

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
 Data File : BM019715.D  
 Acq On : 03 Apr 2019 05:41  
 Operator : JU/SJ  
 Sample : K2168-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF638

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.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         | 5.575       | 434           | 440         | 447          | rVB2     | 257055         | 419571        | 5.65%           | 0.323%        |
| 2         | 6.751       | 635           | 640         | 655          | rVB      | 535072         | 1028990       | 13.86%          | 0.793%        |
| 3         | 6.922       | 664           | 669         | 678          | rBV2     | 496316         | 954199        | 12.85%          | 0.735%        |
| 4         | 7.104       | 694           | 700         | 705          | rVV      | 674575         | 1128310       | 15.20%          | 0.870%        |
| 5         | 7.157       | 705           | 709         | 713          | rVV      | 460654         | 628030        | 8.46%           | 0.484%        |
| 6         | 7.504       | 763           | 768         | 773          | rVB      | 239083         | 339015        | 4.57%           | 0.261%        |
| 7         | 7.569       | 773           | 779         | 789          | rVB      | 588572         | 1032863       | 13.91%          | 0.796%        |
| 8         | 8.187       | 877           | 884         | 890          | rBV2     | 310599         | 507717        | 6.84%           | 0.391%        |
| 9         | 8.281       | 894           | 900         | 908          | rBV2     | 852630         | 1525518       | 20.55%          | 1.176%        |
| 10        | 8.645       | 956           | 962         | 969          | rVV2     | 282167         | 546754        | 7.36%           | 0.421%        |
| 11        | 8.745       | 971           | 979         | 990          | rVB3     | 1207340        | 2592844       | 34.93%          | 1.998%        |
| 12        | 8.981       | 1009          | 1019        | 1026         | rBV6     | 114990         | 348786        | 4.70%           | 0.269%        |
| 13        | 9.163       | 1045          | 1050        | 1056         | rVB2     | 238060         | 357562        | 4.82%           | 0.276%        |
| 14        | 9.451       | 1087          | 1099        | 1105         | rBV2     | 640009         | 1271228       | 17.12%          | 0.980%        |
| 15        | 9.739       | 1141          | 1148        | 1155         | rBV3     | 420848         | 796024        | 10.72%          | 0.614%        |
| 16        | 9.916       | 1170          | 1178        | 1183         | rBV4     | 120281         | 296811        | 4.00%           | 0.229%        |
| 17        | 9.981       | 1183          | 1189        | 1195         | rBV      | 803869         | 1528814       | 20.59%          | 1.178%        |
| 18        | 10.034      | 1195          | 1198        | 1205         | rVB      | 174713         | 283508        | 3.82%           | 0.219%        |
| 19        | 10.292      | 1236          | 1242        | 1247         | rBV      | 1343085        | 2100836       | 28.30%          | 1.619%        |
| 20        | 10.345      | 1247          | 1251        | 1256         | rVB      | 734058         | 1113664       | 15.00%          | 0.858%        |
| 21        | 10.475      | 1263          | 1273        | 1276         | rBV      | 597105         | 1085348       | 14.62%          | 0.837%        |
| 22        | 10.510      | 1276          | 1279        | 1291         | rVB      | 758687         | 1487309       | 20.03%          | 1.146%        |
| 23        | 10.798      | 1324          | 1328        | 1335         | rVB2     | 216622         | 349946        | 4.71%           | 0.270%        |
| 24        | 11.016      | 1355          | 1365        | 1367         | rBV5     | 140928         | 367296        | 4.95%           | 0.283%        |
| 25        | 11.151      | 1383          | 1388        | 1393         | rBV4     | 152639         | 279268        | 3.76%           | 0.215%        |
| 26        | 11.233      | 1397          | 1402        | 1409         | rVB4     | 299177         | 577041        | 7.77%           | 0.445%        |
| 27        | 11.334      | 1414          | 1419        | 1429         | rBV      | 906649         | 1554413       | 20.94%          | 1.198%        |
| 28        | 11.598      | 1459          | 1464        | 1468         | rVB2     | 223679         | 348574        | 4.70%           | 0.269%        |
| 29        | 11.751      | 1483          | 1490        | 1495         | rBV      | 1854365        | 2739841       | 36.91%          | 2.112%        |
| 30        | 11.969      | 1524          | 1527        | 1531         | rVB      | 310629         | 432755        | 5.83%           | 0.334%        |
| 31        | 12.022      | 1531          | 1536        | 1540         | rBV      | 570154         | 827810        | 11.15%          | 0.638%        |
| 32        | 12.245      | 1570          | 1574        | 1581         | rVB      | 321223         | 480429        | 6.47%           | 0.370%        |
| 33        | 12.392      | 1594          | 1599        | 1603         | rBV4     | 303637         | 522198        | 7.03%           | 0.402%        |
| 34        | 12.457      | 1603          | 1610        | 1616         | rVV6     | 177776         | 530255        | 7.14%           | 0.409%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
 Data File : BM019715.D  
 Acq On : 03 Apr 2019 05:41  
 Operator : JU/SJ  
 Sample : K2168-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF638

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 35 | 12.580 | 1627 | 1631 | 1635 | rBV2 | 251452  | 308269  | 4.15%   | 0.238% |
| 36 | 12.669 | 1643 | 1646 | 1650 | rBV2 | 228228  | 330916  | 4.46%   | 0.255% |
| 37 | 12.722 | 1650 | 1655 | 1661 | rVB  | 1283539 | 1675199 | 22.57%  | 1.291% |
| 38 | 13.022 | 1698 | 1706 | 1714 | rVB  | 2775047 | 4233155 | 57.02%  | 3.263% |
| 39 | 13.251 | 1739 | 1745 | 1752 | rBV3 | 307455  | 729088  | 9.82%   | 0.562% |
| 40 | 13.380 | 1762 | 1767 | 1768 | rBV  | 434790  | 578559  | 7.79%   | 0.446% |
| 41 | 13.539 | 1789 | 1794 | 1800 | rBV3 | 721863  | 1392564 | 18.76%  | 1.073% |
| 42 | 13.598 | 1800 | 1804 | 1809 | rVV  | 810435  | 1248851 | 16.82%  | 0.963% |
| 43 | 13.651 | 1809 | 1813 | 1816 | rVV  | 1915203 | 2735193 | 36.84%  | 2.108% |
| 44 | 13.686 | 1816 | 1819 | 1823 | rVV2 | 1652758 | 2146802 | 28.92%  | 1.655% |
| 45 | 13.727 | 1823 | 1826 | 1832 | rVB2 | 408905  | 562686  | 7.58%   | 0.434% |
| 46 | 13.922 | 1854 | 1859 | 1864 | rVV  | 1976346 | 2864371 | 38.58%  | 2.208% |
| 47 | 14.033 | 1874 | 1878 | 1885 | rVB5 | 335471  | 687488  | 9.26%   | 0.530% |
| 48 | 14.110 | 1885 | 1891 | 1896 | rBV  | 3392762 | 4402872 | 59.31%  | 3.393% |
| 49 | 14.175 | 1898 | 1902 | 1906 | rVV4 | 318505  | 577910  | 7.78%   | 0.445% |
| 50 | 14.227 | 1907 | 1911 | 1917 | rVB2 | 1511807 | 2274549 | 30.64%  | 1.753% |
| 51 | 14.439 | 1942 | 1947 | 1951 | rBV3 | 342336  | 688828  | 9.28%   | 0.531% |
| 52 | 14.616 | 1973 | 1977 | 1982 | rBV3 | 496840  | 915618  | 12.33%  | 0.706% |
| 53 | 14.704 | 1989 | 1992 | 1994 | rBV2 | 478772  | 612146  | 8.25%   | 0.472% |
| 54 | 14.733 | 1994 | 1997 | 2001 | rVB3 | 654223  | 899171  | 12.11%  | 0.693% |
| 55 | 14.851 | 2012 | 2017 | 2022 | rBV3 | 397367  | 722645  | 9.73%   | 0.557% |
| 56 | 14.904 | 2022 | 2026 | 2031 | rVV  | 323211  | 510660  | 6.88%   | 0.394% |
| 57 | 15.004 | 2039 | 2043 | 2045 | rBV4 | 222634  | 366263  | 4.93%   | 0.282% |
| 58 | 15.069 | 2050 | 2054 | 2059 | rVB  | 2980791 | 3584148 | 48.28%  | 2.762% |
| 59 | 15.233 | 2077 | 2082 | 2087 | rBV  | 2546217 | 3655090 | 49.24%  | 2.817% |
| 60 | 15.327 | 2094 | 2098 | 2102 | rBV2 | 194941  | 333546  | 4.49%   | 0.257% |
| 61 | 15.380 | 2104 | 2107 | 2113 | rVB  | 584024  | 774683  | 10.44%  | 0.597% |
| 62 | 15.474 | 2117 | 2123 | 2129 | rVB  | 1265262 | 2103449 | 28.33%  | 1.621% |
| 63 | 15.627 | 2146 | 2149 | 2156 | rBV3 | 231566  | 416572  | 5.61%   | 0.321% |
| 64 | 15.692 | 2157 | 2160 | 2164 | rVB2 | 343448  | 410221  | 5.53%   | 0.316% |
| 65 | 15.957 | 2197 | 2205 | 2210 | rBV2 | 3226333 | 7423736 | 100.00% | 5.722% |
| 66 | 16.280 | 2256 | 2260 | 2262 | rBV3 | 220345  | 358274  | 4.83%   | 0.276% |
| 67 | 16.339 | 2268 | 2270 | 2275 | rVB3 | 289696  | 309900  | 4.17%   | 0.239% |
| 68 | 16.445 | 2285 | 2288 | 2293 | rVB4 | 250383  | 305518  | 4.12%   | 0.235% |
| 69 | 16.727 | 2332 | 2336 | 2340 | rBV  | 1716899 | 2137651 | 28.79%  | 1.648% |
| 70 | 16.774 | 2340 | 2344 | 2349 | rVB  | 792549  | 1106905 | 14.91%  | 0.853% |
| 71 | 16.992 | 2375 | 2381 | 2385 | rBV  | 1895316 | 2681030 | 36.11%  | 2.066% |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
 Data File : BM019715.D  
 Acq On : 03 Apr 2019 05:41  
 Operator : JU/SJ  
 Sample : K2168-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF638

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 17.092 | 2392 | 2398 | 2405 | rVB  | 2963715 | 4434077 | 59.73% | 3.417% |
| 73  | 17.268 | 2424 | 2428 | 2433 | rVB5 | 215743  | 355258  | 4.79%  | 0.274% |
| 74  | 17.386 | 2444 | 2448 | 2454 | rBV3 | 218694  | 352696  | 4.75%  | 0.272% |
| 75  | 17.468 | 2458 | 2462 | 2467 | rVB  | 1278662 | 1485277 | 20.01% | 1.145% |
| 76  | 17.880 | 2526 | 2532 | 2536 | rBV2 | 644417  | 1076286 | 14.50% | 0.830% |
| 77  | 18.162 | 2576 | 2580 | 2585 | rBV  | 772771  | 896556  | 12.08% | 0.691% |
| 78  | 18.804 | 2683 | 2689 | 2694 | rBV2 | 532514  | 728971  | 9.82%  | 0.562% |
| 79  | 19.174 | 2748 | 2752 | 2760 | rBV  | 335677  | 536441  | 7.23%  | 0.413% |
| 80  | 19.398 | 2784 | 2790 | 2795 | rBV  | 3867366 | 5134127 | 69.16% | 3.957% |
| 81  | 19.733 | 2843 | 2847 | 2850 | rBV  | 1661443 | 1695593 | 22.84% | 1.307% |
| 82  | 19.774 | 2850 | 2854 | 2860 | rVB  | 2827495 | 3256778 | 43.87% | 2.510% |
| 83  | 19.927 | 2877 | 2880 | 2886 | rVB  | 253733  | 301033  | 4.06%  | 0.232% |
| 84  | 20.709 | 3009 | 3013 | 3016 | rBV  | 728798  | 809892  | 10.91% | 0.624% |
| 85  | 20.745 | 3016 | 3019 | 3026 | rVB  | 1494453 | 1560591 | 21.02% | 1.203% |
| 86  | 20.898 | 3041 | 3045 | 3055 | rVB  | 1204151 | 1493845 | 20.12% | 1.151% |
| 87  | 21.198 | 3091 | 3096 | 3108 | rBV  | 2416826 | 3197386 | 43.07% | 2.464% |
| 88  | 21.327 | 3115 | 3118 | 3122 | rBV  | 248426  | 302215  | 4.07%  | 0.233% |
| 89  | 21.574 | 3157 | 3160 | 3163 | rBV  | 373961  | 439517  | 5.92%  | 0.339% |
| 90  | 21.609 | 3163 | 3166 | 3170 | rVB  | 790435  | 834388  | 11.24% | 0.643% |
| 91  | 21.756 | 3187 | 3191 | 3194 | rBV  | 350851  | 387104  | 5.21%  | 0.298% |
| 92  | 22.203 | 3264 | 3267 | 3274 | rVB  | 411020  | 538063  | 7.25%  | 0.415% |
| 93  | 22.480 | 3311 | 3314 | 3317 | rBV  | 285545  | 370061  | 4.98%  | 0.285% |
| 94  | 22.521 | 3317 | 3321 | 3326 | rVB  | 615052  | 817630  | 11.01% | 0.630% |
| 95  | 22.697 | 3347 | 3351 | 3356 | rVB  | 517846  | 701970  | 9.46%  | 0.541% |
| 96  | 23.262 | 3439 | 3447 | 3462 | rBV2 | 2852795 | 5795308 | 78.06% | 4.467% |
| 97  | 23.403 | 3464 | 3471 | 3480 | rVB  | 1585349 | 3047071 | 41.04% | 2.348% |
| 98  | 23.868 | 3545 | 3550 | 3559 | rVB  | 388269  | 733067  | 9.87%  | 0.565% |
| 99  | 23.974 | 3564 | 3568 | 3577 | rVB3 | 213122  | 447346  | 6.03%  | 0.345% |
| 100 | 24.591 | 3668 | 3673 | 3680 | rVB  | 281992  | 570469  | 7.68%  | 0.440% |

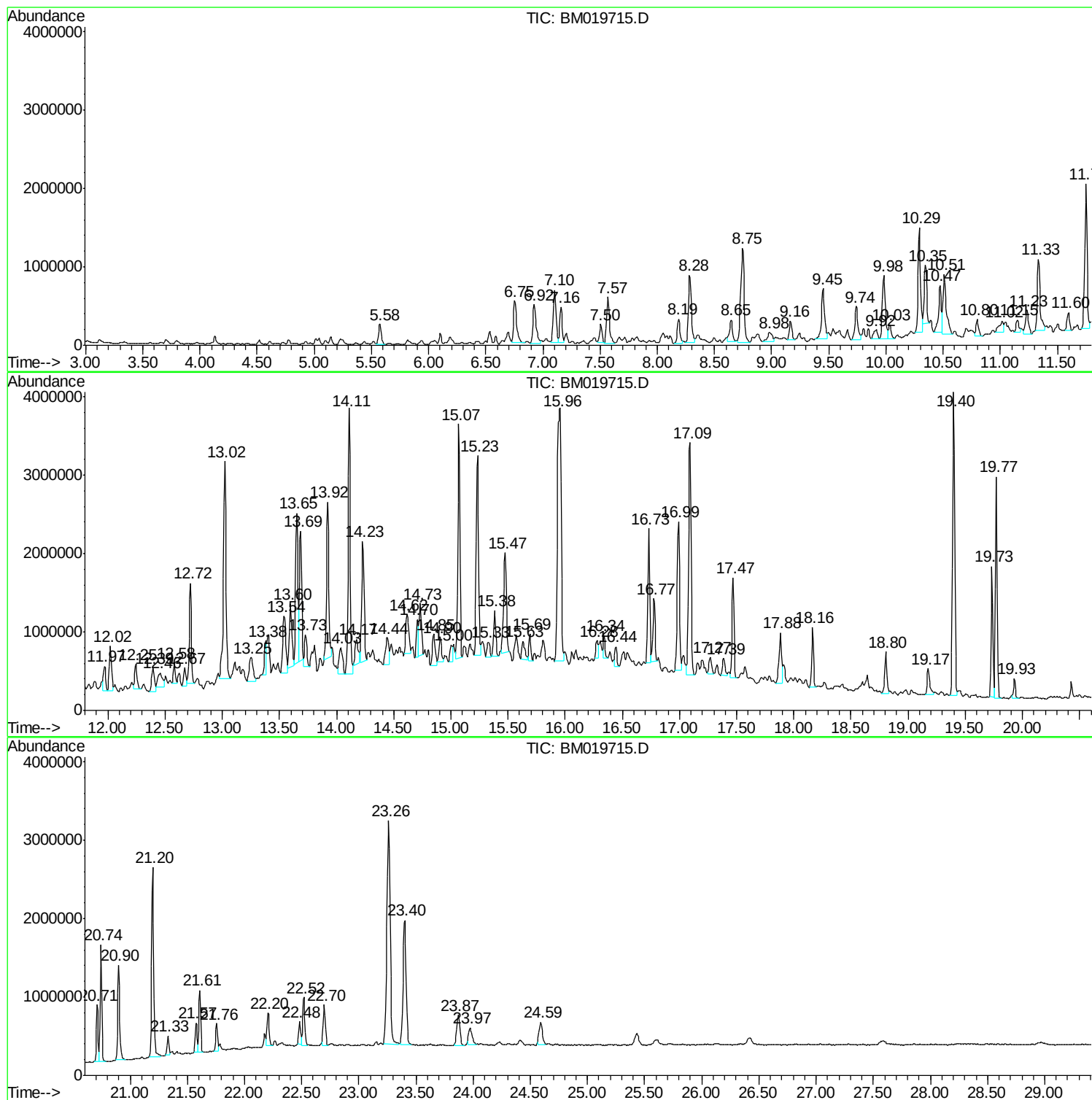
Sum of corrected areas: 129747069

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

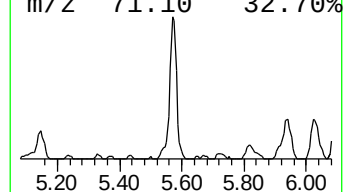
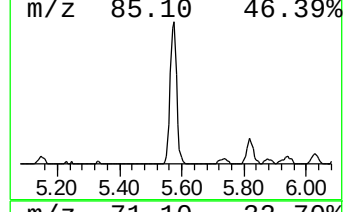
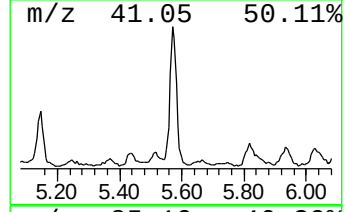
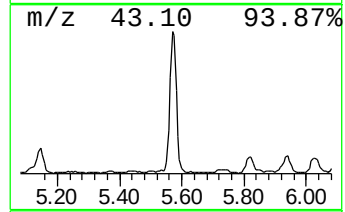
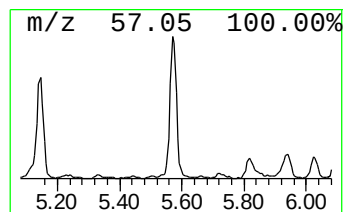
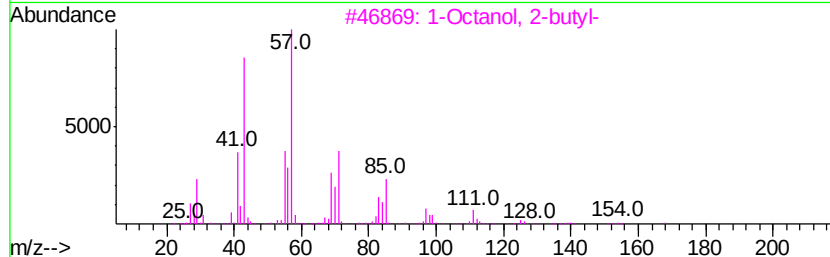
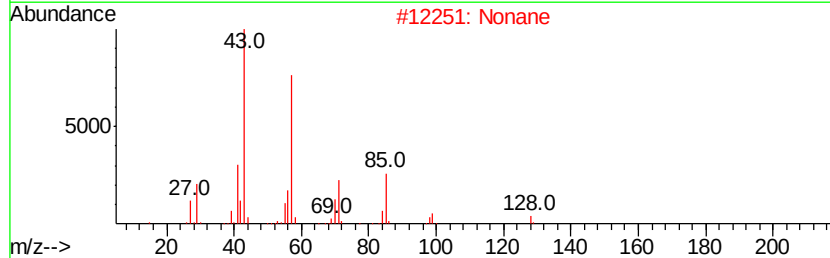
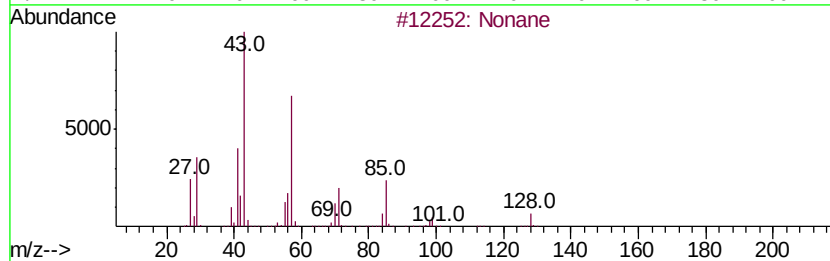
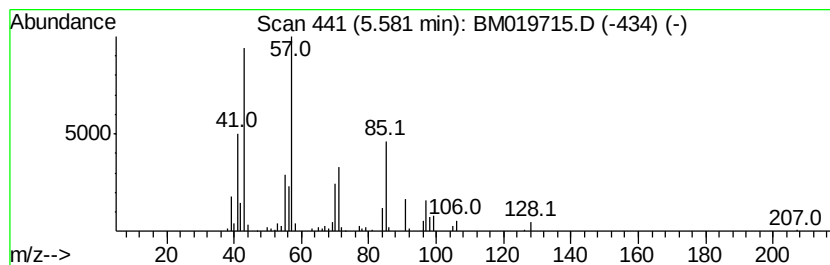
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.58 | 8.12 ng/ul | 419571 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of                  | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|---------------------|--------------|-----|---------|-------------|------|
| 1    | Nonane              |              | 128 | C9H20   | 000111-84-2 | 93   |
| 2    | Nonane              |              | 128 | C9H20   | 000111-84-2 | 90   |
| 3    | 1-Octanol, 2-butyl- |              | 186 | C12H26O | 003913-02-8 | 64   |
| 4    | Octane              |              | 114 | C8H18   | 000111-65-9 | 64   |
| 5    | Decane              |              | 142 | C10H22  | 000124-18-5 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

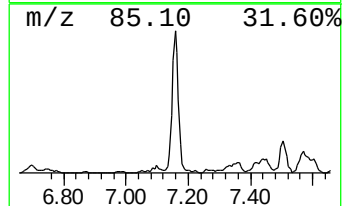
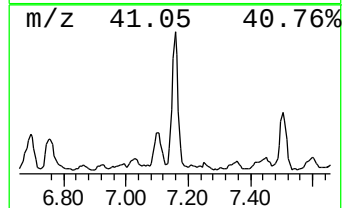
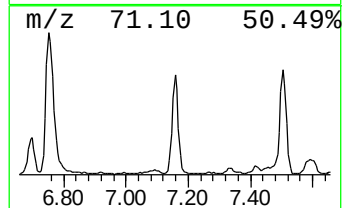
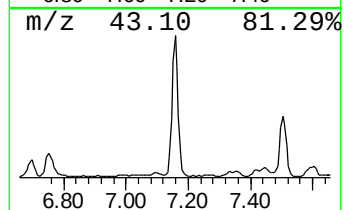
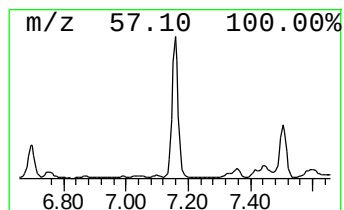
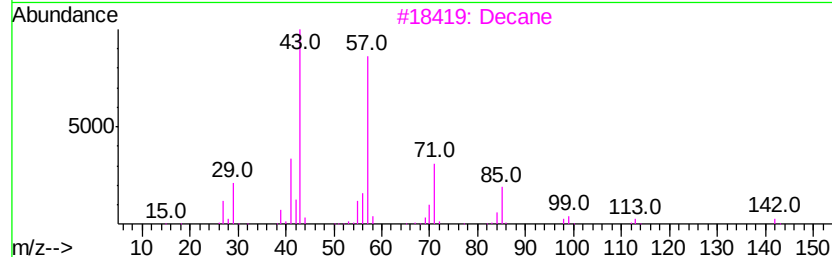
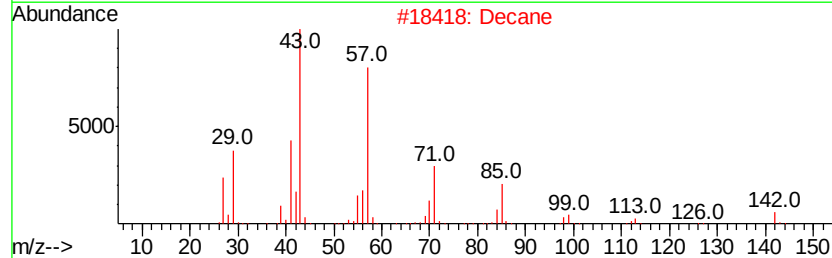
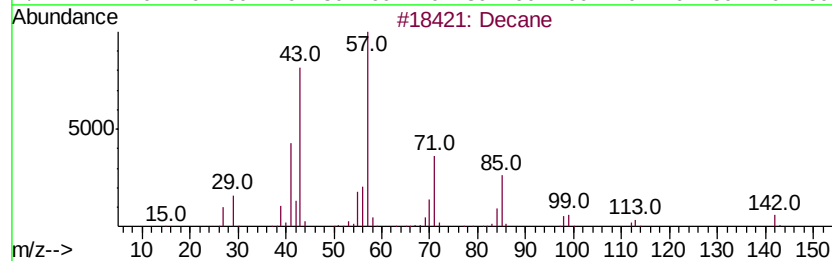
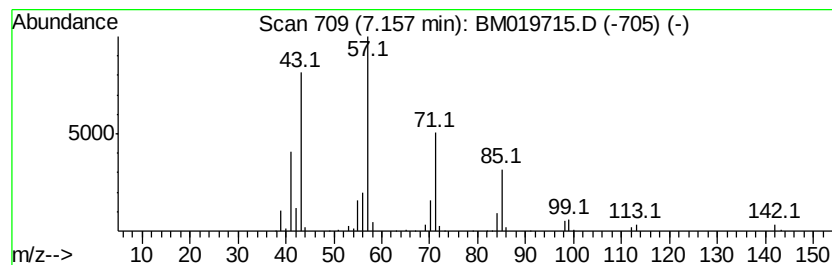
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.16 | 12.16 ng/ul | 628030 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Decane       | 142 | C10H22  | 000124-18-5 | 95   |
| 2    |    | Decane       | 142 | C10H22  | 000124-18-5 | 95   |
| 3    |    | Decane       | 142 | C10H22  | 000124-18-5 | 94   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 5    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

