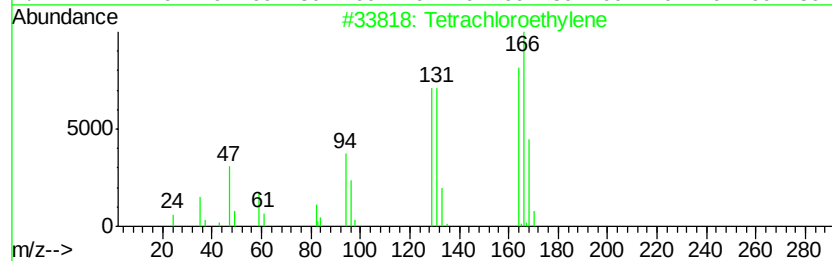
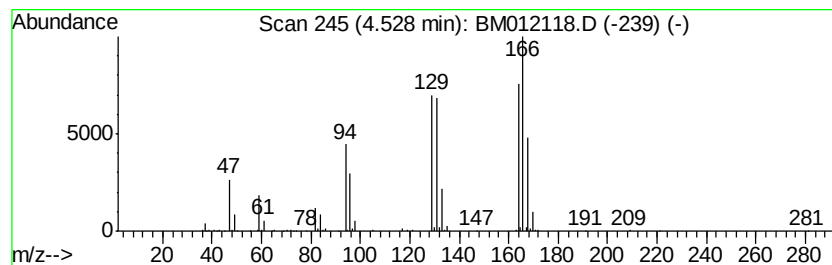
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Tetrachloroethylene Concentration Rank 14

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 4.53 | 10.33 ng/ul | 1017060 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 98   |
| 2    |    | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 98   |
| 3    |    | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 97   |
| 4    |    | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 97   |
| 5    |    | Pyrimidine, 5-fluoro-2,4-dichloro- | 166 | C4HCl2FN2 | 002927-71-1 | 38   |



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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

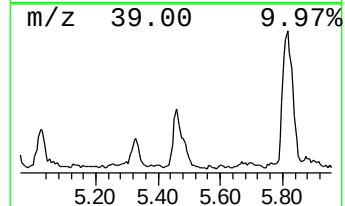
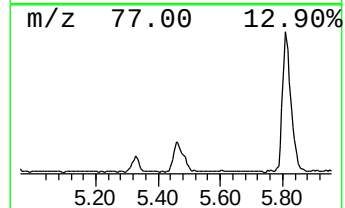
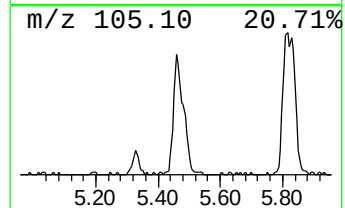
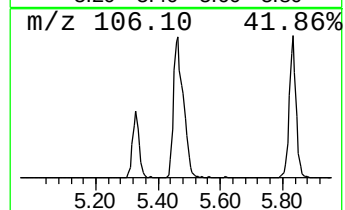
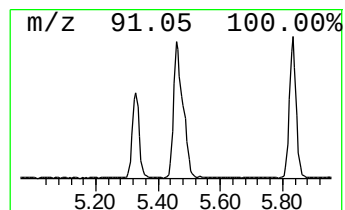
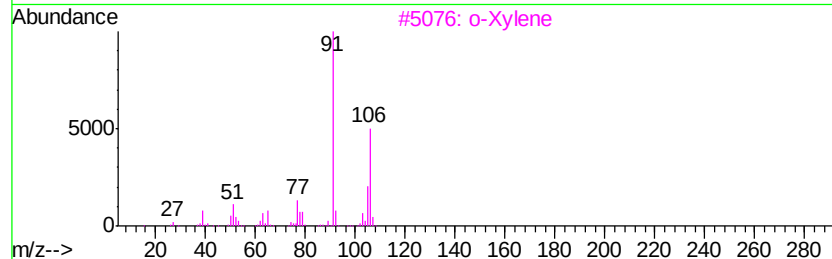
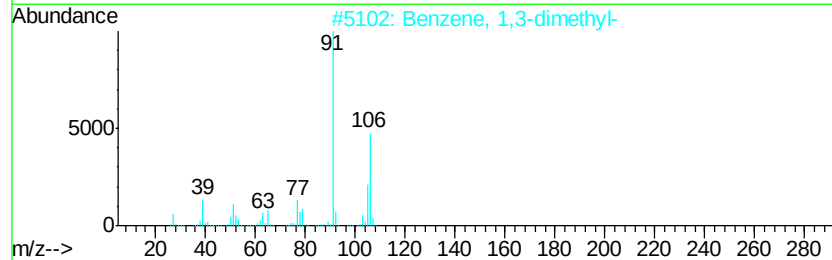
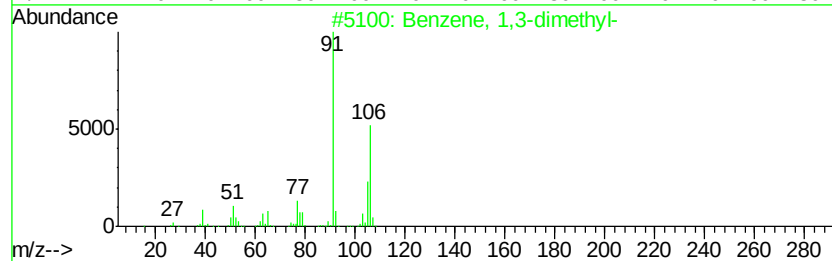
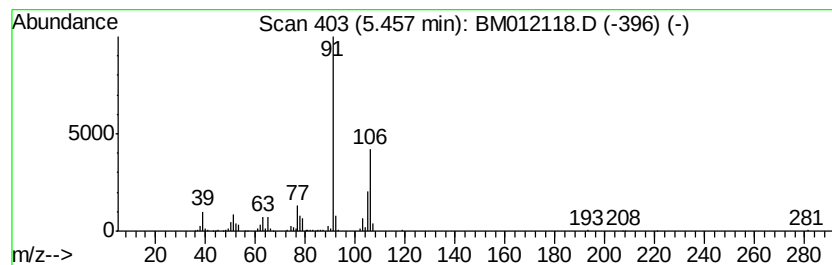
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 24

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.46 | 5.61 ng/ul | 551618 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 95   |
| 2    |    | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 95   |
| 3    |    | o-Xylene               | 106 | C8H10   | 000095-47-6 | 95   |
| 4    |    | p-Xylene               | 106 | C8H10   | 000106-42-3 | 95   |
| 5    |    | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 93   |



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Quant Title : SVOA CALIBRATION

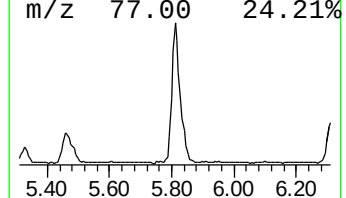
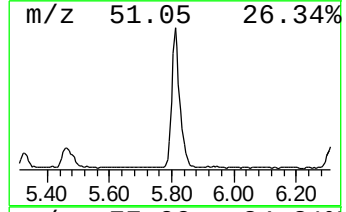
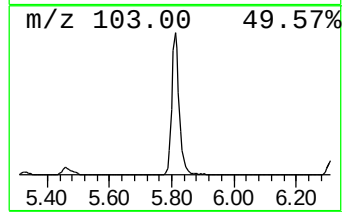
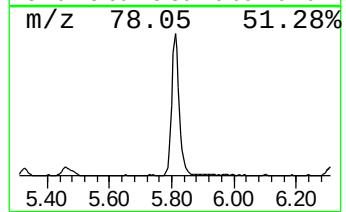
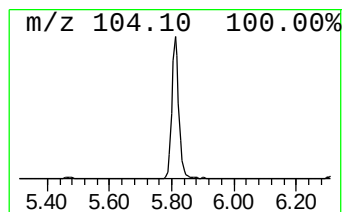
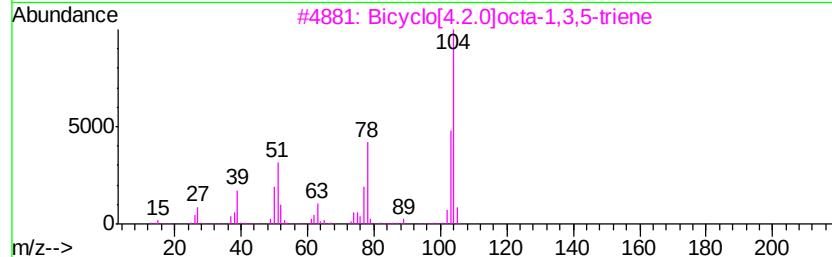
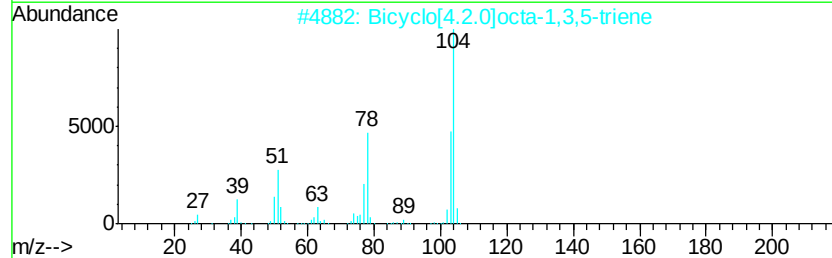
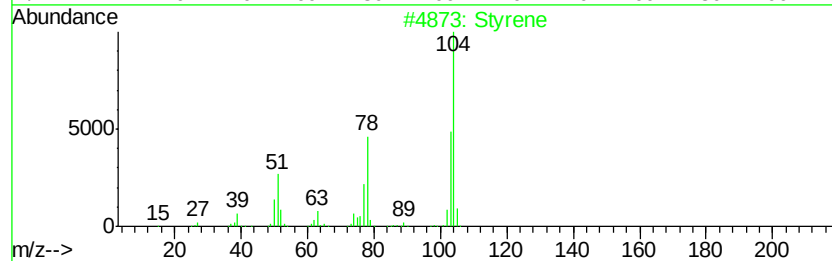
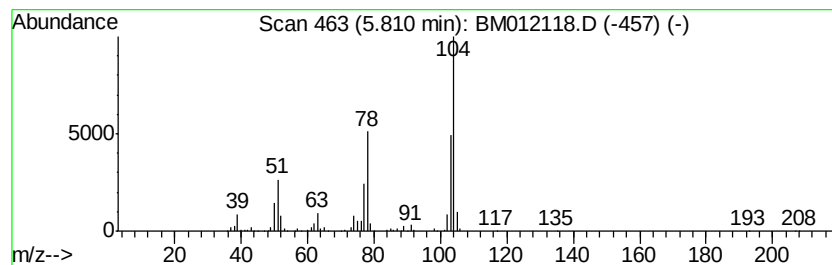
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Styrene Concentration Rank 11

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 5.81 | 16.38 ng/ul | 1611550 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Styrene                         | 104 | C8H8    | 000100-42-5 | 96   |
| 2    |    | Bicyclo[4.2.0]octa-1,3,5-triene | 104 | C8H8    | 000694-87-1 | 96   |
| 3    |    | Bicyclo[4.2.0]octa-1,3,5-triene | 104 | C8H8    | 000694-87-1 | 95   |
| 4    |    | Styrene                         | 104 | C8H8    | 000100-42-5 | 94   |
| 5    |    | Styrene                         | 104 | C8H8    | 000100-42-5 | 93   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

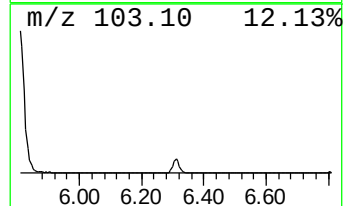
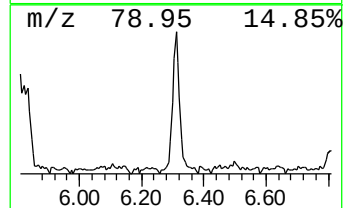
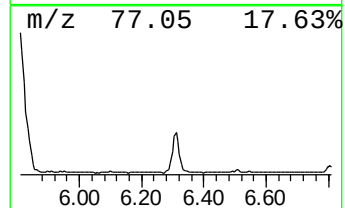
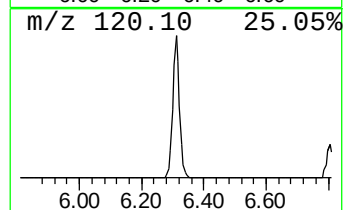
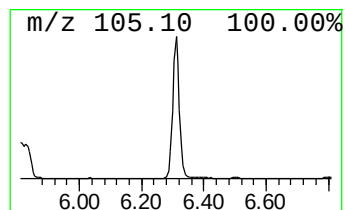
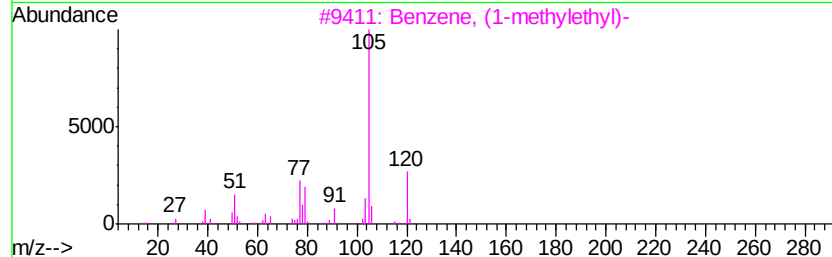
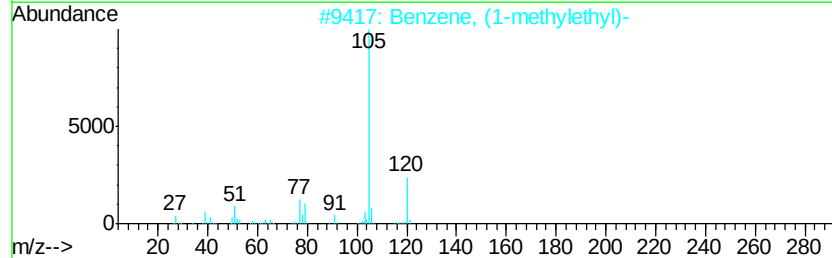
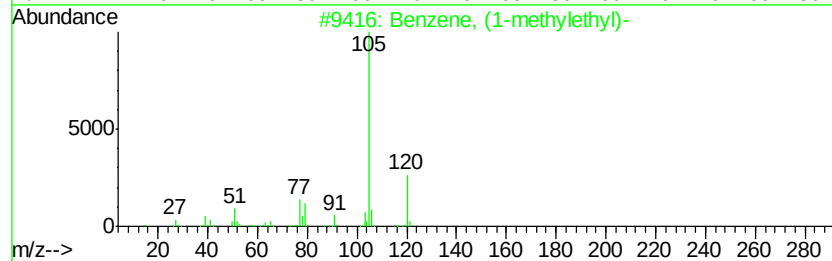
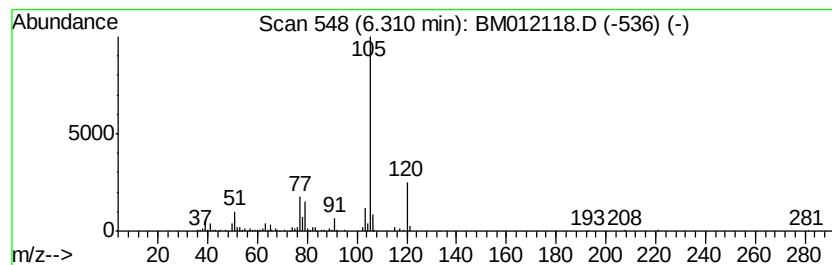
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Quant Title : SVOA CALIBRATION

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

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Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 30

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.31 | 4.11 ng/ul | 404630 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 94   |
| 2    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 94   |
| 3    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 94   |
| 4    |    |   | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 5    |    |   | Benzene, (1-methylethyl)- | 120 | C9H12   | 000098-82-8 | 90   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

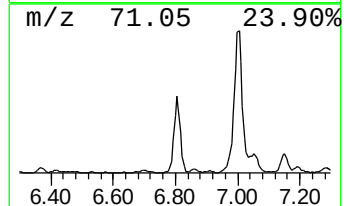
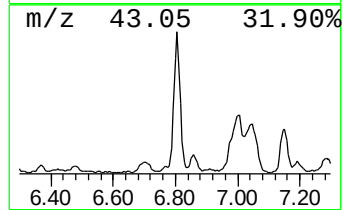
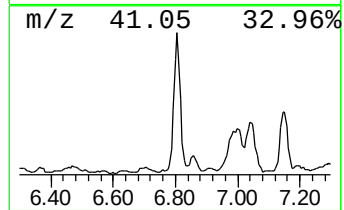
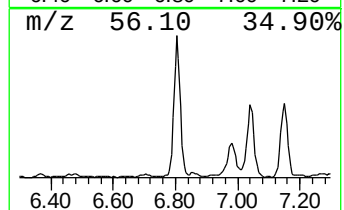
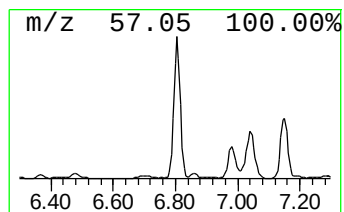
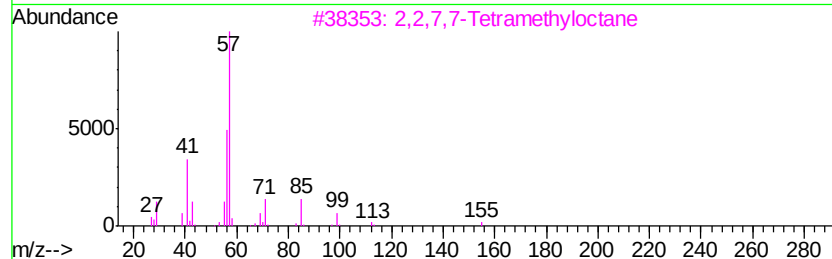
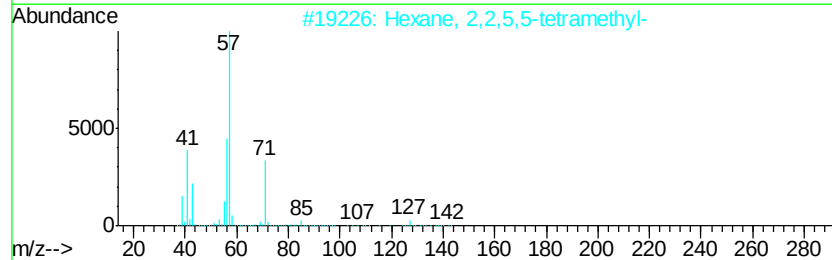
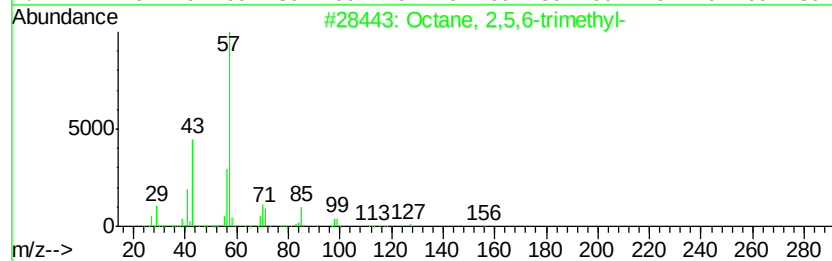
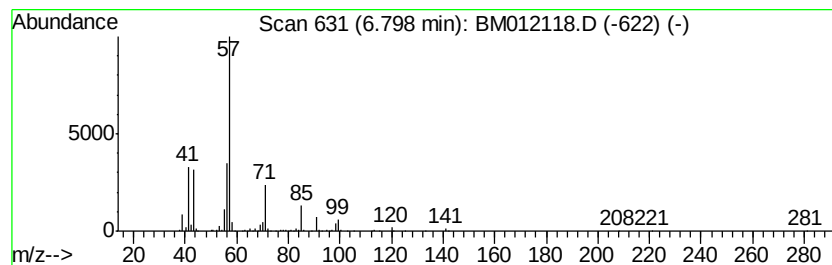
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 6.80 | 13.08 ng/ul | 1287040 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2,5,6-trimethyl-     | 156 | C11H24  | 062016-14-2 | 53   |
| 2    |    | Hexane, 2,2,5,5-tetramethyl- | 142 | C10H22  | 001071-81-4 | 53   |
| 3    |    | 2,2,7,7-Tetramethyloctane    | 170 | C12H26  | 001071-31-4 | 53   |
| 4    |    | Hexane, 2,2,5,5-tetramethyl- | 142 | C10H22  | 001071-81-4 | 47   |
| 5    |    | Decane, 2,6,8-trimethyl-     | 184 | C13H28  | 062108-26-3 | 45   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

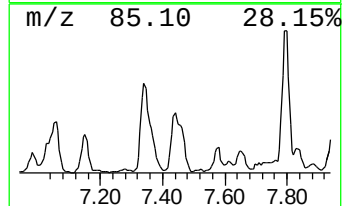
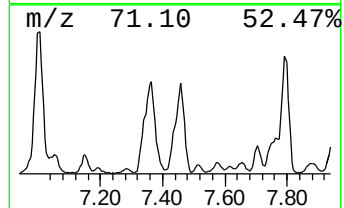
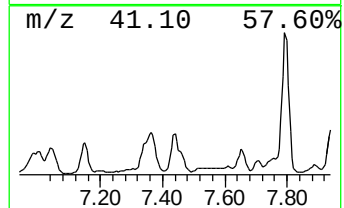
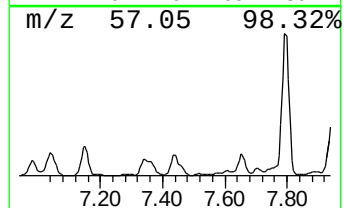
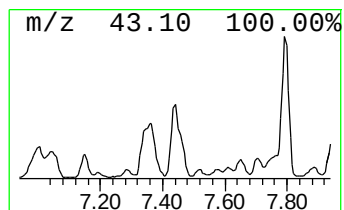
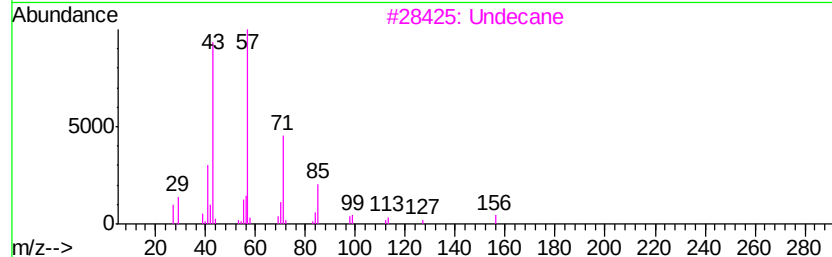
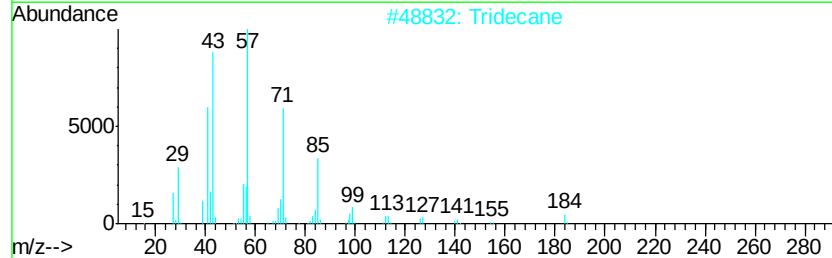
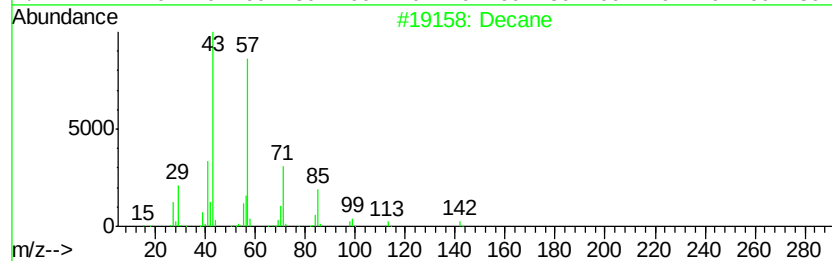
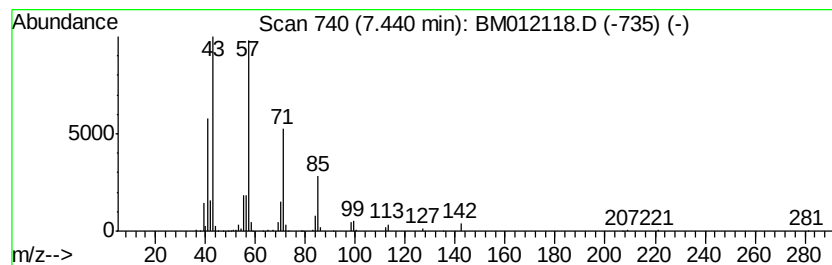
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.44 | 6.04 ng/ul | 594861 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Decane       | 142 | C10H22  | 000124-18-5 | 93   |
| 2    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 90   |
| 3    |    |   | Undecane     | 156 | C11H24  | 001120-21-4 | 83   |
| 4    |    |   | Tridecane    | 184 | C13H28  | 000629-50-5 | 83   |
| 5    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 83   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

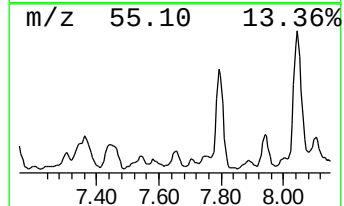
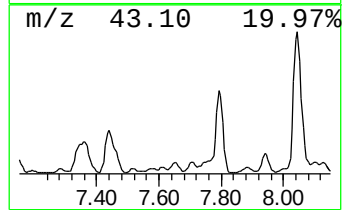
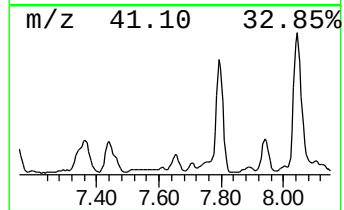
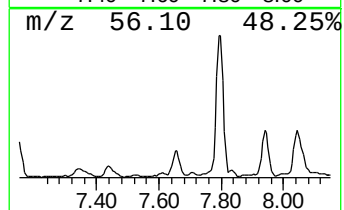
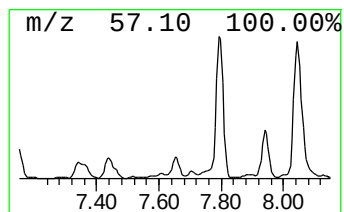
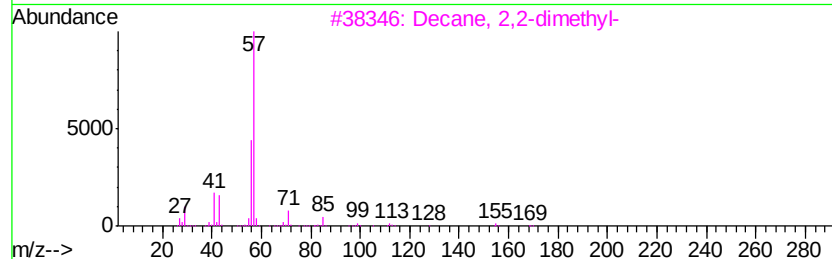
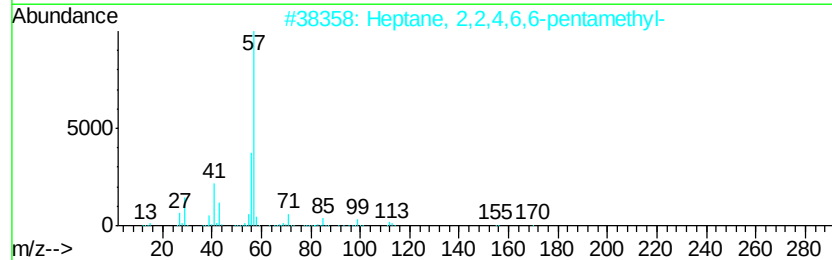
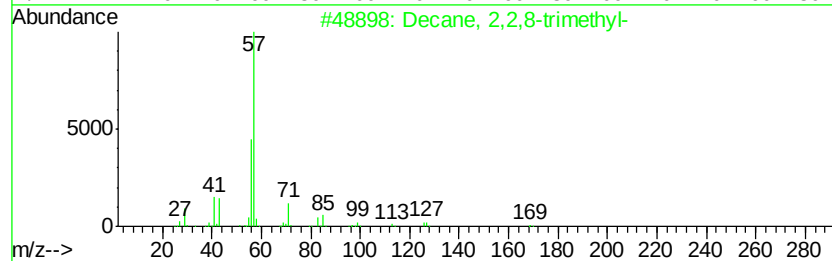
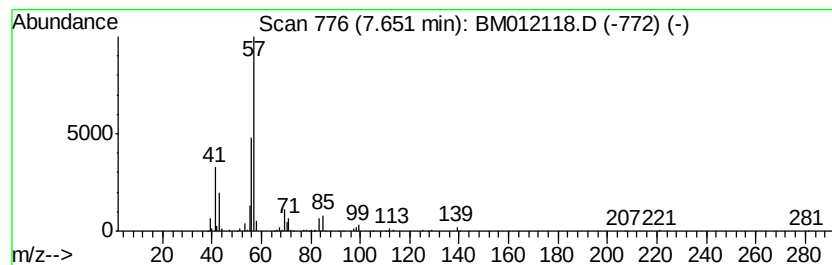
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.65 | 3.69 ng/ul | 363279 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,2,8-trimethyl-        | 184 | C13H28  | 062238-01-1 | 64   |
| 2    |    |   | Heptane, 2,2,4,6,6-pentamethyl- | 170 | C12H26  | 013475-82-6 | 59   |
| 3    |    |   | Decane, 2,2-dimethyl-           | 170 | C12H26  | 017302-37-3 | 59   |
| 4    |    |   | Heptane, 2,2,4,6,6-pentamethyl- | 170 | C12H26  | 013475-82-6 | 50   |
| 5    |    |   | Pentane, 2,2,4,4-tetramethyl-   | 128 | C9H20   | 001070-87-7 | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

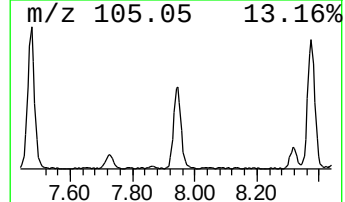
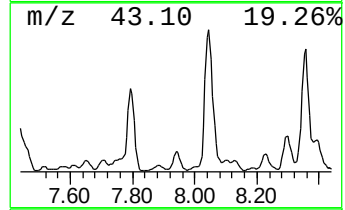
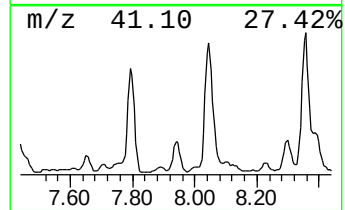
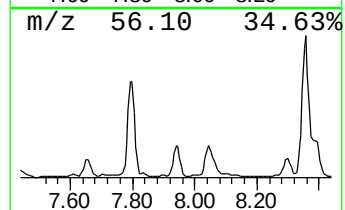
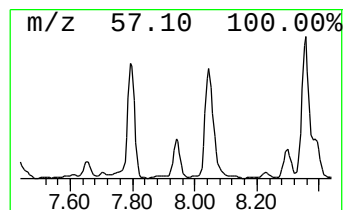
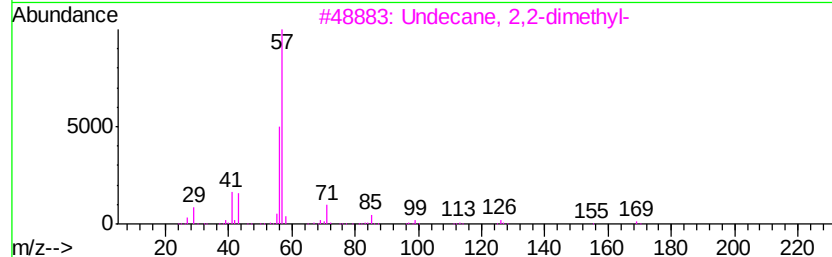
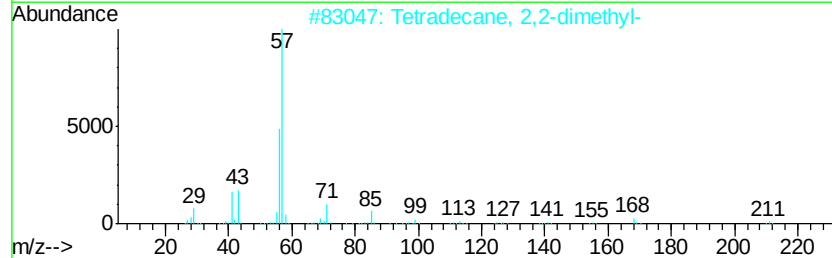
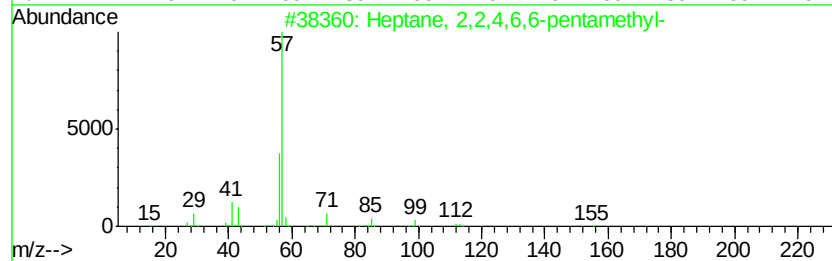
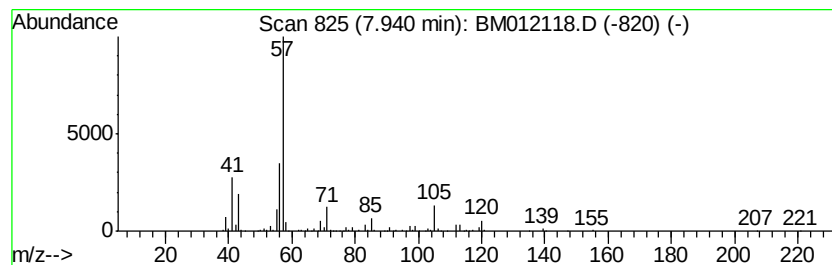
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.94 | 8.47 ng/ul | 833122 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 2,2,4,6,6-pentamethyl- | 170 | C12H26  | 013475-82-6 | 53   |
| 2    |    | Tetradecane, 2,2-dimethyl-      | 226 | C16H34  | 059222-86-5 | 53   |
| 3    |    | Undecane, 2,2-dimethyl-         | 184 | C13H28  | 017312-64-0 | 53   |
| 4    |    | Decane, 2,2,6-trimethyl-        | 184 | C13H28  | 062237-97-2 | 50   |
| 5    |    | Hexane, 2,2,5-trimethyl-        | 128 | C9H20   | 003522-94-9 | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

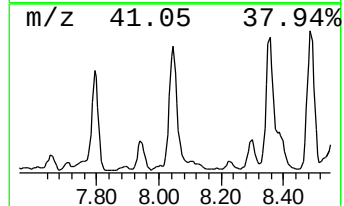
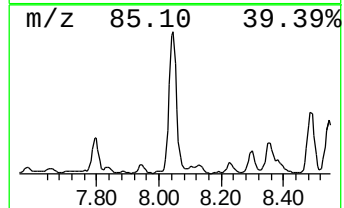
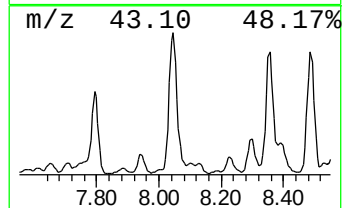
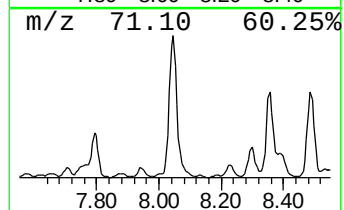
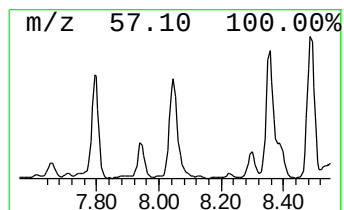
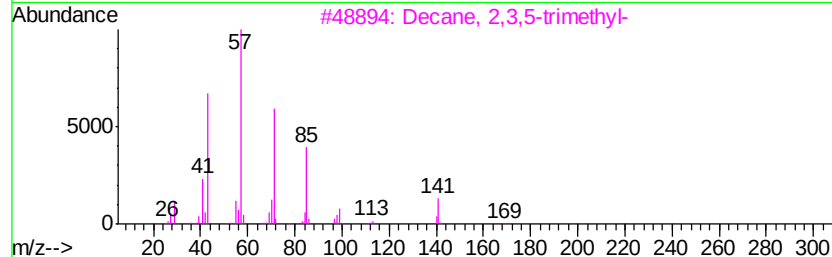
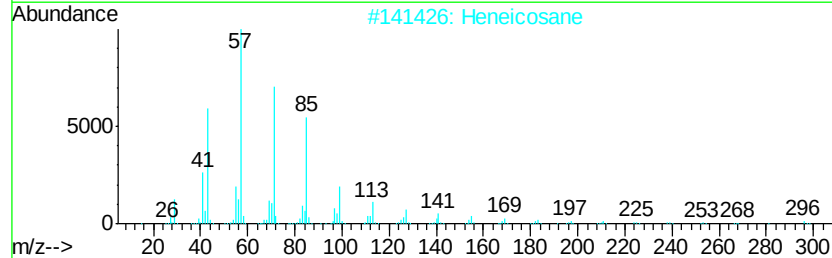
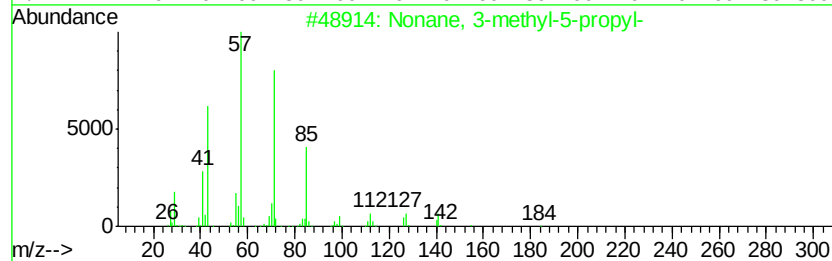
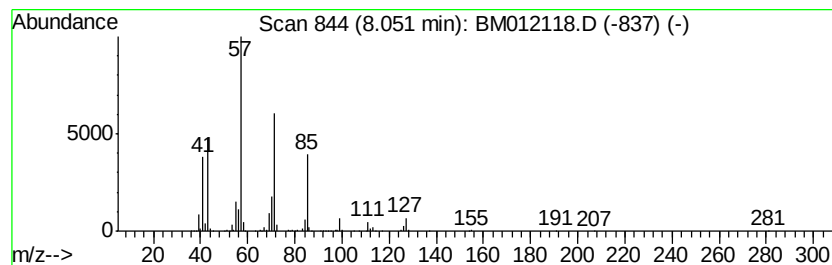
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.05 | 38.27 ng/ul | 3765680 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 3-methyl-5-propyl- | 184 | C13H28  | 031081-18-2 | 78   |
| 2    |    | Heneicosane                | 296 | C21H44  | 000629-94-7 | 78   |
| 3    |    | Decane, 2,3,5-trimethyl-   | 184 | C13H28  | 062238-11-3 | 72   |
| 4    |    | Undecane, 4,7-dimethyl-    | 184 | C13H28  | 017301-32-5 | 64   |
| 5    |    | Nonane, 5-butyl-           | 184 | C13H28  | 017312-63-9 | 59   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

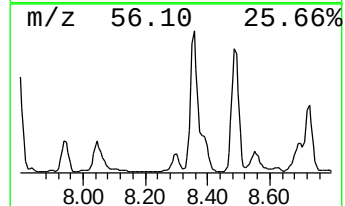
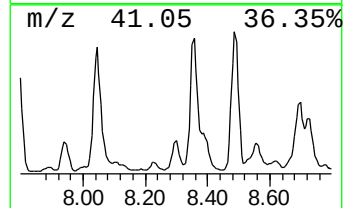
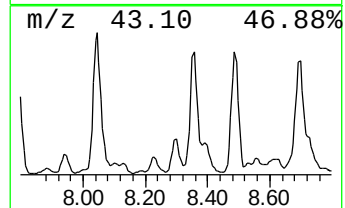
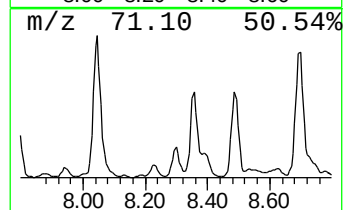
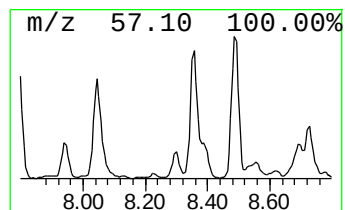
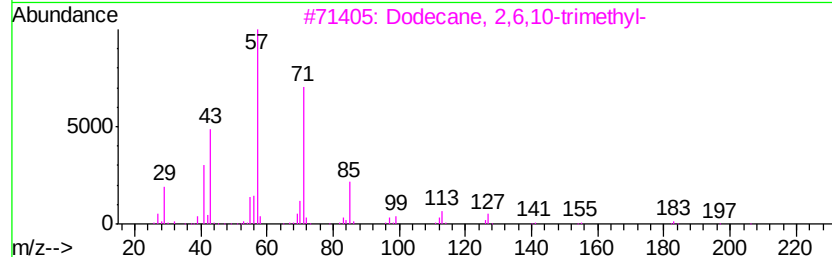
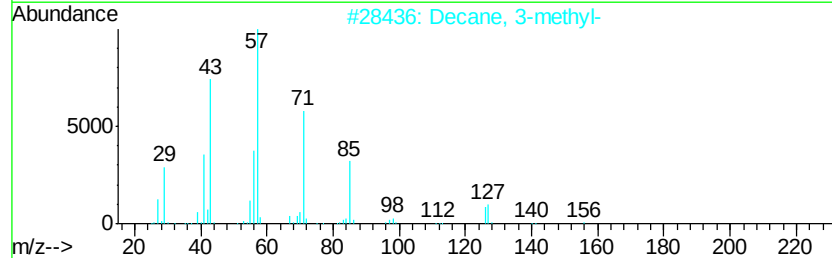
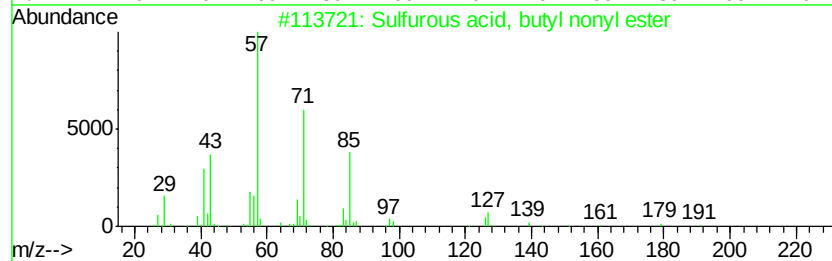
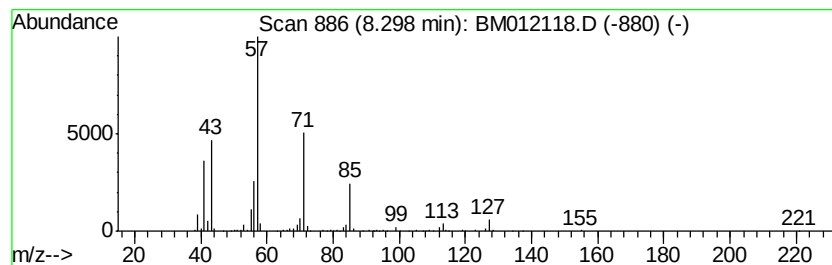
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Sulfurous acid, butyl nonyl... Concentration Rank 18

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.30 | 8.60 ng/ul | 846153 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|-----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Sulfurous acid, butyl nonyl ester | 264 | C13H28O3S | 1000309-17-6 | 64   |
| 2    |    | Decane, 3-methyl-                 | 156 | C11H24    | 013151-34-3  | 64   |
| 3    |    | Dodecane, 2,6,10-trimethyl-       | 212 | C15H32    | 003891-98-3  | 64   |
| 4    |    | Nonane, 3,7-dimethyl-             | 156 | C11H24    | 017302-32-8  | 64   |
| 5    |    | Octane, 2,6-dimethyl-             | 142 | C10H22    | 002051-30-1  | 64   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

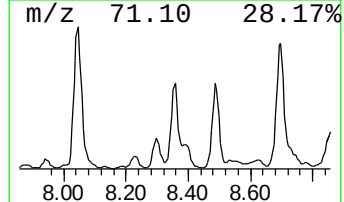
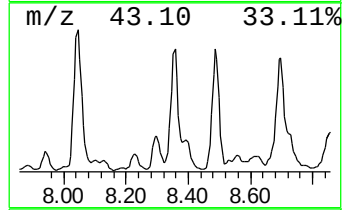
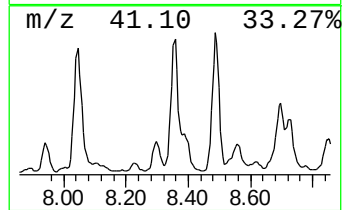
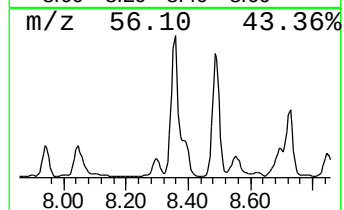
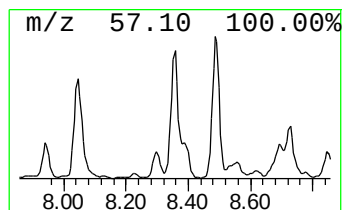
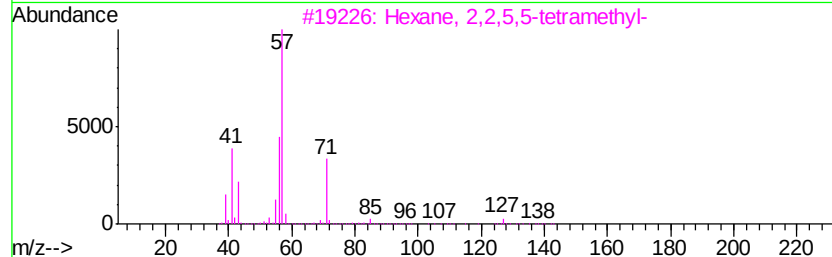
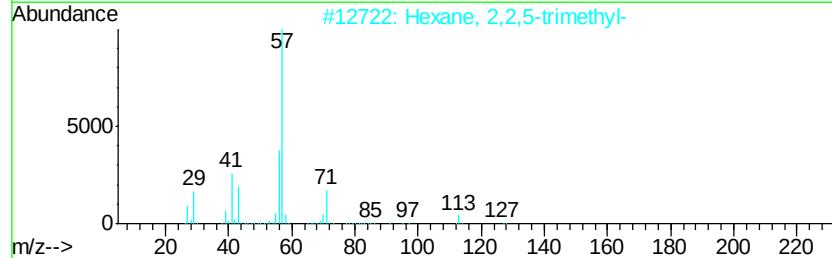
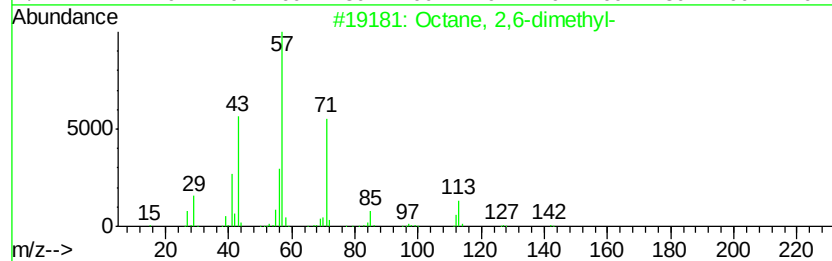
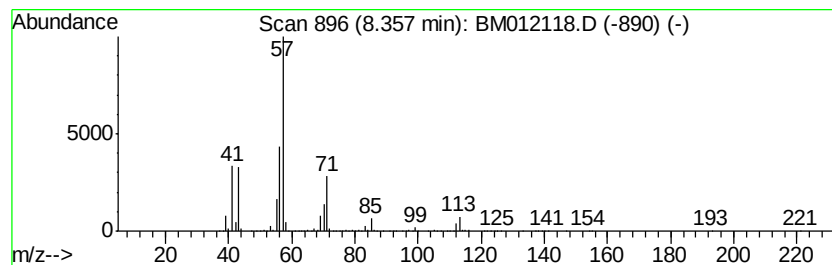
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.36 | 50.55 ng/ul | 4975060 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-             | 142 | C10H22  | 002051-30-1 | 72   |
| 2    |    | Hexane, 2,2,5-trimethyl-          | 128 | C9H20   | 003522-94-9 | 59   |
| 3    |    | Hexane, 2,2,5,5-tetramethyl-      | 142 | C10H22  | 001071-81-4 | 59   |
| 4    |    | Heptane, 5-ethyl-2,2,3-trimethyl- | 170 | C12H26  | 062199-06-8 | 56   |
| 5    |    | Heptane, 2,2-dimethyl-            | 128 | C9H20   | 001071-26-7 | 53   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

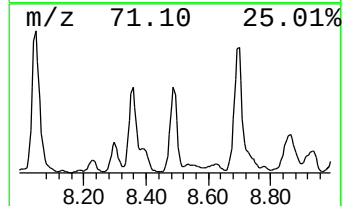
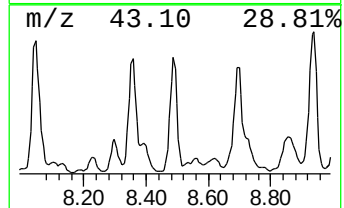
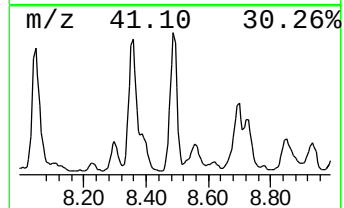
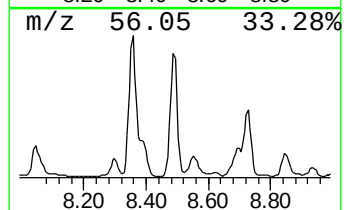
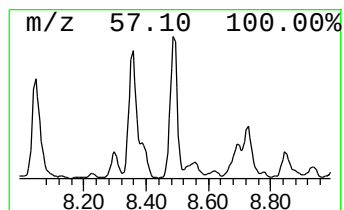
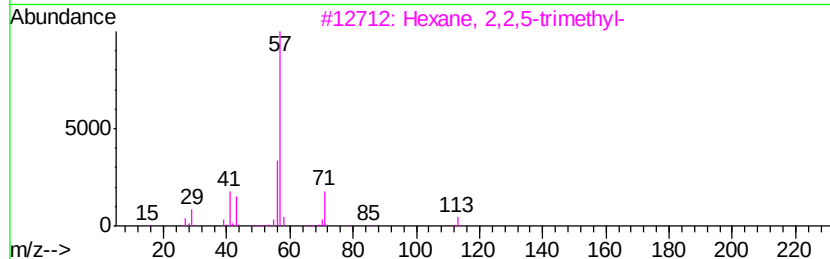
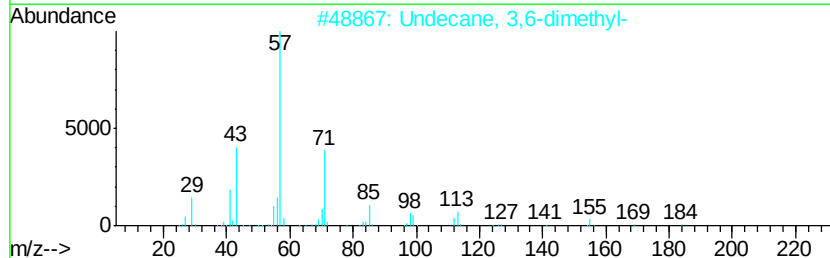
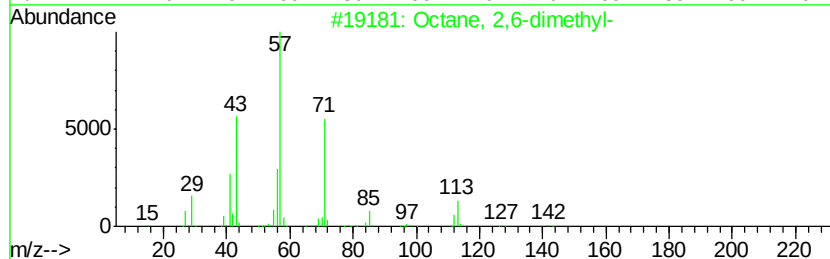
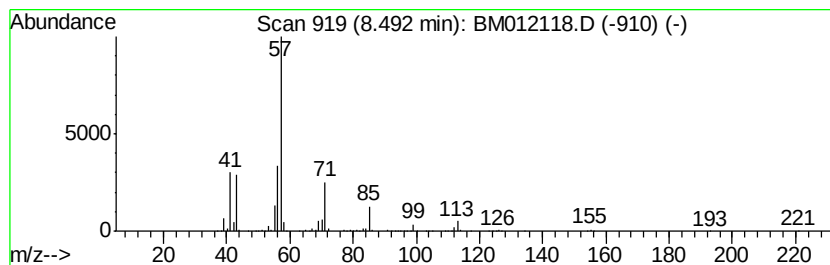
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.49 | 39.87 ng/ul | 3923370 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-    | 142 | C10H22  | 002051-30-1 | 72   |
| 2    |    | Undecane, 3,6-dimethyl-  | 184 | C13H28  | 017301-28-9 | 64   |
| 3    |    | Hexane, 2,2,5-trimethyl- | 128 | C9H20   | 003522-94-9 | 64   |
| 4    |    | Octane, 2,4,6-trimethyl- | 156 | C11H24  | 062016-37-9 | 59   |
| 5    |    | Octane, 2,5,6-trimethyl- | 156 | C11H24  | 062016-14-2 | 53   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