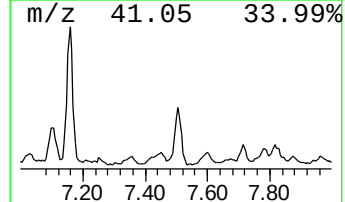
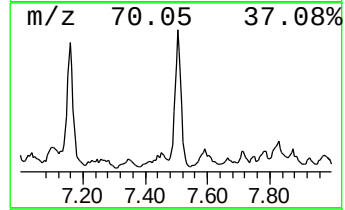
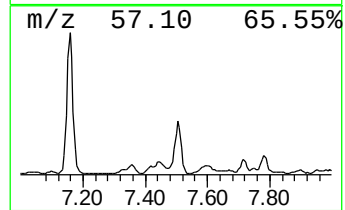
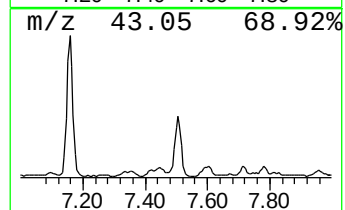
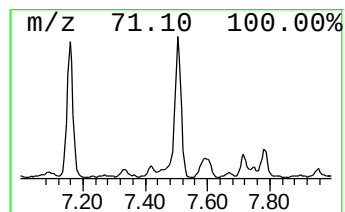
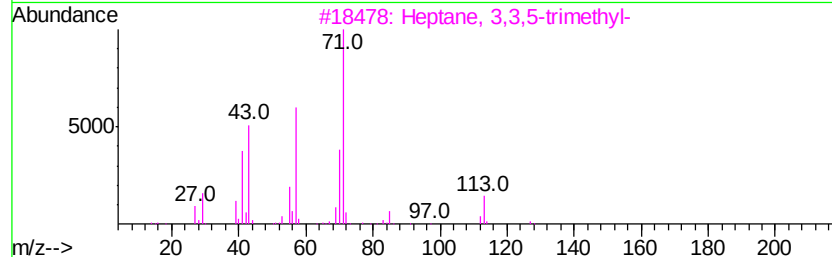
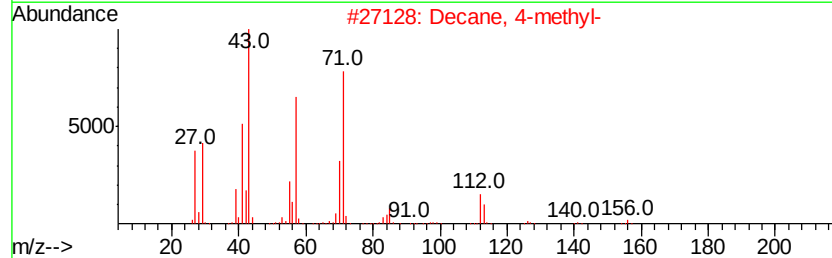
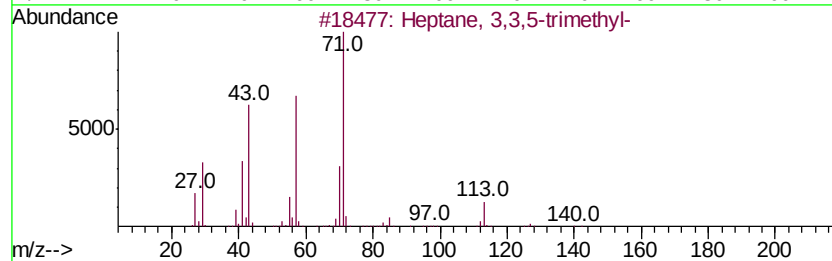
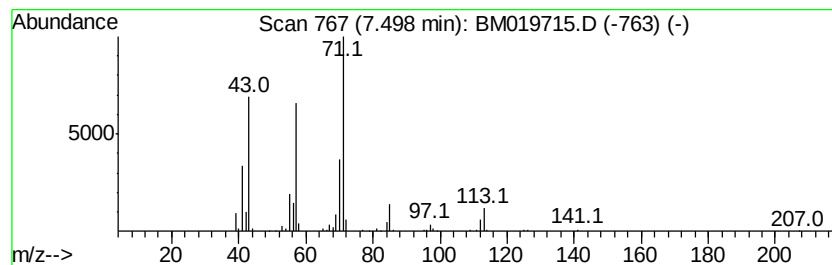
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.50 | 6.56 ng/ul | 339015 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 78   |
| 2    |    |   | Decane, 4-methyl-         | 156 | C11H24  | 002847-72-5 | 78   |
| 3    |    |   | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 78   |
| 4    |    |   | Octane, 3,3-dimethyl-     | 142 | C10H22  | 004110-44-5 | 72   |
| 5    |    |   | Octane, 2,6-dimethyl-     | 142 | C10H22  | 002051-30-1 | 72   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

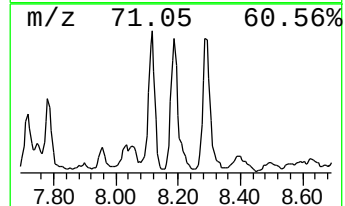
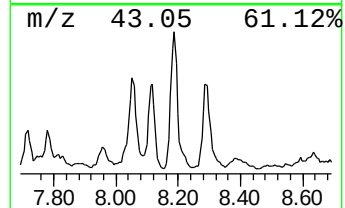
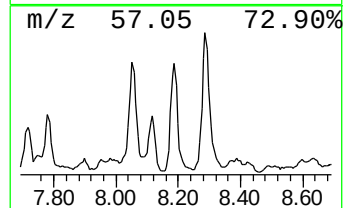
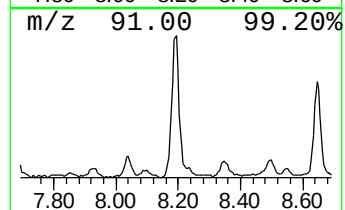
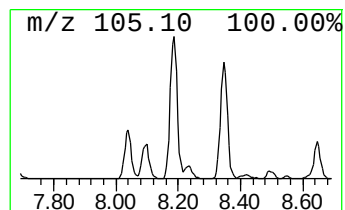
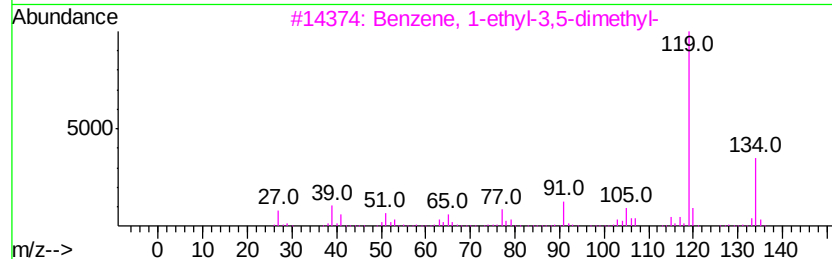
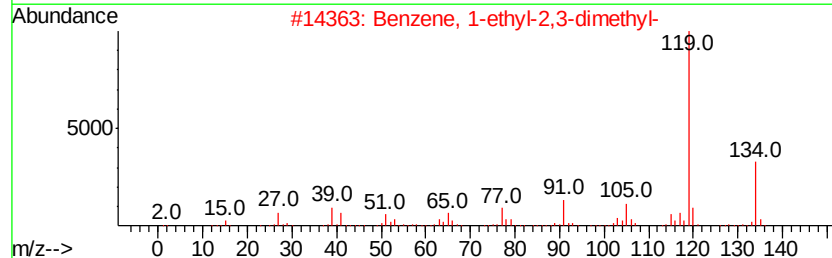
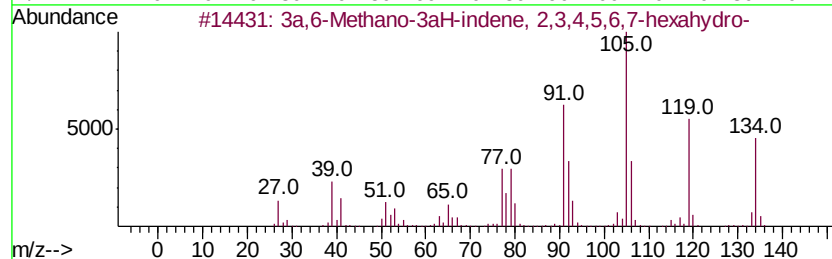
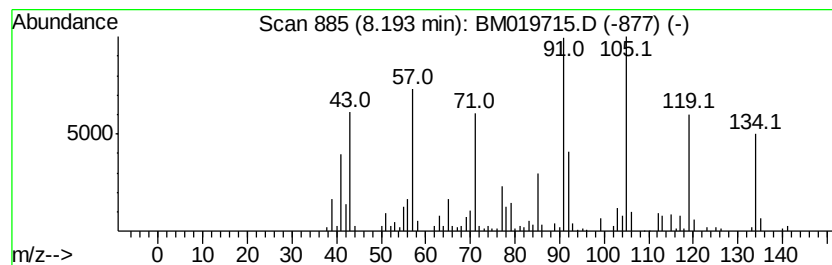
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 3a,6-Methano-3aH-indene, 2,... Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.19 | 9.83 ng/ul | 507717 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------------|-----|---------|-------------|------|
| 1       |   | 3a,6-Methano-3aH-indene, 2,3,4,5... | 134 | C10H14  | 098640-10-9 | 55   |
| 2       |   | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 50   |
| 3       |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 50   |
| 4       |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 46   |
| 5       |   | Benzene, 1,2,3,4-tetramethyl-       | 134 | C10H14  | 000488-23-3 | 46   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

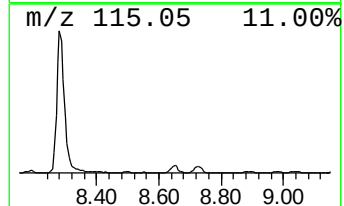
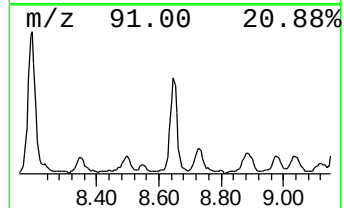
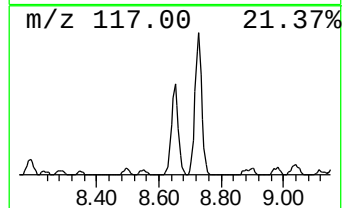
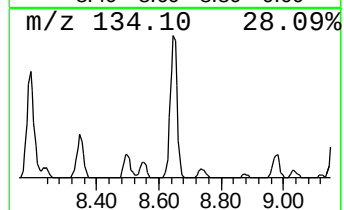
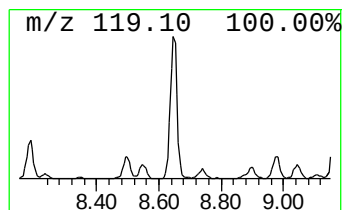
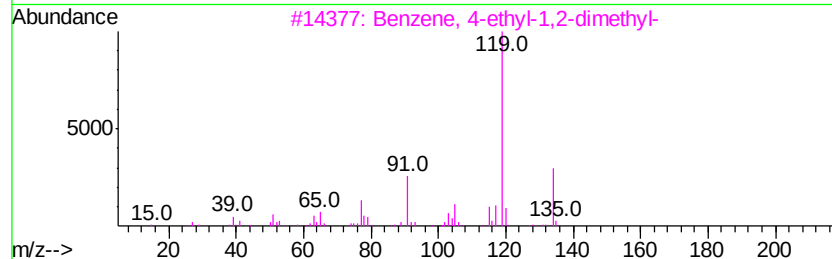
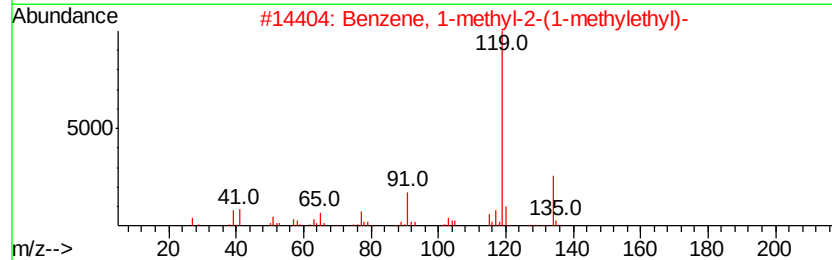
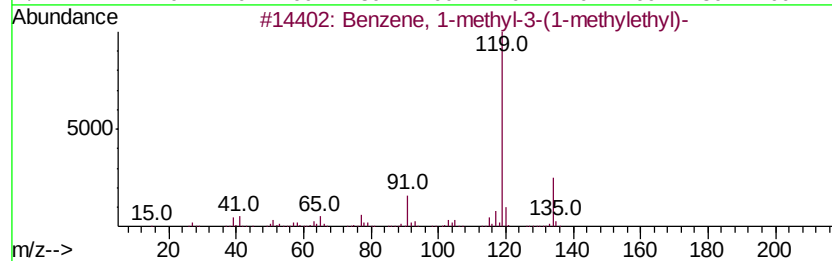
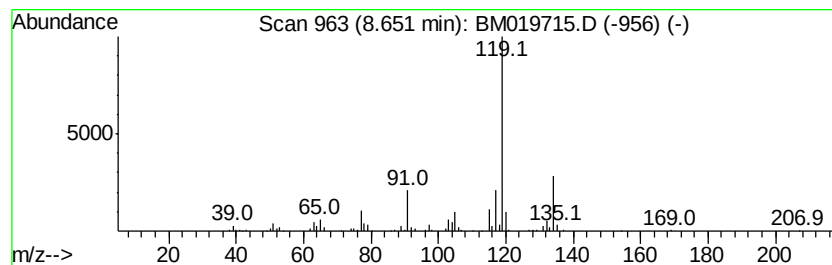
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Benzene, 1-methyl-3-(1-meth... Concentration Rank 18

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 8.65 | 10.59 ng/ul | 546754 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 95   |
| 2    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 95   |
| 3    |    | Benzene, 4-ethyl-1,2-dimethyl-       | 134 | C10H14  | 000934-80-5 | 94   |
| 4    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 94   |
| 5    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

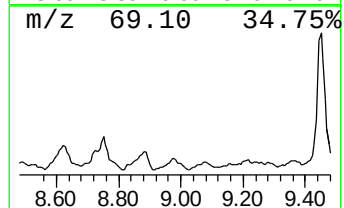
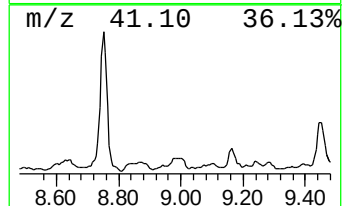
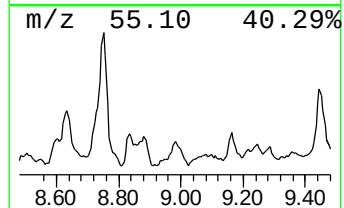
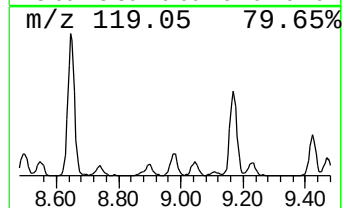
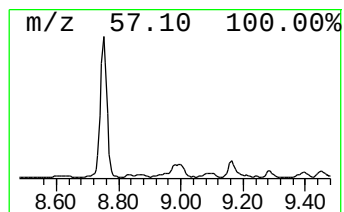
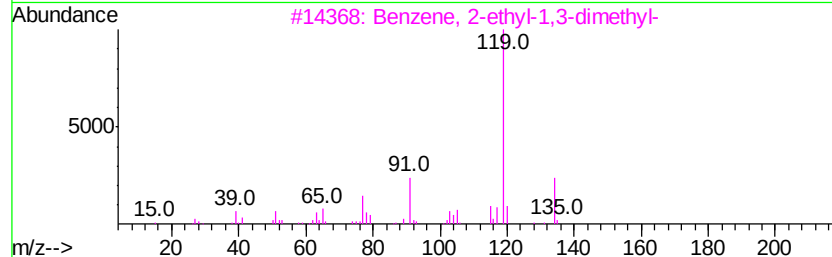
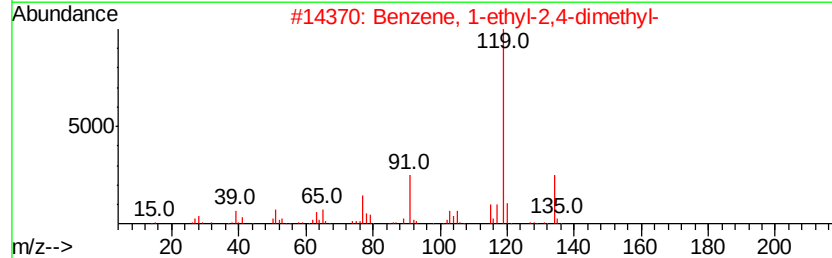
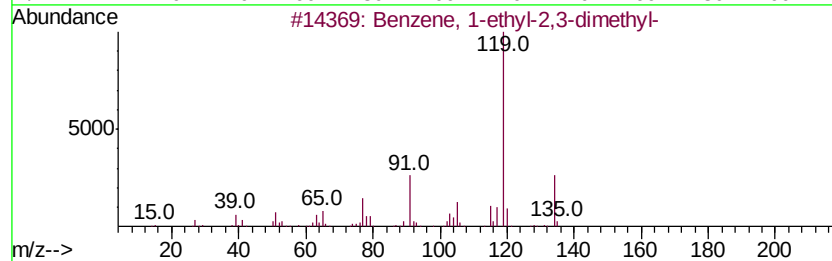
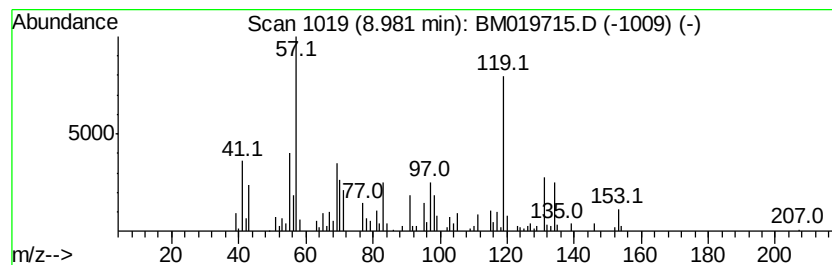
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 37

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 8.98 | 6.26 ng/ul | 348786 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 89   |
| 2    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 86   |
| 3    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 86   |
| 4    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 86   |
| 5    |    | 1,4-Cyclohexadiene, 3-ethenyl-1,... | 134 | C10H14  | 062338-57-2 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

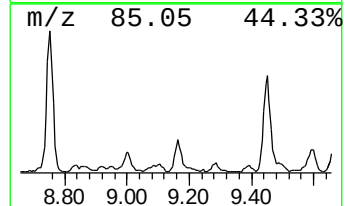
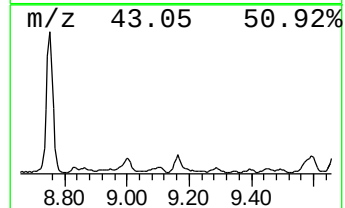
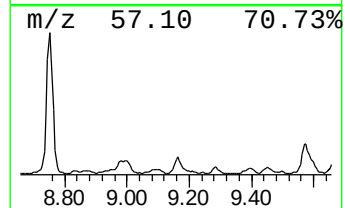
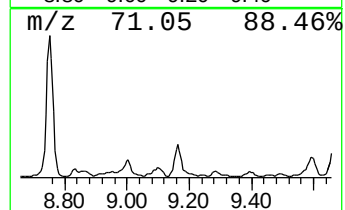
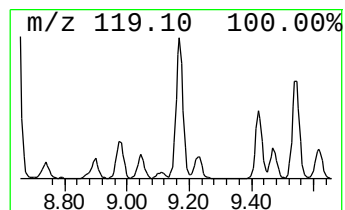
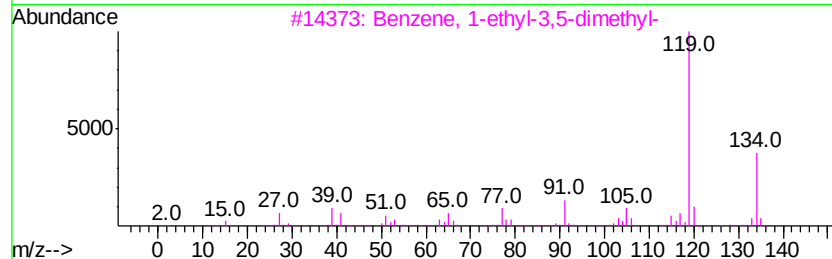
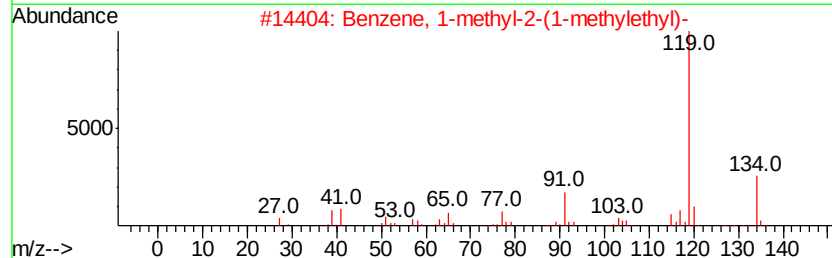
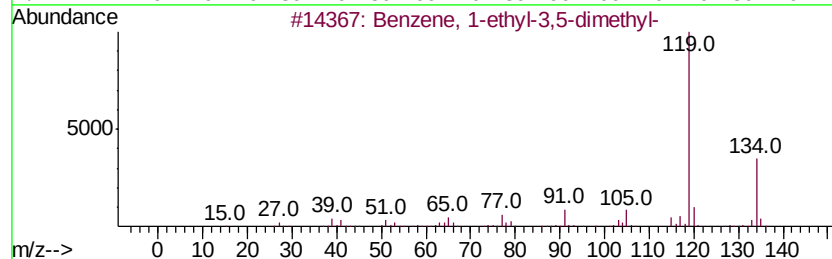
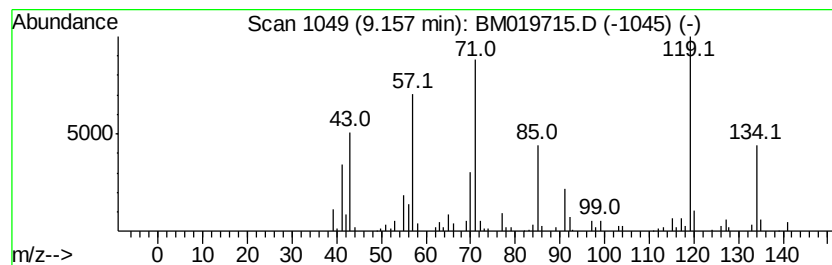
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 33

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.16 | 6.42 ng/ul | 357562 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 90   |
| 2    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 87   |
| 3    |    | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 60   |
| 4    |    | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 60   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

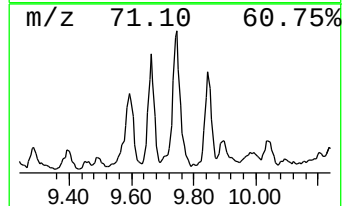
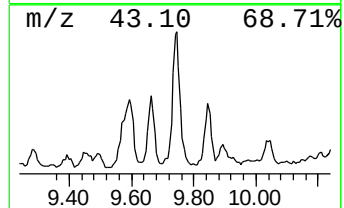
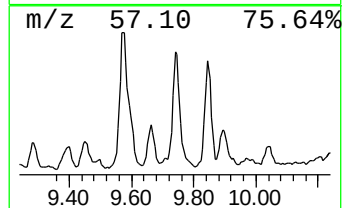
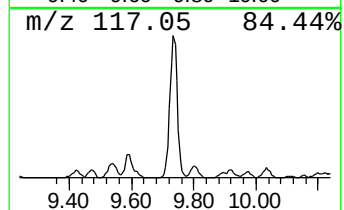
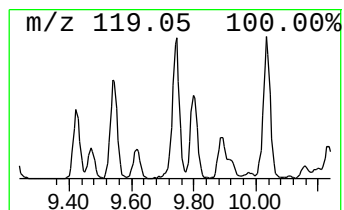
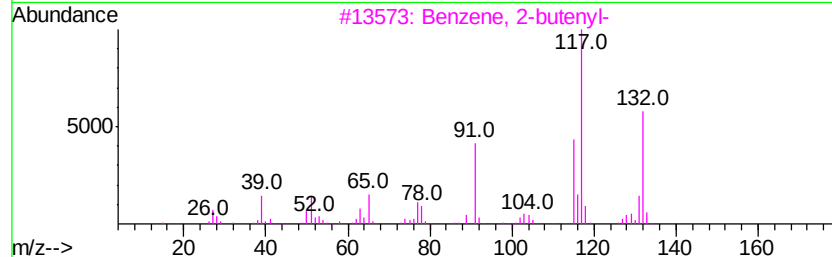
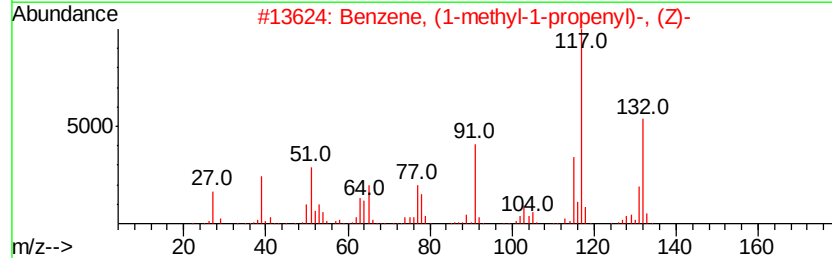
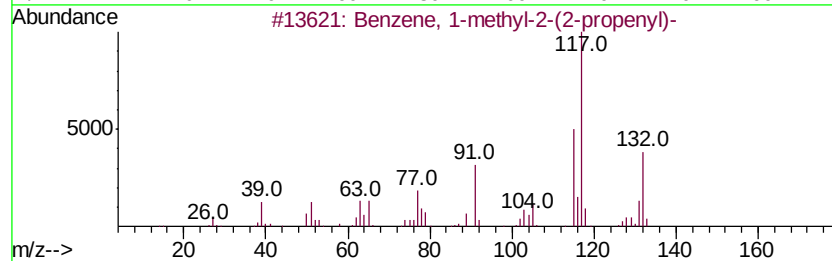
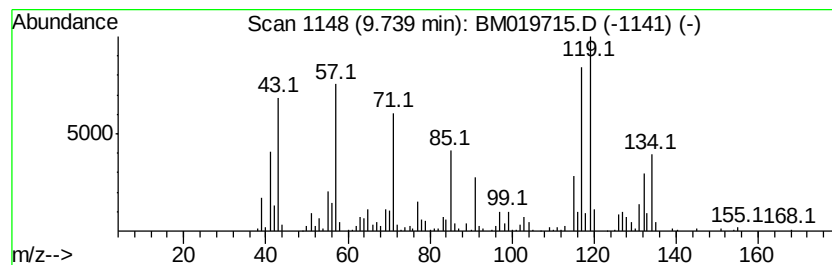
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Benzene, 1-methyl-2-(2-prop... Concentration Rank 12

| R.T. | EstConc     | Area   | Relative to ISTD | R.T.  |
|------|-------------|--------|------------------|-------|
| 9.74 | 14.30 ng/ul | 796024 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 53   |
| 2    |    | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000767-99-7 | 43   |
| 3    |    | Benzene, 2-butenyl-                 | 132 | C10H12  | 001560-06-1 | 43   |
| 4    |    | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 43   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 42   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