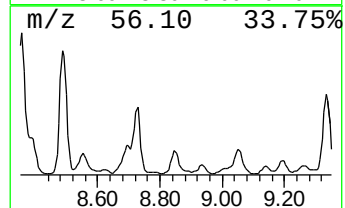
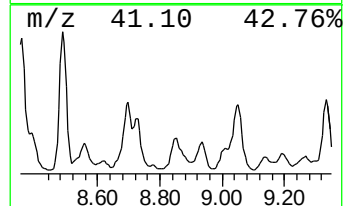
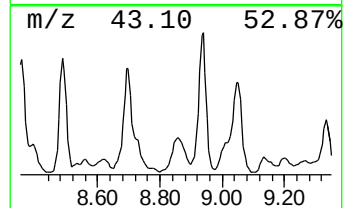
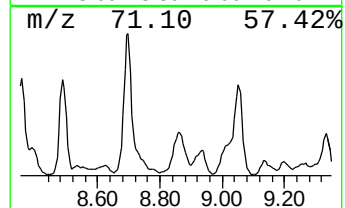
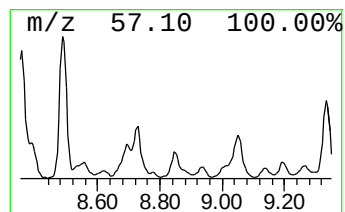
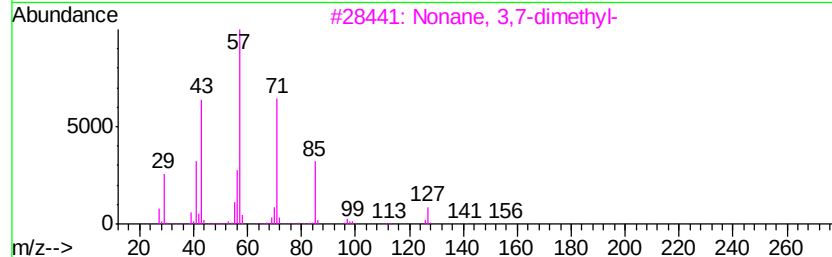
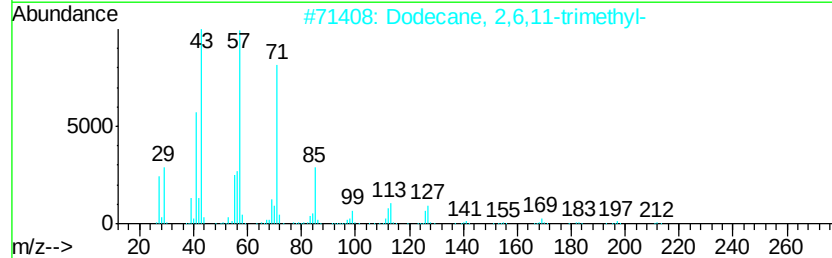
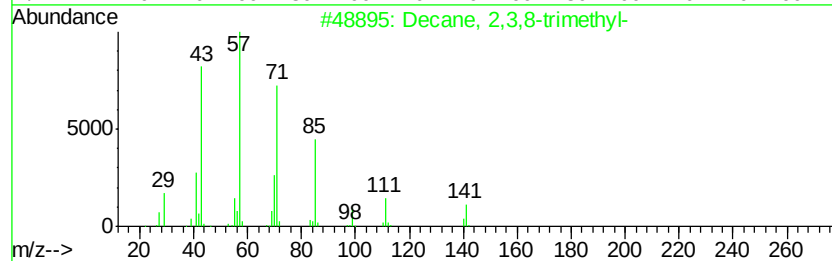
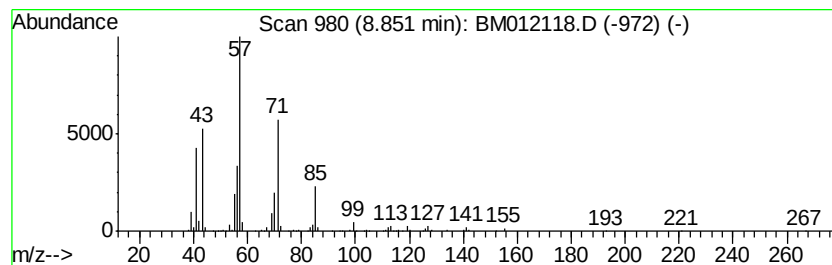
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.85 | 14.72 ng/ul | 1448830 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | Tentative ID                       | MW  | MolForm  | CAS#         | Qual |
|------|----|------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Decane, 2,3,8-trimethyl-           | 184 | C13H28   | 062238-14-6  | 59   |
| 2    |    | Dodecane, 2,6,11-trimethyl-        | 212 | C15H32   | 031295-56-4  | 53   |
| 3    |    | Nonane, 3,7-dimethyl-              | 156 | C11H24   | 017302-32-8  | 53   |
| 4    |    | Undecane, 6-ethyl-                 | 184 | C13H28   | 017312-60-6  | 50   |
| 5    |    | Oxalic acid, isobutyl pentyl ester | 216 | C11H20O4 | 1000309-37-0 | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

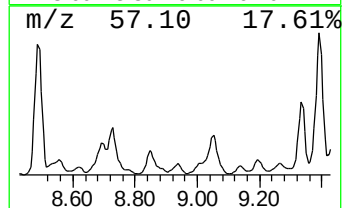
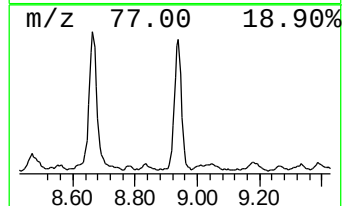
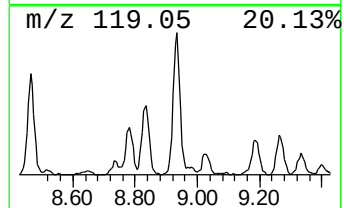
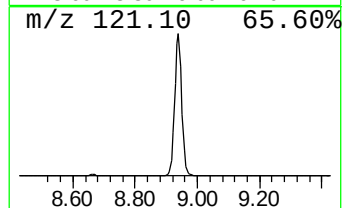
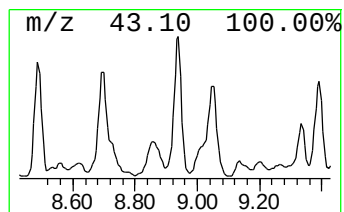
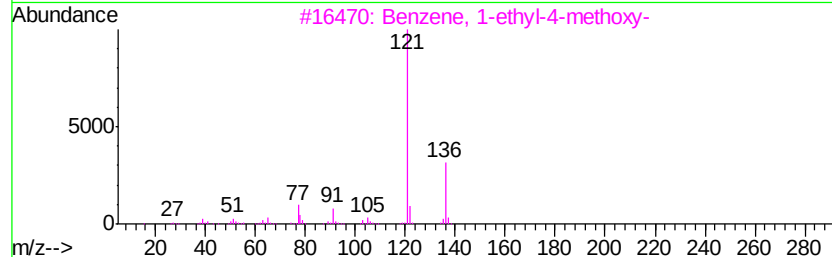
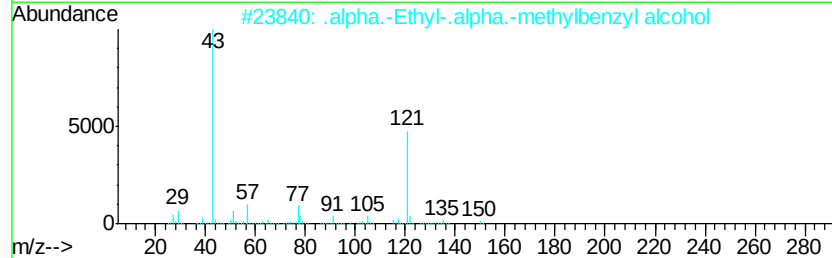
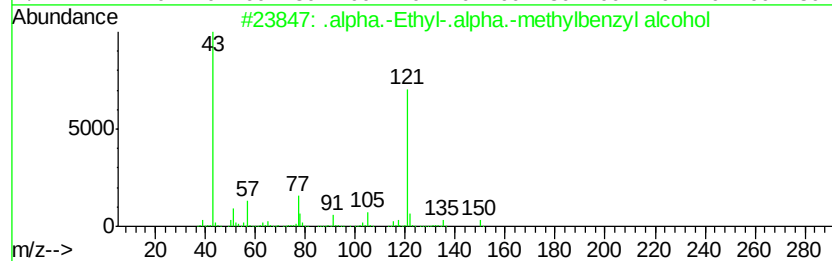
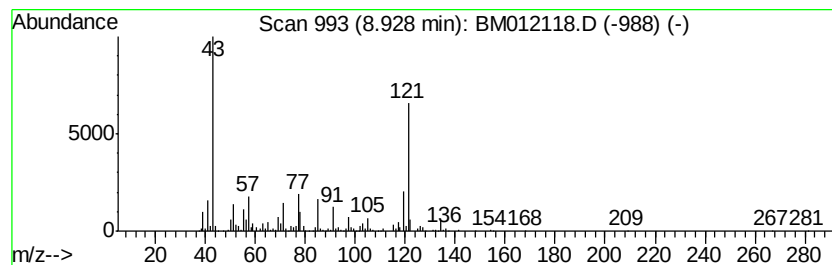
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 unknown-01 Concentration Rank 7

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.93 | 22.39 ng/ul | 2203770 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of                          | 5                                  | Tentative ID | MW      | MolForm     | CAS# | Qual |
|------|-----------------------------|------------------------------------|--------------|---------|-------------|------|------|
| 1    | .                           | alpha.-Ethyl-.alpha.-methylbenz... | 150          | C10H14O | 001565-75-9 | 47   |      |
| 2    | .                           | alpha.-Ethyl-.alpha.-methylbenz... | 150          | C10H14O | 001565-75-9 | 47   |      |
| 3    | Benzene, 1-ethyl-4-methoxy- |                                    | 136          | C9H12O  | 001515-95-3 | 46   |      |
| 4    | Benzene, 1-ethyl-4-methoxy- |                                    | 136          | C9H12O  | 001515-95-3 | 43   |      |
| 5    | dl-2-Phenyl-1,2-propanediol |                                    | 152          | C9H12O2 | 004217-66-7 | 43   |      |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

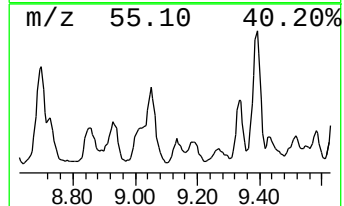
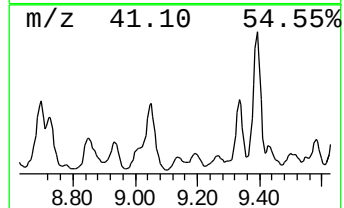
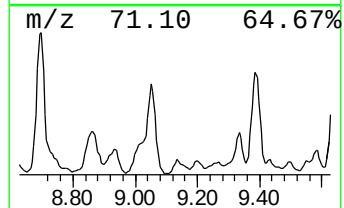
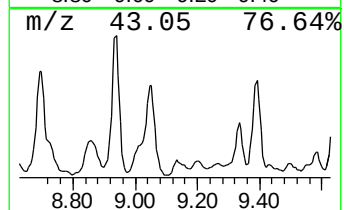
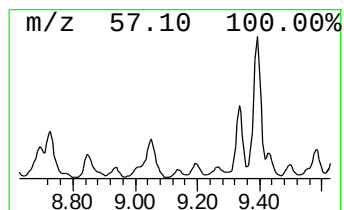
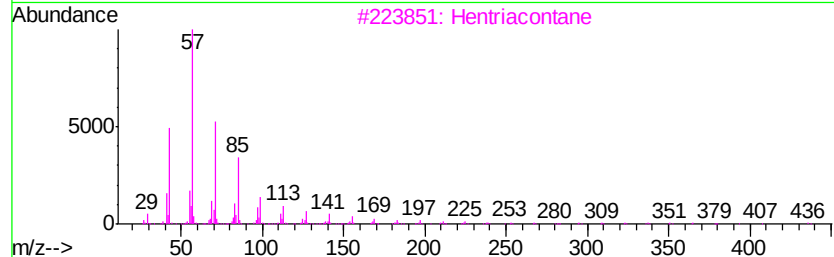
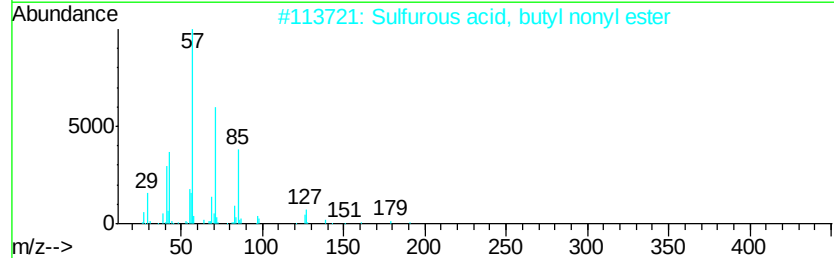
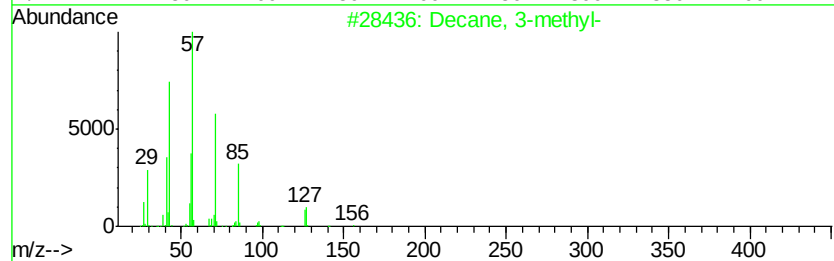
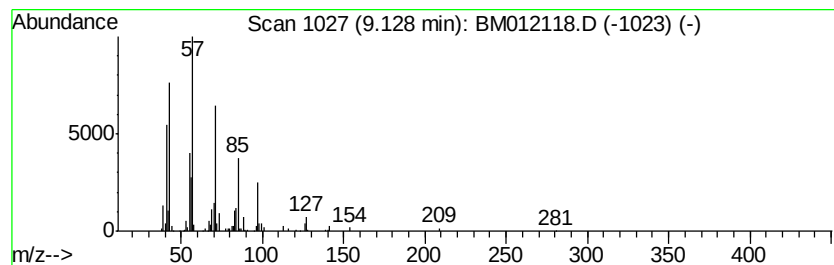
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.13 | 5.89 ng/ul | 579570 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-----------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 3-methyl-                 | 156 | C11H24    | 013151-34-3  | 64   |
| 2    |    |   | Sulfurous acid, butyl nonyl ester | 264 | C13H28O3S | 1000309-17-6 | 59   |
| 3    |    |   | Hentriacontane                    | 437 | C31H64    | 000630-04-6  | 53   |
| 4    |    |   | Nonane, 3,7-dimethyl-             | 156 | C11H24    | 017302-32-8  | 53   |
| 5    |    |   | Decane, 2,4-dimethyl-             | 170 | C12H26    | 002801-84-5  | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

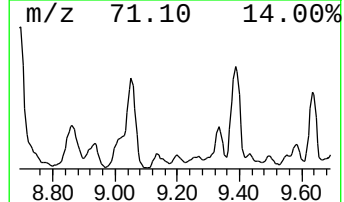
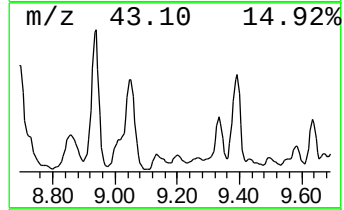
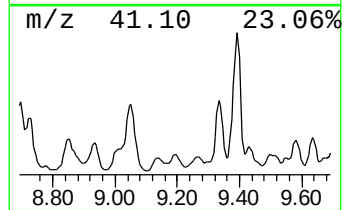
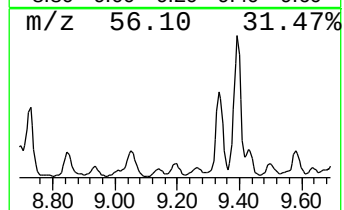
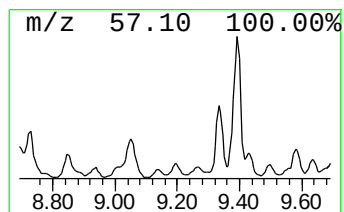
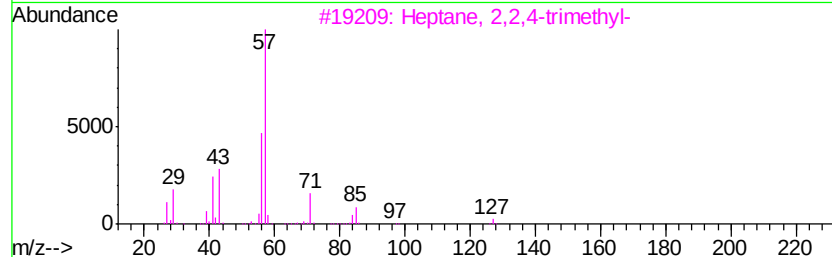
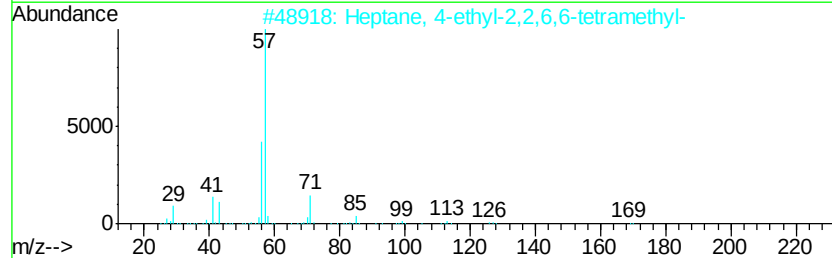
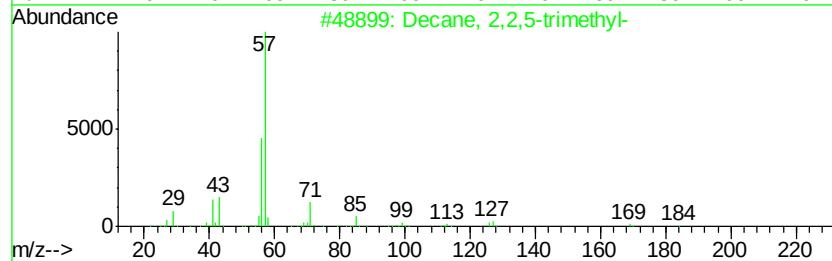
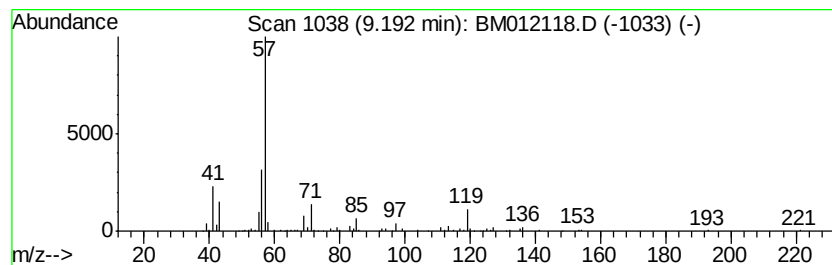
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.19 | 5.41 ng/ul | 532554 | 1,4-Dichlorobenzene-d4 | 7.83 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,2,5-trimethyl-            | 184 | C13H28  | 062237-96-1 | 53   |
| 2    |    |   | Heptane, 4-ethyl-2,2,6,6-tetrame... | 184 | C13H28  | 062108-31-0 | 50   |
| 3    |    |   | Heptane, 2,2,4-trimethyl-           | 142 | C10H22  | 014720-74-2 | 50   |
| 4    |    |   | Octane, 2,2-dimethyl-               | 142 | C10H22  | 015869-87-1 | 50   |
| 5    |    |   | Decane, 2,2,7-trimethyl-            | 184 | C13H28  | 062237-99-4 | 47   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

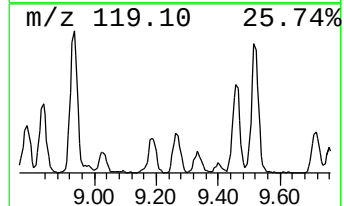
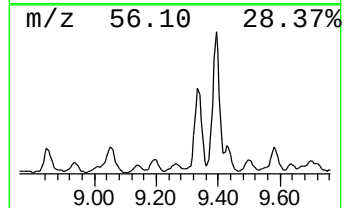
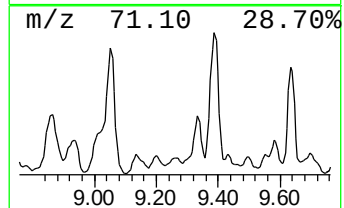
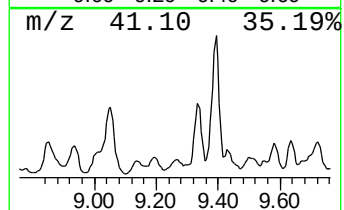
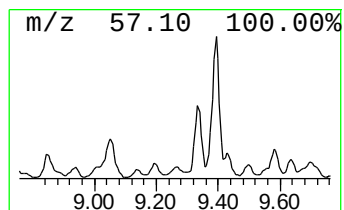
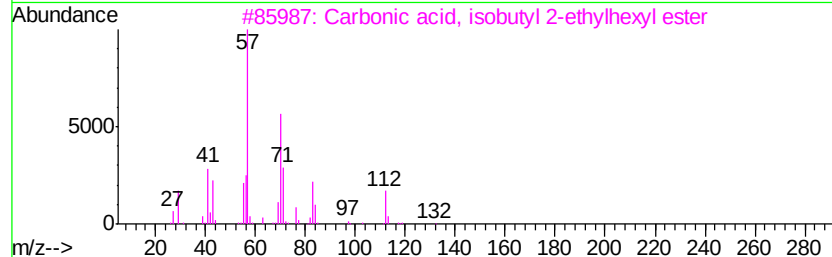
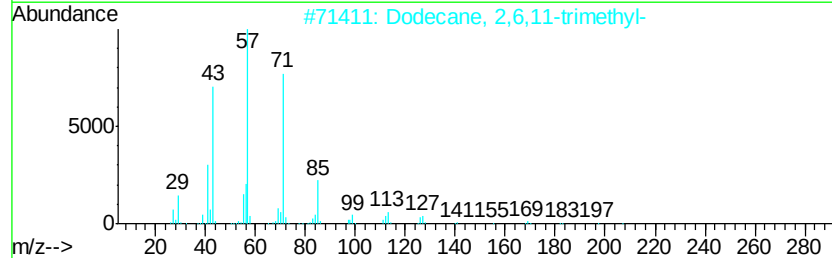
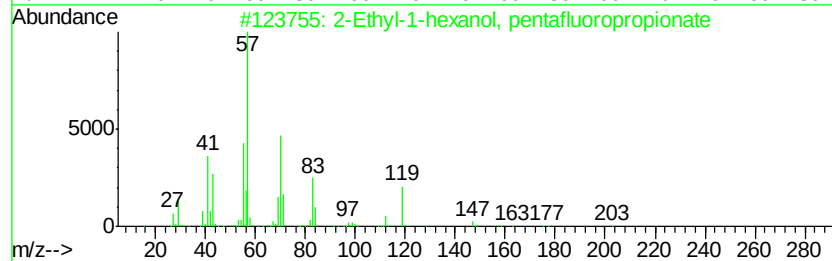
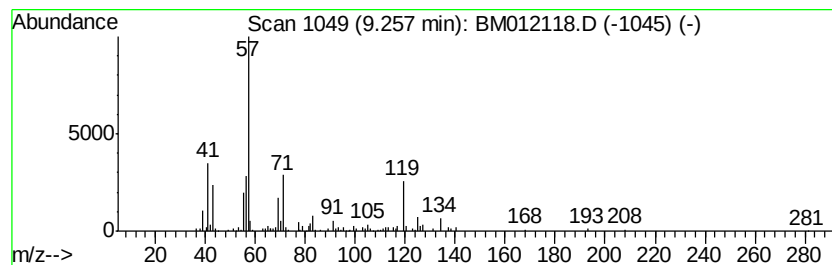
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 17 2-Ethyl-1-hexanol, pentaflu... Concentration Rank 43

| R.T.    | EstConc                             | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------------|----------------|------------------|-----------|
| 9.26    | 2.00 ng/ul                          | 344808         | Napthalene-d8    | 10.64     |
| Hit# of | 5                                   | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 2-Ethyl-1-hexanol, pentafluoropr... | 276 C11H17F5O2 | 1000365-51-1     | 72        |
| 2       | Dodecane, 2,6,11-trimethyl-         | 212 C15H32     | 031295-56-4      | 43        |
| 3       | Carbonic acid, isobutyl 2-ethylh... | 230 C13H26O3   | 1000357-82-9     | 38        |
| 4       | 2-Ethyl-1-hexanol, heptafluorobu... | 326 C12H17F7O2 | 1000365-16-0     | 28        |
| 5       | 1-Pentanol, 4-methyl-2-propyl-      | 144 C9H20O     | 054004-41-0      | 27        |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

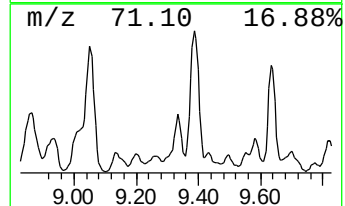
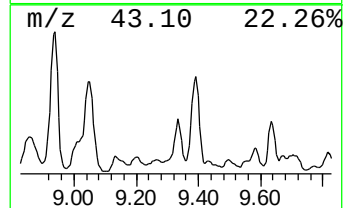
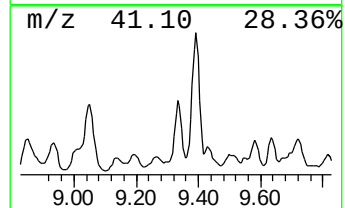
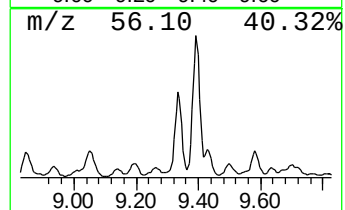
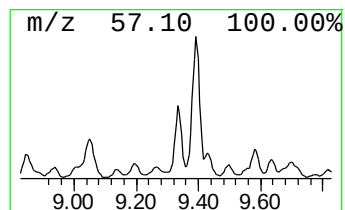
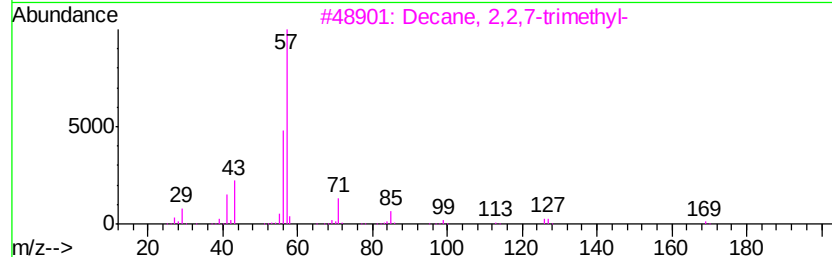
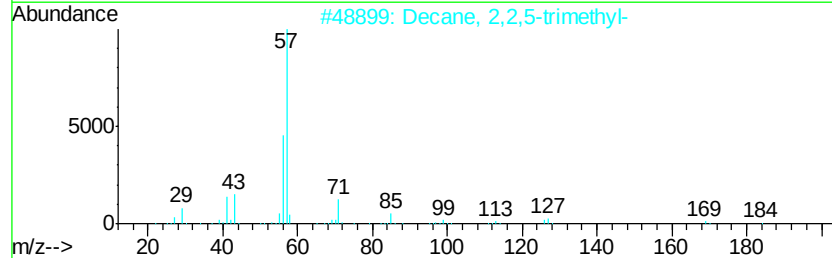
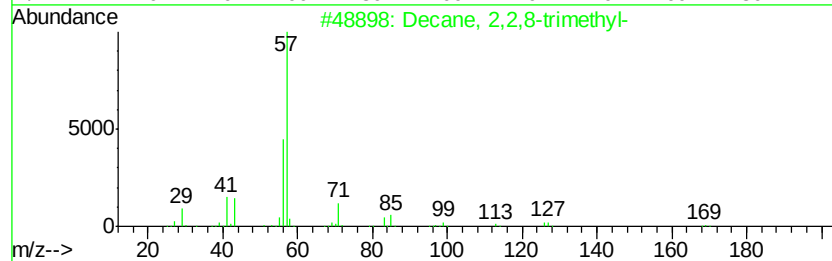
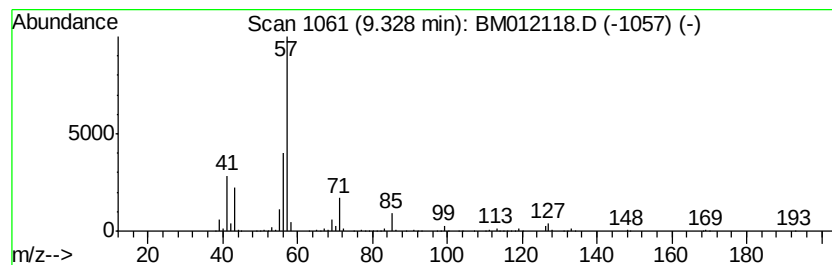
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T. | EstConc    | Area    | Relative to ISTD | R.T.  |
|------|------------|---------|------------------|-------|
| 9.33 | 9.12 ng/ul | 1569940 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,2,8-trimethyl-   | 184 | C13H28  | 062238-01-1 | 83   |
| 2    |    |   | Decane, 2,2,5-trimethyl-   | 184 | C13H28  | 062237-96-1 | 83   |
| 3    |    |   | Decane, 2,2,7-trimethyl-   | 184 | C13H28  | 062237-99-4 | 78   |
| 4    |    |   | Tetradecane, 2,2-dimethyl- | 226 | C16H34  | 059222-86-5 | 78   |
| 5    |    |   | Decane, 2,2-dimethyl-      | 170 | C12H26  | 017302-37-3 | 78   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

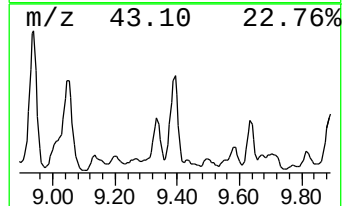
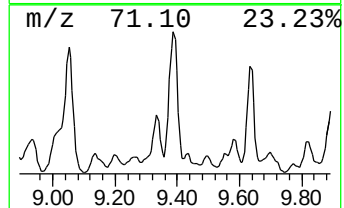
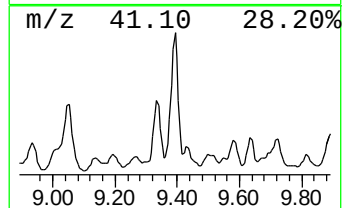
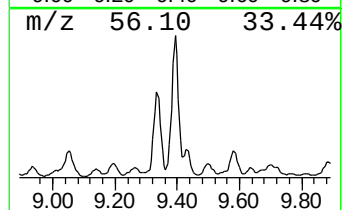
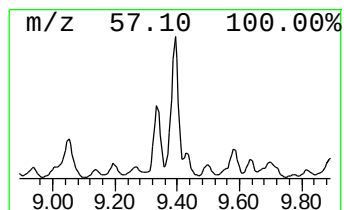
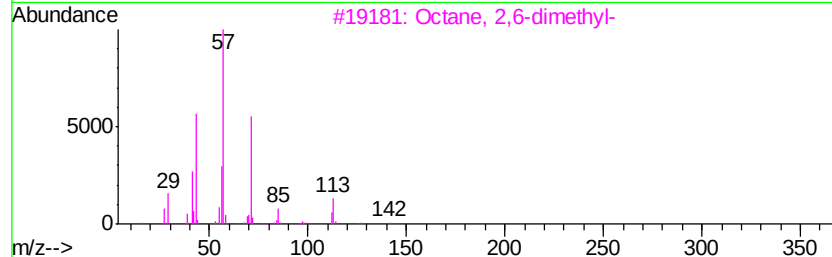
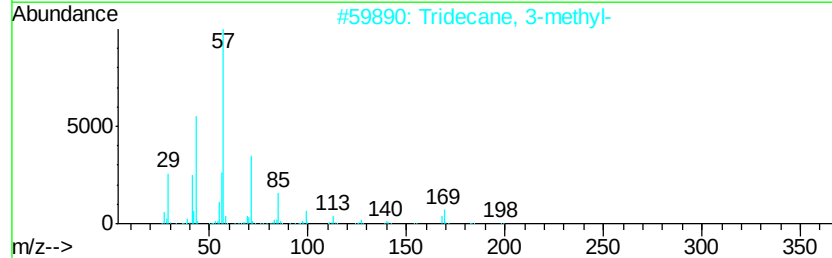
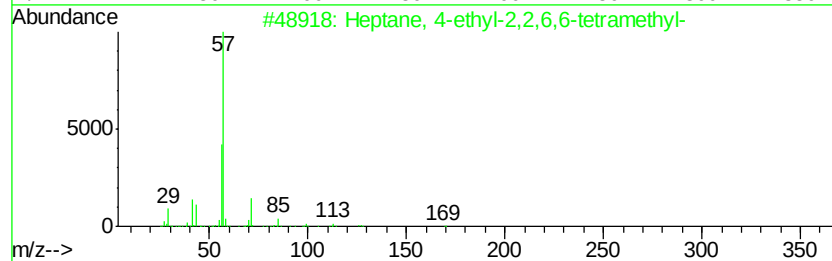
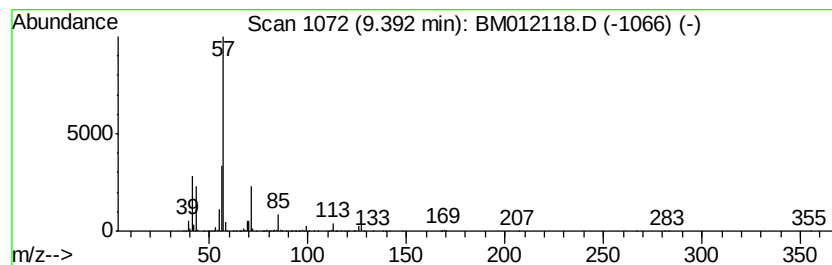
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T. | EstConc     | Area    | Relative to ISTD | R.T.  |
|------|-------------|---------|------------------|-------|
| 9.39 | 20.71 ng/ul | 3565930 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane, 4-ethyl-2,2,6,6-tetrame... | 184 | C13H28  | 062108-31-0 | 72   |
| 2    |    | Tridecane, 3-methyl-                | 198 | C14H30  | 006418-41-3 | 72   |
| 3    |    | Octane, 2,6-dimethyl-               | 142 | C10H22  | 002051-30-1 | 64   |
| 4    |    | Decane, 2,6,8-trimethyl-            | 184 | C13H28  | 062108-26-3 | 64   |
| 5    |    | Decane, 2,2,5-trimethyl-            | 184 | C13H28  | 062237-96-1 | 64   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

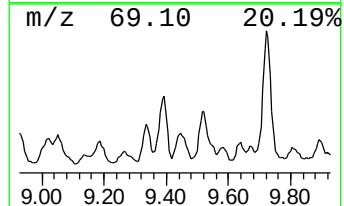
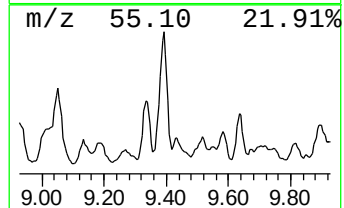
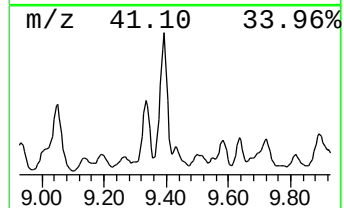
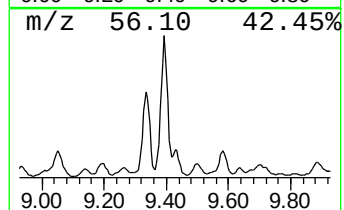
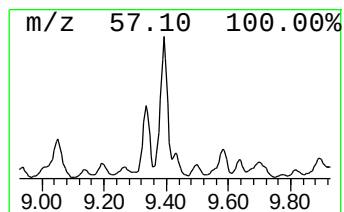
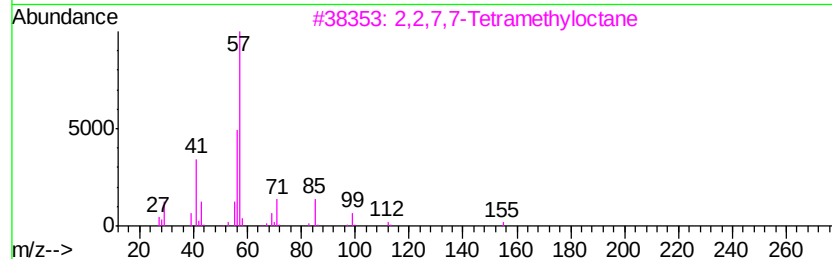
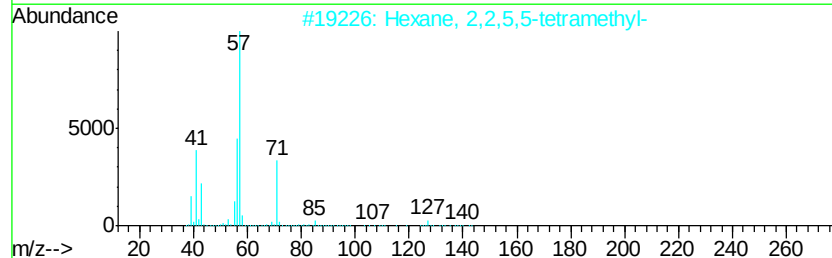
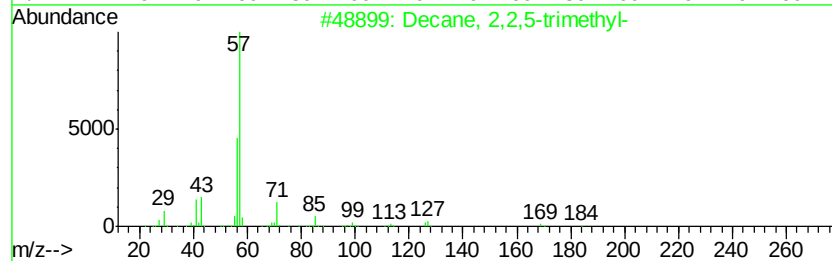
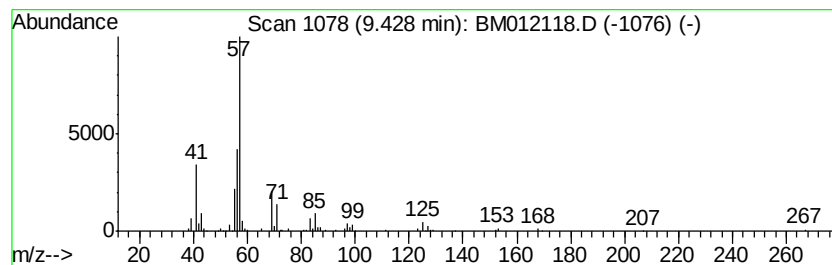
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.43 | 3.16 ng/ul | 543987 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,2,5-trimethyl-     | 184 | C13H28  | 062237-96-1 | 53   |
| 2    |    |   | Hexane, 2,2,5,5-tetramethyl- | 142 | C10H22  | 001071-81-4 | 50   |
| 3    |    |   | 2,2,7,7-Tetramethyloctane    | 170 | C12H26  | 001071-31-4 | 50   |
| 4    |    |   | Decanal                      | 156 | C10H20O | 000112-31-2 | 47   |
| 5    |    |   | 1-Octanol, 2-butyl-          | 186 | C12H26O | 003913-02-8 | 45   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

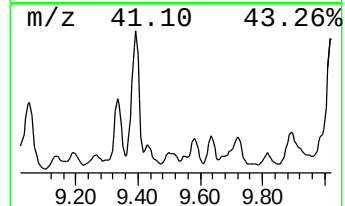
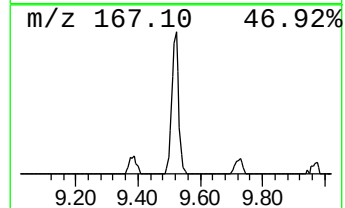
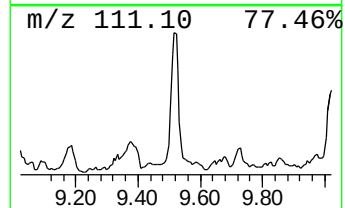
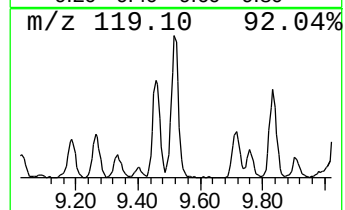
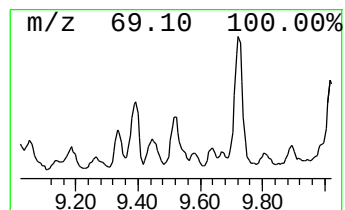
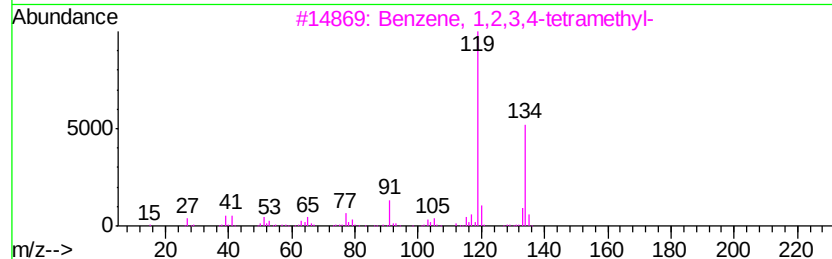
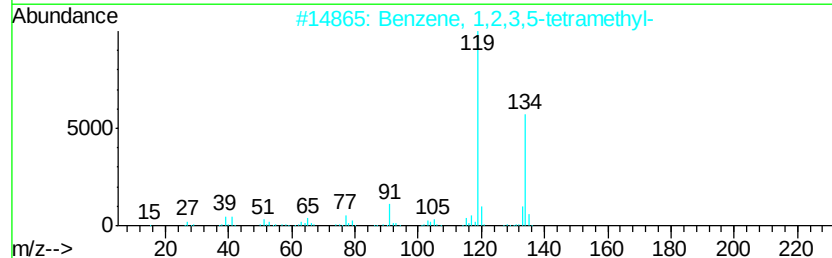
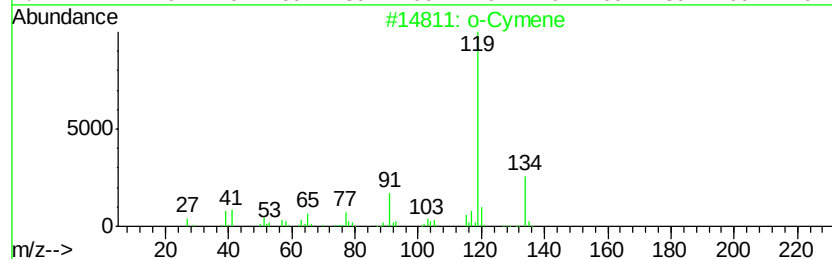
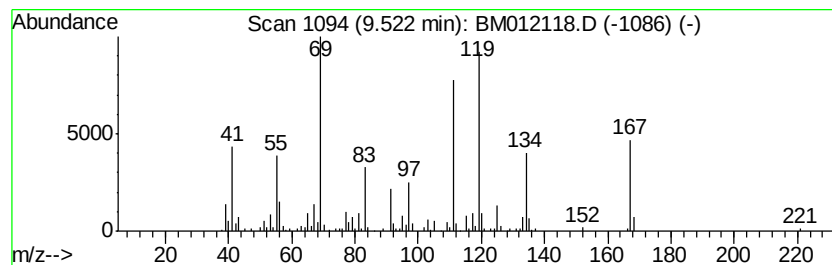
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 o-Cymene Concentration Rank 35

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.52 | 3.43 ng/ul | 590182 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | o-Cymene                      | 134 | C10H14  | 000527-84-4 | 70   |
| 2    |    | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14  | 000527-53-7 | 38   |
| 3    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 38   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14  | 000488-23-3 | 38   |
| 5    |    | p-Cymene                      | 134 | C10H14  | 000099-87-6 | 38   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

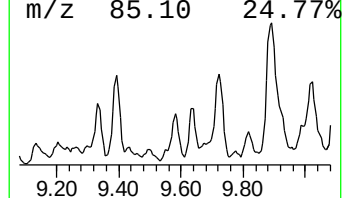
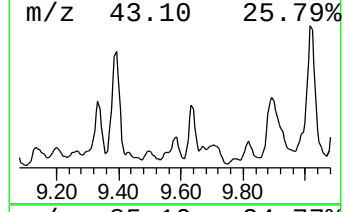
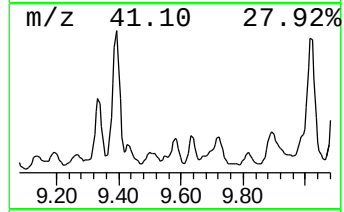
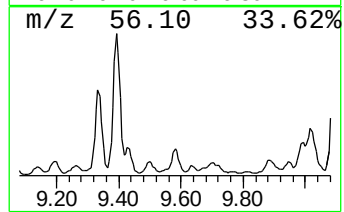
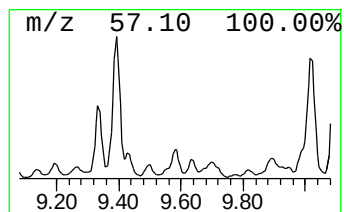
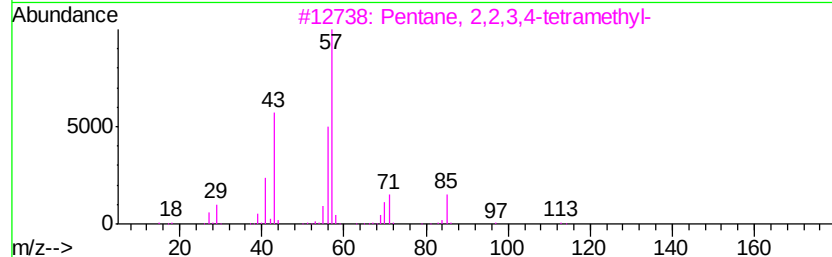
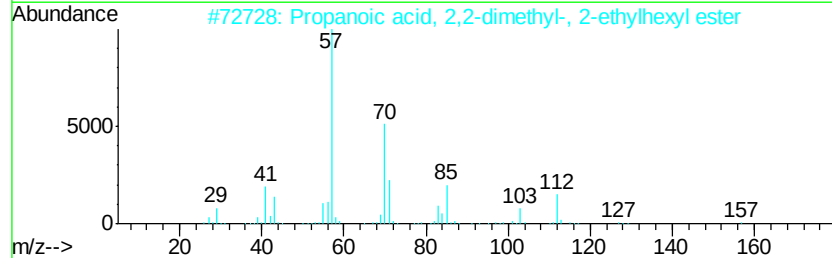
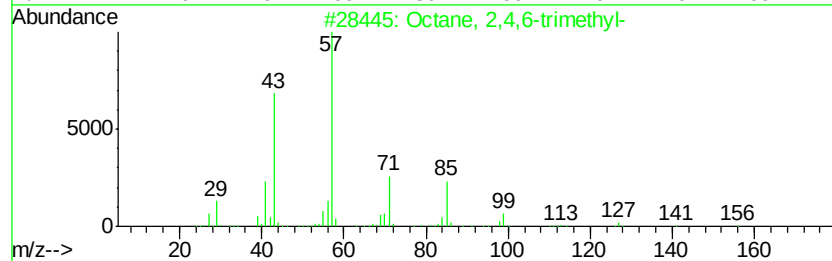
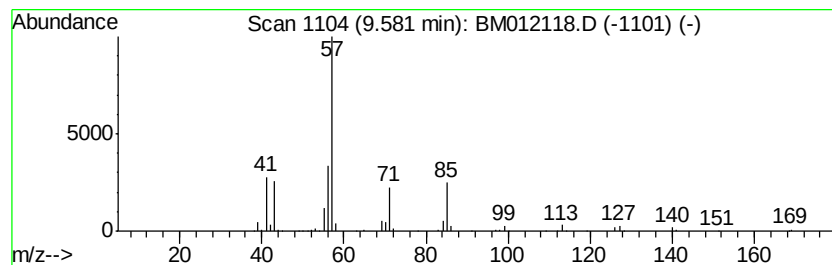
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.58 | 3.98 ng/ul | 685737 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Octane, 2,4,6-trimethyl-            | 156 | C11H24   | 062016-37-9 | 64   |
| 2    |    | Propanoic acid, 2,2-dimethyl-, 2... | 214 | C13H26O2 | 016387-18-1 | 64   |
| 3    |    | Pentane, 2,2,3,4-tetramethyl-       | 128 | C9H20    | 001186-53-4 | 59   |
| 4    |    | Undecane, 5,6-dimethyl-             | 184 | C13H28   | 017615-91-7 | 56   |
| 5    |    | Heptane, 2,2,4-trimethyl-           | 142 | C10H22   | 014720-74-2 | 50   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