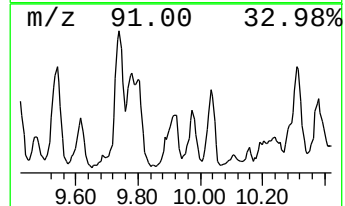
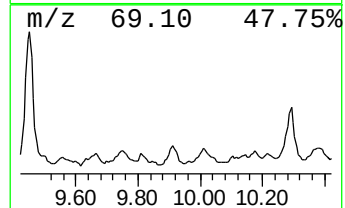
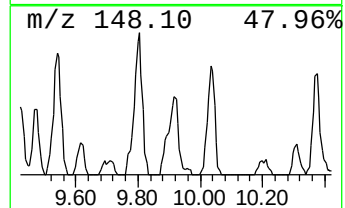
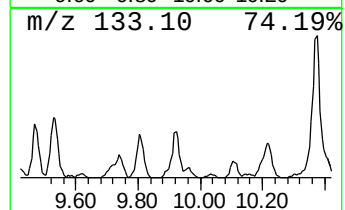
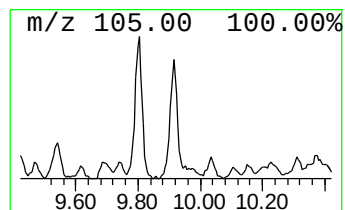
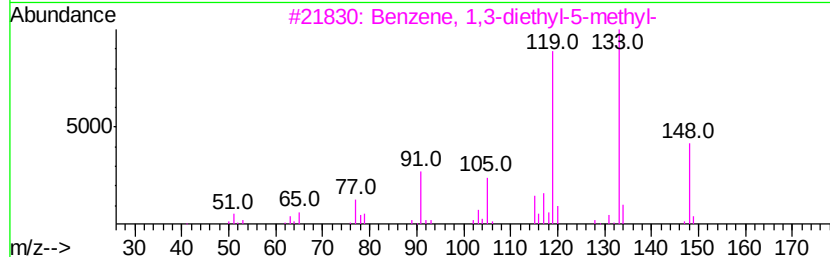
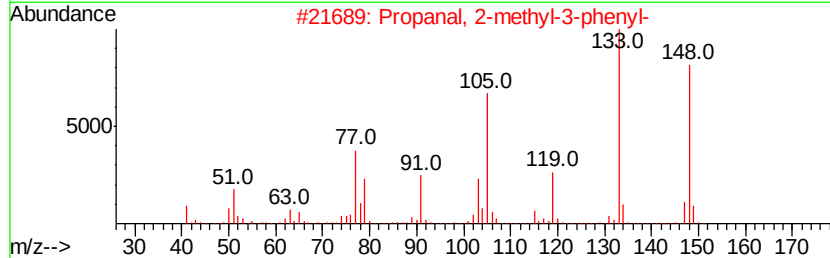
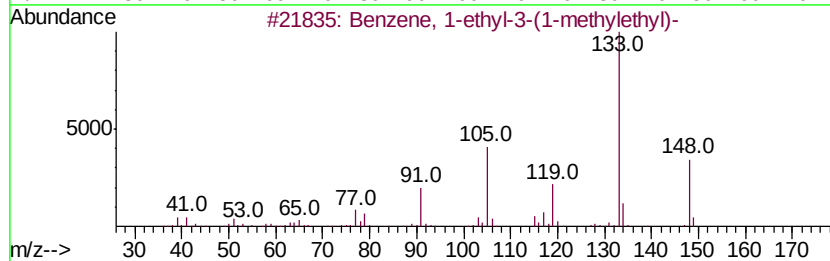
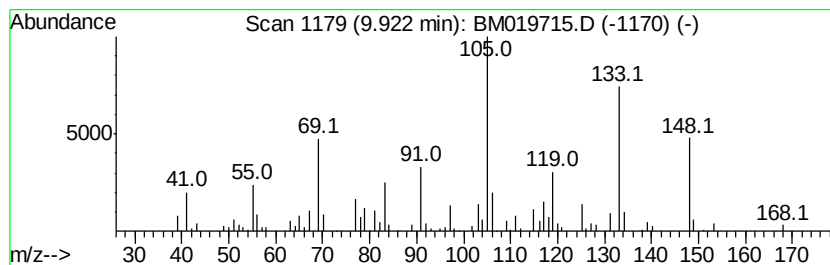
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Benzene, 1-ethyl-3-(1-methylethyl) Concentration Rank 43

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.92 | 5.33 ng/ul | 296811 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16  | 004920-99-4  | 70   |
| 2    |    | Propanal, 2-methyl-3-phenyl-        | 148 | C10H12O | 1000131-87-6 | 53   |
| 3    |    | Benzene, 1,3-diethyl-5-methyl-      | 148 | C11H16  | 002050-24-0  | 49   |
| 4    |    | Benzaldehyde, 4-(1-methylethyl)-    | 148 | C10H12O | 000122-03-2  | 47   |
| 5    |    | 1-(2,3-Dimethylphenyl)ethanone      | 148 | C10H12O | 002142-71-4  | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

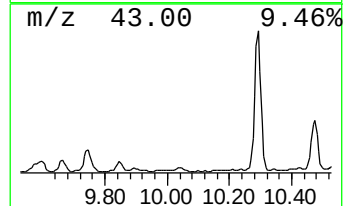
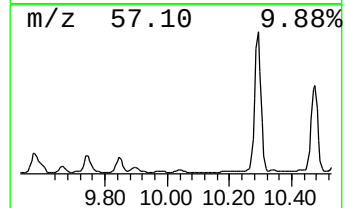
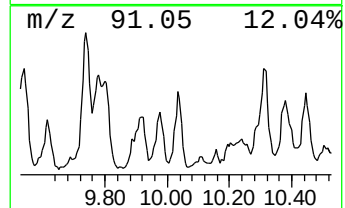
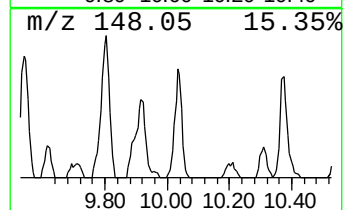
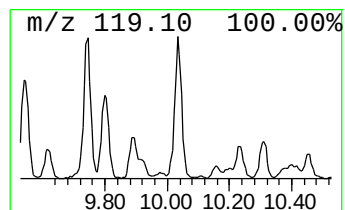
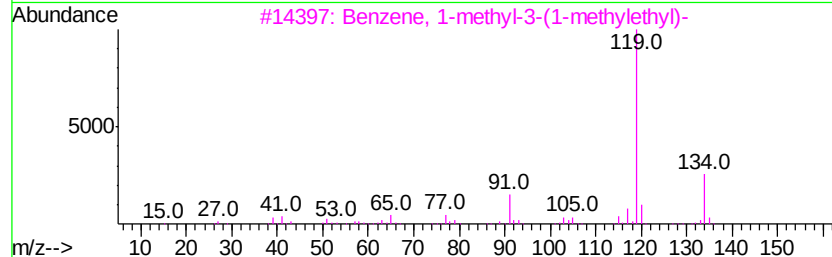
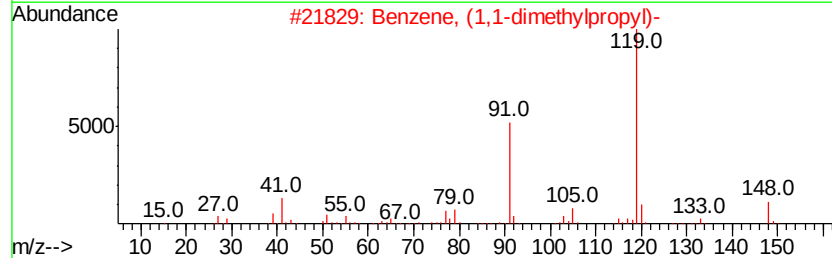
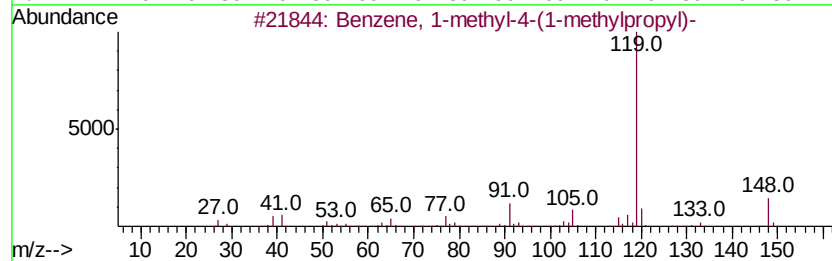
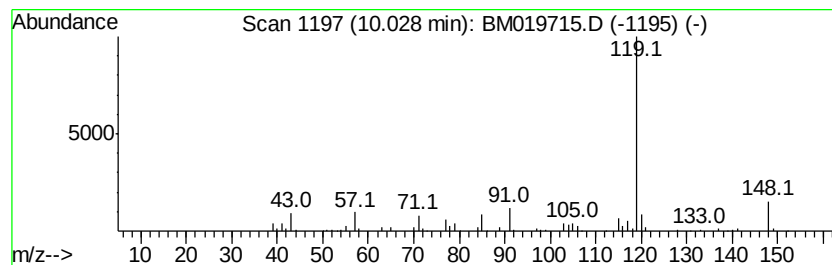
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzene, 1-methyl-4-(1-meth... Concentration Rank 45

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.03 | 5.09 ng/ul | 283508 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-(1-methylpro... | 148 | C11H16  | 001595-16-0 | 68   |
| 2    |    | Benzene, (1,1-dimethylpropyl)-      | 148 | C11H16  | 002049-95-8 | 59   |
| 3    |    | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 53   |
| 4    |    | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14  | 000099-87-6 | 53   |
| 5    |    | Pentadecane, 2-methyl-2-phenyl-     | 302 | C22H38  | 029138-94-1 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

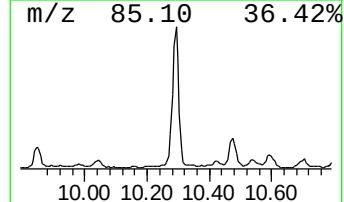
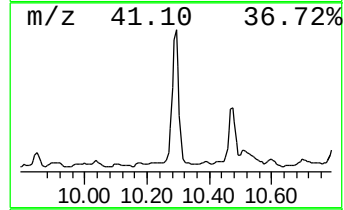
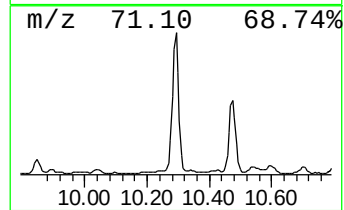
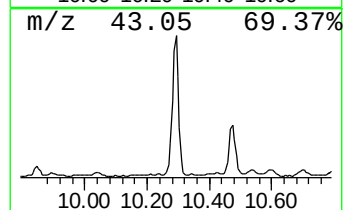
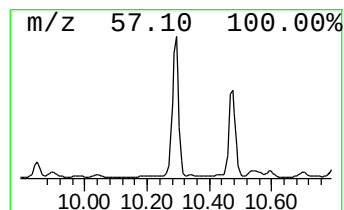
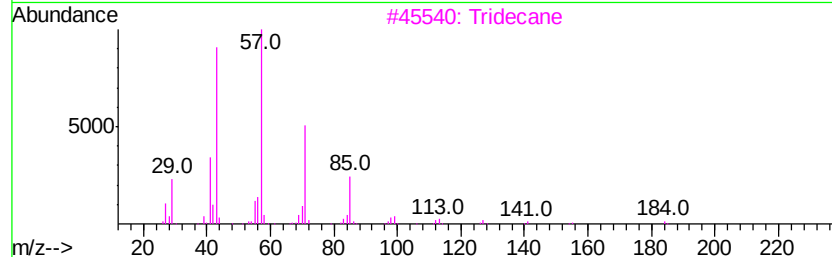
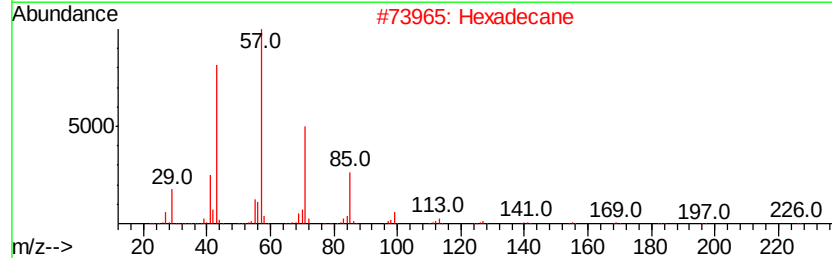
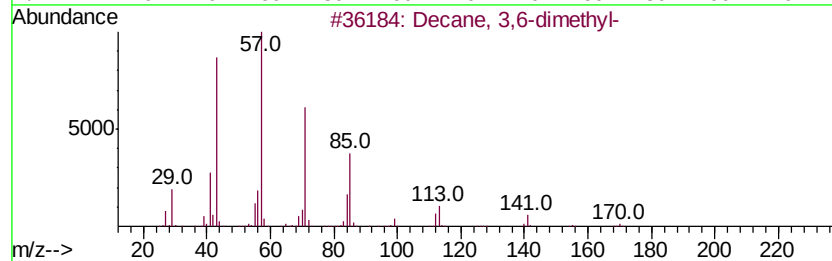
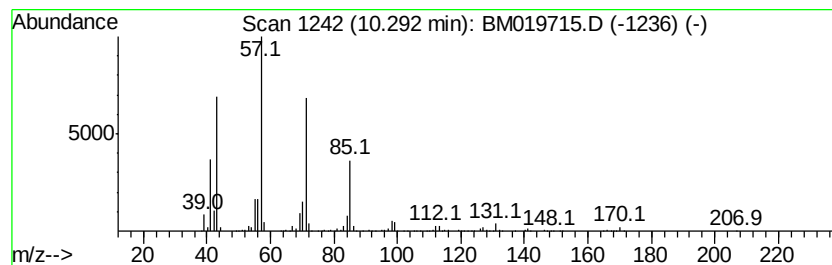
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 10.29 | 37.73 ng/ul | 2100840 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 3,6-dimethyl-   | 170 | C12H26  | 017312-53-7 | 72   |
| 2    |    | Hexadecane              | 226 | C16H34  | 000544-76-3 | 72   |
| 3    |    | Tridecane               | 184 | C13H28  | 000629-50-5 | 64   |
| 4    |    | Dodecane, 4,6-dimethyl- | 198 | C14H30  | 061141-72-8 | 64   |
| 5    |    | Undecane, 5-methyl-     | 170 | C12H26  | 001632-70-8 | 64   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

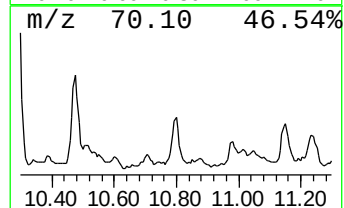
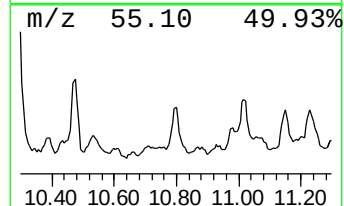
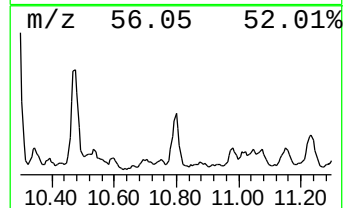
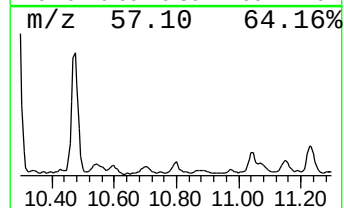
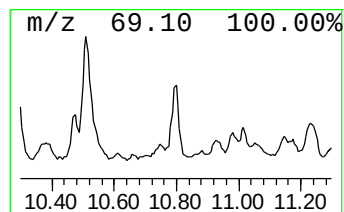
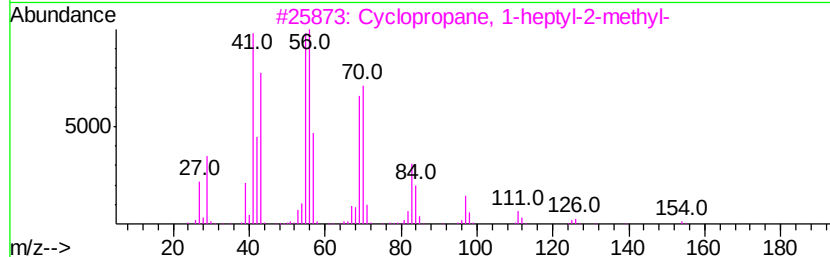
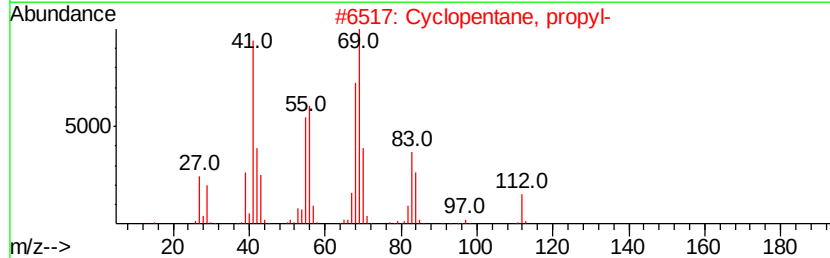
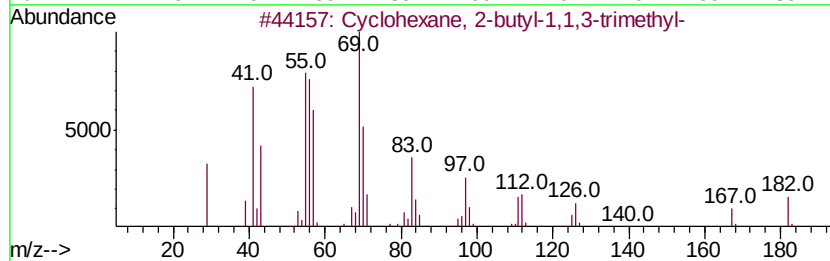
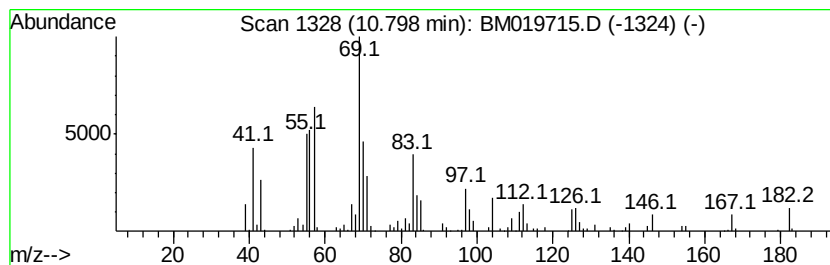
Instrument :  
BNA\_M  
ClientSampled :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Cyclic10.80 Concentration Rank 36

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 10.80   | 6.28 ng/ul | 349946                              | Napthalene-d8    | 10.35          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Cyclohexane, 2-butyl-1,1,3-trime... | 182 C13H26       | 054676-39-0 83 |
| 2       |            | Cyclopentane, propyl-               | 112 C8H16        | 002040-96-2 58 |
| 3       |            | Cyclopropane, 1-heptyl-2-methyl-    | 154 C11H22       | 074663-91-5 53 |
| 4       |            | 1.alpha.,2.beta.,3.alpha.,4.beta... | 126 C9H18        | 002532-67-4 52 |
| 5       |            | 2-Dodecene, 2-methyl-               | 182 C13H26       | 055103-82-7 49 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

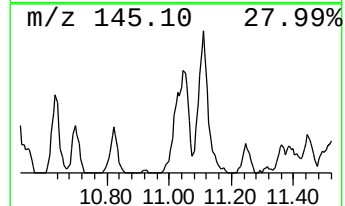
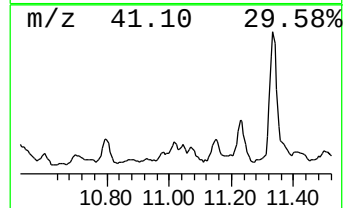
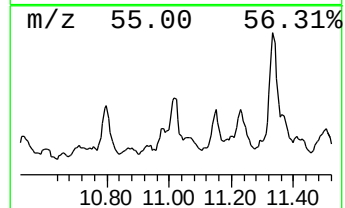
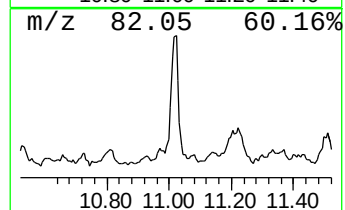
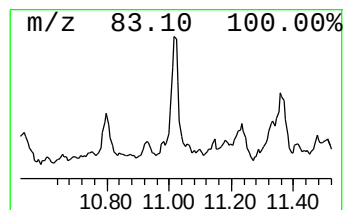
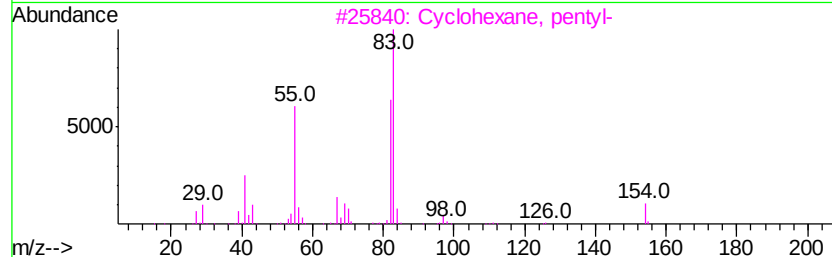
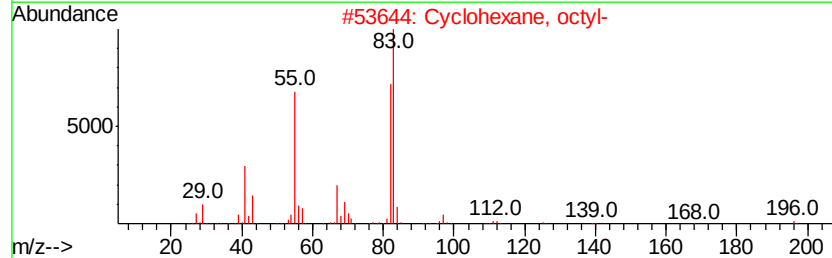
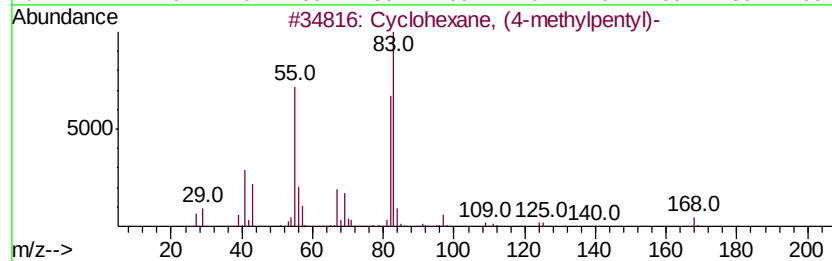
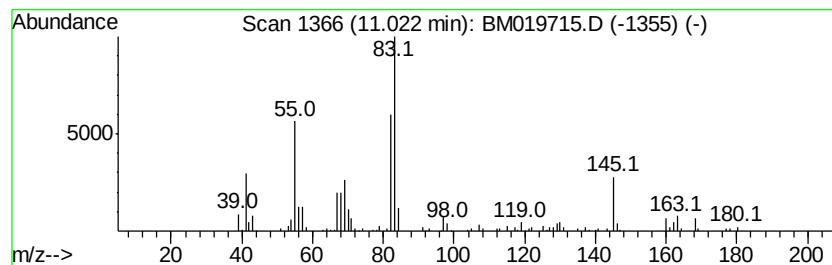
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Cyclic11.02 Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.02 | 6.60 ng/ul | 367296 | Napthalene-d8    | 10.35 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 58   |
| 2    |    |   | Cyclohexane, octyl-            | 196 | C14H28  | 001795-15-9 | 53   |
| 3    |    |   | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 50   |
| 4    |    |   | Cyclohexane, propyl-           | 126 | C9H18   | 001678-92-8 | 50   |
| 5    |    |   | Cyclohexane, (1-methylethyl)-  | 126 | C9H18   | 000696-29-7 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

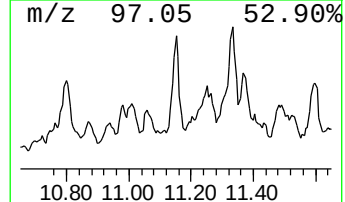
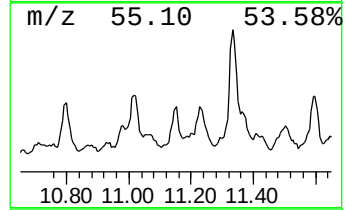
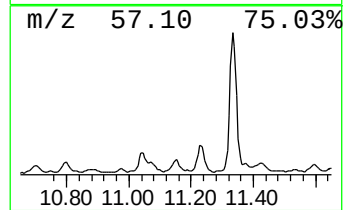
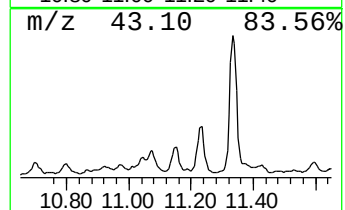
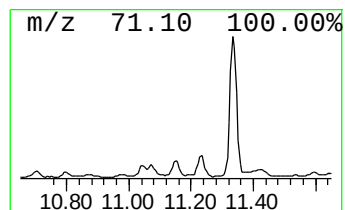
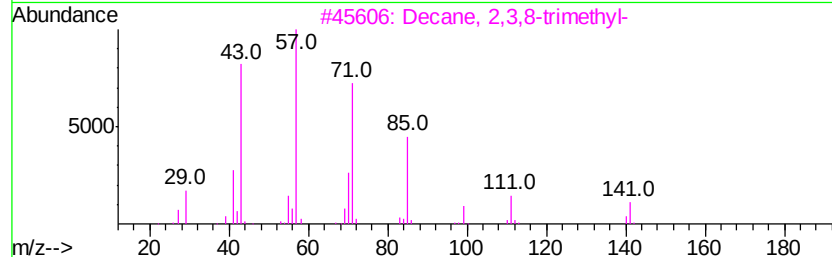
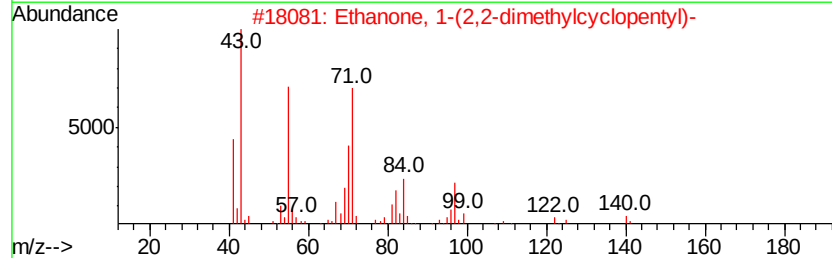
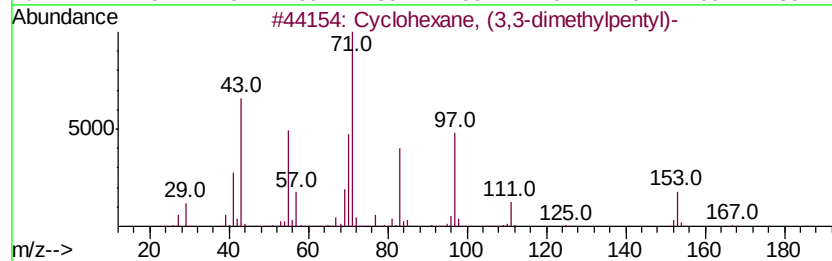
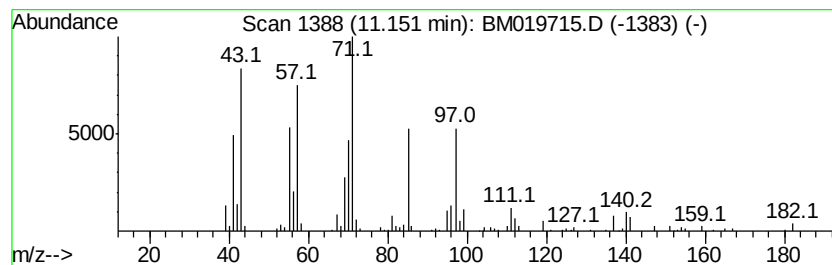
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Cyclic11.15 Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.15 | 5.02 ng/ul | 279268 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, (3,3-dimethylpentyl)-  | 182 | C13H26  | 061142-22-1 | 53   |
| 2    |    | Ethanone, 1-(2,2-dimethylcyclope... | 140 | C9H16O  | 003664-75-3 | 50   |
| 3    |    | Decane, 2,3,8-trimethyl-            | 184 | C13H28  | 062238-14-6 | 50   |
| 4    |    | Decane, 3,3,8-trimethyl-            | 184 | C13H28  | 062338-16-3 | 47   |
| 5    |    | Tritetracontane                     | 605 | C43H88  | 007098-21-7 | 47   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