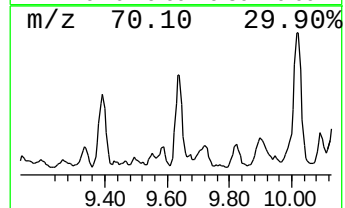
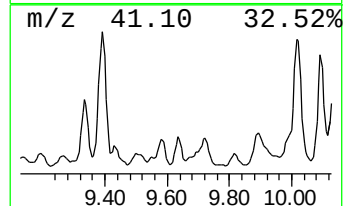
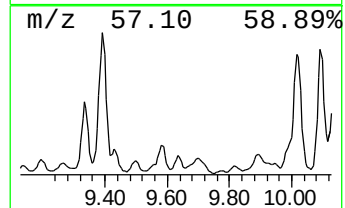
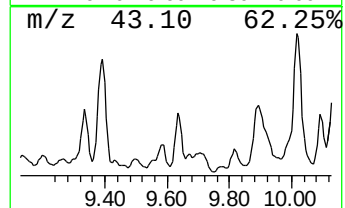
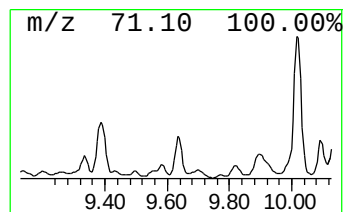
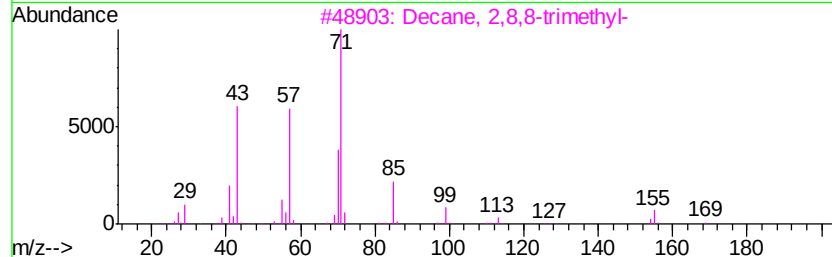
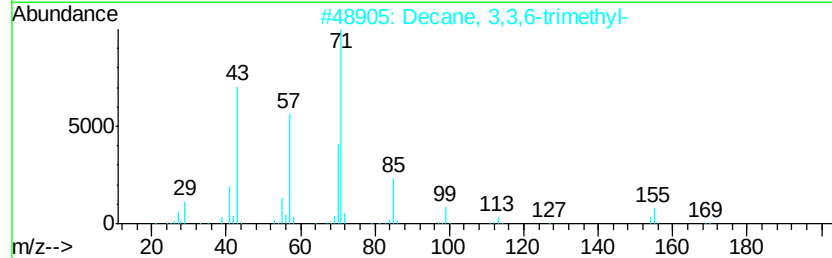
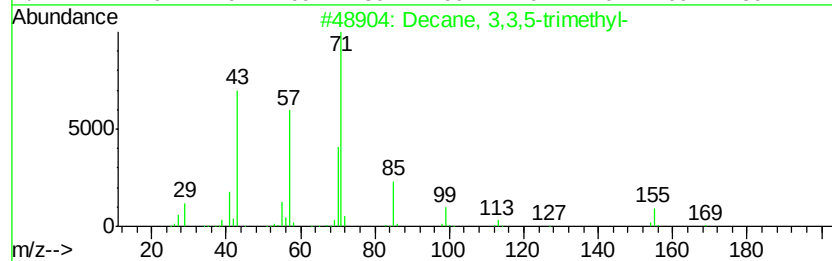
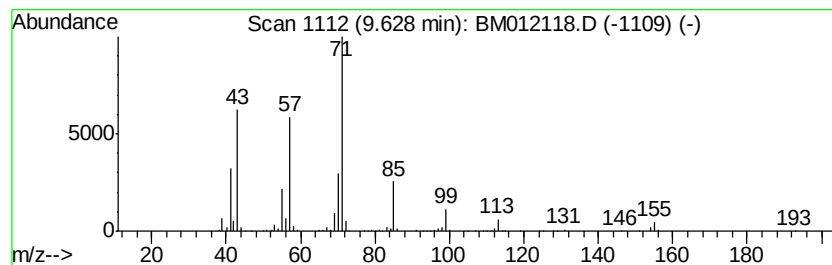
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.63 | 5.42 ng/ul | 932566 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------------|-----|---------|--------------|------|
| 1    | 5  | Decane, 3,3,5-trimethyl- | 184 | C13H28  | 062338-13-0  | 90   |
| 2    |    | Decane, 3,3,6-trimethyl- | 184 | C13H28  | 062338-14-1  | 83   |
| 3    |    | Decane, 2,8,8-trimethyl- | 184 | C13H28  | 1000060-81-2 | 83   |
| 4    |    | Decane, 3,3,8-trimethyl- | 184 | C13H28  | 062338-16-3  | 78   |
| 5    |    | Undecane, 4-methyl-      | 170 | C12H26  | 002980-69-0  | 72   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

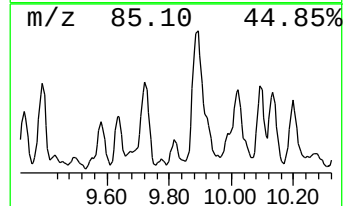
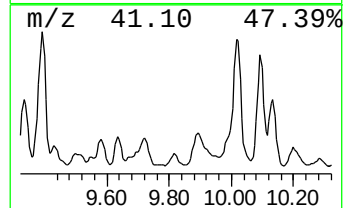
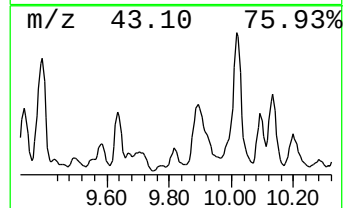
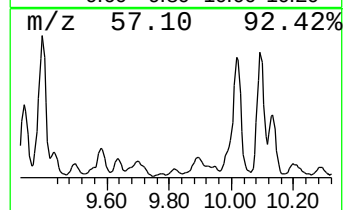
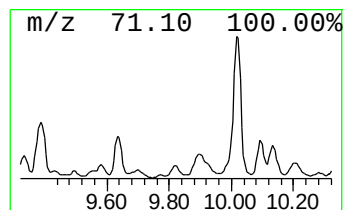
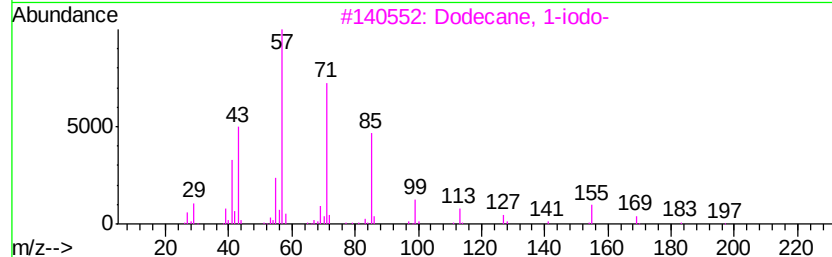
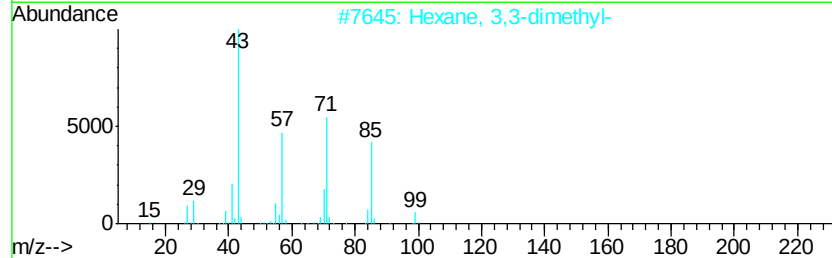
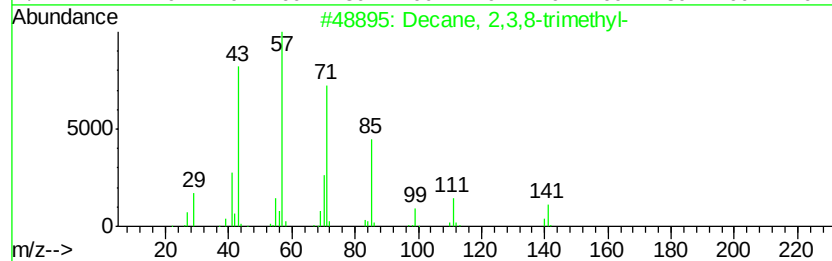
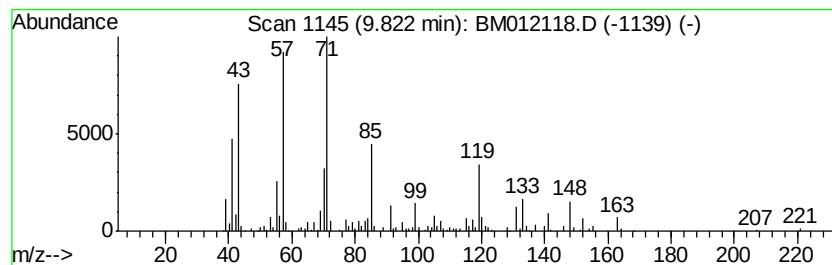
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.82 | 3.56 ng/ul | 613473 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 2,3,8-trimethyl- | 184 | C13H28  | 062238-14-6 | 64   |
| 2    |    |   | Hexane, 3,3-dimethyl-    | 114 | C8H18   | 000563-16-6 | 58   |
| 3    |    |   | Dodecane, 1-iodo-        | 296 | C12H25I | 004292-19-7 | 50   |
| 4    |    |   | Octane, 3-ethyl-         | 142 | C10H22  | 005881-17-4 | 43   |
| 5    |    |   | Heptane, 3,3-dimethyl-   | 128 | C9H20   | 004032-86-4 | 38   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

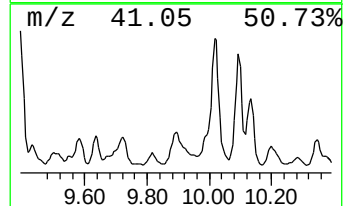
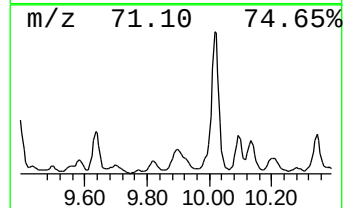
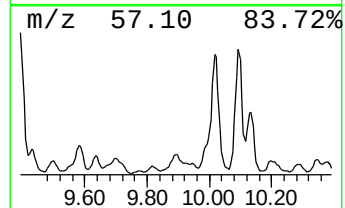
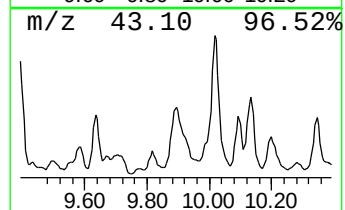
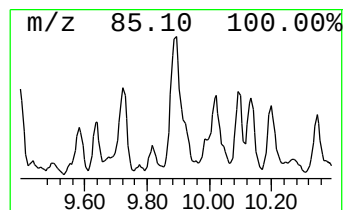
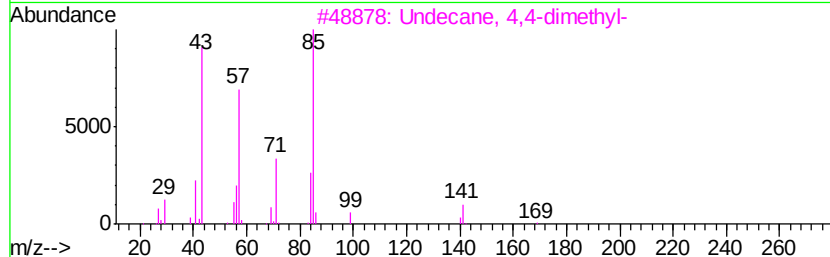
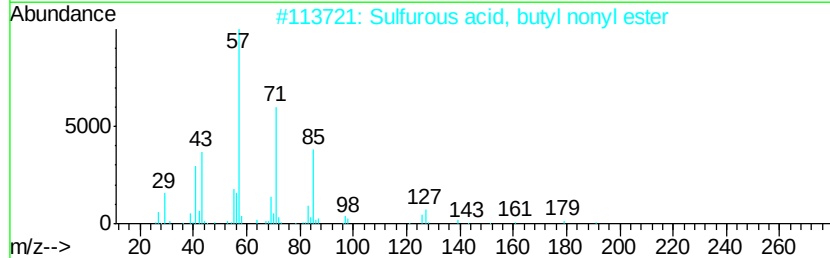
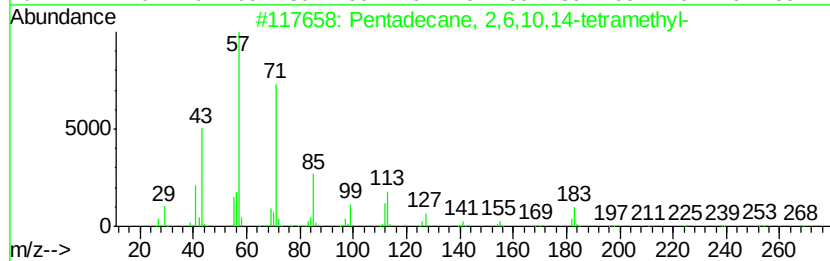
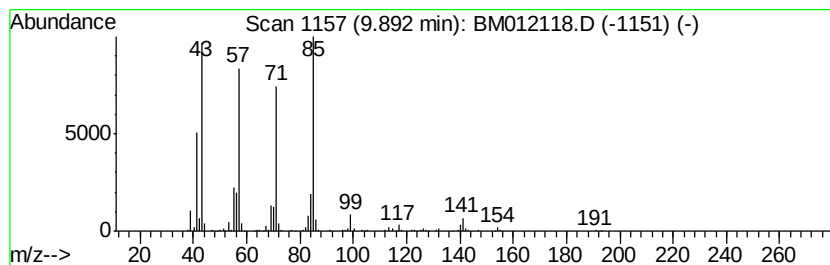
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T. | EstConc    | Area    | Relative to ISTD | R.T.  |
|------|------------|---------|------------------|-------|
| 9.89 | 9.94 ng/ul | 1711510 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40    | 001921-70-6  | 50   |
| 2    |    |   | Sulfurous acid, butyl nonyl ester   | 264 | C13H28O3S | 1000309-17-6 | 50   |
| 3    |    |   | Undecane, 4,4-dimethyl-             | 184 | C13H28    | 017312-68-4  | 47   |
| 4    |    |   | Oxalic acid, octyl propyl ester     | 244 | C13H24O4  | 1000309-26-1 | 47   |
| 5    |    |   | Tetrahydropyran 12-tetradecyn-1-... | 294 | C19H34O2  | 096249-40-0  | 43   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

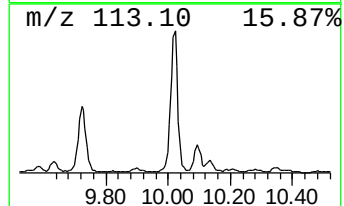
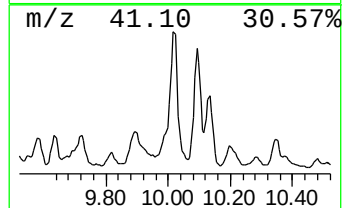
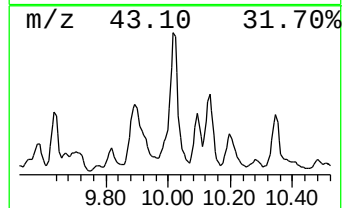
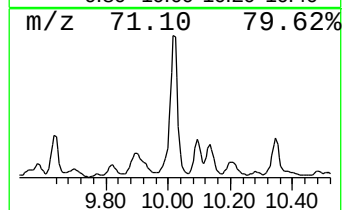
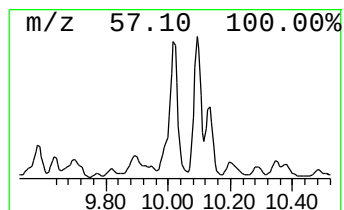
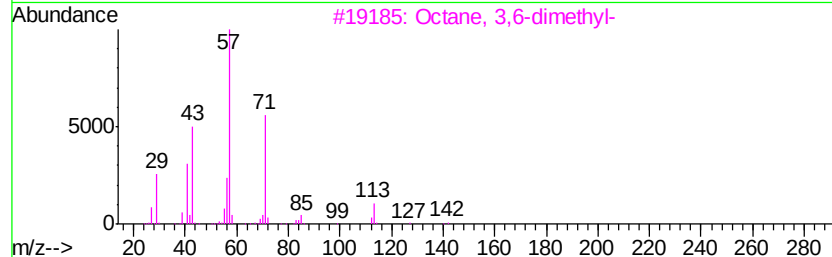
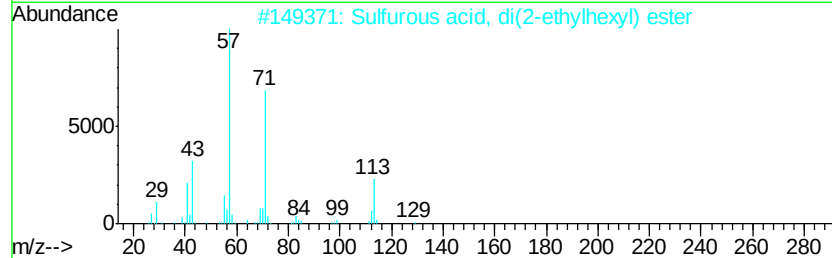
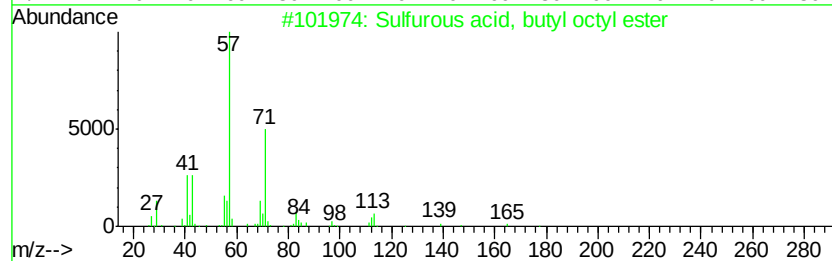
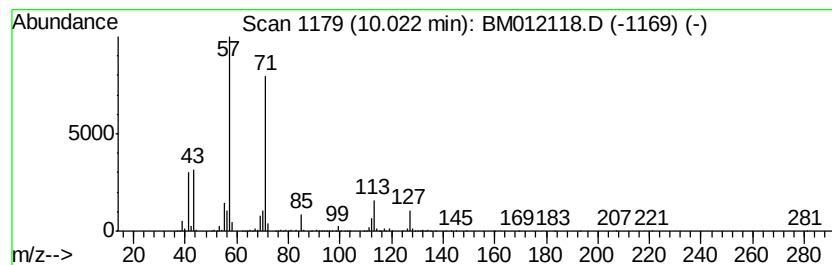
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 26 Sulfurous acid, butyl octyl... Concentration Rank 6

| R.T.    | EstConc     | Area                                | Relative to ISTD | R.T.            |
|---------|-------------|-------------------------------------|------------------|-----------------|
| 10.02   | 30.49 ng/ul | 5250810                             | Napthalene-d8    | 10.64           |
| Hit# of | 5           | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |             | Sulfurous acid, butyl octyl ester   | 250 C12H26O3S    | 1000309-17-5 78 |
| 2       |             | Sulfurous acid, di(2-ethylhexyl)... | 306 C16H34O3S    | 1000309-19-1 78 |
| 3       |             | Octane, 3,6-dimethyl-               | 142 C10H22       | 015869-94-0 72  |
| 4       |             | Tridecane, 7-methyl-                | 198 C14H30       | 026730-14-3 72  |
| 5       |             | Nonane, 3-methyl-                   | 142 C10H22       | 005911-04-6 72  |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

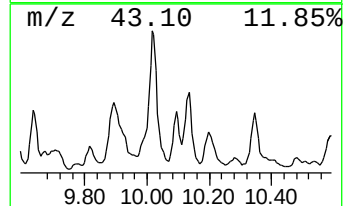
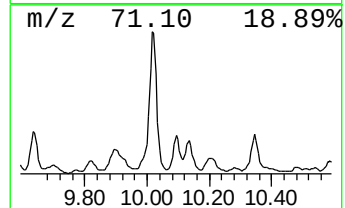
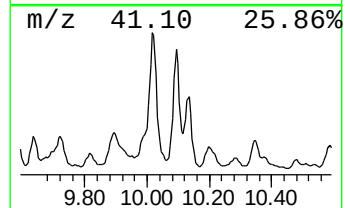
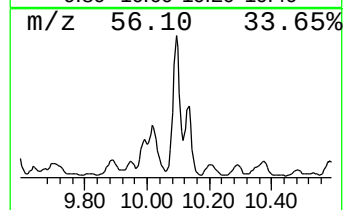
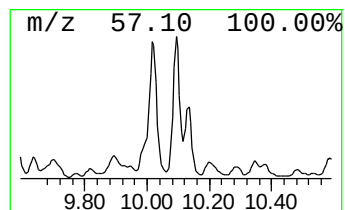
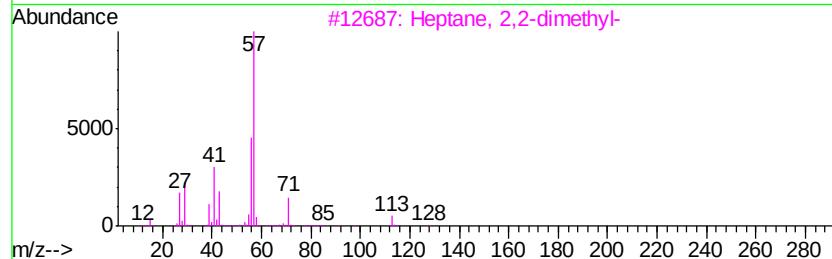
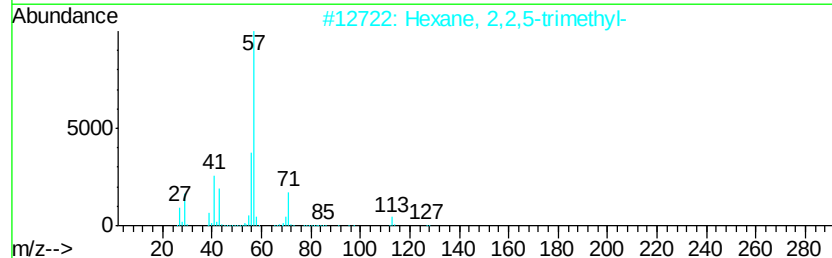
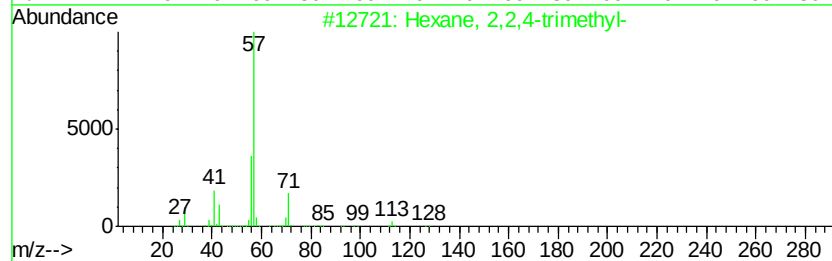
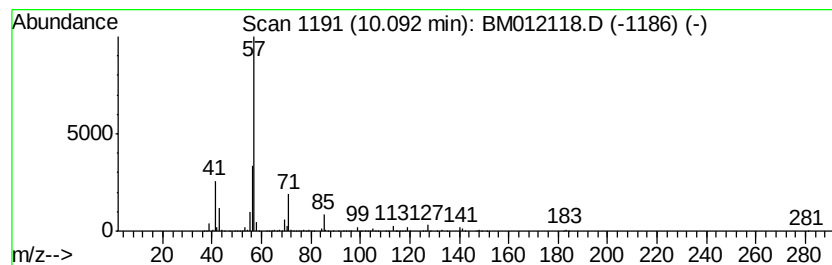
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.09 | 17.59 ng/ul | 3029550 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexane, 2,2,4-trimethyl- | 128 | C9H20   | 016747-26-5 | 59   |
| 2    |    | Hexane, 2,2,5-trimethyl- | 128 | C9H20   | 003522-94-9 | 59   |
| 3    |    | Heptane, 2,2-dimethyl-   | 128 | C9H20   | 001071-26-7 | 59   |
| 4    |    | Hexane, 2,2,4-trimethyl- | 128 | C9H20   | 016747-26-5 | 59   |
| 5    |    | Decane, 2,2,5-trimethyl- | 184 | C13H28  | 062237-96-1 | 56   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

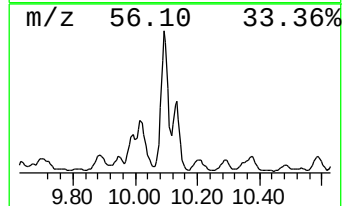
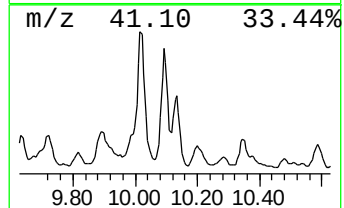
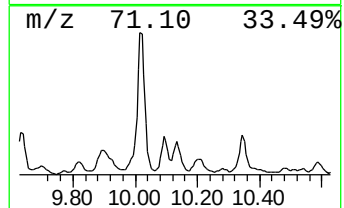
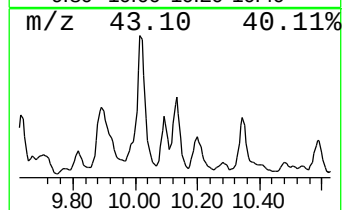
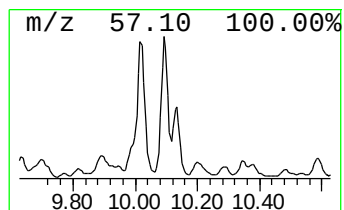
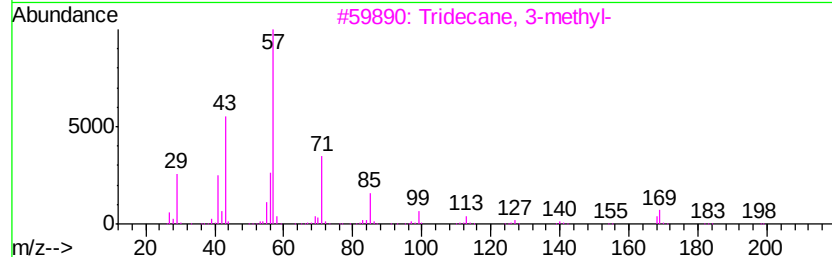
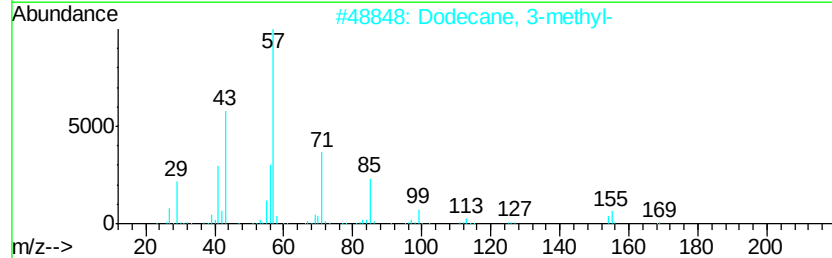
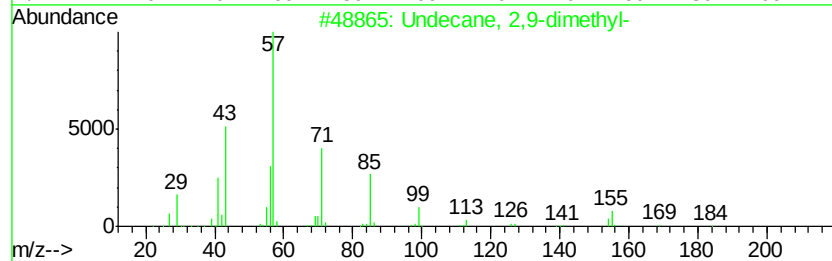
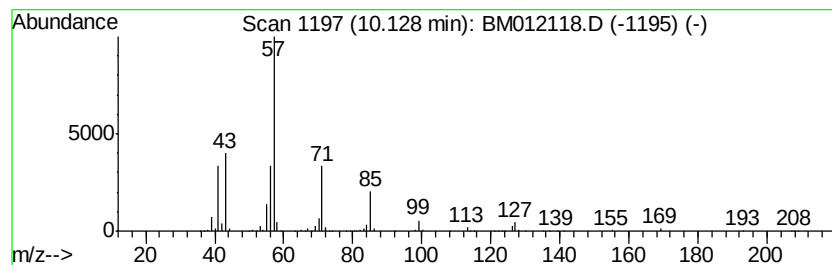
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 10.13 | 9.59 ng/ul | 1651440 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 2,9-dimethyl- | 184 | C13H28  | 017301-26-7 | 78   |
| 2    |    |   | Dodecane, 3-methyl-     | 184 | C13H28  | 017312-57-1 | 78   |
| 3    |    |   | Tridecane, 3-methyl-    | 198 | C14H30  | 006418-41-3 | 78   |
| 4    |    |   | Dodecane, 3-methyl-     | 184 | C13H28  | 017312-57-1 | 78   |
| 5    |    |   | Undecane, 3-methyl-     | 170 | C12H26  | 001002-43-3 | 78   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