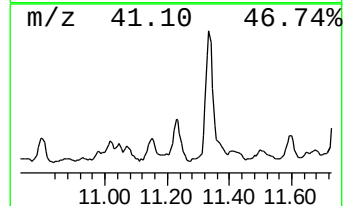
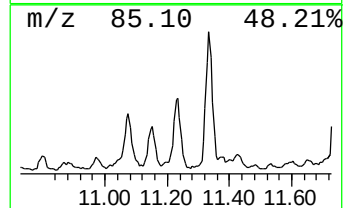
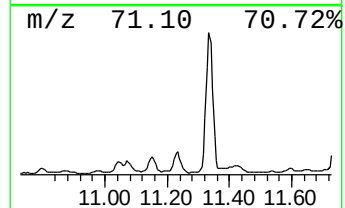
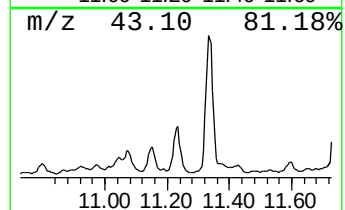
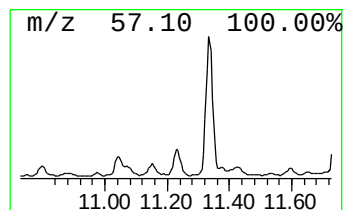
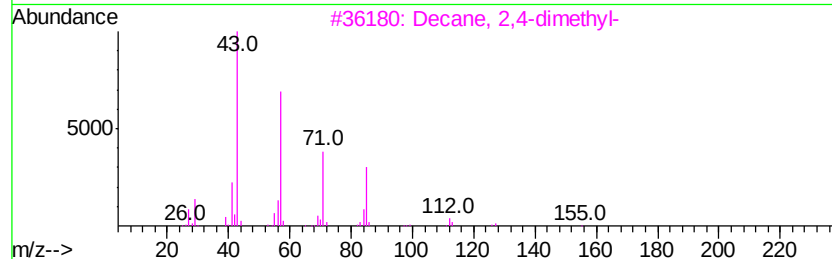
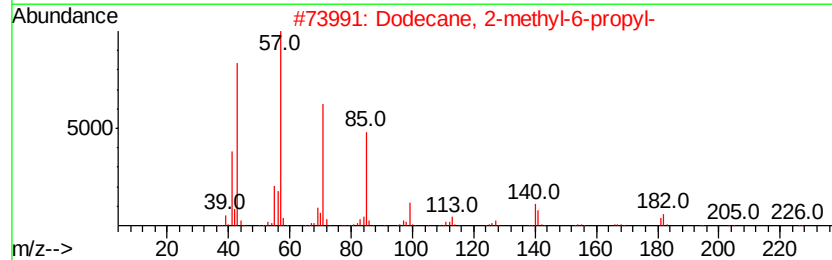
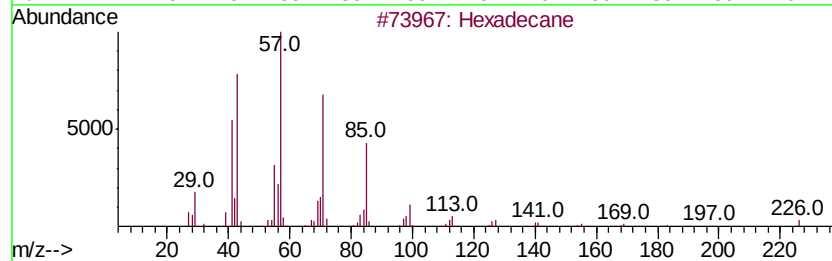
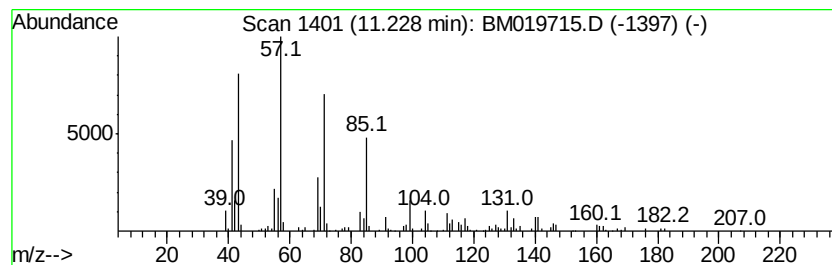
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 11.23 | 10.36 ng/ul | 577041 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane                   | 226 | C16H34  | 000544-76-3 | 74   |
| 2    |    | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 72   |
| 3    |    | Decane, 2,4-dimethyl-        | 170 | C12H26  | 002801-84-5 | 59   |
| 4    |    | Nonadecane, 9-methyl-        | 282 | C20H42  | 013287-24-6 | 53   |
| 5    |    | Heptane, 3,4-dimethyl-       | 128 | C9H20   | 000922-28-1 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

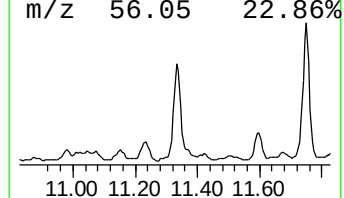
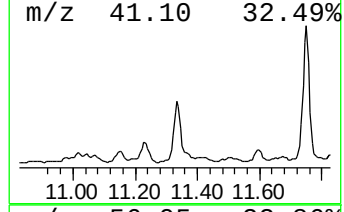
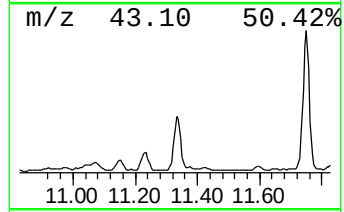
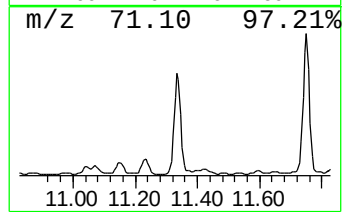
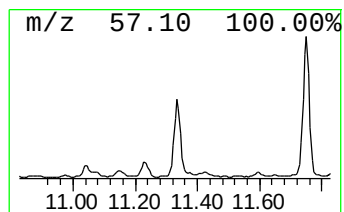
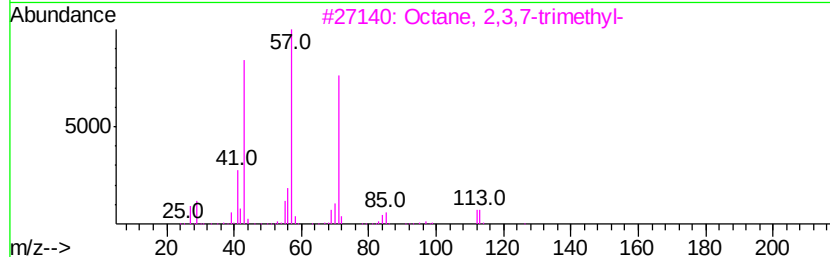
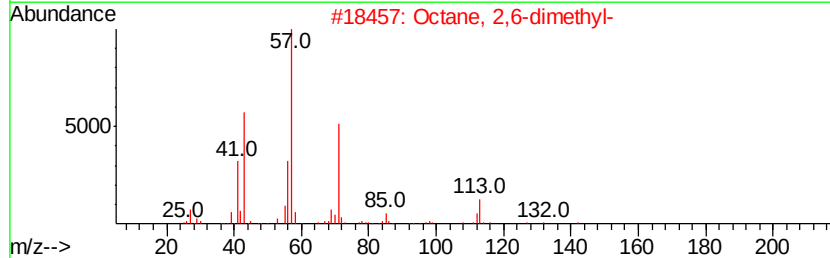
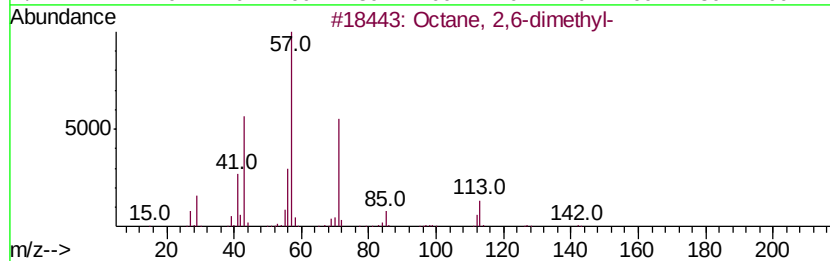
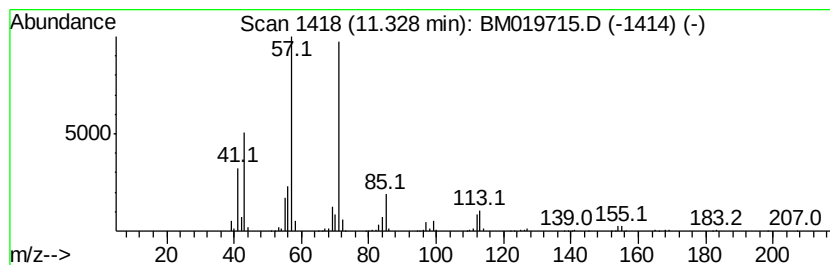
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.33 | 27.92 ng/ul | 1554410 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-    | 142 | C10H22  | 002051-30-1 | 90   |
| 2    |    | Octane, 2,6-dimethyl-    | 142 | C10H22  | 002051-30-1 | 86   |
| 3    |    | Octane, 2,3,7-trimethyl- | 156 | C11H24  | 062016-34-6 | 83   |
| 4    |    | Octane, 3,6-dimethyl-    | 142 | C10H22  | 015869-94-0 | 78   |
| 5    |    | Octane, 2-bromo-         | 192 | C8H17Br | 000557-35-7 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

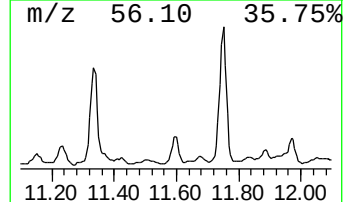
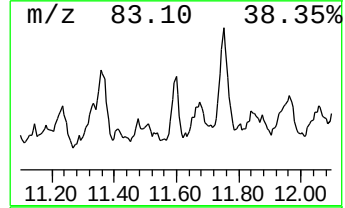
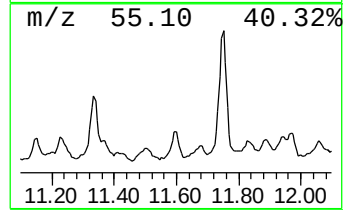
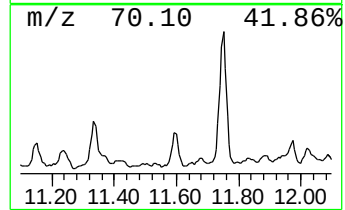
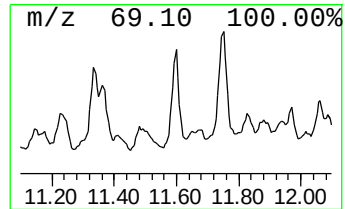
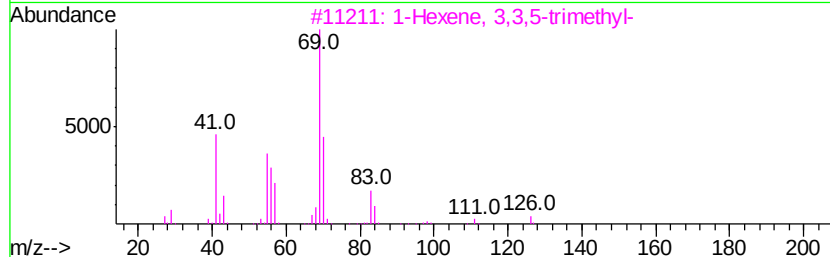
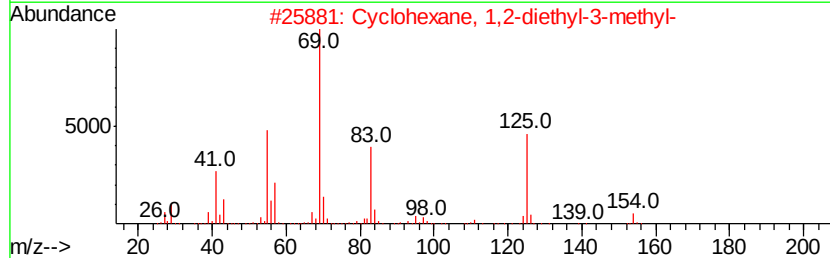
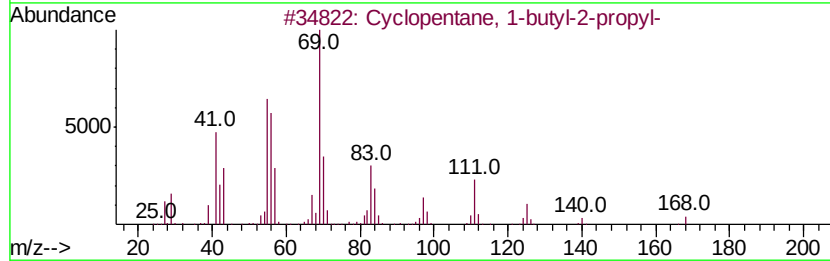
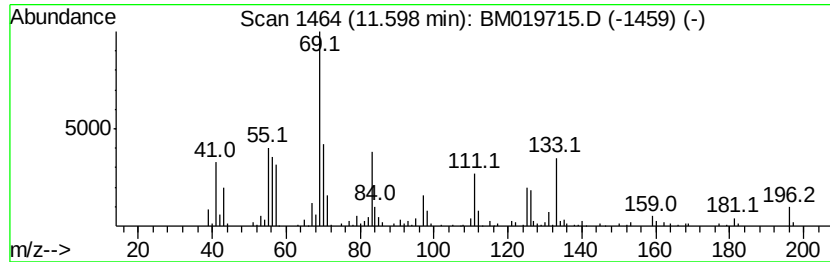
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Cyclic11.60 Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.60 | 6.26 ng/ul | 348574 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1-butyl-2-propyl-    | 168 | C12H24  | 062199-50-2 | 76   |
| 2    |    | Cyclohexane, 1,2-diethyl-3-methyl- | 154 | C11H22  | 061141-80-8 | 50   |
| 3    |    | 1-Hexene, 3,3,5-trimethyl-         | 126 | C9H18   | 013427-43-5 | 47   |
| 4    |    | 3-Heptene, 2,2,3,5,6-pentamethyl-  | 168 | C12H24  | 116164-06-8 | 47   |
| 5    |    | 1-Hexene, 3,3-dimethyl-            | 112 | C8H16   | 003404-77-1 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

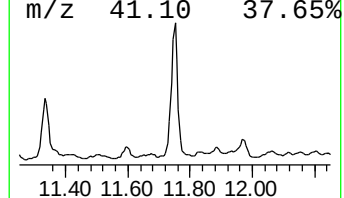
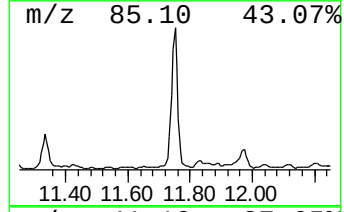
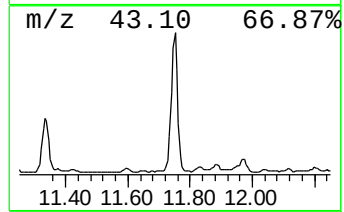
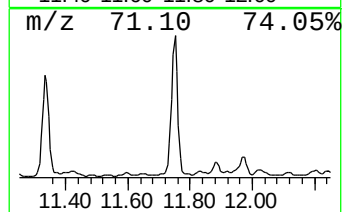
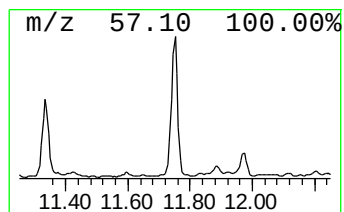
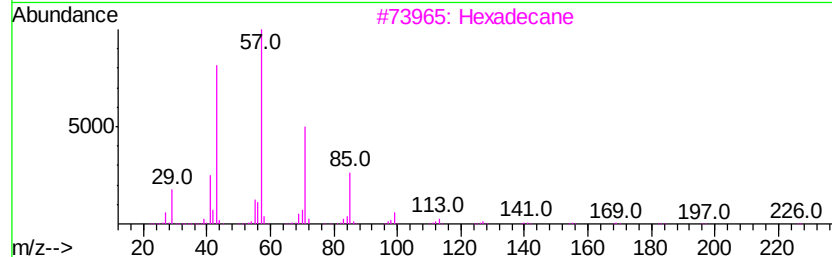
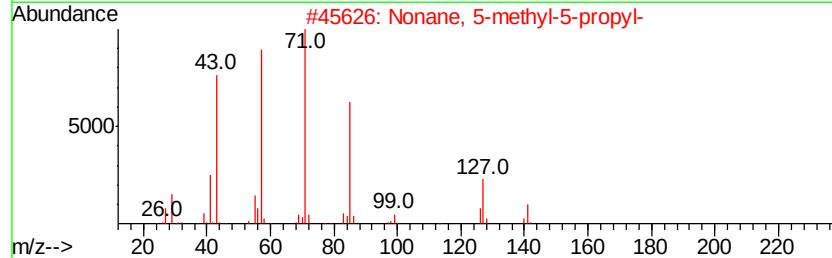
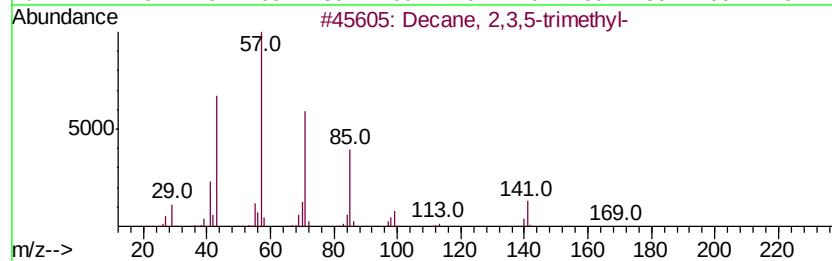
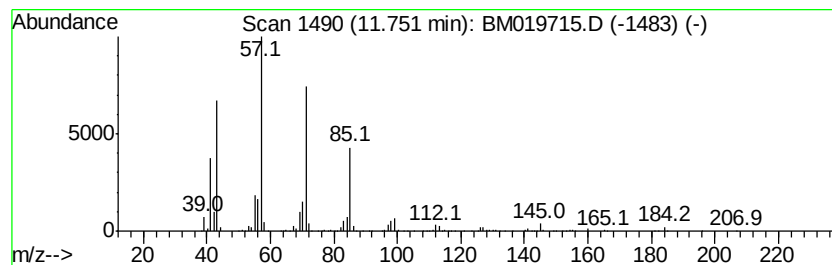
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.75 | 49.20 ng/ul | 2739840 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,3,5-trimethyl-   | 184 | C13H28  | 062238-11-3 | 83   |
| 2    |    | Nonane, 5-methyl-5-propyl- | 184 | C13H28  | 017312-75-3 | 72   |
| 3    |    | Hexadecane                 | 226 | C16H34  | 000544-76-3 | 72   |
| 4    |    | Undecane, 2-methyl-        | 170 | C12H26  | 007045-71-8 | 64   |
| 5    |    | Tridecane                  | 184 | C13H28  | 000629-50-5 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

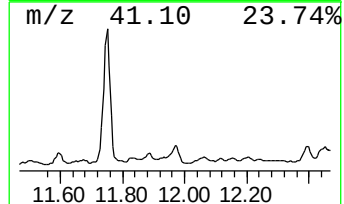
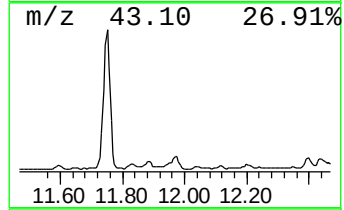
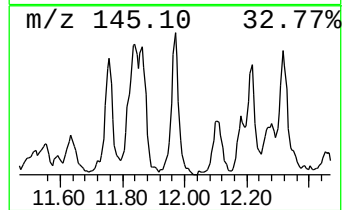
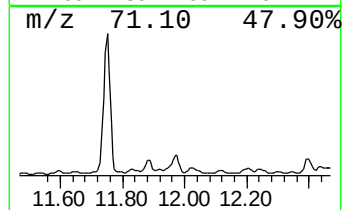
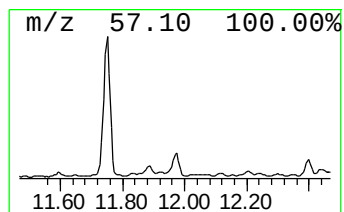
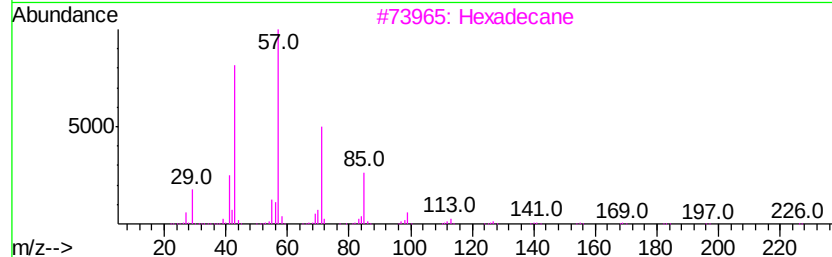
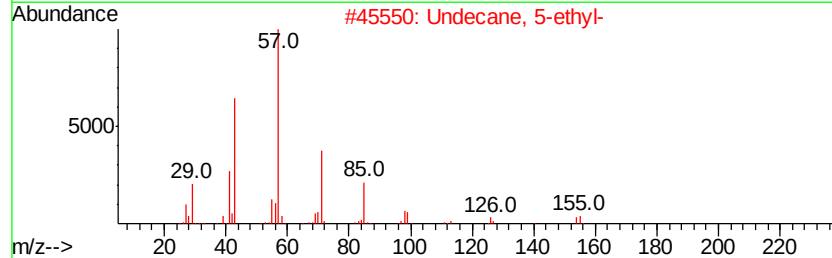
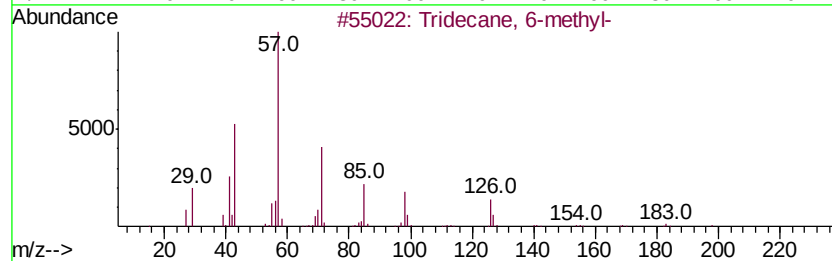
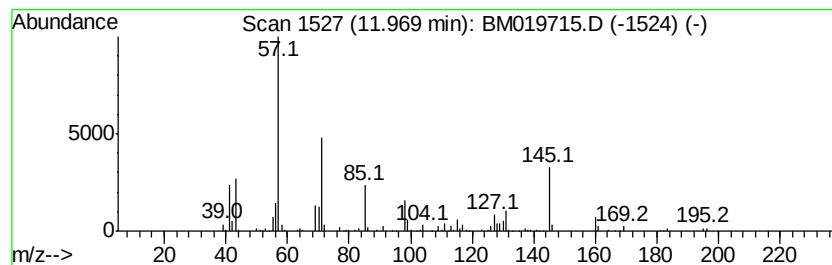
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.97 | 7.77 ng/ul | 432755 | Napthalene-d8    | 10.35 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 6-methyl- | 198 | C14H30  | 013287-21-3 | 50   |
| 2    |    | Undecane, 5-ethyl-   | 184 | C13H28  | 017453-94-0 | 50   |
| 3    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 47   |
| 4    |    | Pentacosane          | 352 | C25H52  | 000629-99-2 | 43   |
| 5    |    | Undecane             | 156 | C11H24  | 001120-21-4 | 43   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

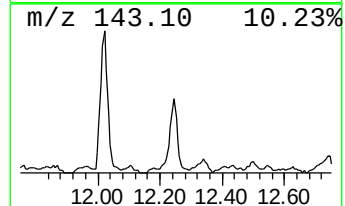
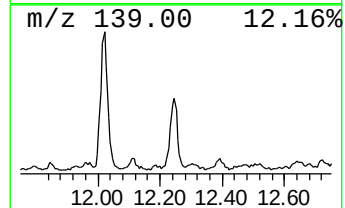
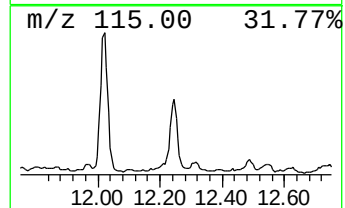
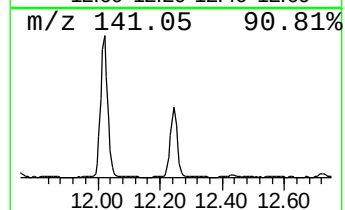
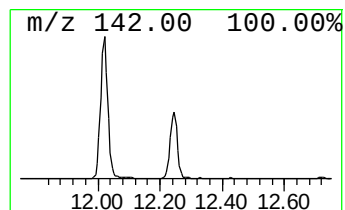
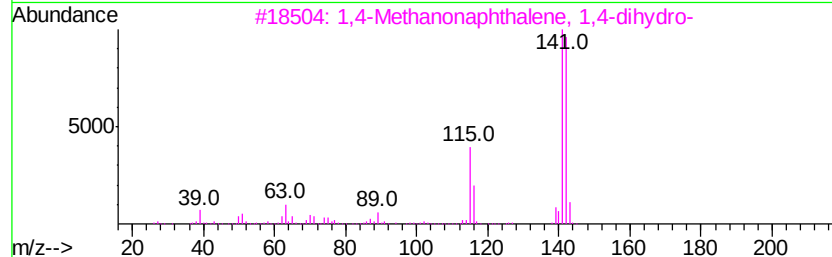
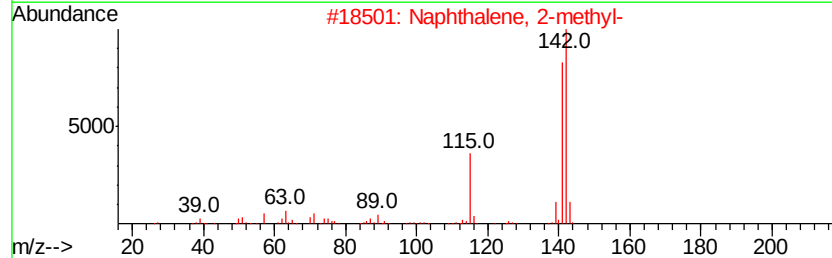
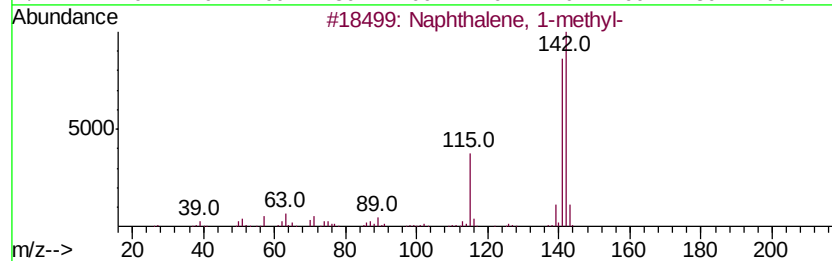
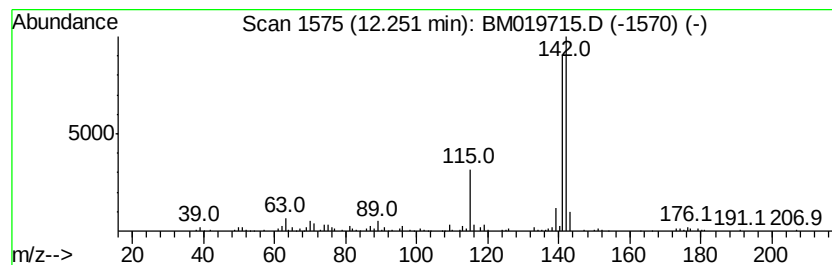
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Naphthalene, 1-methyl- Concentration Rank 23

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.25 | 8.63 ng/ul | 480429 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 97   |
| 2    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 97   |
| 3    |    | 1,4-Methanonaphthalene, 1,4-dihv... | 142 | C11H10  | 004453-90-1 | 91   |
| 4    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 91   |
| 5    |    | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

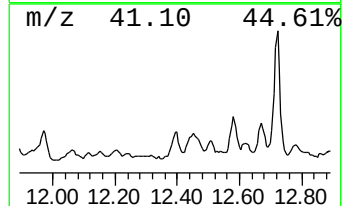
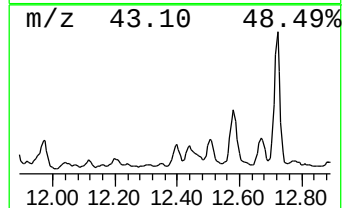
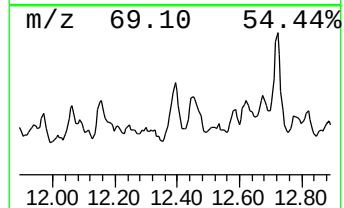
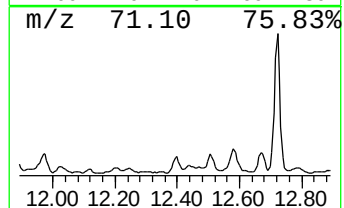
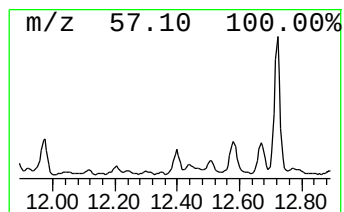
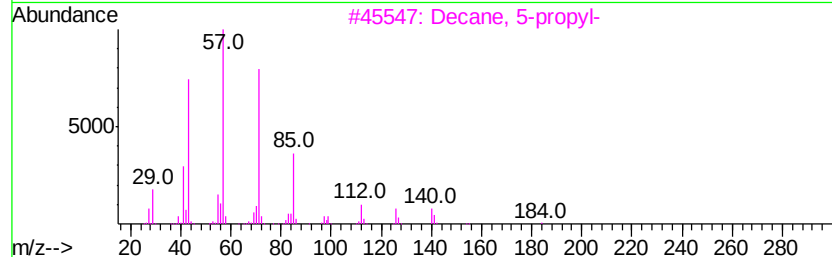
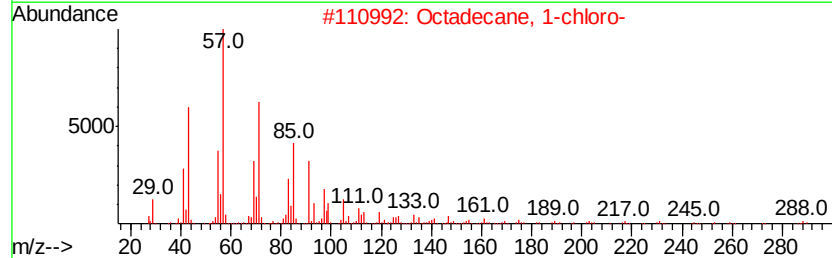
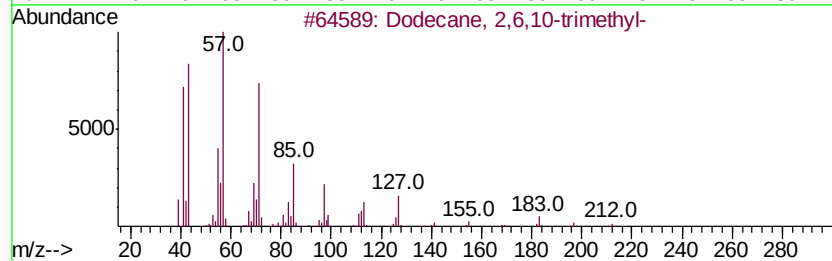
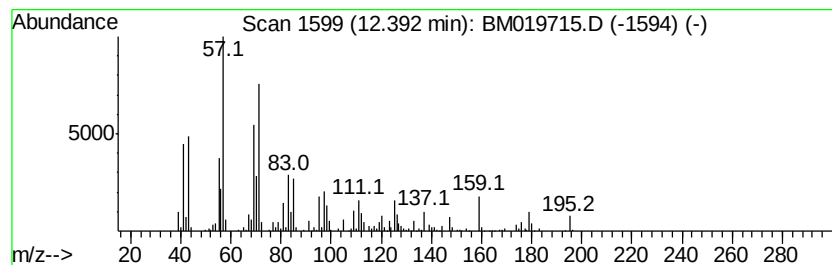
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 53

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.39 | 4.59 ng/ul | 522198 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------------------|-----|----------|-------------|------|
| 1    |    |   | Dodecane, 2,6,10-trimethyl- | 212 | C15H32   | 003891-98-3 | 53   |
| 2    |    |   | Octadecane, 1-chloro-       | 288 | C18H37Cl | 003386-33-2 | 47   |
| 3    |    |   | Decane, 5-propyl-           | 184 | C13H28   | 017312-62-8 | 47   |
| 4    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32   | 031295-56-4 | 43   |
| 5    |    |   | Tridecane, 7-methyl-        | 198 | C14H30   | 026730-14-3 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

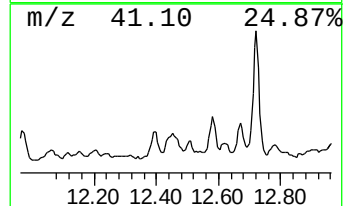
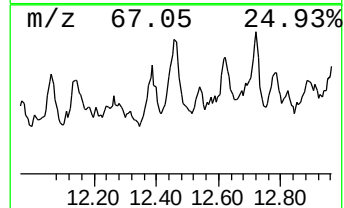
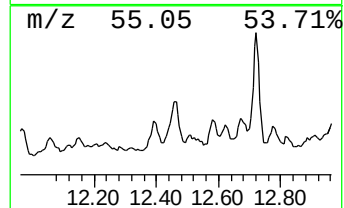
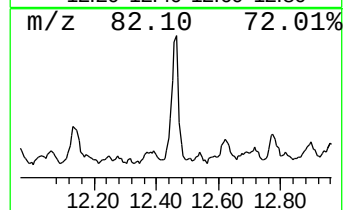
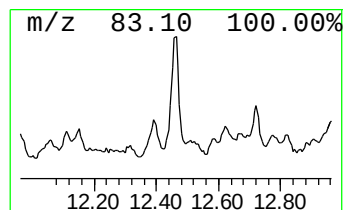
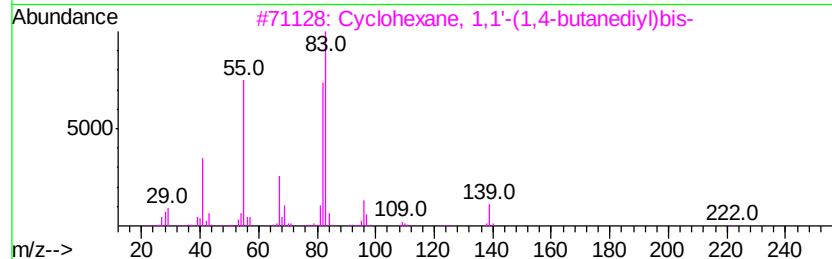
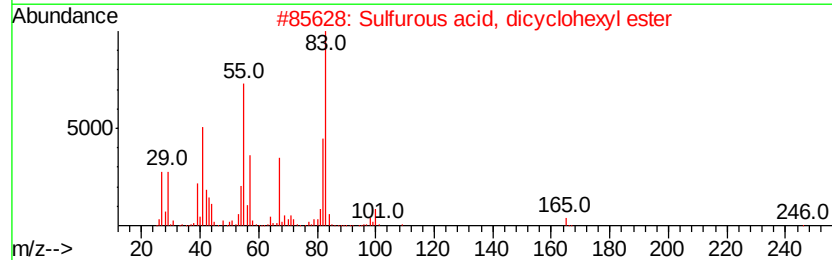
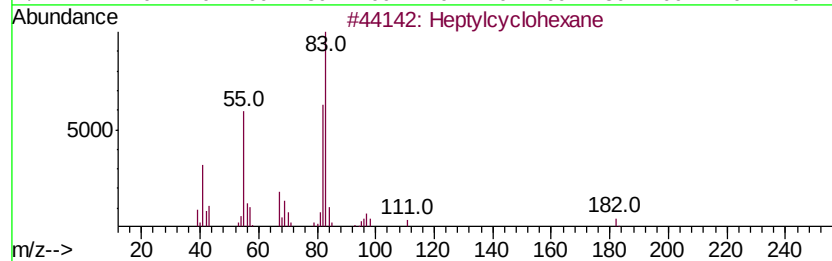
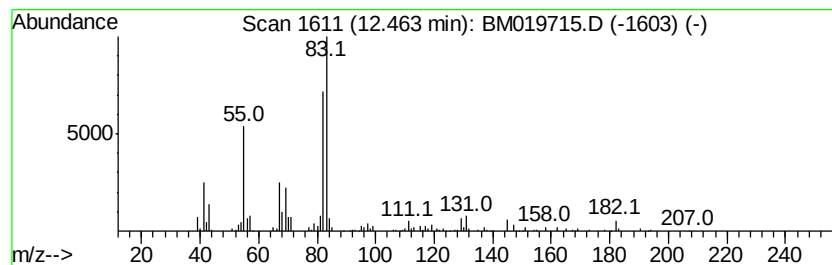
Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Cyclic12.46 Concentration Rank 51

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 12.46   | 4.66 ng/ul | 530255                              | Acenaphthene-d10 | 14.23          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Heptylcyclohexane                   | 182 C13H26       | 005617-41-4 81 |
| 2       |            | Sulfurous acid, dicyclohexyl ester  | 246 C12H22O3S    | 006214-17-1 64 |
| 3       |            | Cyclohexane, 1,1'-(1,4-butanediyl)- | 222 C16H30       | 006165-44-2 64 |
| 4       |            | Heptylcyclohexane                   | 182 C13H26       | 005617-41-4 64 |
| 5       |            | Cyclohexane, (1-methylethyl)-       | 126 C9H18        | 000696-29-7 64 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

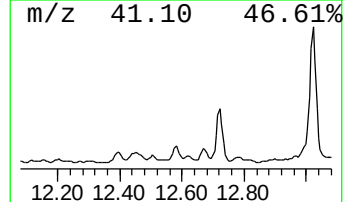
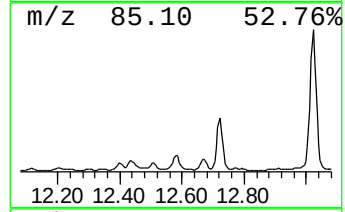
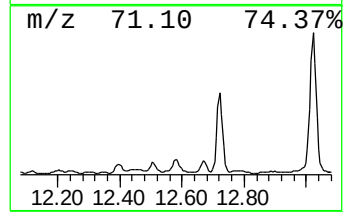
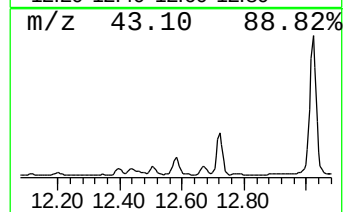
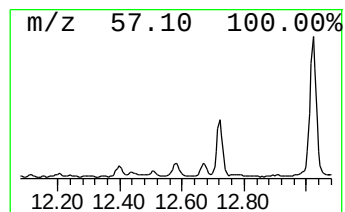
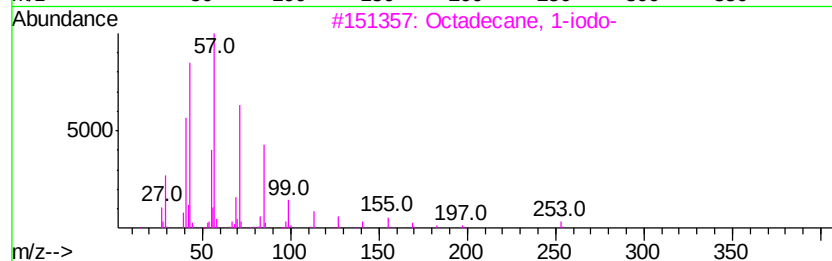
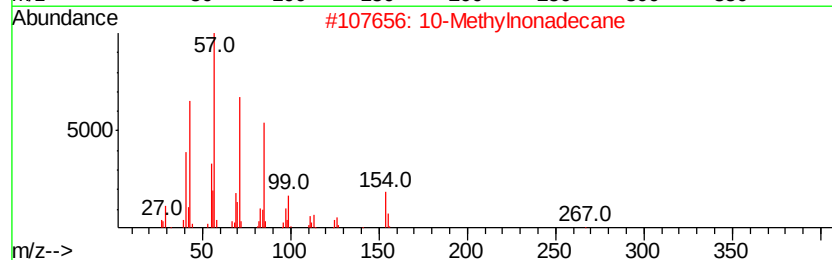
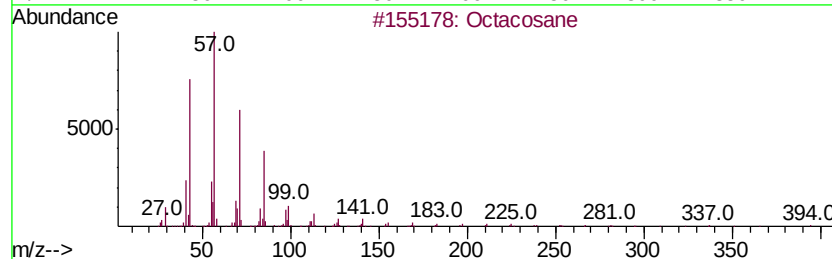
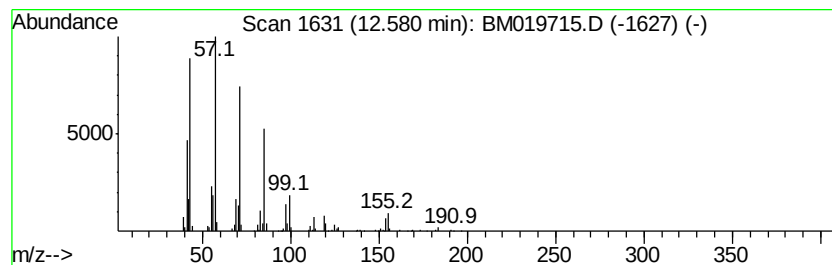
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 65

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.58 | 2.71 ng/ul | 308269 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Octacosane          | 394 | C28H58   | 000630-02-4 | 87   |
| 2    |    |   | 10-Methylnonadecane | 282 | C20H42   | 056862-62-5 | 87   |
| 3    |    |   | Octadecane, 1-iodo- | 380 | C18H37I  | 000629-93-6 | 80   |
| 4    |    |   | Tridecane, 1-iodo-  | 310 | C13H27I  | 035599-77-0 | 72   |
| 5    |    |   | 2-Bromo dodecane    | 248 | C12H25Br | 013187-99-0 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

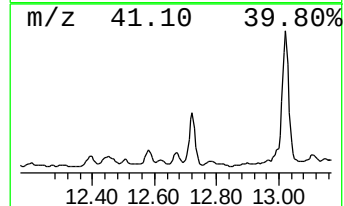
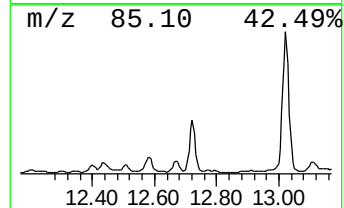
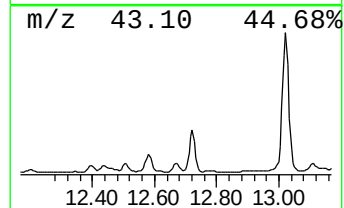
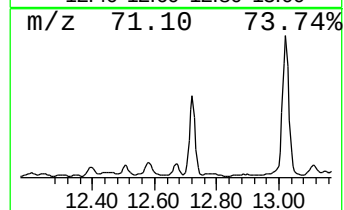
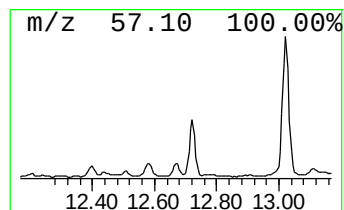
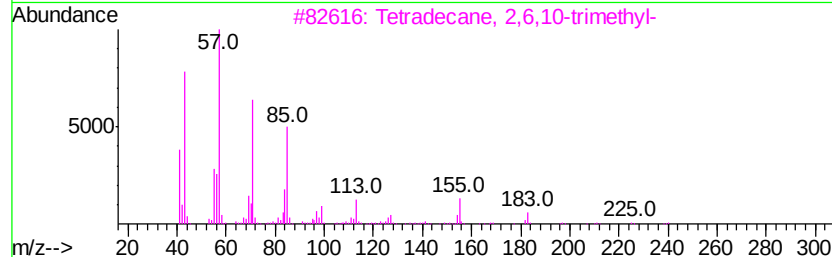
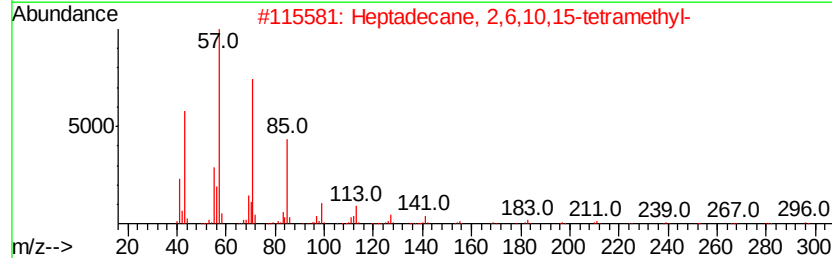
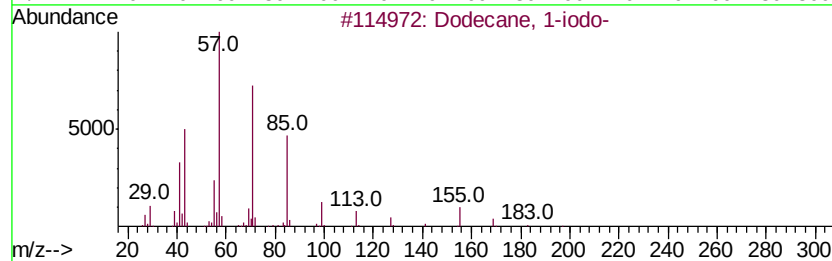
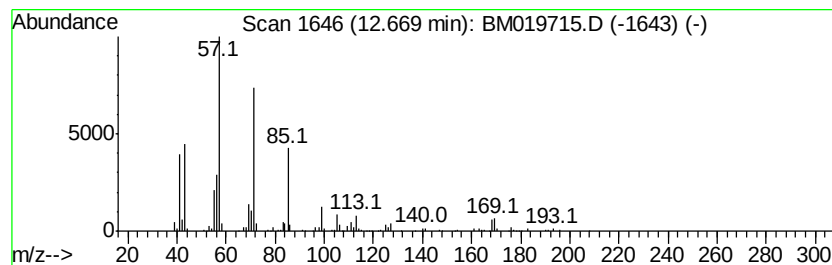
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 Dodecane, 1-iodo- Concentration Rank 63

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.67 | 2.91 ng/ul | 330916 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 1-iodo-                   | 296 | C12H25I | 004292-19-7 | 83   |
| 2    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 80   |
| 3    |    |   | Tetradecane, 2,6,10-trimethyl-      | 240 | C17H36  | 014905-56-7 | 78   |
| 4    |    |   | Decane, 3,8-dimethyl-               | 170 | C12H26  | 017312-55-9 | 78   |
| 5    |    |   | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

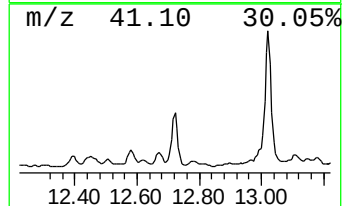
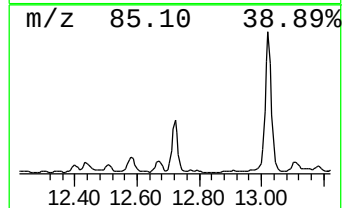
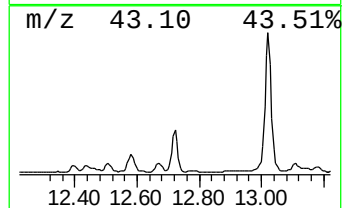
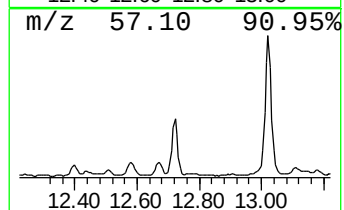
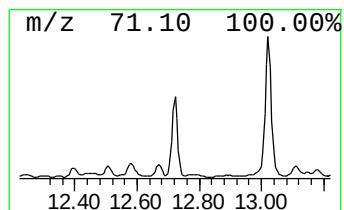
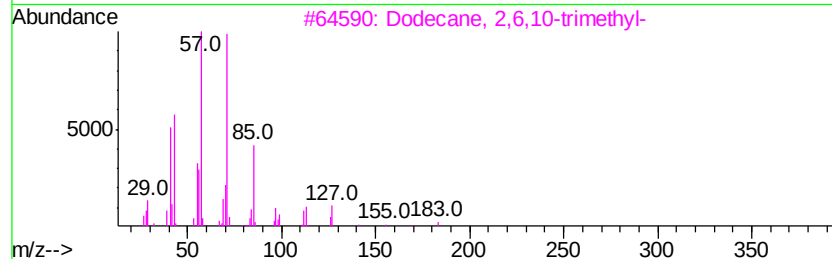
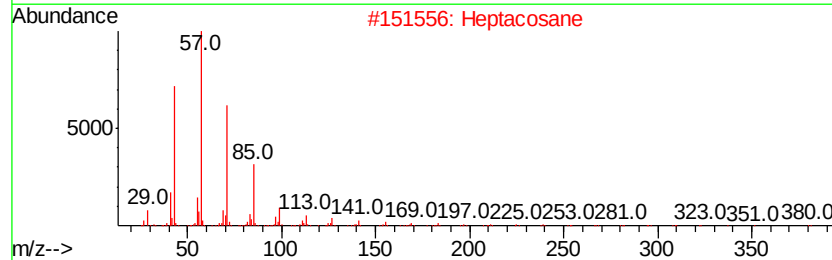
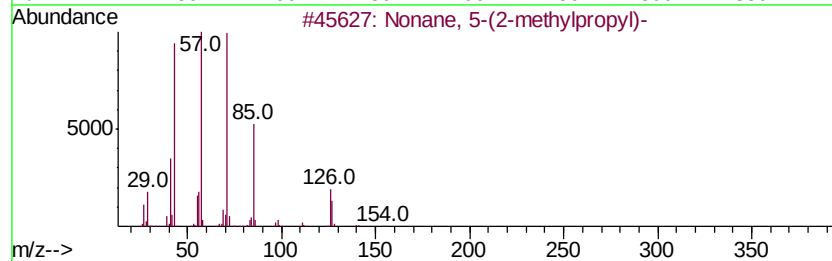
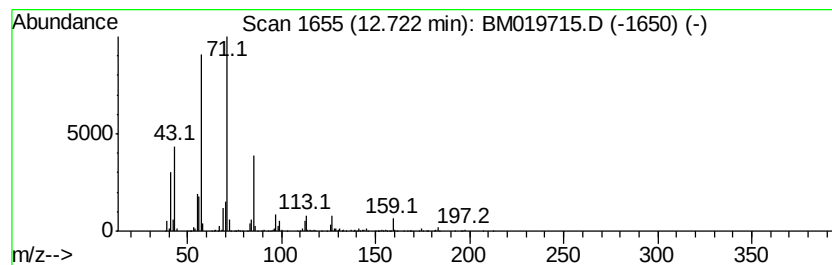
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.72 | 14.73 ng/ul | 1675200 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 72   |
| 2    |    | Heptacosane                 | 380 | C27H56  | 000593-49-7 | 64   |
| 3    |    | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 59   |
| 4    |    | Octacosane                  | 394 | C28H58  | 000630-02-4 | 53   |
| 5    |    | Decane, 5-propyl-           | 184 | C13H28  | 017312-62-8 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