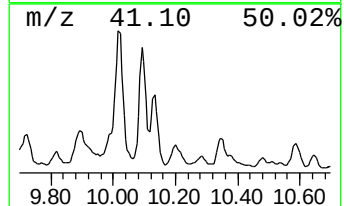
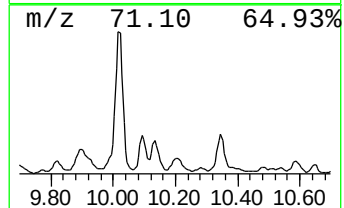
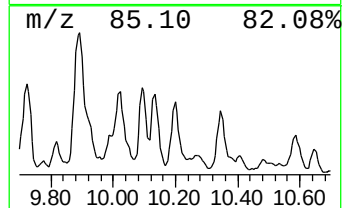
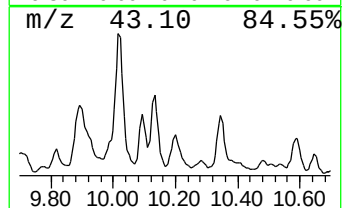
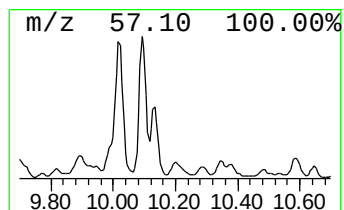
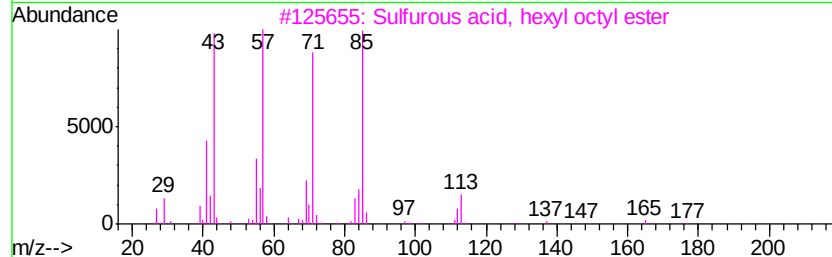
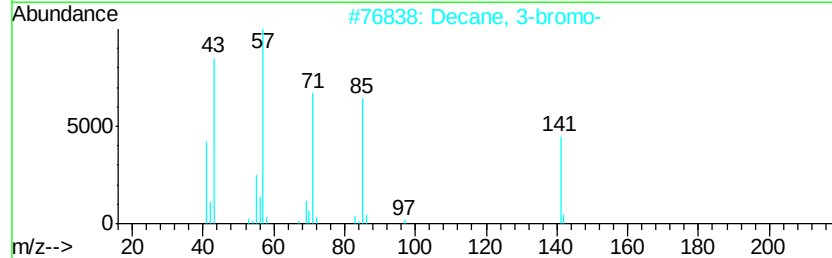
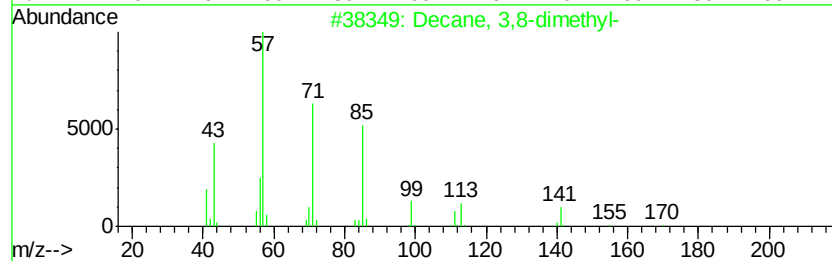
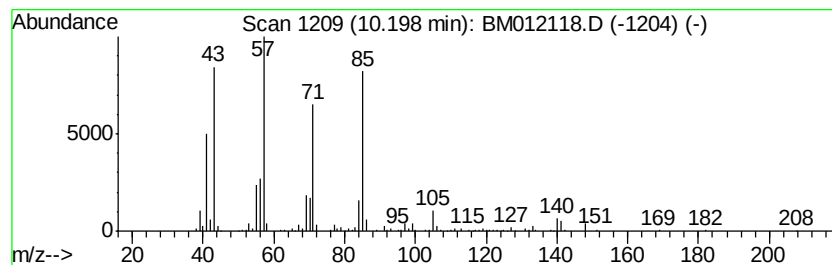
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.20 | 4.84 ng/ul | 833522 | Napthalene-d8    | 10.64 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 3,8-dimethyl-               | 170 | C12H26    | 017312-55-9  | 64   |
| 2    |    |   | Decane, 3-bromo-                    | 220 | C10H21Br  | 030571-71-2  | 59   |
| 3    |    |   | Sulfurous acid, hexyl octyl ester   | 278 | C14H30O3S | 1000309-13-0 | 59   |
| 4    |    |   | Hexane, 2,4-dimethyl-               | 114 | C8H18     | 000589-43-5  | 53   |
| 5    |    |   | 3,3-Dimethylbutan-2-yl isobutyl ... | 202 | C11H22O3  | 1000372-76-5 | 47   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

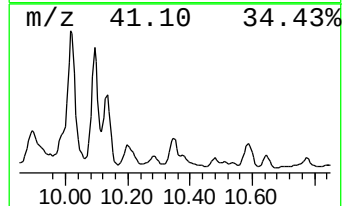
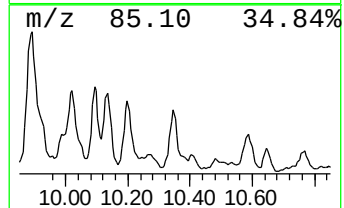
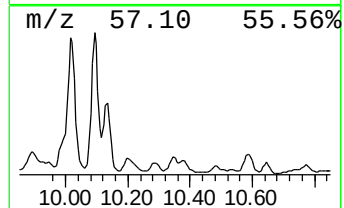
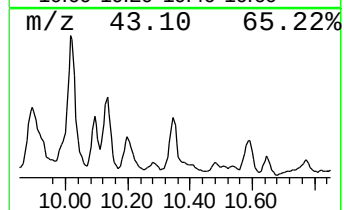
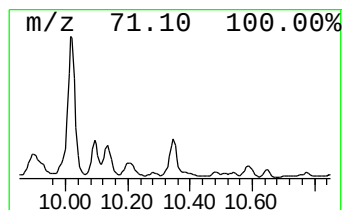
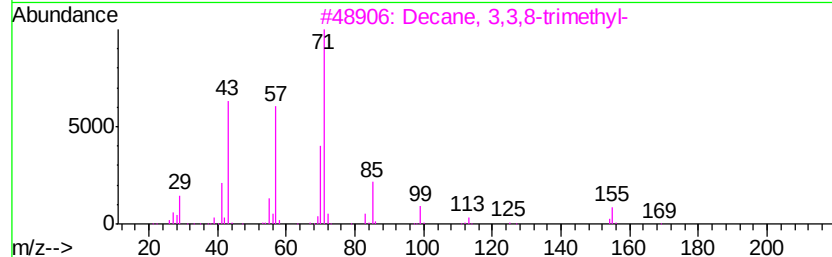
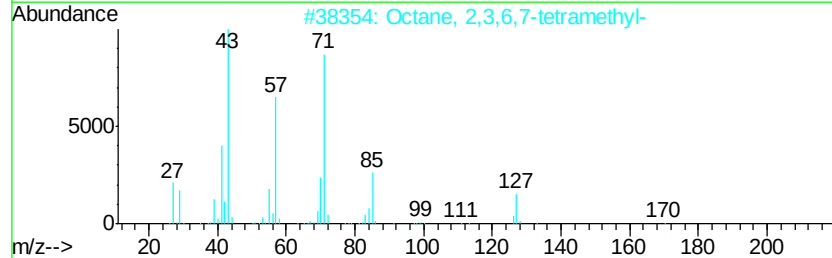
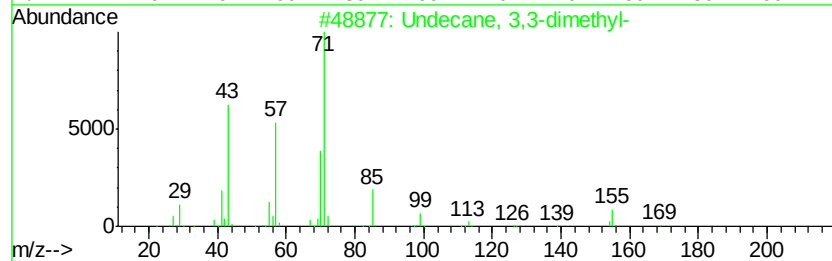
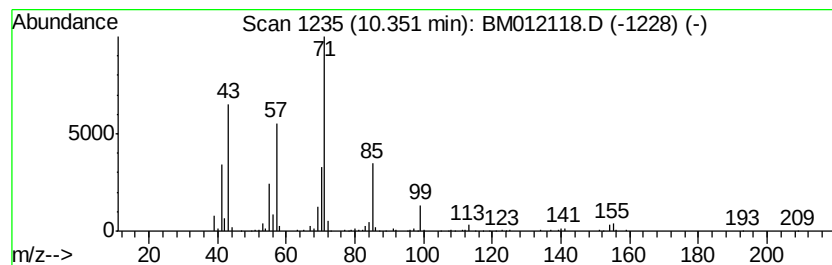
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 10.35 | 6.54 ng/ul | 1126240 | Naphthalene-d8   | 10.64 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 3,3-dimethyl-      | 184 | C13H28  | 017312-65-1 | 64   |
| 2    |    |   | Octane, 2,3,6,7-tetramethyl- | 170 | C12H26  | 052670-34-5 | 64   |
| 3    |    |   | Decane, 3,3,8-trimethyl-     | 184 | C13H28  | 062338-16-3 | 64   |
| 4    |    |   | Heptane, 3,3,5-trimethyl-    | 142 | C10H22  | 007154-80-5 | 59   |
| 5    |    |   | Octane, 2,6,6-trimethyl-     | 156 | C11H24  | 054166-32-4 | 59   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

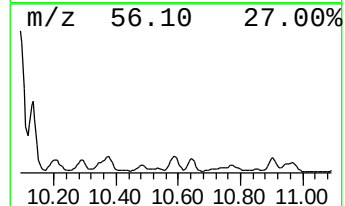
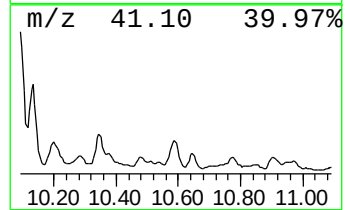
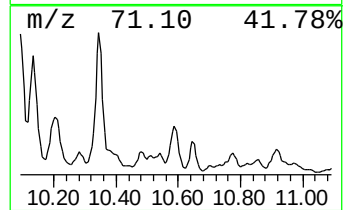
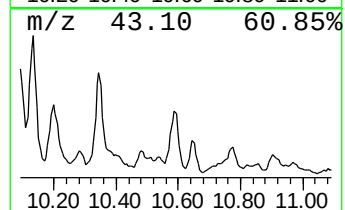
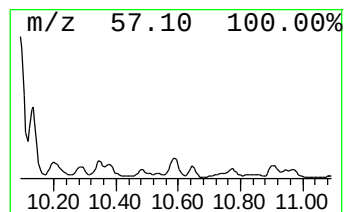
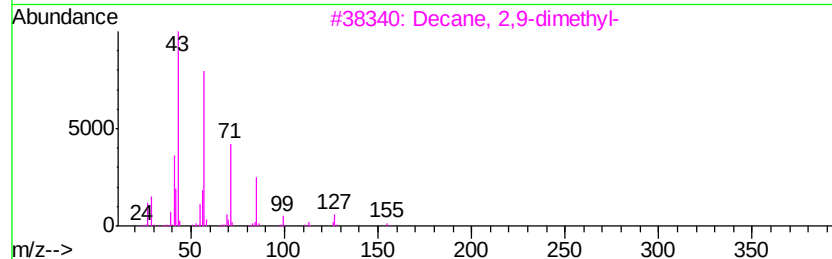
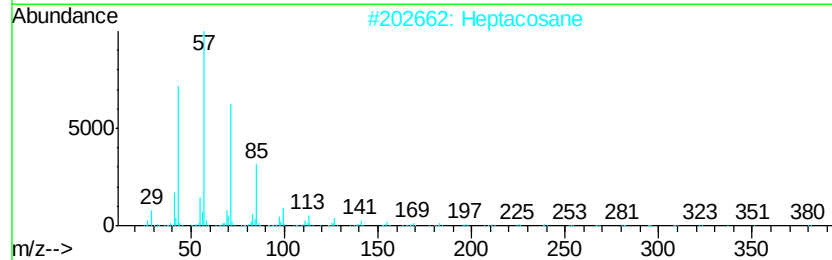
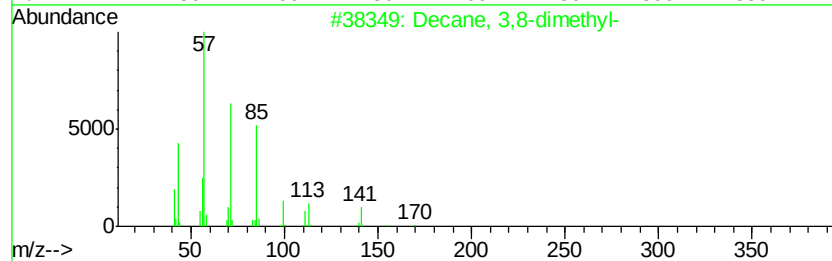
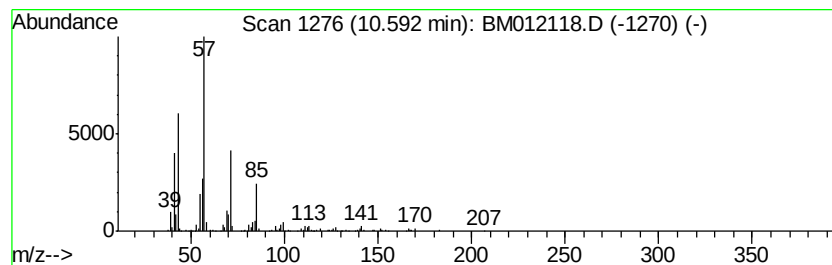
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.59 | 4.00 ng/ul | 689555 | Napthalene-d8    | 10.64 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 3,8-dimethyl-    | 170 | C12H26  | 017312-55-9 | 80   |
| 2    |    | Heptacosane              | 380 | C27H56  | 000593-49-7 | 59   |
| 3    |    | Decane, 2,9-dimethyl-    | 170 | C12H26  | 001002-17-1 | 59   |
| 4    |    | Octane, 2,3,3-trimethyl- | 156 | C11H24  | 062016-30-2 | 59   |
| 5    |    | Hexadecane               | 226 | C16H34  | 000544-76-3 | 58   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

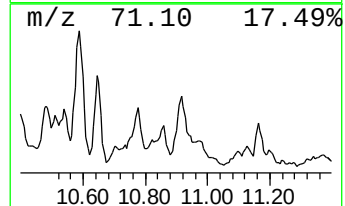
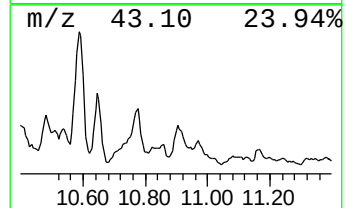
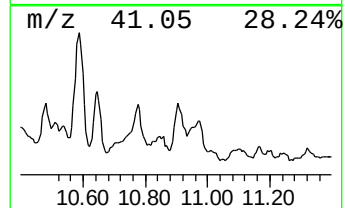
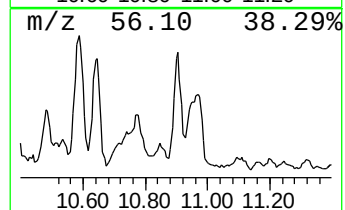
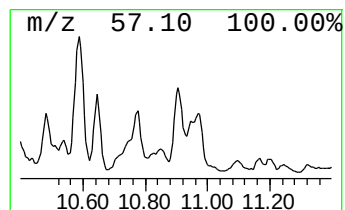
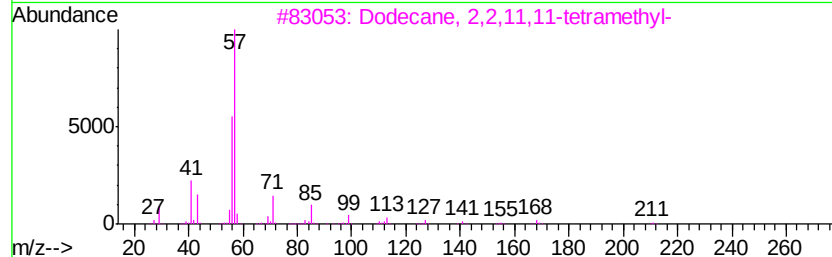
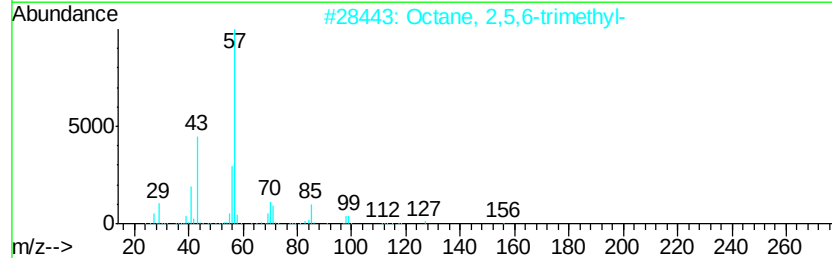
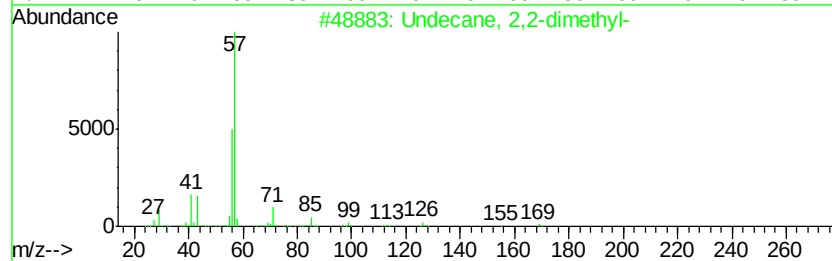
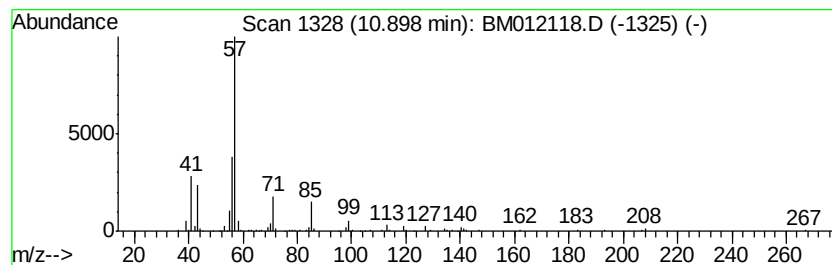
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 41

| R.T.    | EstConc    | Area                             | Relative to ISTD | R.T.           |
|---------|------------|----------------------------------|------------------|----------------|
| 10.90   | 2.03 ng/ul | 349349                           | Napthalene-d8    | 10.64          |
| Hit# of | 5          | Tentative ID                     | MW MolForm       | CAS# Qual      |
| 1       |            | Undecane, 2,2-dimethyl-          | 184 C13H28       | 017312-64-0 59 |
| 2       |            | Octane, 2,5,6-trimethyl-         | 156 C11H24       | 062016-14-2 59 |
| 3       |            | Dodecane, 2,2,11,11-tetramethyl- | 226 C16H34       | 127204-12-0 53 |
| 4       |            | Dodecane, 3-methyl-              | 184 C13H28       | 017312-57-1 47 |
| 5       |            | Heptane, 2,2,4,6,6-pentamethyl-  | 170 C12H26       | 013475-82-6 47 |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

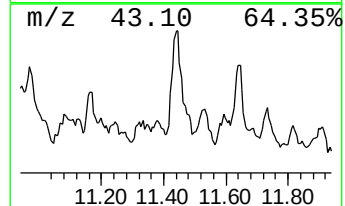
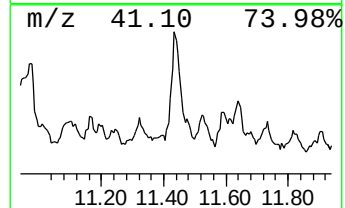
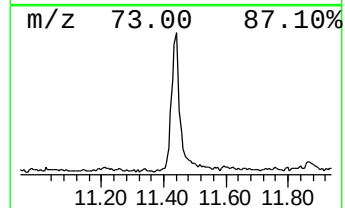
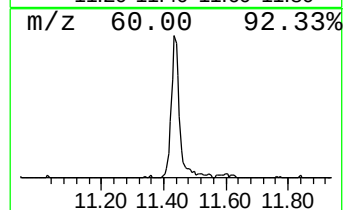
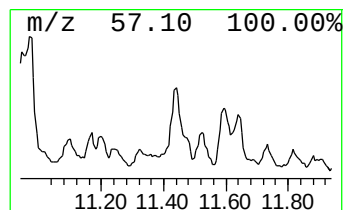
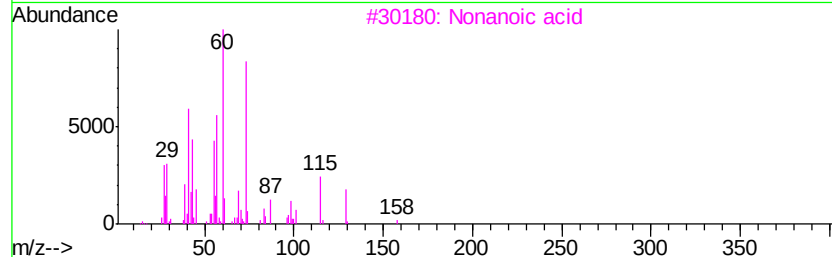
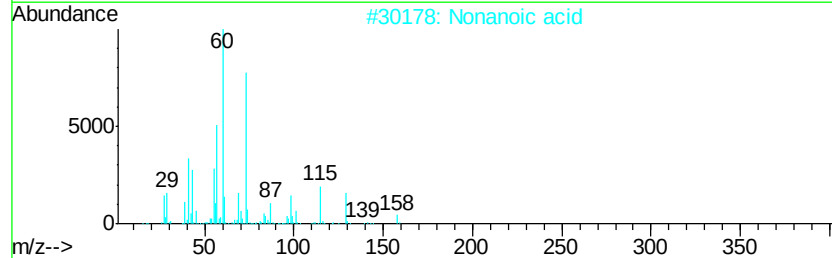
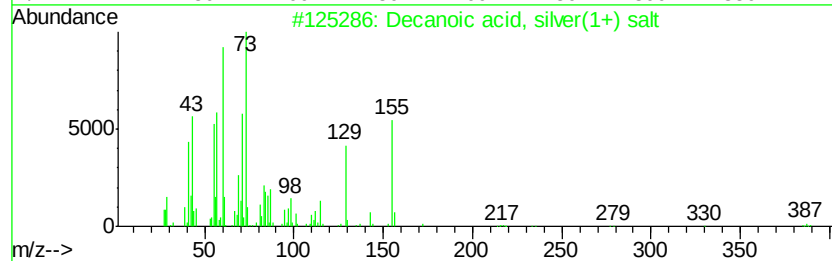
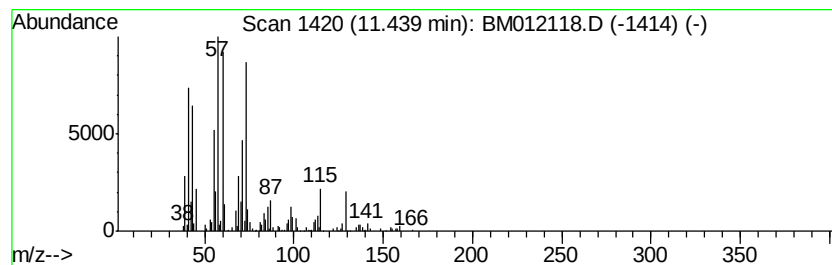
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 33 Decanoic acid, silver(1+) salt Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.44 | 2.64 ng/ul | 454811 | Naphthalene-d8   | 10.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Decanoic acid, silver(1+) salt      | 278 | C10H19AgO2 | 013126-67-5 | 64   |
| 2    |    | Nonanoic acid                       | 158 | C9H18O2    | 000112-05-0 | 59   |
| 3    |    | Nonanoic acid                       | 158 | C9H18O2    | 000112-05-0 | 52   |
| 4    |    | Isoamyl nitrite                     | 117 | C5H11NO2   | 000110-46-3 | 47   |
| 5    |    | .beta.-D-Glucopyranose, 1,6-anhy... | 162 | C6H10O5    | 000498-07-7 | 38   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