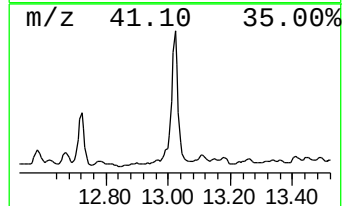
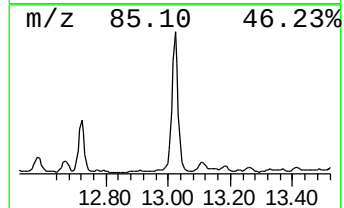
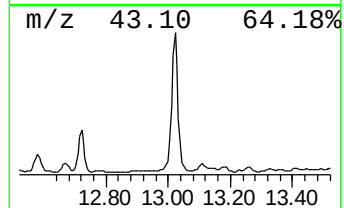
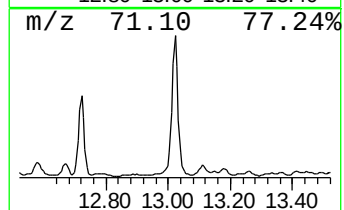
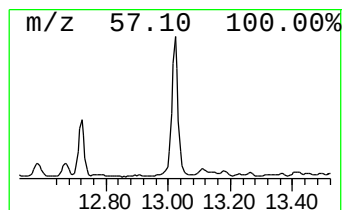
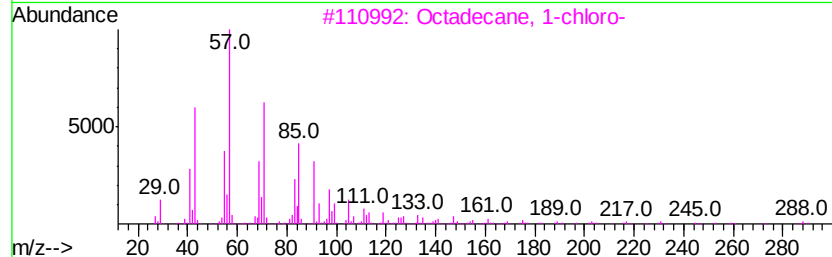
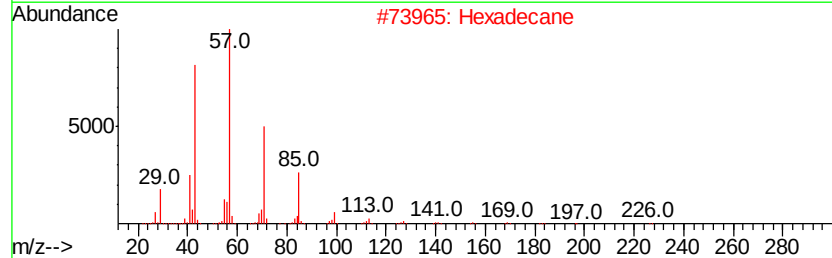
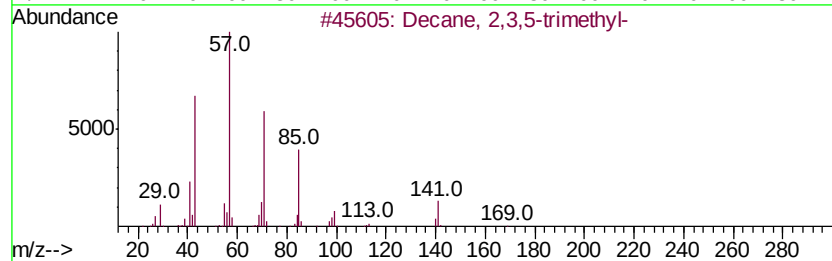
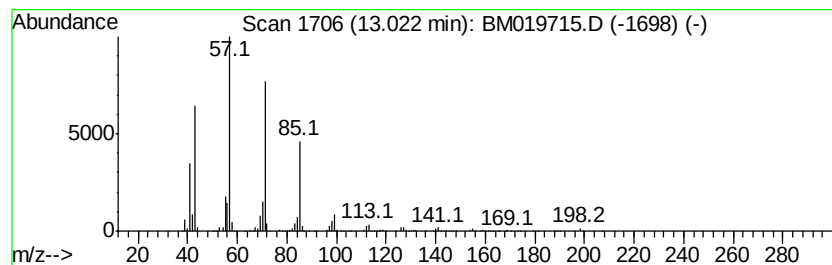
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.02 | 37.22 ng/ul | 4233160 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------|-----|----------|-------------|------|
| 1    | 5  | Decane, 2,3,5-trimethyl-    | 184 | C13H28   | 062238-11-3 | 83   |
| 2    |    | Hexadecane                  | 226 | C16H34   | 000544-76-3 | 80   |
| 3    |    | Octadecane, 1-chloro-       | 288 | C18H37Cl | 003386-33-2 | 78   |
| 4    |    | Nonane, 5-(2-methylpropyl)- | 184 | C13H28   | 062185-53-9 | 72   |
| 5    |    | Nonane, 3,7-dimethyl-       | 156 | C11H24   | 017302-32-8 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

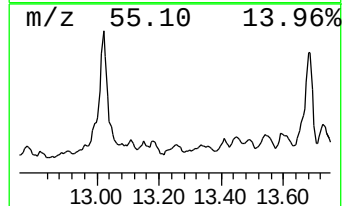
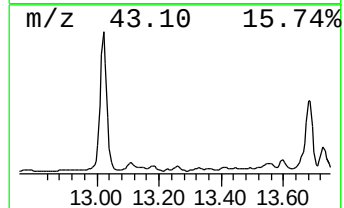
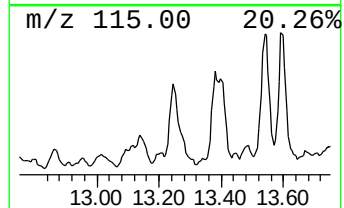
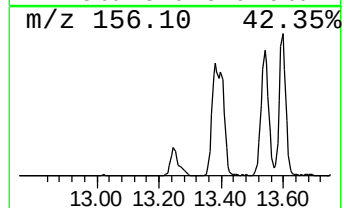
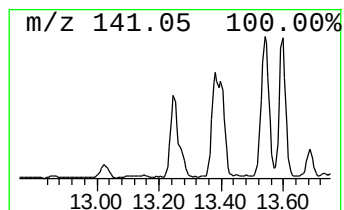
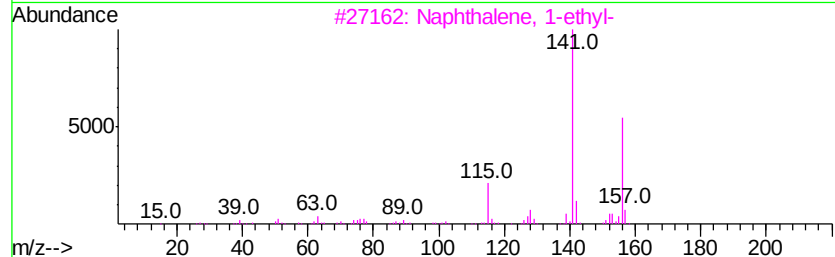
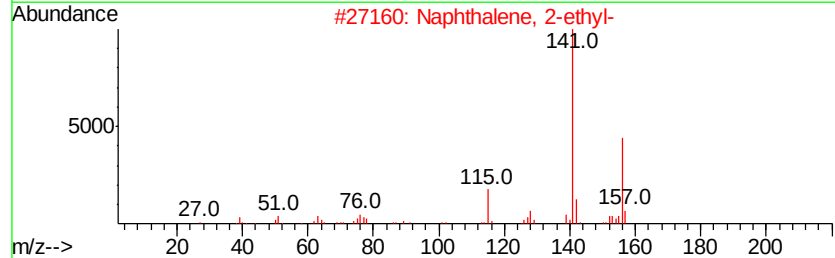
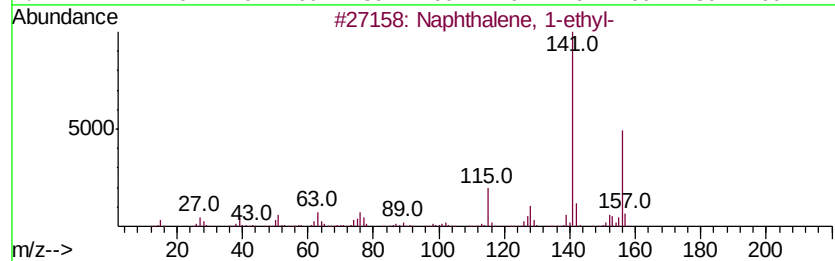
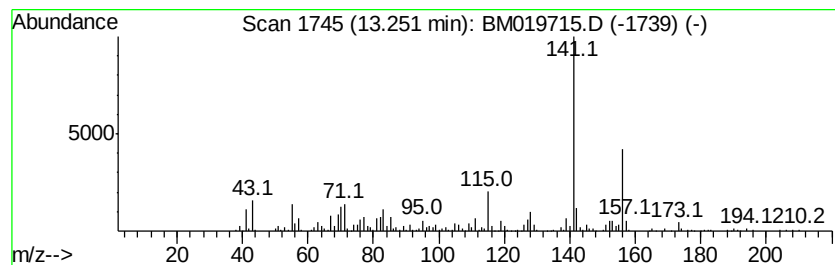
Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 27 unknown-01 Concentration Rank 34

| R.T.    | EstConc    | Area                  | Relative to ISTD | R.T.           |
|---------|------------|-----------------------|------------------|----------------|
| 13.25   | 6.41 ng/ul | 729088                | Acenaphthene-d10 | 14.23          |
| Hit# of | 5          | Tentative ID          | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 1-ethyl- | 156 C12H12       | 001127-76-0 95 |
| 2       |            | Naphthalene, 2-ethyl- | 156 C12H12       | 000939-27-5 95 |
| 3       |            | Naphthalene, 1-ethyl- | 156 C12H12       | 001127-76-0 95 |
| 4       |            | Naphthalene, 2-ethyl- | 156 C12H12       | 000939-27-5 93 |
| 5       |            | Naphthalene, 1-ethyl- | 156 C12H12       | 001127-76-0 93 |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

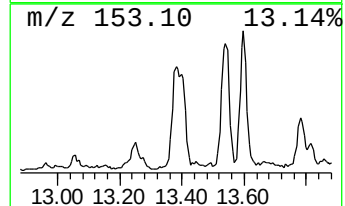
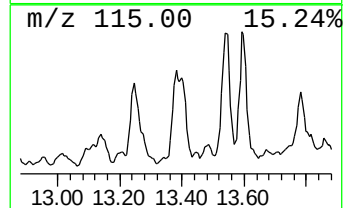
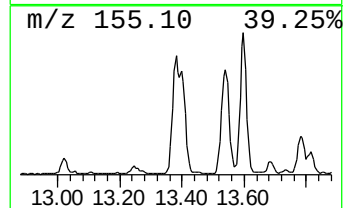
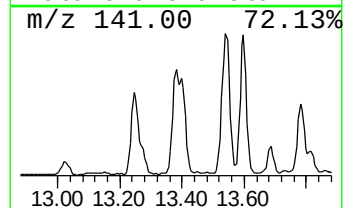
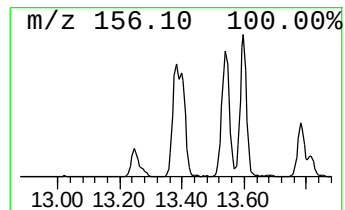
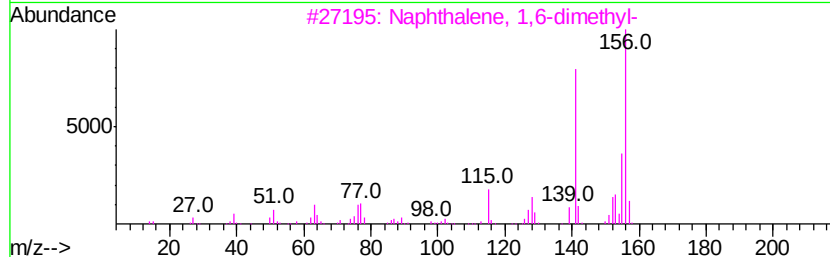
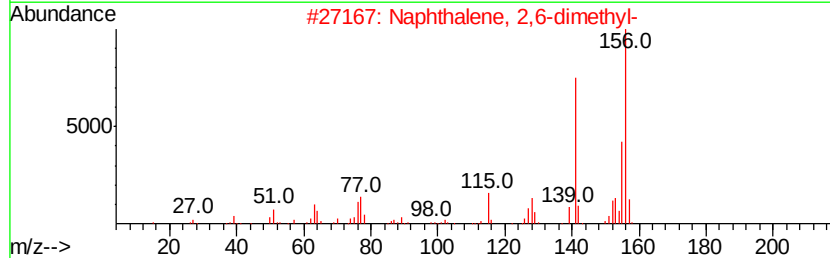
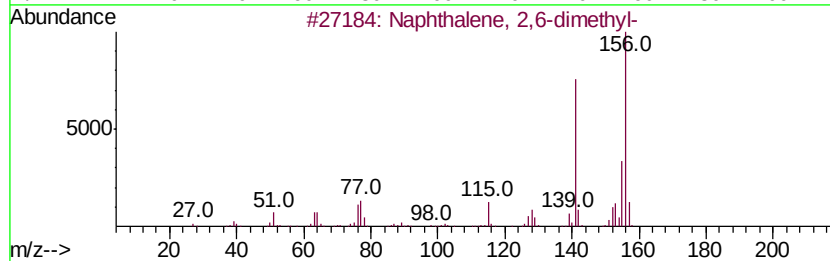
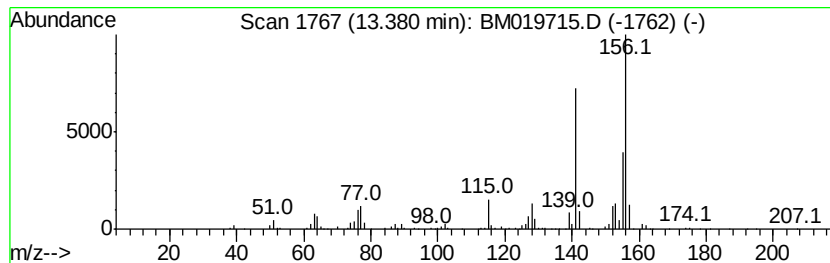
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 Naphthalene, 2,6-dimethyl- Concentration Rank 46

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.38 | 5.09 ng/ul | 578559 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 2    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 3    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 4    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 98   |
| 5    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

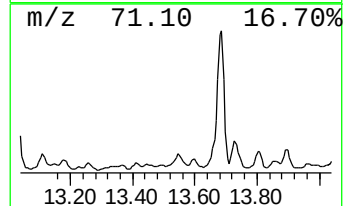
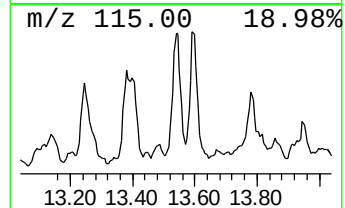
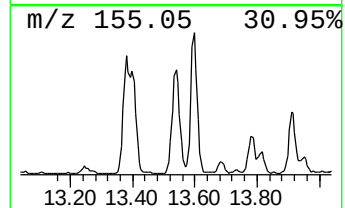
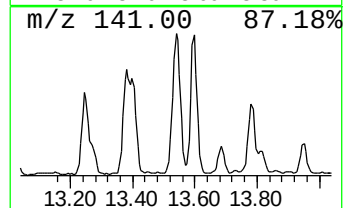
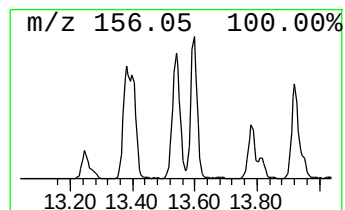
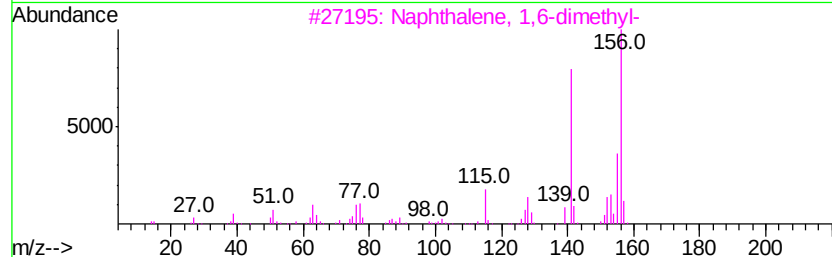
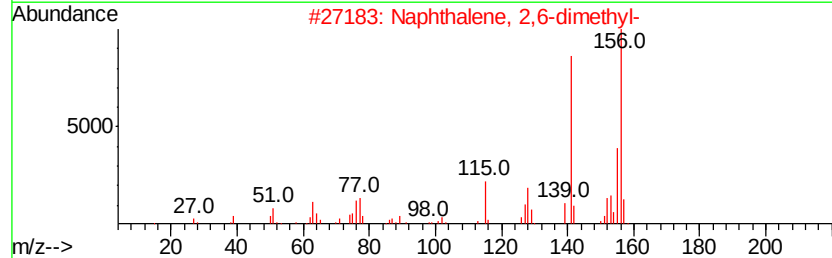
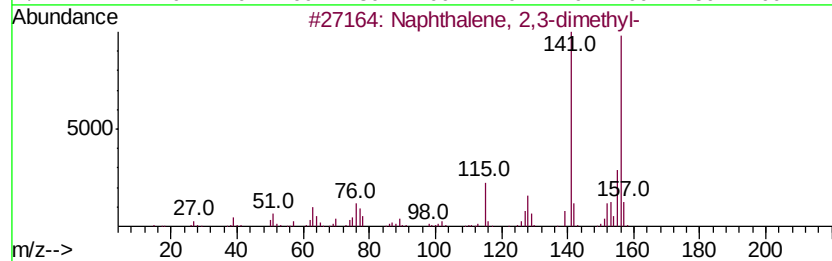
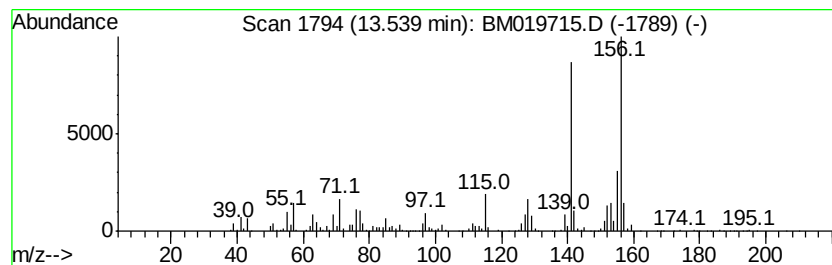
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Naphthalene, 2,3-dimethyl- Concentration Rank 13

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.54 | 12.24 ng/ul | 1392560 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 98   |
| 2    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 4    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 98   |
| 5    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

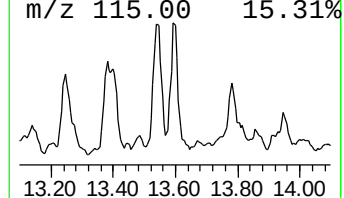
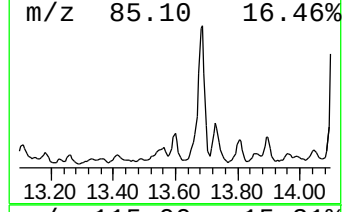
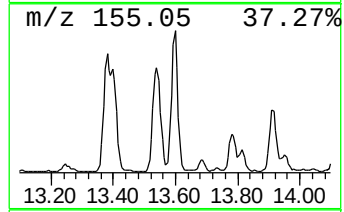
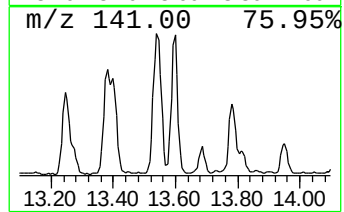
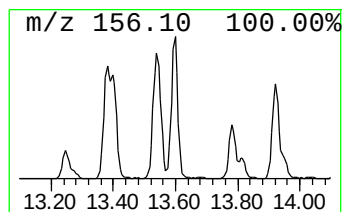
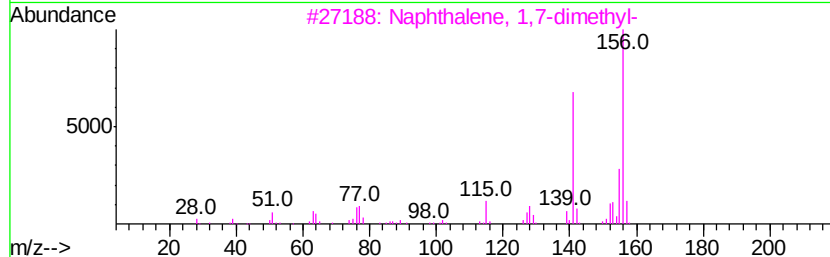
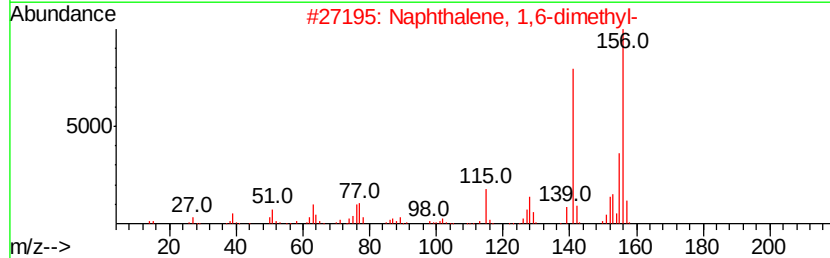
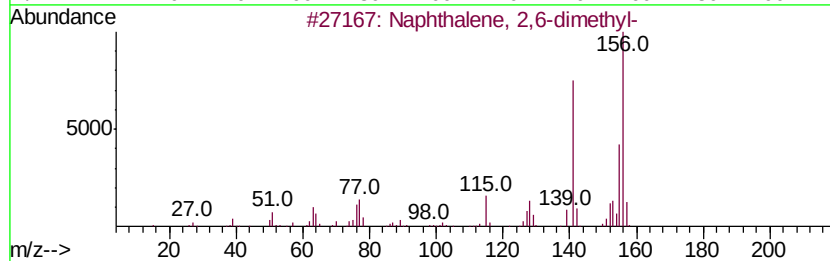
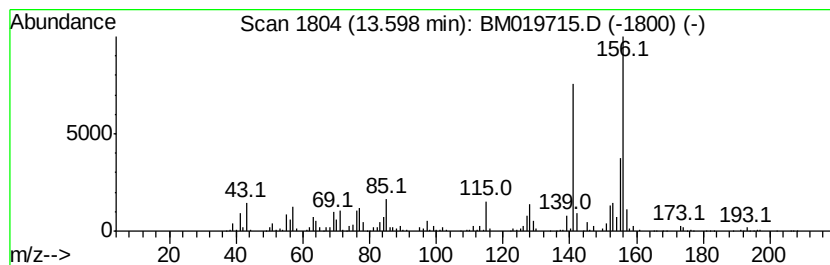
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 Naphthalene, 1,6-dimethyl- Concentration Rank 16

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 13.60 | 10.98 ng/ul | 1248850 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 2    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 3    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |    | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 96   |
| 5    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

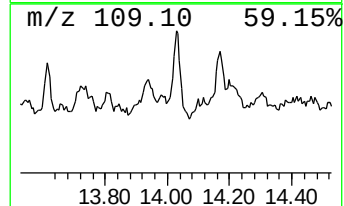
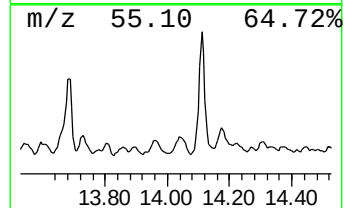
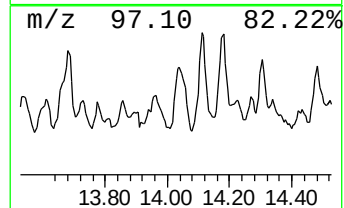
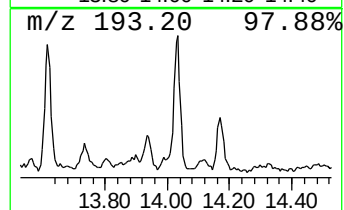
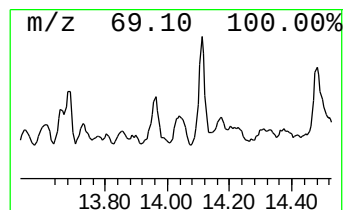
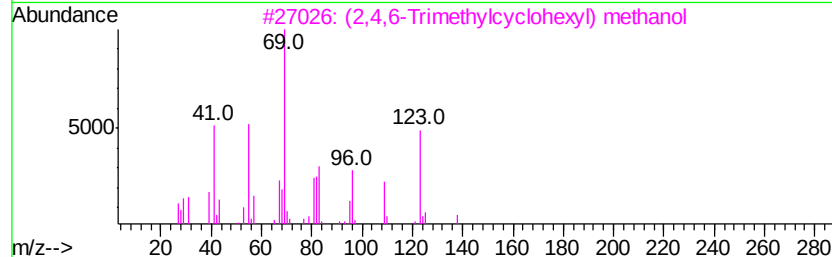
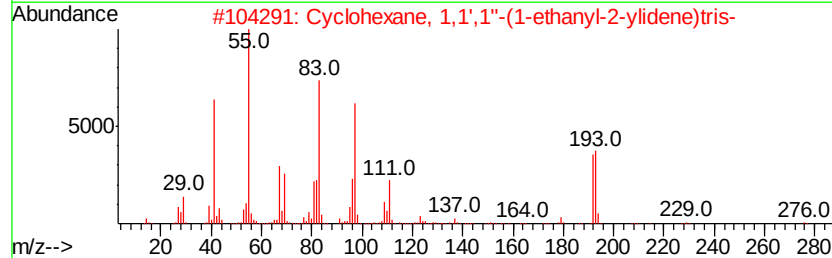
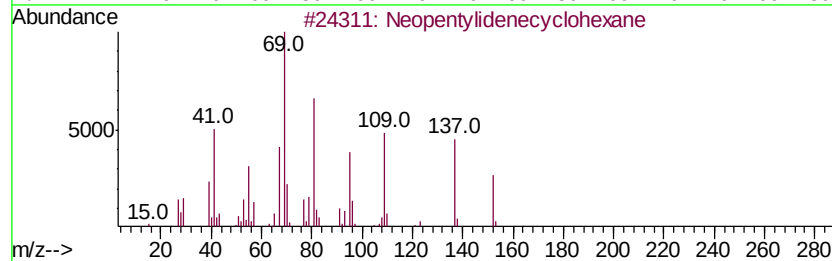
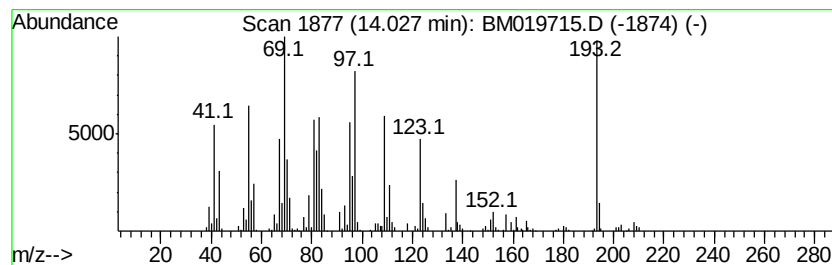
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 unknown-02 Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.03 | 6.05 ng/ul | 687488 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Neopentylidenecyclohexane           | 152 | C11H20  | 039546-80-0 | 38   |
| 2    |    | Cyclohexane, 1,1',1''-(1-ethanyl... | 276 | C20H36  | 055682-86-5 | 35   |
| 3    |    | (2,4,6-Trimethylcyclohexyl) meth... | 156 | C10H20O | 013702-56-2 | 27   |
| 4    |    | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28  | 080655-44-3 | 27   |
| 5    |    | 1-Hexyne, 3-ethoxy-3,4-dimethyl-    | 154 | C10H18O | 054244-91-6 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