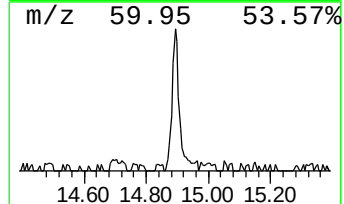
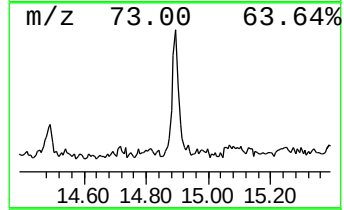
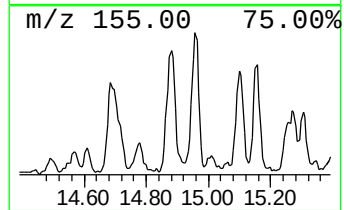
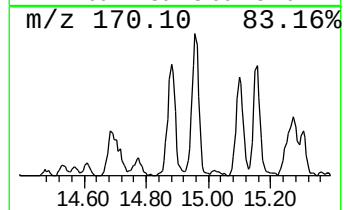
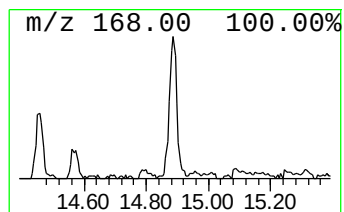
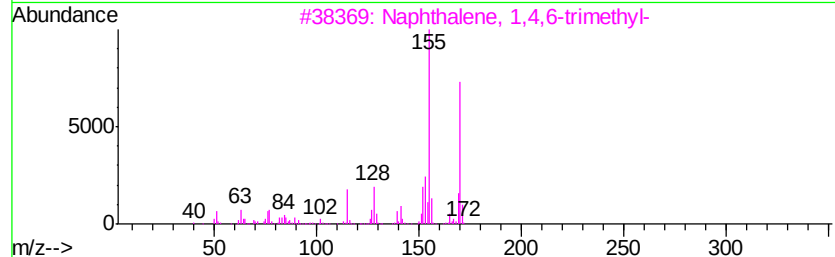
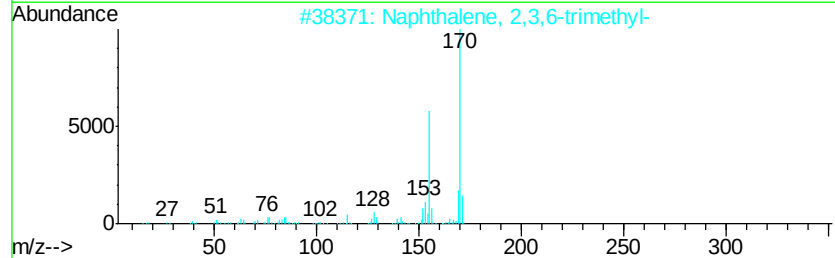
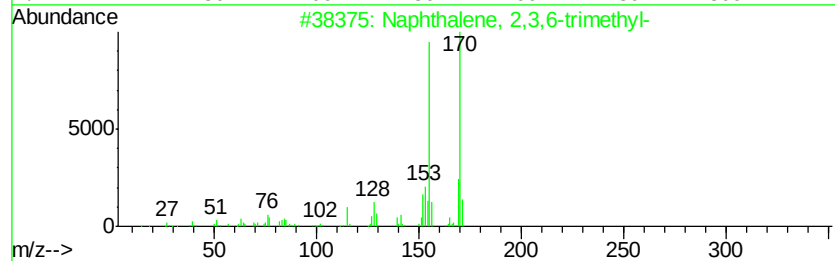
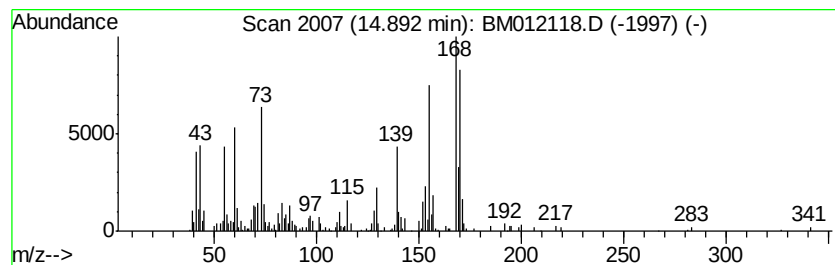
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 34 Naphthalene, 2,3,6-trimethyl- Concentration Rank 40

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.89 | 2.21 ng/ul | 419194 | Acenaphthene-d10 | 14.49 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |
| 2    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 89   |
| 3    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 89   |
| 4    |    | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 87   |
| 5    |    | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 80   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

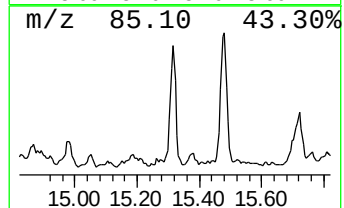
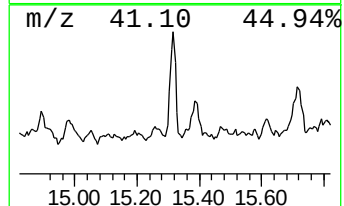
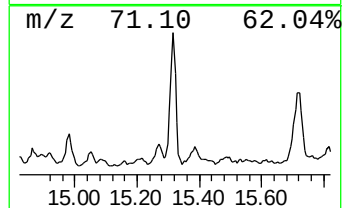
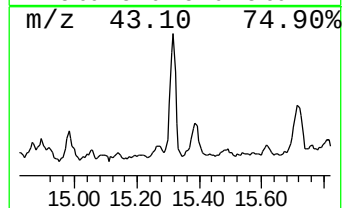
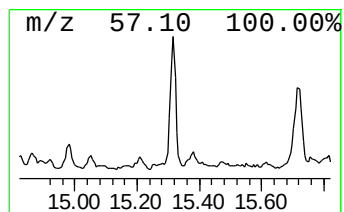
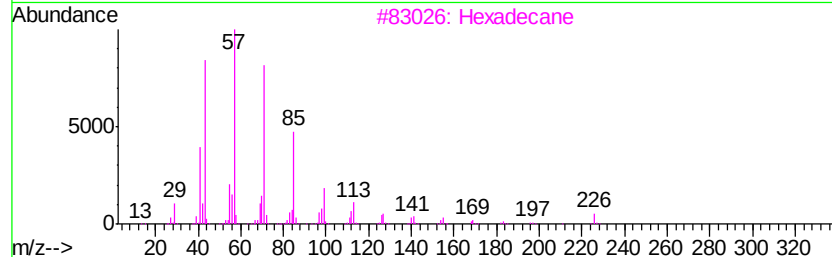
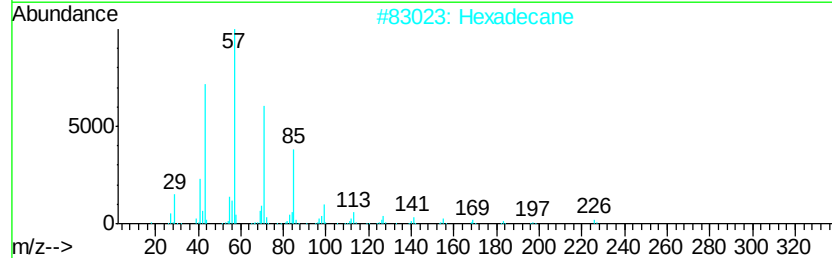
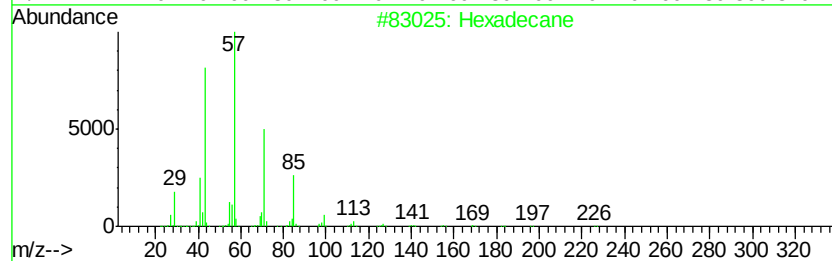
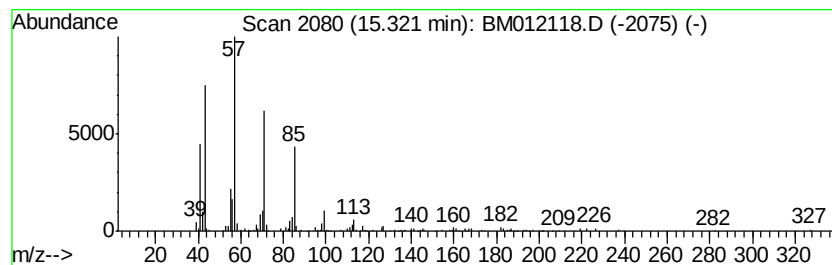
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.32 | 2.01 ng/ul | 382108 | Acenaphthene-d10 | 14.49 |

| Hit# | of          | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|-------------|---|--------------|-----|---------|-------------|------|
| 1    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 96   |
| 2    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 94   |
| 3    | Hexadecane  |   |              | 226 | C16H34  | 000544-76-3 | 93   |
| 4    | Nonadecane  |   |              | 268 | C19H40  | 000629-92-5 | 90   |
| 5    | Tetradecane |   |              | 198 | C14H30  | 000629-59-4 | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

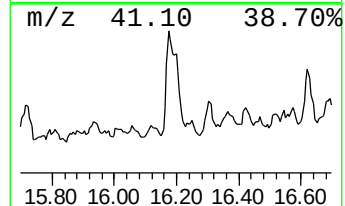
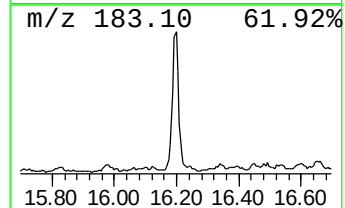
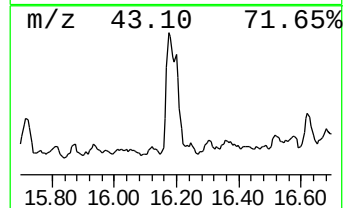
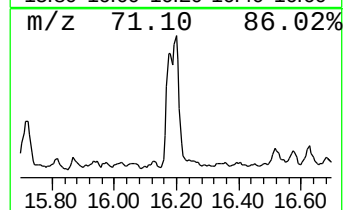
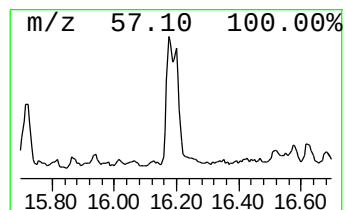
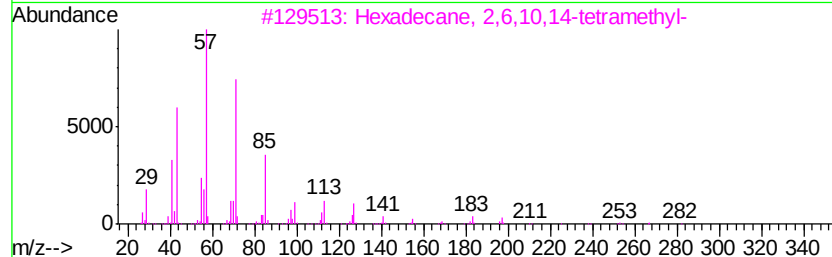
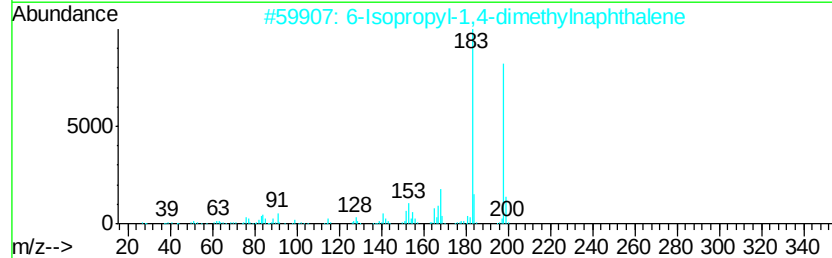
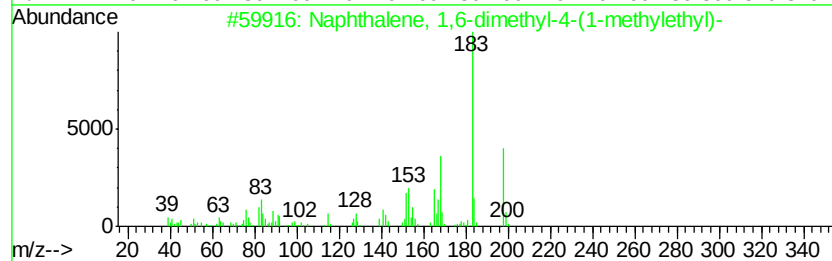
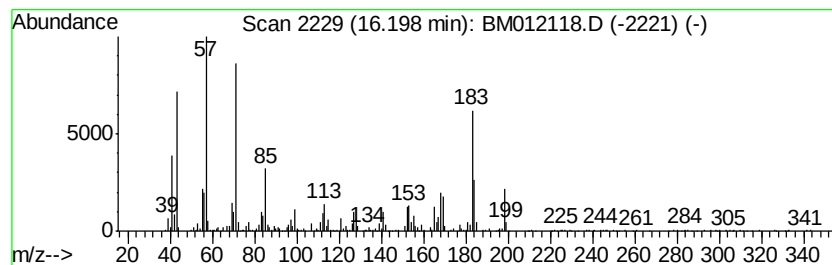
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 Naphthalene, 1,6-dimethyl-4... Concentration Rank 20

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.20 | 7.78 ng/ul | 1450610 | Phenanthrene-d10 | 17.23 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18  | 000483-78-3 | 96   |
| 2    |    |   | 6-Isopropyl-1,4-dimethylnaphthalene | 198 | C15H18  | 000489-77-0 | 94   |
| 3    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 94   |
| 4    |    |   | Azulene, 1,4-dimethyl-7-(1-methy... | 198 | C15H18  | 000489-84-9 | 83   |
| 5    |    |   | Naphthalene, 1,6-dimethyl-4-(1-m... | 198 | C15H18  | 000483-78-3 | 64   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

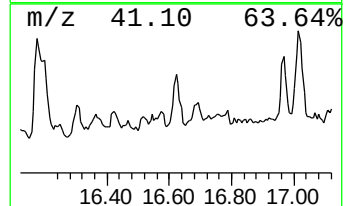
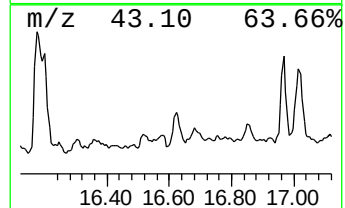
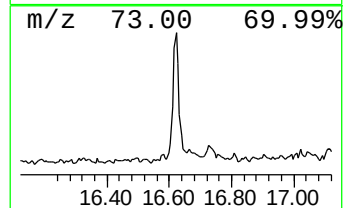
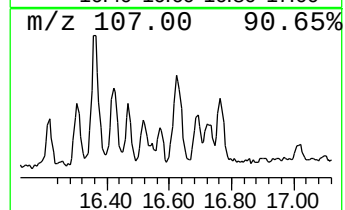
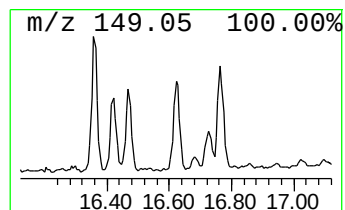
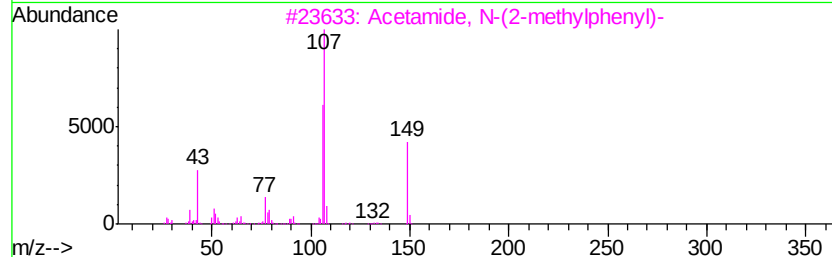
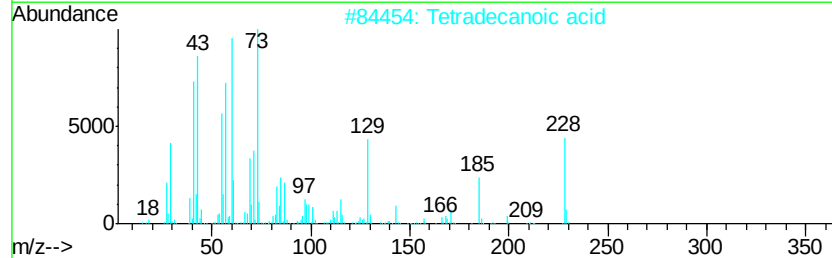
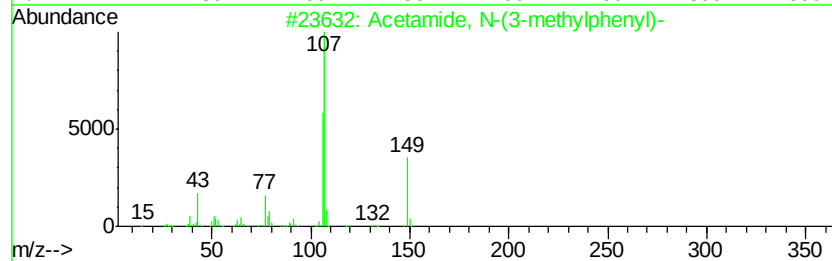
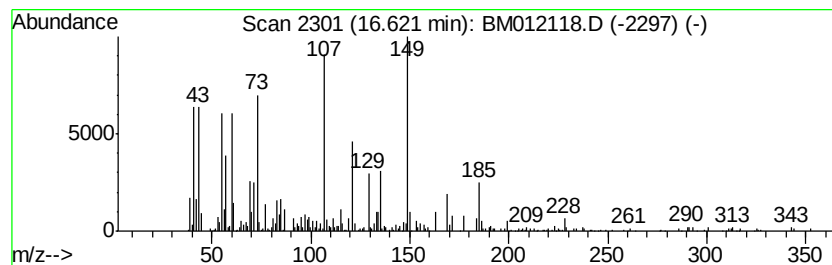
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 37 unknown-02 Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD               |     | R.T.       |             |      |
|-------|------------|--------|--------------------------------|-----|------------|-------------|------|
| 16.62 | 2.59 ng/ul | 482570 | Phenanthrene-d10               |     | 17.23      |             |      |
| Hit#  | of         | 5      | Tentative ID                   | MW  | MolForm    | CAS#        | Qual |
| 1     |            |        | Acetamide, N-(3-methylphenyl)- | 149 | C9H11NO    | 000537-92-8 | 46   |
| 2     |            |        | Tetradecanoic acid             | 228 | C14H28O2   | 000544-63-8 | 41   |
| 3     |            |        | Acetamide, N-(2-methylphenyl)- | 149 | C9H11NO    | 000120-66-1 | 27   |
| 4     |            |        | Undecanoic acid                | 186 | C11H22O2   | 000112-37-8 | 25   |
| 5     |            |        | 11-Bromoundecanoic acid        | 264 | C11H21BrO2 | 002834-05-1 | 15   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

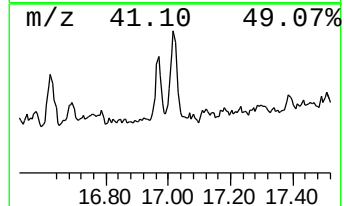
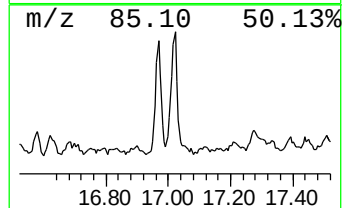
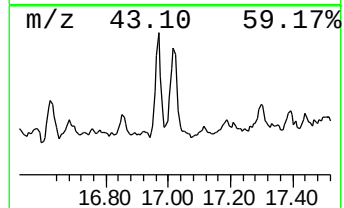
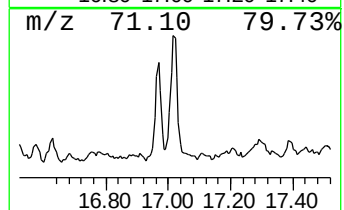
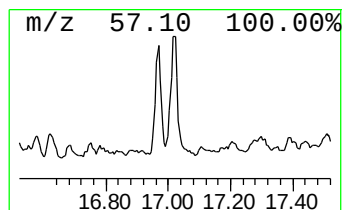
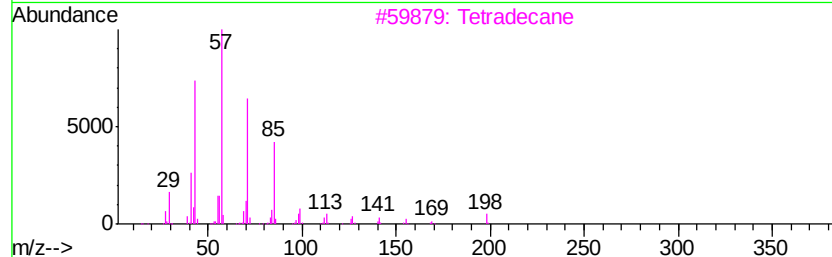
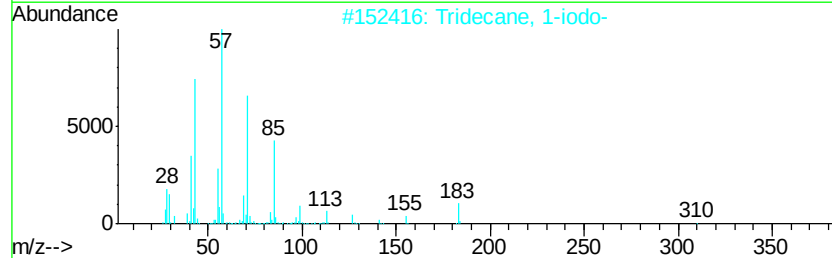
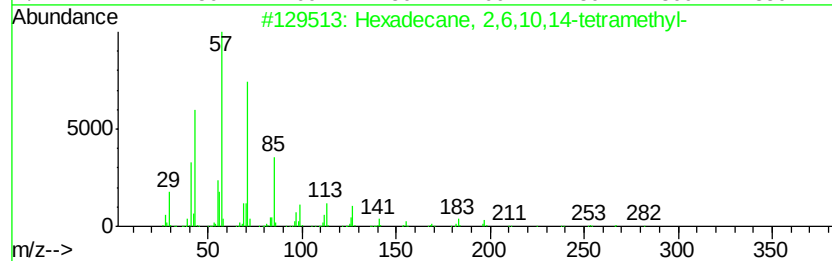
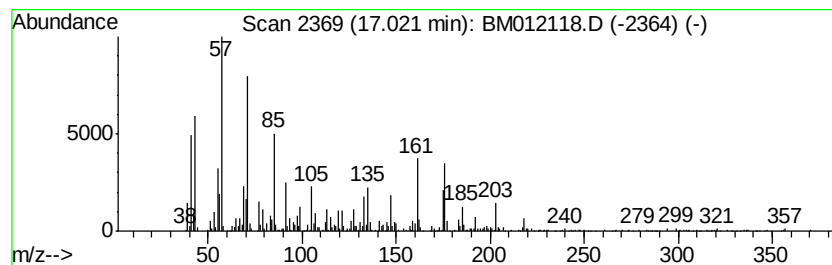
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.02 | 4.16 ng/ul | 775569 | Phenanthrene-d10 | 17.23 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 46   |
| 2    |    |   | Tridecane, 1-iodo-                 | 310 | C13H27I | 035599-77-0 | 46   |
| 3    |    |   | Tetradecane                        | 198 | C14H30  | 000629-59-4 | 41   |
| 4    |    |   | Tridecane                          | 184 | C13H28  | 000629-50-5 | 41   |
| 5    |    |   | Hexadecane                         | 226 | C16H34  | 000544-76-3 | 38   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

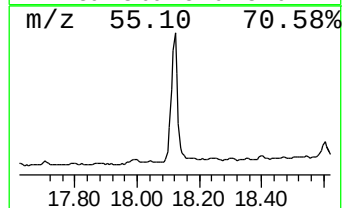
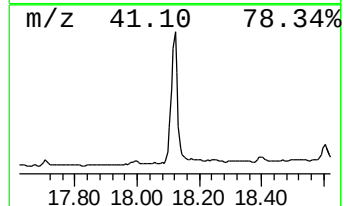
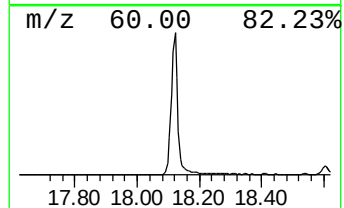
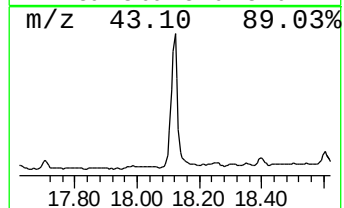
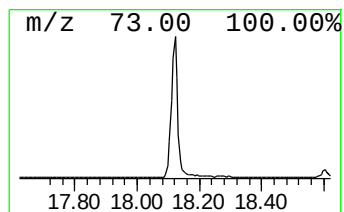
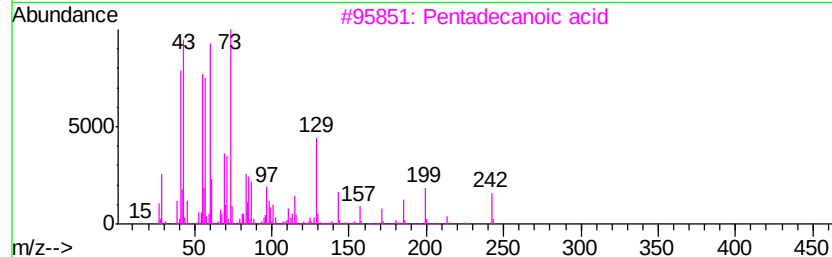
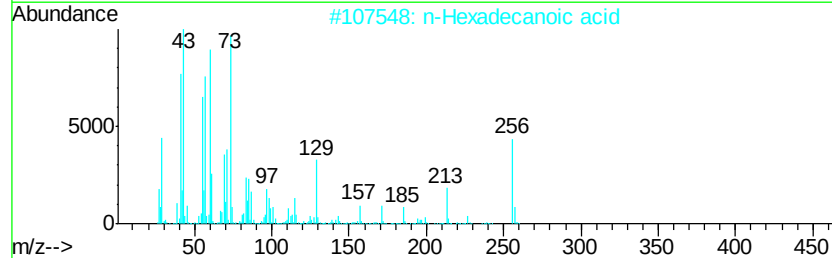
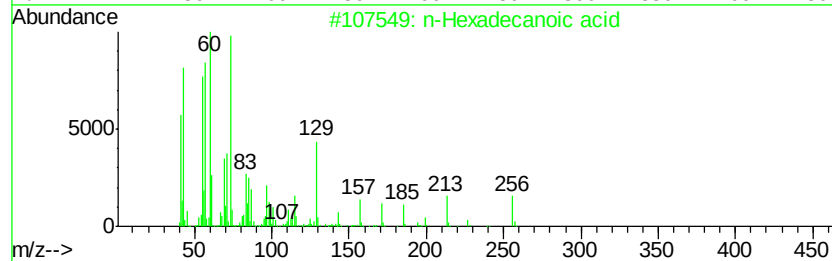
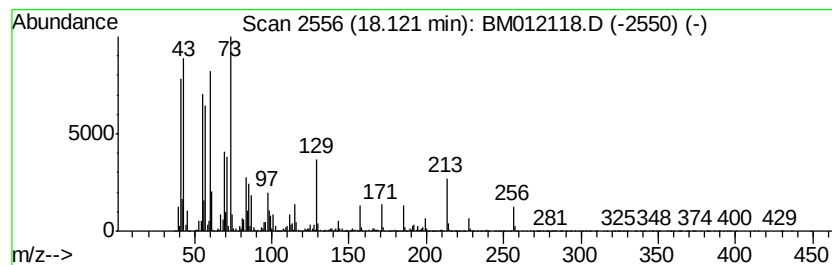
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 n-Hexadecanoic acid Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.12 | 46.25 ng/ul | 8628340 | Phenanthrene-d10 | 17.23 |

| 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    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 90   |
| 4    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 90   |
| 5    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 86   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

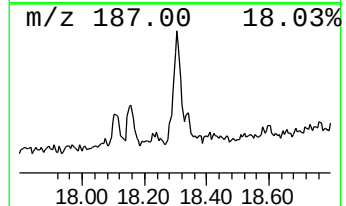
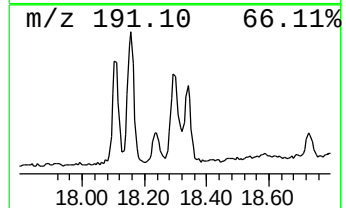
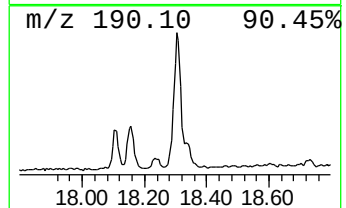
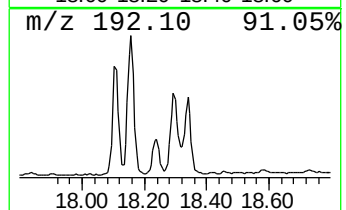
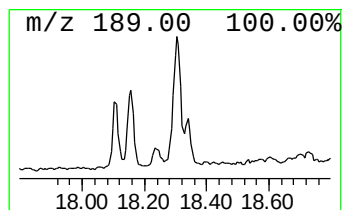
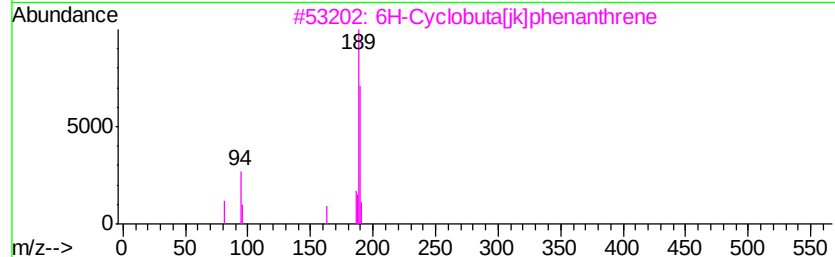
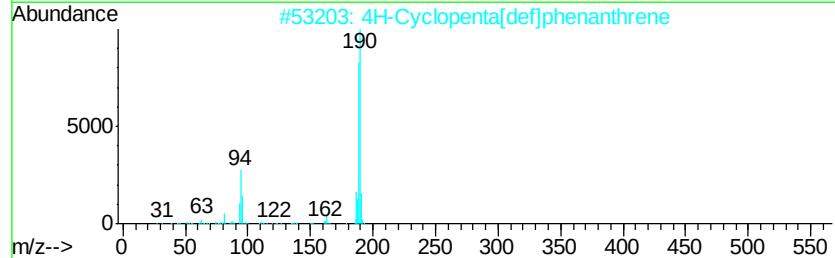
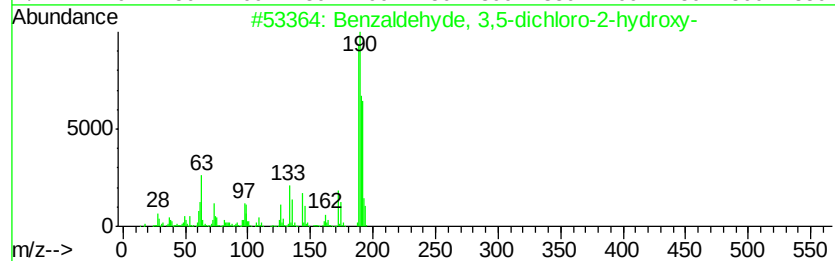
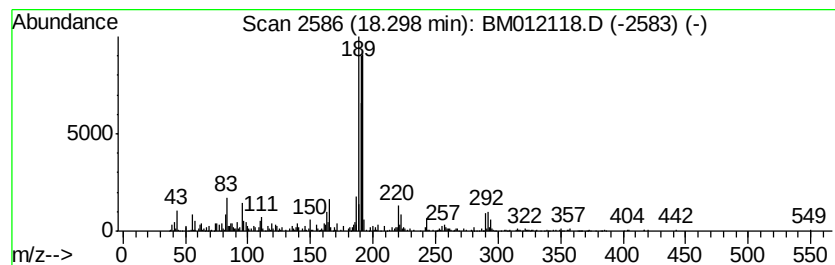
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 40 Benzaldehyde, 3,5-dichloro-... Concentration Rank 37

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.      |             |      |
|---------|------------|-------------------------------------|------------------|-----------|-------------|------|
| 18.30   | 2.76 ng/ul | 514865                              | Phenanthrene-d10 | 17.23     |             |      |
| Hit# of | 5          | Tentative ID                        | MW               | MolForm   | CAS#        | Qual |
| 1       |            | Benzaldehyde, 3,5-dichloro-2-hyd... | 190              | C7H4Cl2O2 | 000090-60-8 | 58   |
| 2       |            | 4H-Cyclopenta[def]phenanthrene      | 190              | C15H10    | 000203-64-5 | 46   |
| 3       |            | 6H-Cyclobuta[jk]phenanthrene        | 190              | C15H10    | 083469-43-6 | 46   |
| 4       |            | 4H-Cyclopenta[def]phenanthrene      | 190              | C15H10    | 000203-64-5 | 46   |
| 5       |            | 2,2'-Bis(4,5-dimethylimidazole)     | 190              | C10H14N4  | 069286-06-2 | 43   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

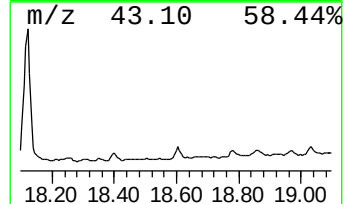
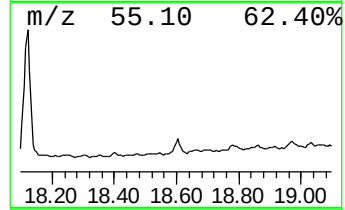
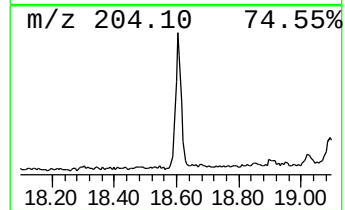
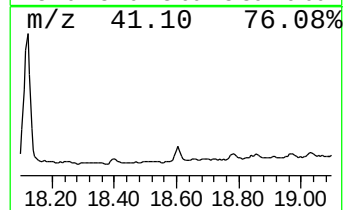
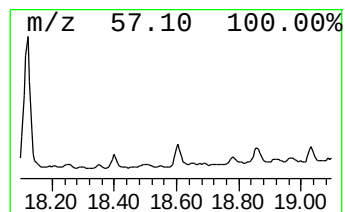
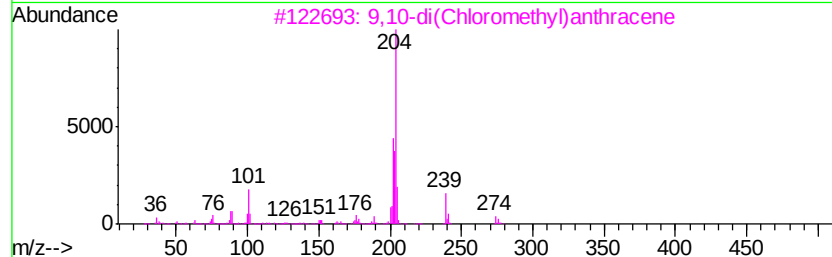
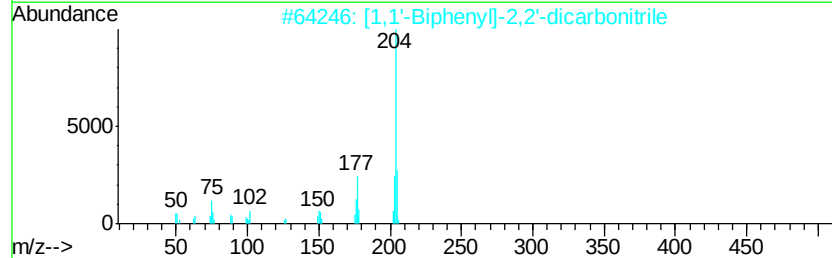
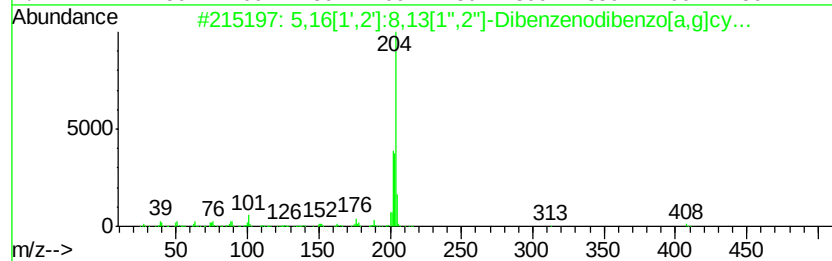
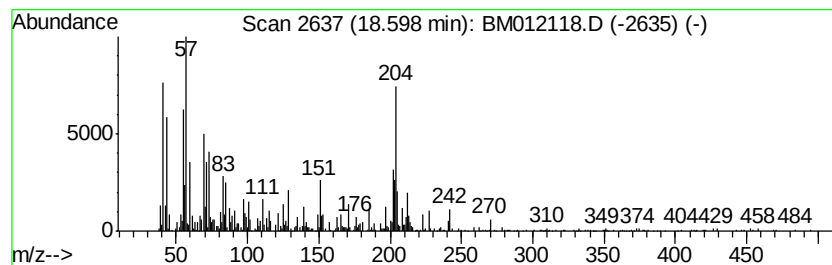
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 41 unknown-03 Concentration Rank 27

| R.T.    | EstConc                             | Area                    | Relative to ISTD | R.T.           |
|---------|-------------------------------------|-------------------------|------------------|----------------|
| 18.60   | 5.38 ng/ul                          | 1004150                 | Phenanthrene-d10 | 17.23          |
| Hit# of | 5                                   | Tentative ID            | MW MolForm       | CAS# Qual      |
| 1       | 5,16[1',2']:                        | 8,13[1'',2'']-Dibenz... | 408 C32H24       | 005672-97-9 43 |
| 2       | [1,1'-Biphenyl]-2,2'-dicarbonitrile |                         | 204 C14H8N2      | 004341-02-0 38 |
| 3       | 9,10-di(Chloromethyl)anthracene     |                         | 274 C16H12Cl2    | 010387-13-0 38 |
| 4       | [1,1'-Biphenyl]-4-ol, 4'-chloro-    |                         | 204 C12H9ClO     | 028034-99-3 35 |
| 5       | Naphthalene, 2-phenyl-              |                         | 204 C16H12       | 000612-94-2 30 |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

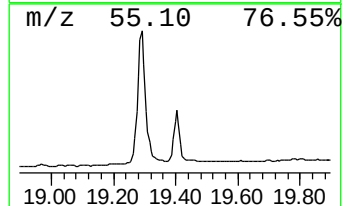
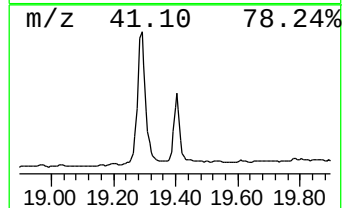
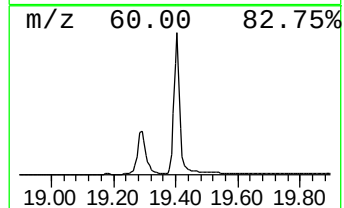
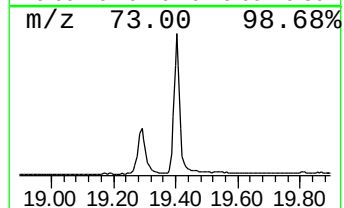
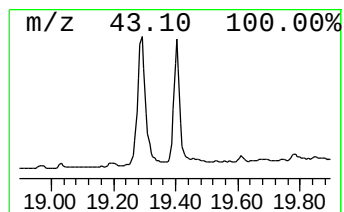
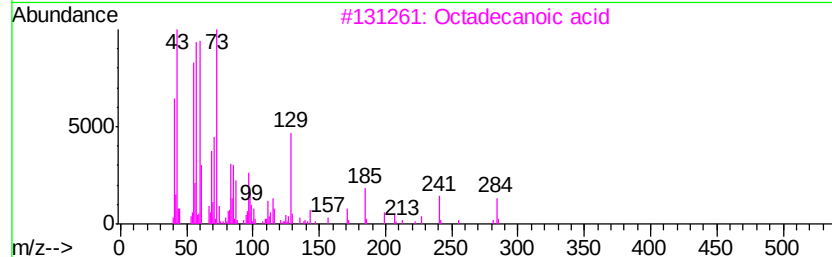
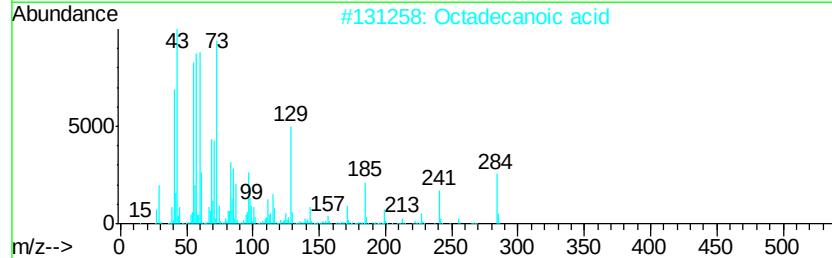
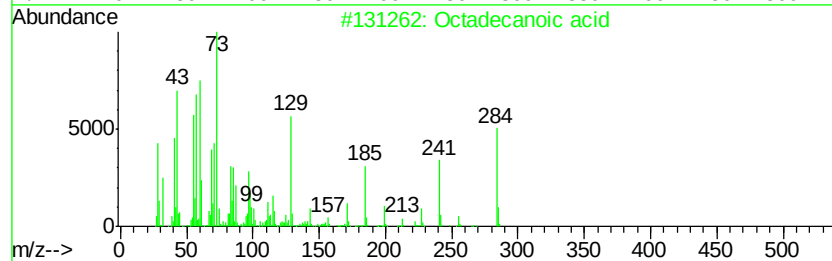
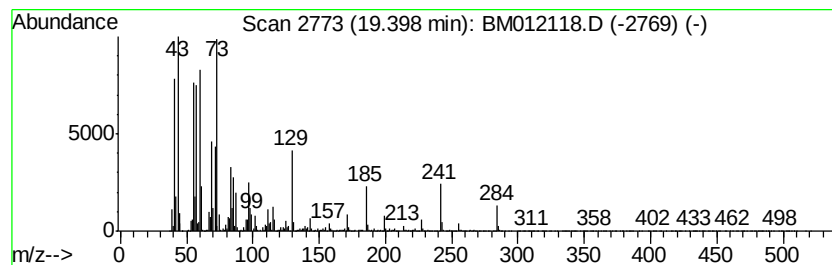
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 42 Octadecanoic acid Concentration Rank 1

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 19.40 | 77.60 ng/ul | 11788100 | Chrysene-d12     | 21.42 |

| Hit# | of | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 3    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 4    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 98   |
| 5    |    | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
Data File : BM012118.D  
Acq On : 13 Oct 2017 02:52  
Operator : SJ/JU  
Sample : I5490-06  
Misc :  
ALS Vial : 18 Sample Multiplier: 1

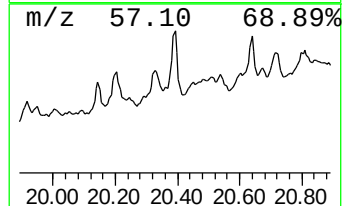
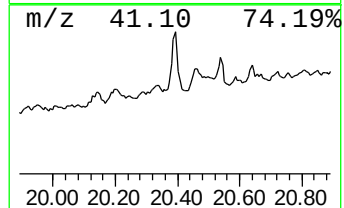
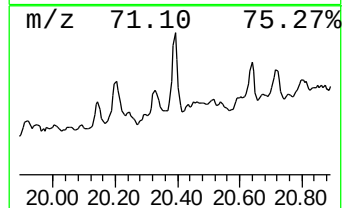
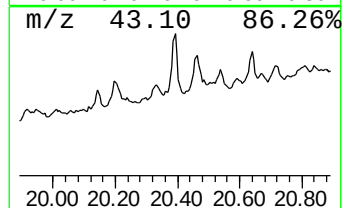
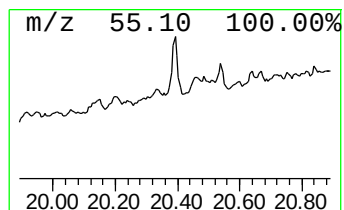
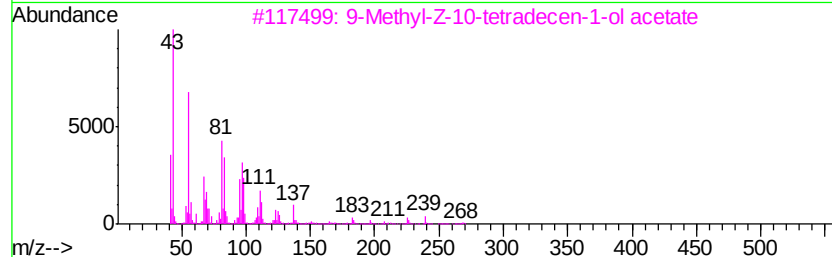
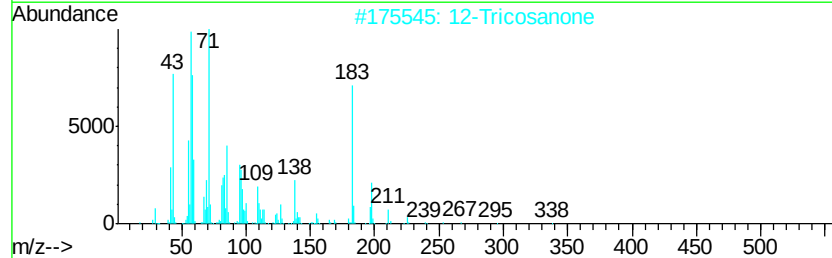
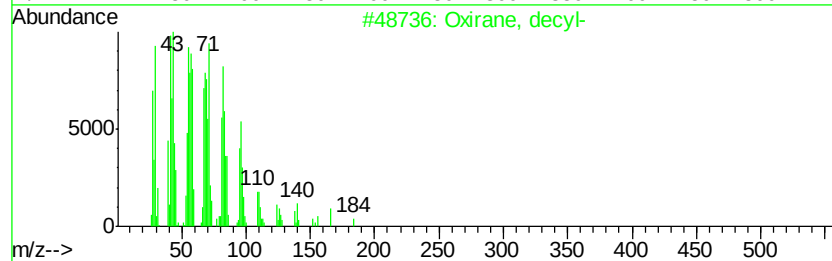
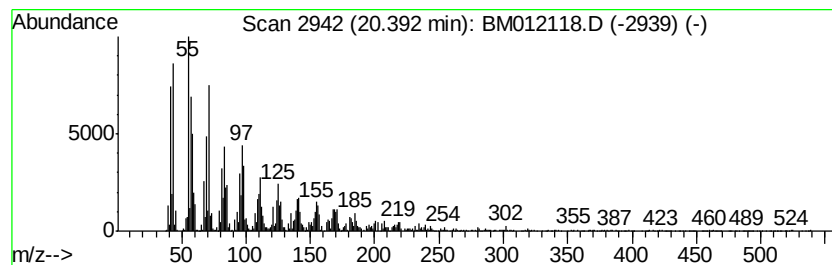
Instrument :  
BNA\_M  
ClientSampleId :  
B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 43 Oxirane, decyl- Concentration Rank 10

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 20.39   | 17.18 ng/ul                         | 2609830      | Chrysene-d12     | 21.42     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Oxirane, decyl-                     | 184 C12H24O  | 002855-19-8      | 64        |
| 2       | 12-Tricosanone                      | 338 C23H46O  | 000540-09-0      | 59        |
| 3       | 9-Methyl-Z-10-tetradecen-1-ol ac... | 268 C17H32O2 | 1000130-99-4     | 58        |
| 4       | Oxirane, tetradecyl-                | 240 C16H32O  | 007320-37-8      | 58        |
| 5       | 9-Heptadecanone                     | 254 C17H34O  | 000540-08-9      | 55        |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM101217\  
 Data File : BM012118.D  
 Acq On : 13 Oct 2017 02:52  
 Operator : SJ/JU  
 Sample : I5490-06  
 Misc :  
 ALS Vial : 18 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 B-69(1-3)-092717

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM101217.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 |
| Tetrachloroethylene  | 4.53  | 10.3    | ng/ul | 1017060  | 1                     | 7.83  | 1968210 | 20.0 |
| Benzene, 1,3-dime... | 5.46  | 5.6     | ng/ul | 551618   | 1                     | 7.83  | 1968210 | 20.0 |
| Styrene              | 5.81  | 16.4    | ng/ul | 1611550  | 1                     | 7.83  | 1968210 | 20.0 |
| Benzene, (1-methy... | 6.31  | 4.1     | ng/ul | 404630   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 6.80  | 13.1    | ng/ul | 1287040  | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 7.44  | 6.0     | ng/ul | 594861   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 7.65  | 3.7     | ng/ul | 363279   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 7.94  | 8.5     | ng/ul | 833122   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 8.05  | 38.3    | ng/ul | 3765680  | 1                     | 7.83  | 1968210 | 20.0 |
| Sulfurous acid, b... | 8.30  | 8.6     | ng/ul | 846153   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 8.36  | 50.5    | ng/ul | 4975060  | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 8.49  | 39.9    | ng/ul | 3923370  | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 8.85  | 14.7    | ng/ul | 1448830  | 1                     | 7.83  | 1968210 | 20.0 |
| unknown-01           | 8.93  | 22.4    | ng/ul | 2203770  | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 9.13  | 5.9     | ng/ul | 579570   | 1                     | 7.83  | 1968210 | 20.0 |
| (DEL) Alkane: Str... | 9.19  | 5.4     | ng/ul | 532554   | 1                     | 7.83  | 1968210 | 20.0 |
| 2-Ethyl-1-hexanol... | 9.26  | 2.0     | ng/ul | 344808   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.33  | 9.1     | ng/ul | 1569940  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.39  | 20.7    | ng/ul | 3565930  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.43  | 3.2     | ng/ul | 543987   | 2                     | 10.64 | 3443740 | 20.0 |
| o-Cymene             | 9.52  | 3.4     | ng/ul | 590182   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.58  | 4.0     | ng/ul | 685737   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.63  | 5.4     | ng/ul | 932566   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.82  | 3.6     | ng/ul | 613473   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 9.89  | 9.9     | ng/ul | 1711510  | 2                     | 10.64 | 3443740 | 20.0 |
| Sulfurous acid, b... | 10.02 | 30.5    | ng/ul | 5250810  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.09 | 17.6    | ng/ul | 3029550  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.13 | 9.6     | ng/ul | 1651440  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.20 | 4.8     | ng/ul | 833522   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.35 | 6.5     | ng/ul | 1126240  | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.59 | 4.0     | ng/ul | 689555   | 2                     | 10.64 | 3443740 | 20.0 |
| (DEL) Alkane: Str... | 10.90 | 2.0     | ng/ul | 349349   | 2                     | 10.64 | 3443740 | 20.0 |
| Decanoic acid, si... | 11.44 | 2.6     | ng/ul | 454811   | 2                     | 10.64 | 3443740 | 20.0 |
| Naphthalene, 2,3,... | 14.89 | 2.2     | ng/ul | 419194   | 3                     | 14.49 | 3792690 | 20.0 |
| (DEL) Alkane: Str... | 15.32 | 2.0     | ng/ul | 382108   | 3                     | 14.49 | 3792690 | 20.0 |
| Naphthalene, 1,6-... | 16.20 | 7.8     | ng/ul | 1450610  | 4                     | 17.23 | 3731340 | 20.0 |
| unknown-02           | 16.62 | 2.6     | ng/ul | 482570   | 4                     | 17.23 | 3731340 | 20.0 |
| (DEL) Alkane: Str... | 17.02 | 4.2     | ng/ul | 775569   | 4                     | 17.23 | 3731340 | 20.0 |
| n-Hexadecanoic acid  | 18.12 | 46.3    | ng/ul | 8628340  | 4                     | 17.23 | 3731340 | 20.0 |
| Benzaldehyde, 3,5... | 18.30 | 2.8     | ng/ul | 514865   | 4                     | 17.23 | 3731340 | 20.0 |
| unknown-03           | 18.60 | 5.4     | ng/ul | 1004150  | 4                     | 17.23 | 3731340 | 20.0 |
| Octadecanoic acid    | 19.40 | 77.6    | ng/ul | 11788100 | 5                     | 21.42 | 3038280 | 20.0 |
| Oxirane, decyl-      | 20.39 | 17.2    | ng/ul | 2609830  | 5                     | 21.42 | 3038280 | 20.0 |