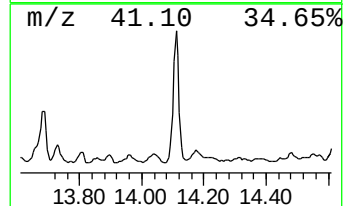
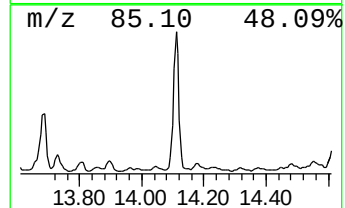
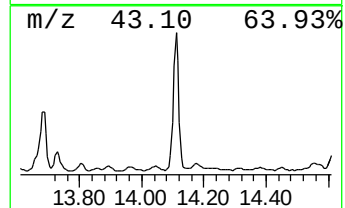
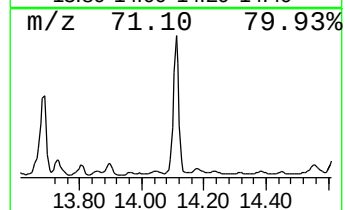
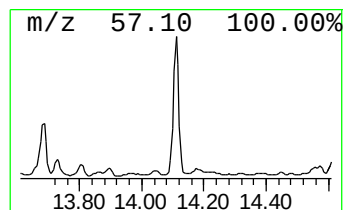
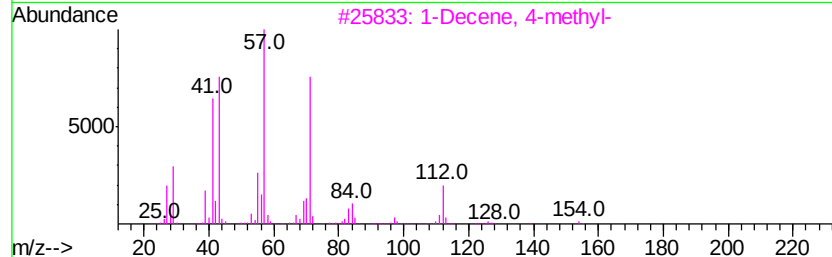
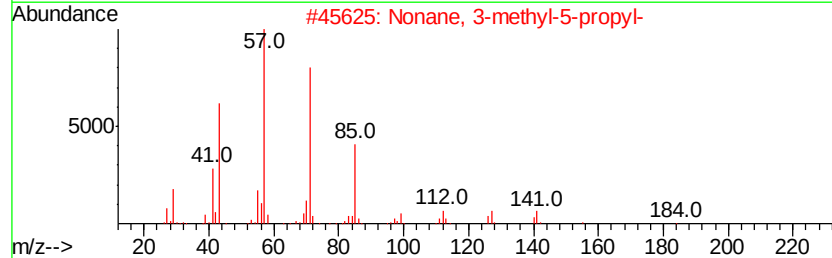
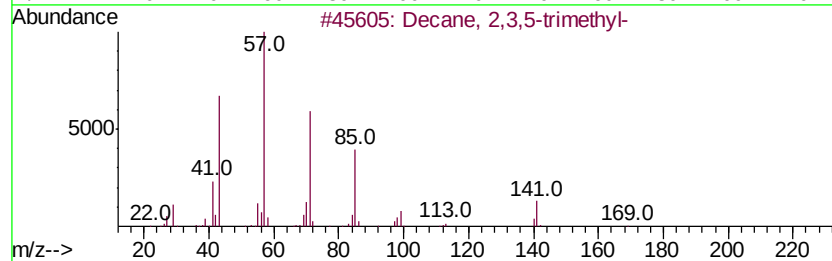
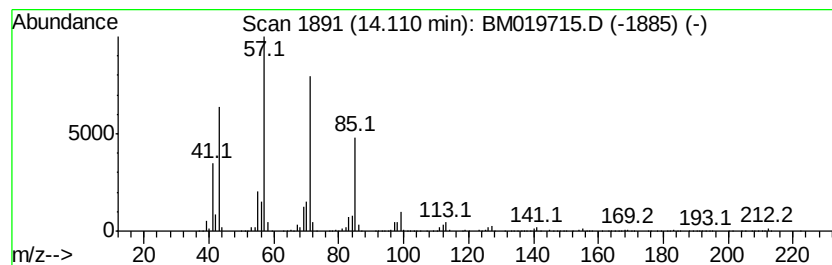
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.11 | 38.71 ng/ul | 4402870 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,3,5-trimethyl-   | 184 | C13H28  | 062238-11-3 | 90   |
| 2    |    | Nonane, 3-methyl-5-propyl- | 184 | C13H28  | 031081-18-2 | 83   |
| 3    |    | 1-Decene, 4-methyl-        | 154 | C11H22  | 013151-29-6 | 64   |
| 4    |    | Nonane, 3,7-dimethyl-      | 156 | C11H24  | 017302-32-8 | 64   |
| 5    |    | Nonadecane                 | 268 | C19H40  | 000629-92-5 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

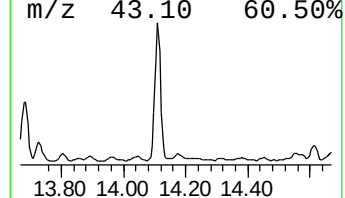
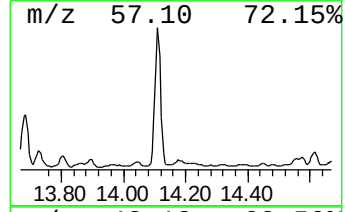
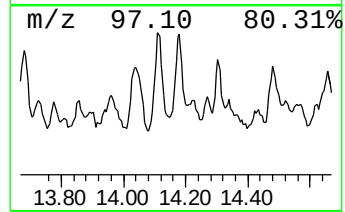
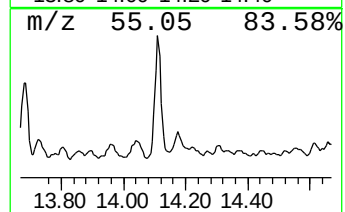
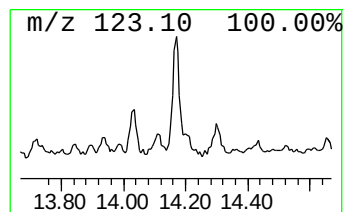
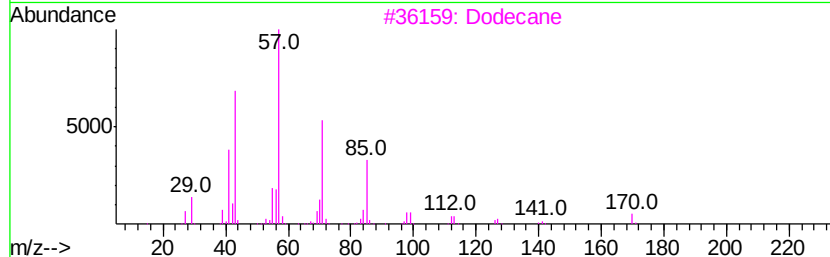
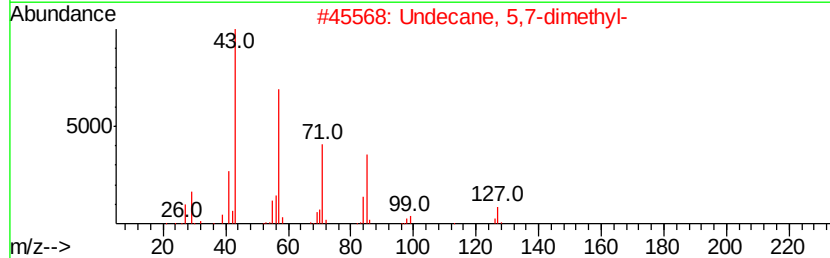
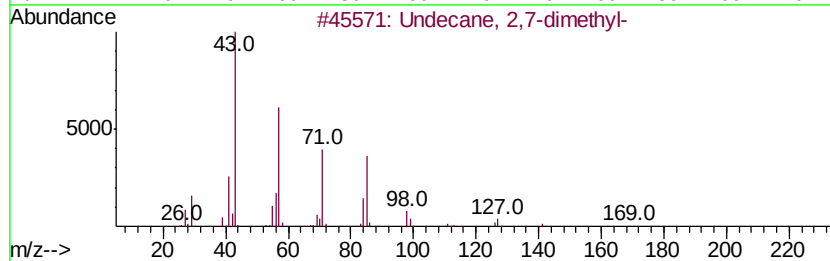
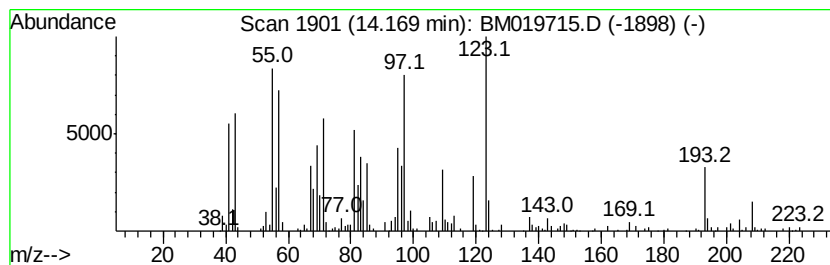
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.17 | 5.08 ng/ul | 577910 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Undecane, 2,7-dimethyl-  | 184 | C13H28  | 017301-24-5 | 25   |
| 2    |    |   | Undecane, 5,7-dimethyl-  | 184 | C13H28  | 017312-83-3 | 25   |
| 3    |    |   | Dodecane                 | 170 | C12H26  | 000112-40-3 | 20   |
| 4    |    |   | 5-Octen-4-one, 7-methyl- | 140 | C9H16O  | 032064-78-1 | 20   |
| 5    |    |   | Decane, 3-methyl-        | 156 | C11H24  | 013151-34-3 | 18   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

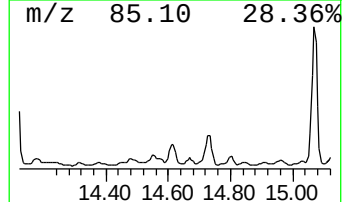
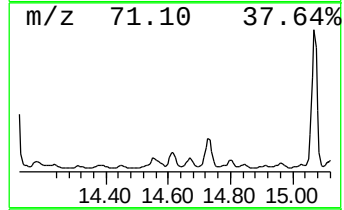
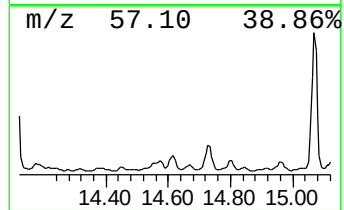
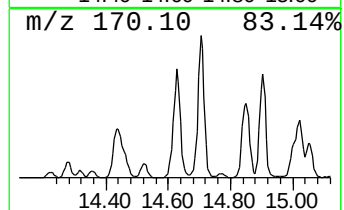
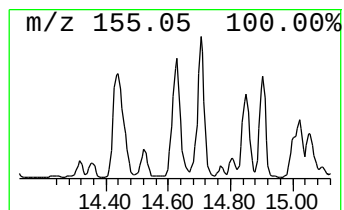
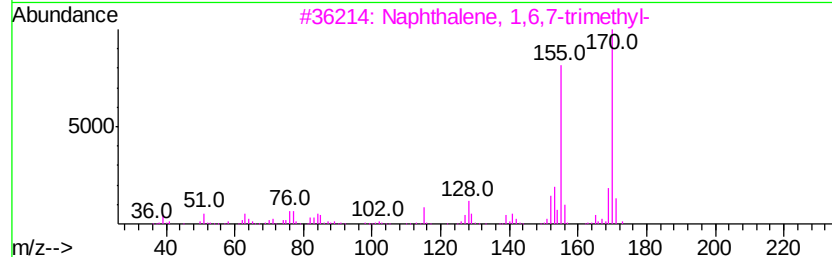
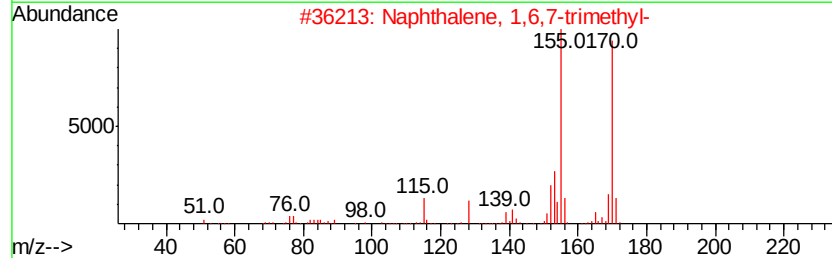
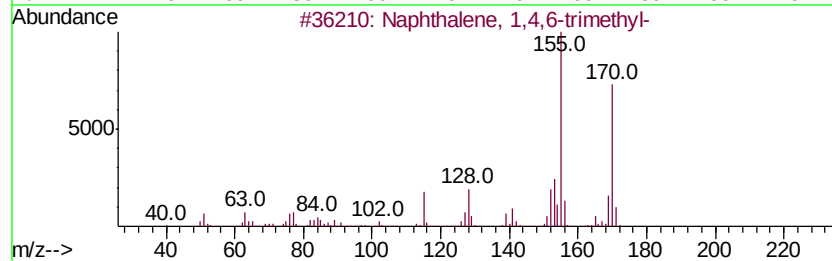
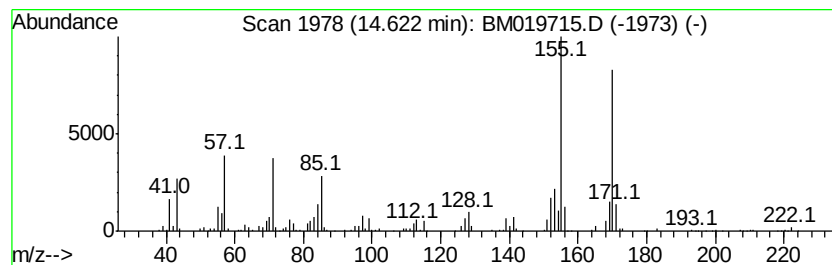
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 Naphthalene, 1,4,6-trimethyl- Concentration Rank 26

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.62 | 8.05 ng/ul | 915618 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 89   |
| 2    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 89   |
| 3    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 78   |
| 4    |    | Azulene, 4,6,8-trimethyl-     | 170 | C13H14  | 000941-81-1 | 70   |
| 5    |    | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

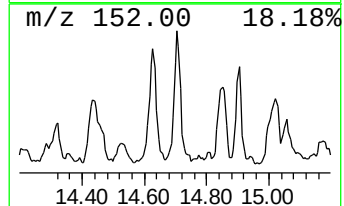
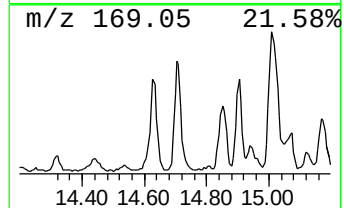
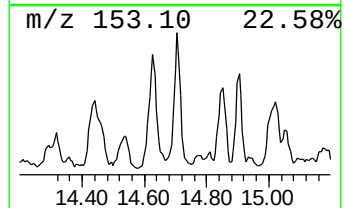
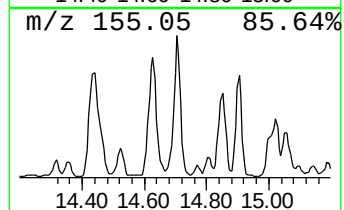
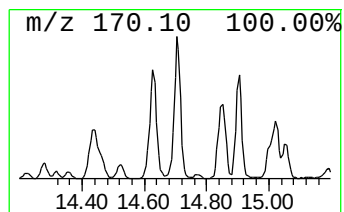
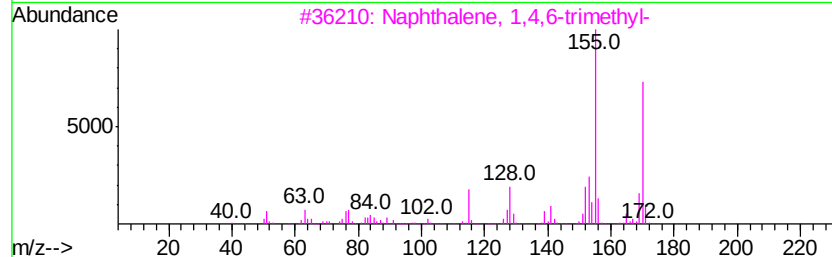
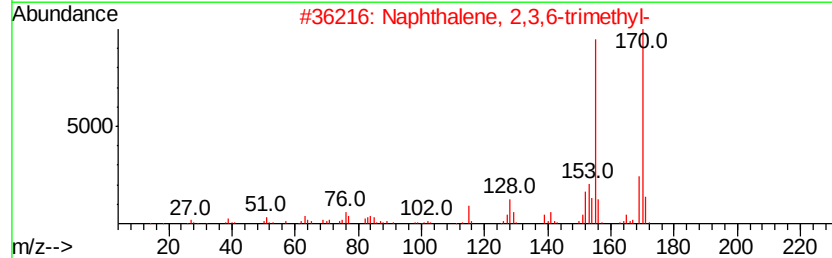
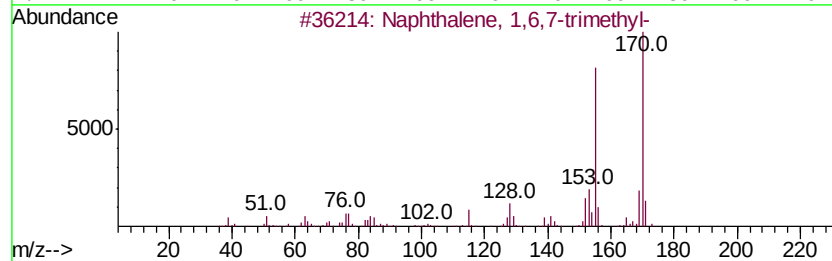
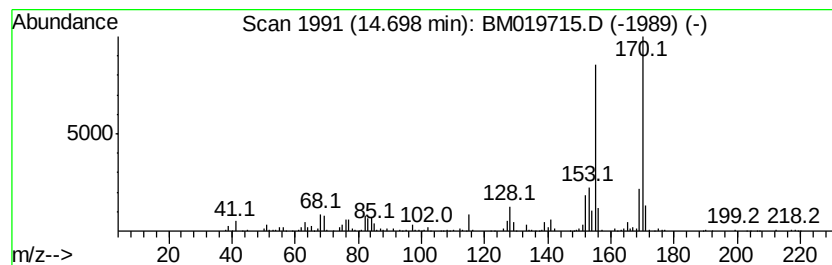
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Naphthalene, 1,6,7-trimethyl- Concentration Rank 41

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.70 | 5.38 ng/ul | 612146 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 98   |
| 2    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 3    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 96   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 95   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

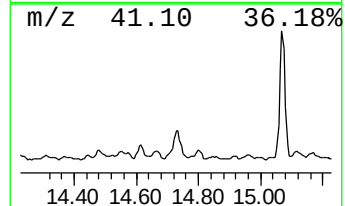
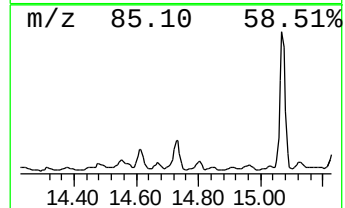
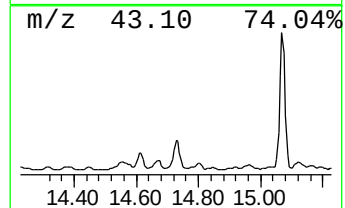
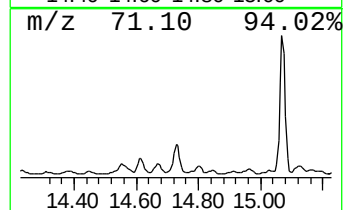
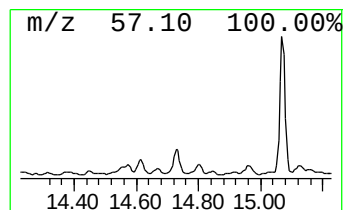
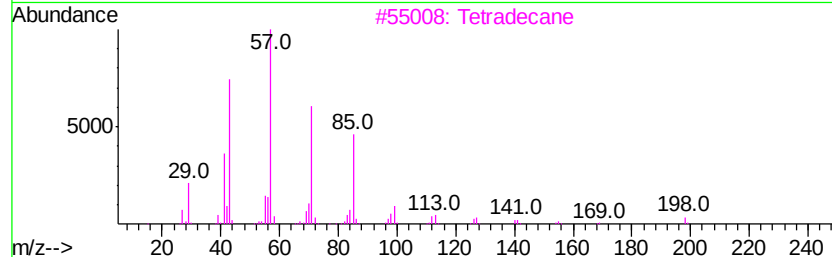
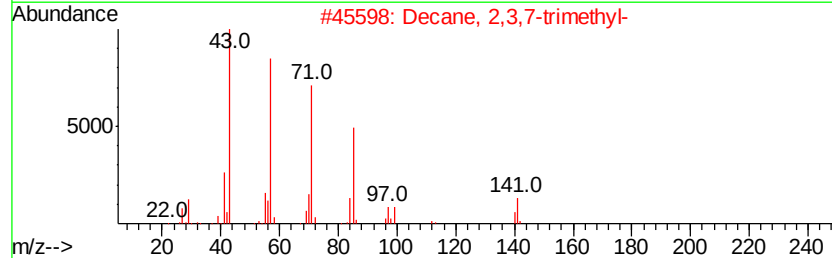
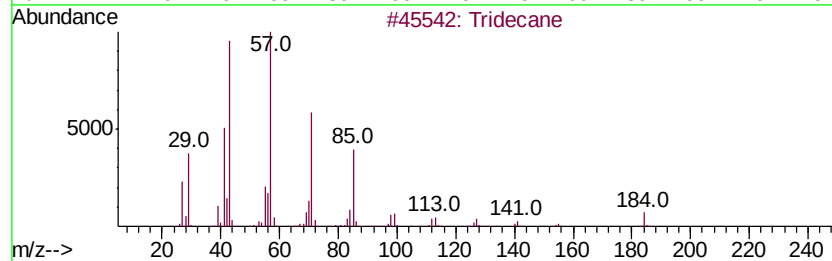
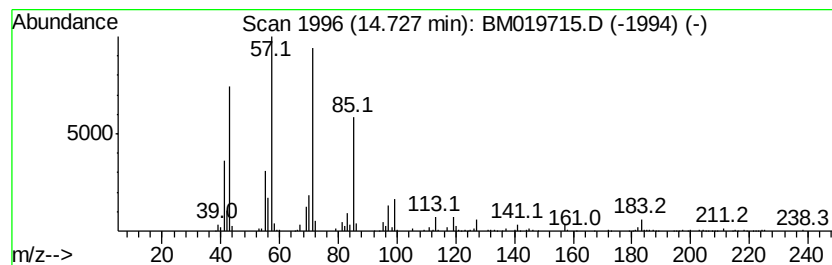
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.73 | 7.91 ng/ul | 899171 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane                           | 184 | C13H28  | 000629-50-5 | 64   |
| 2    |    | Decane, 2,3,7-trimethyl-            | 184 | C13H28  | 062238-13-5 | 64   |
| 3    |    | Tetradecane                         | 198 | C14H30  | 000629-59-4 | 58   |
| 4    |    | Butanal, 4-[(tetrahydro-2H-pyran... | 172 | C9H16O3 | 054911-85-2 | 47   |
| 5    |    | Decane, 5-propyl-                   | 184 | C13H28  | 017312-62-8 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

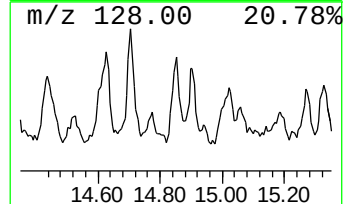
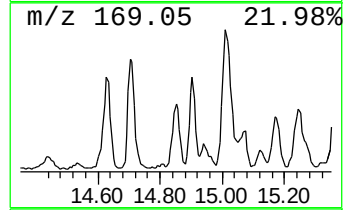
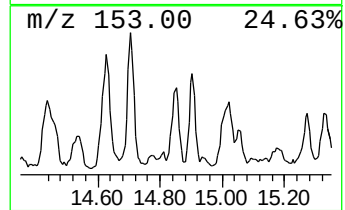
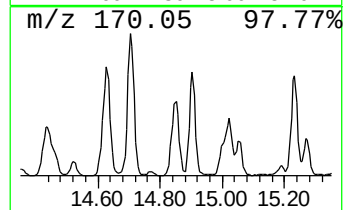
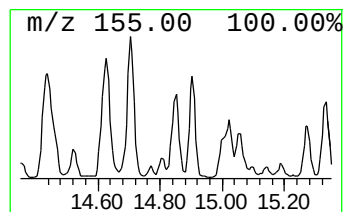
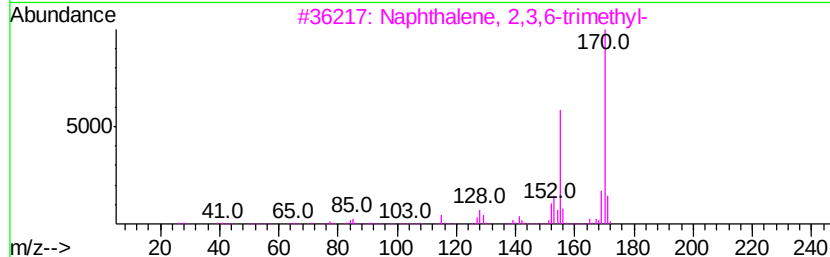
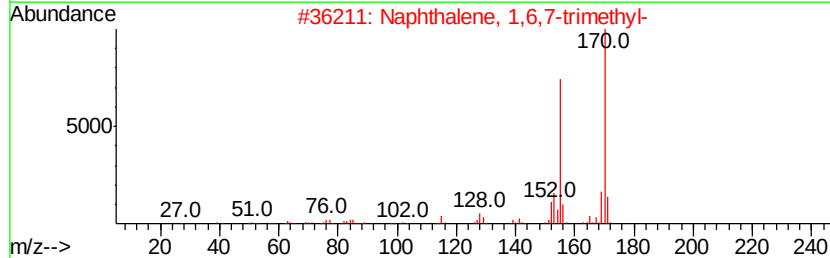
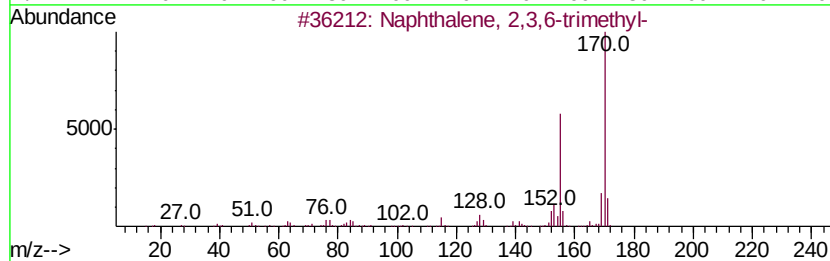
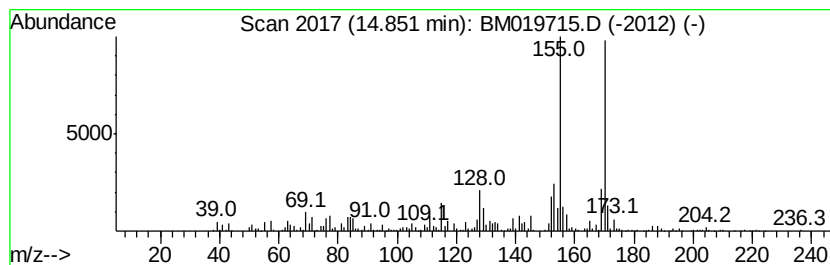
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 Naphthalene, 2,3,6-trimethyl- Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.85 | 6.35 ng/ul | 722645 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 5    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

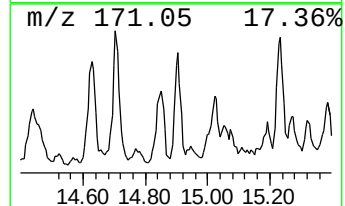
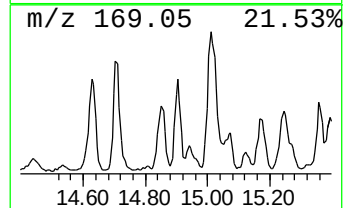
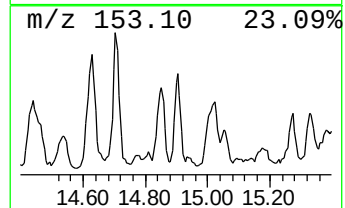
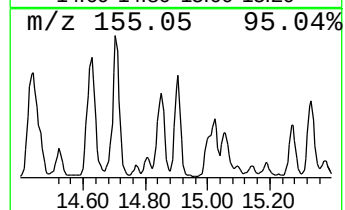
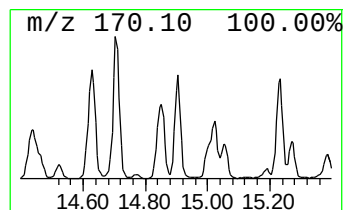
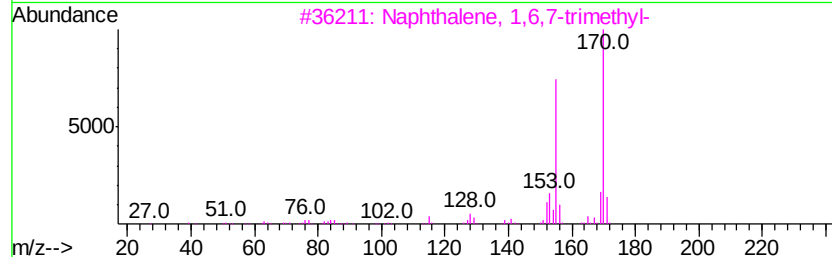
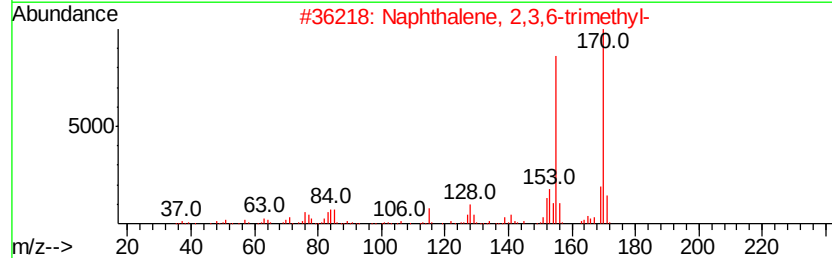
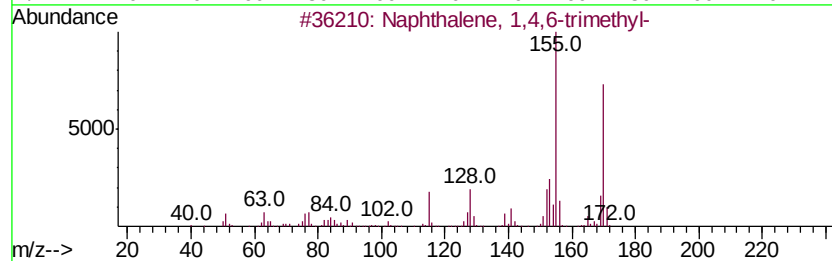
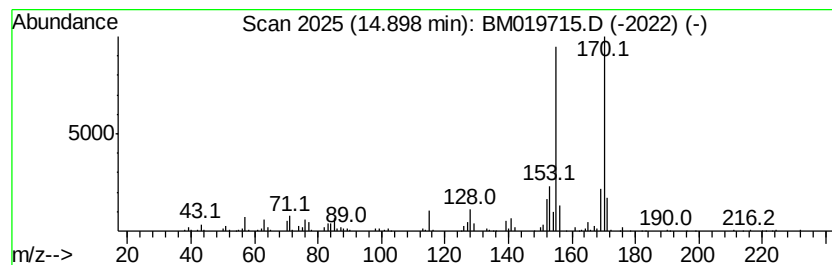
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 unknown-03 Concentration Rank 54

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.90 | 4.49 ng/ul | 510660 | Acenaphthene-d10 | 14.23 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 97   |
| 2    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 97   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |
| 5    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

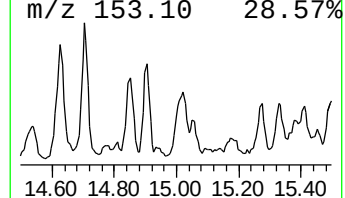
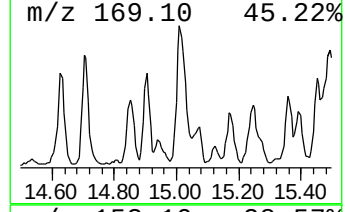
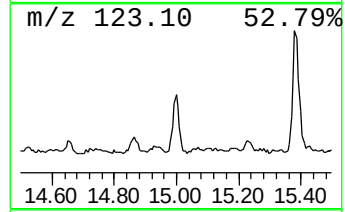
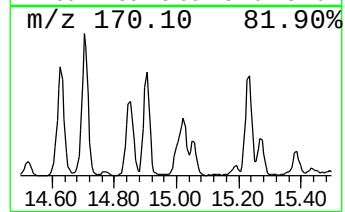
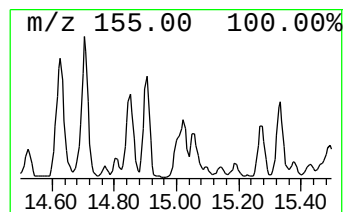
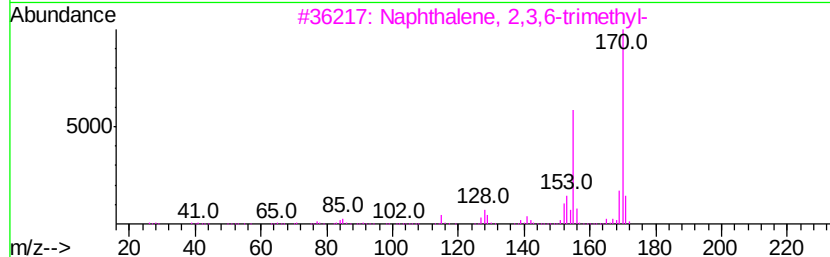
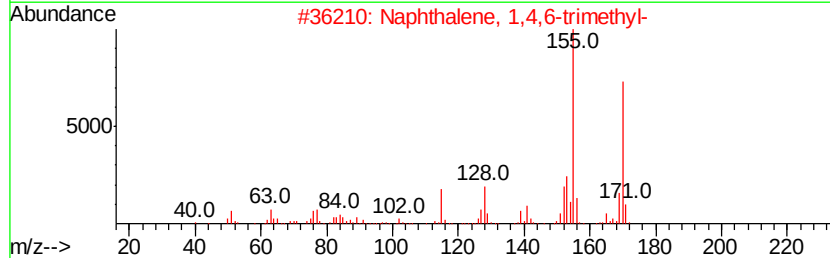
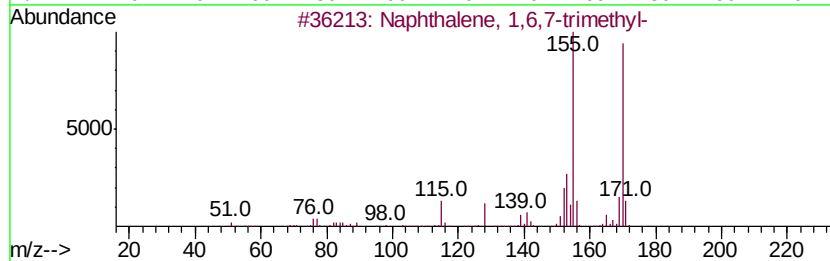
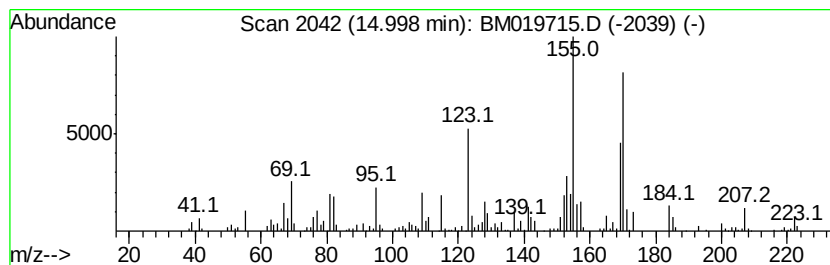
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 Azulene, 4,6,8-trimethyl- Concentration Rank 58

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.00 | 3.22 ng/ul | 366263 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 91   |
| 2    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 83   |
| 3    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 83   |
| 4    |    | Azulene, 4,6,8-trimethyl-     | 170 | C13H14  | 000941-81-1 | 64   |
| 5    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

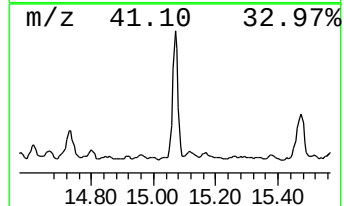
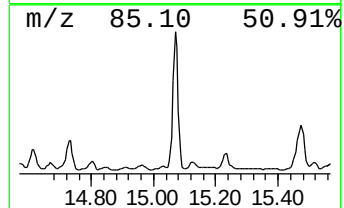
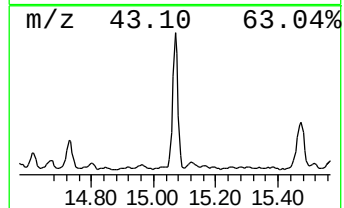
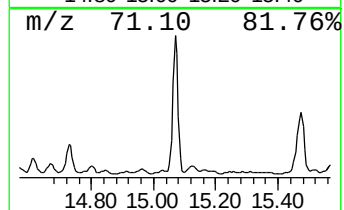
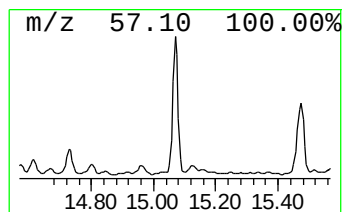
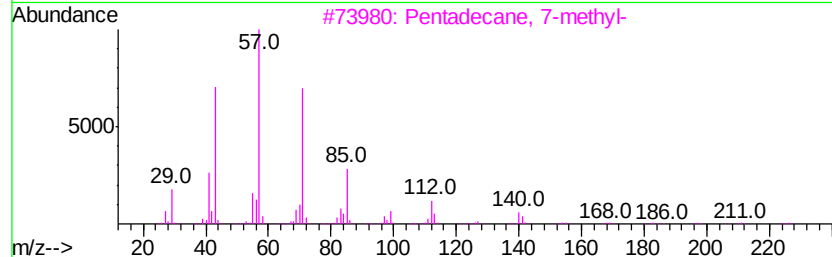
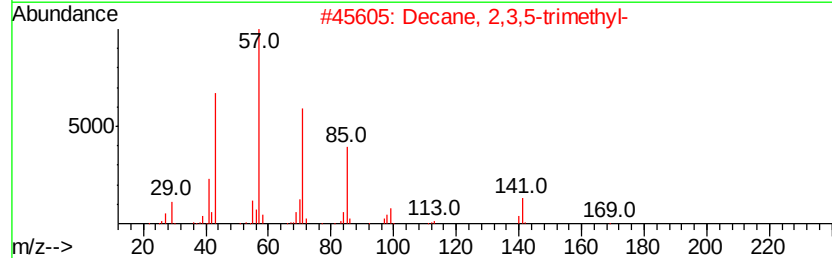
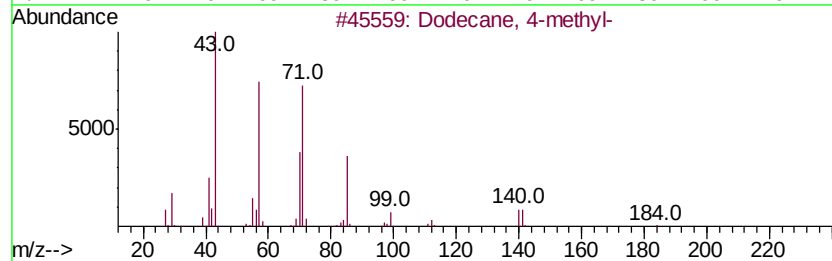
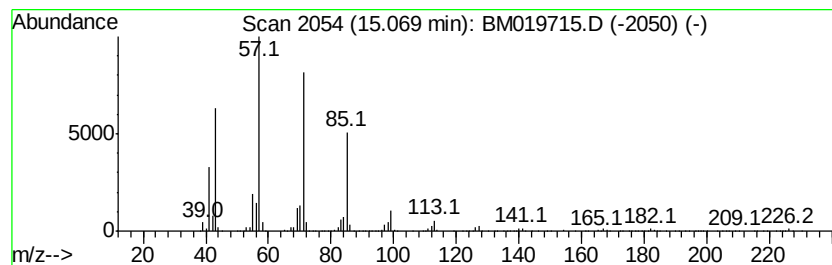
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.07 | 31.52 ng/ul | 3584150 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 4-methyl-      | 184 | C13H28  | 006117-97-1 | 90   |
| 2    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 90   |
| 3    |    | Pentadecane, 7-methyl-   | 226 | C16H34  | 006165-40-8 | 87   |
| 4    |    | Decane, 5-propyl-        | 184 | C13H28  | 017312-62-8 | 87   |
| 5    |    | Nonane, 5-butyl-         | 184 | C13H28  | 017312-63-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

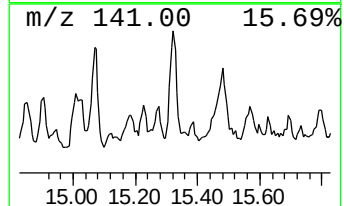
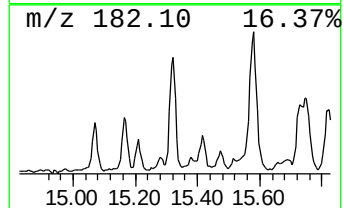
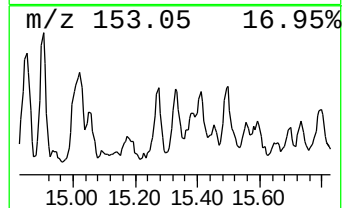
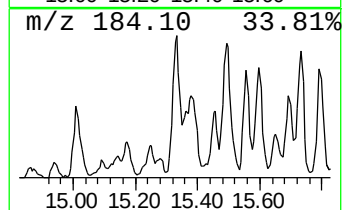
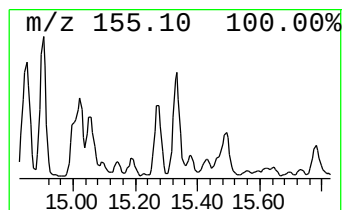
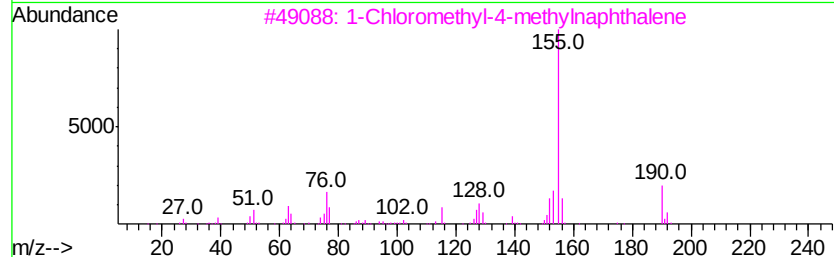
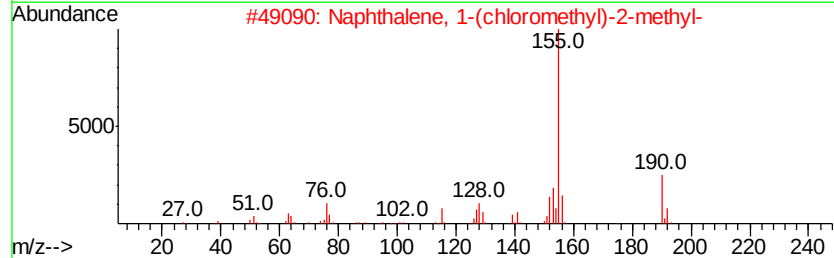
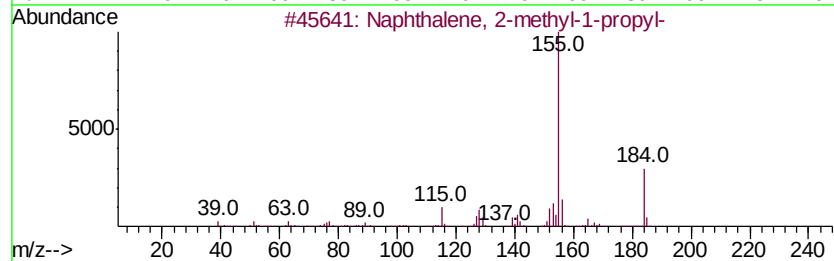
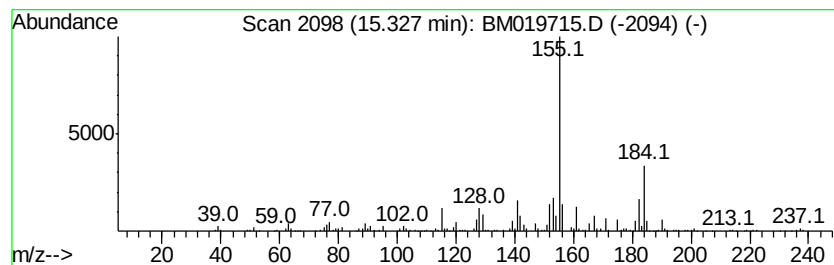
Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 41 Naphthalene, 2-methyl-1-pro... Concentration Rank 62

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 15.33   | 2.93 ng/ul | 333546                              | Acenaphthene-d10 | 14.23          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 2-methyl-1-propyl-     | 184 C14H16       | 054774-89-9 96 |
| 2       |            | Naphthalene, 1-(chloromethyl)-2-... | 190 C12H11Cl     | 006626-23-9 93 |
| 3       |            | 1-Chloromethyl-4-methylnaphthalene  | 190 C12H11Cl     | 005261-50-7 58 |
| 4       |            | Naphthalene, 2-(1-methylethyl)-     | 170 C13H14       | 002027-17-0 53 |
| 5       |            | Pyridine, 4-phenyl-                 | 155 C11H9N       | 000939-23-1 47 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

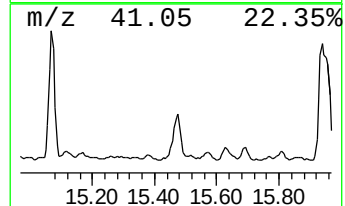
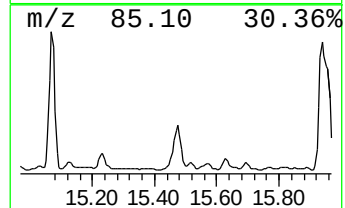
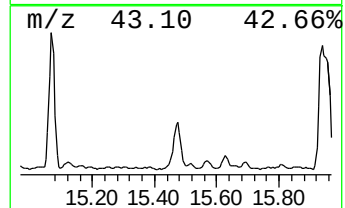
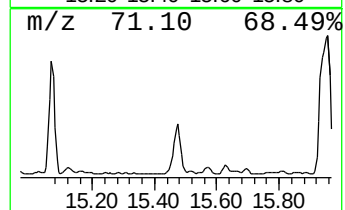
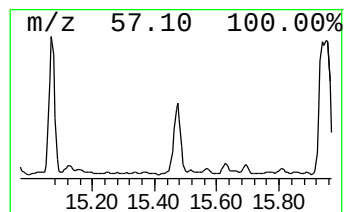
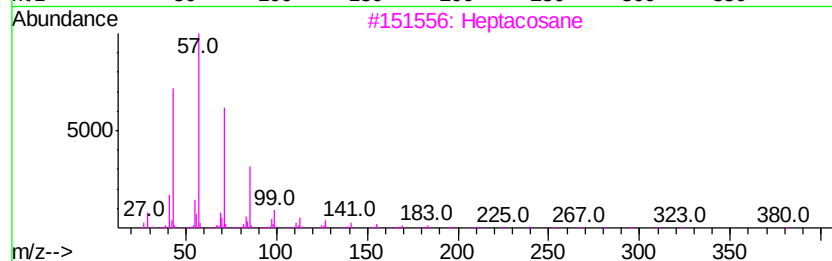
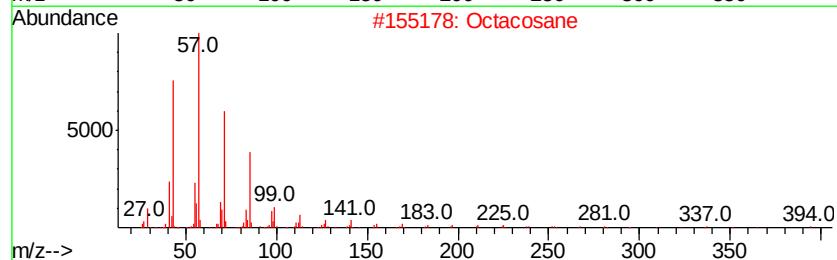
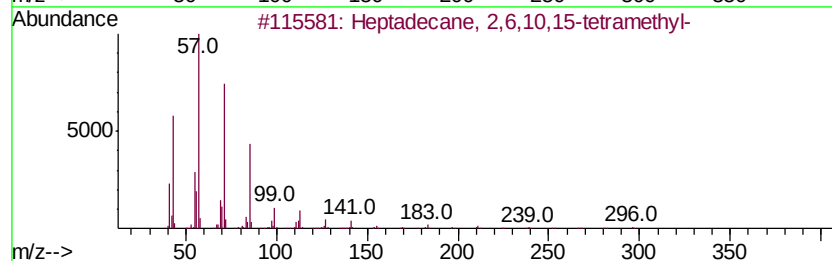
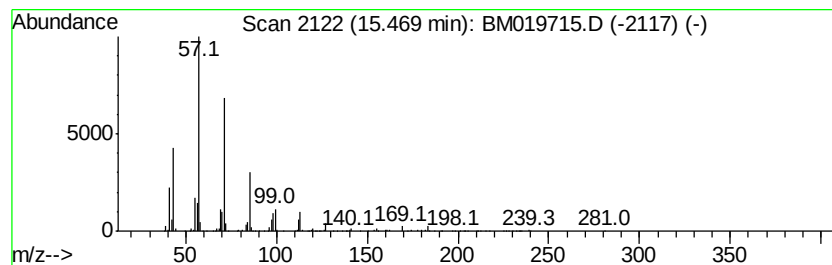
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.47 | 18.50 ng/ul | 2103450 | Acenaphthene-d10 | 14.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 87   |
| 2    |    | Octacosane                          | 394 | C28H58   | 000630-02-4 | 86   |
| 3    |    | Heptacosane                         | 380 | C27H56   | 000593-49-7 | 86   |
| 4    |    | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42   | 000638-36-8 | 80   |
| 5    |    | 2-Bromo dodecane                    | 248 | C12H25Br | 013187-99-0 | 78   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

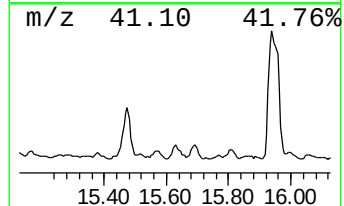
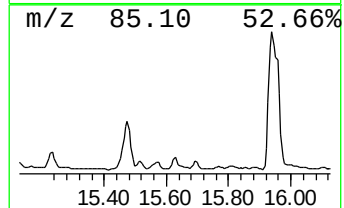
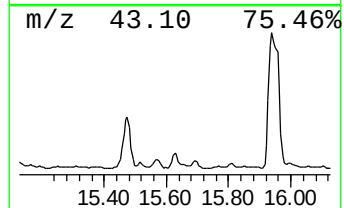
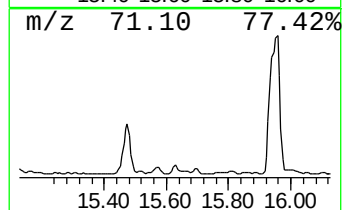
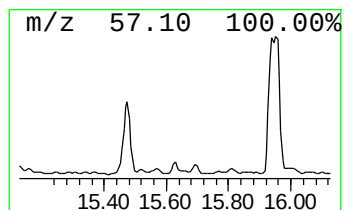
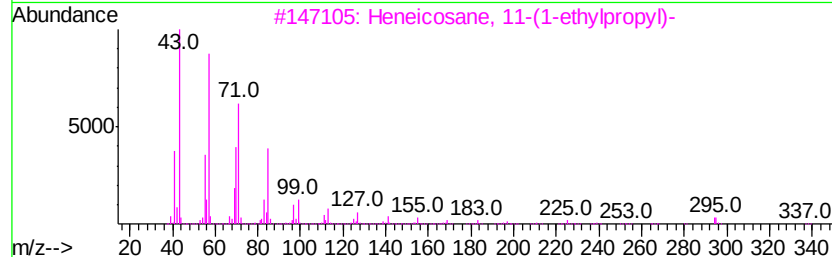
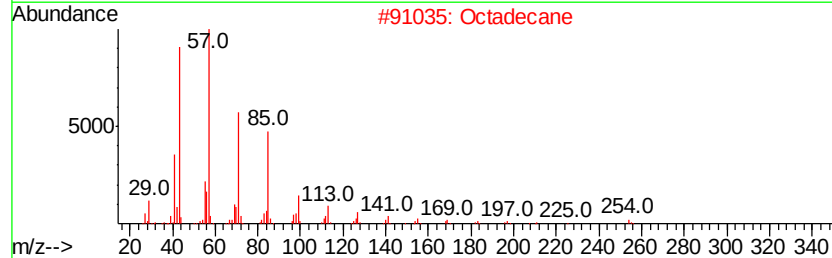
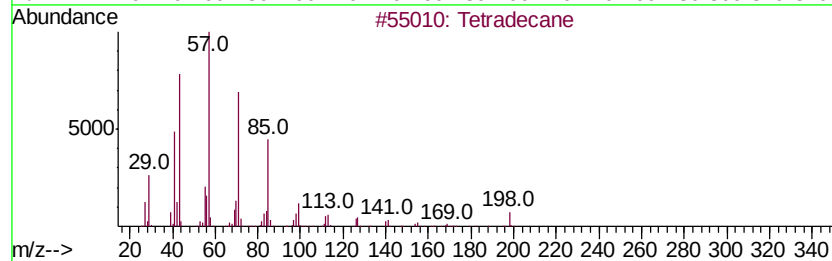
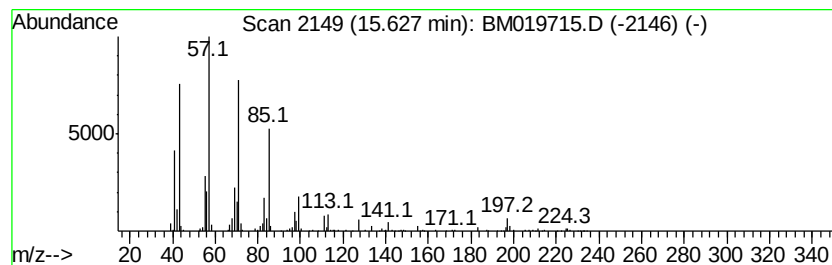
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 59

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.63 | 3.11 ng/ul | 416572 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane                         | 198 | C14H30  | 000629-59-4 | 90   |
| 2    |    |   | Octadecane                          | 254 | C18H38  | 000593-45-3 | 90   |
| 3    |    |   | Heneicosane, 11-(1-ethylpropyl)-    | 366 | C26H54  | 055282-11-6 | 87   |
| 4    |    |   | 10-Methylnonadecane                 | 282 | C20H42  | 056862-62-5 | 86   |
| 5    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 86   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

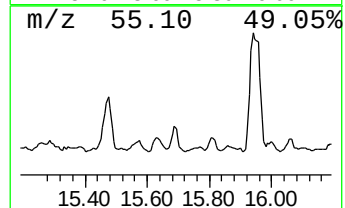
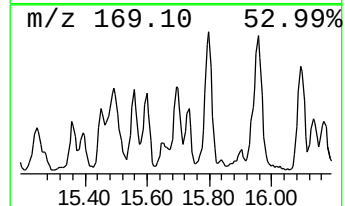
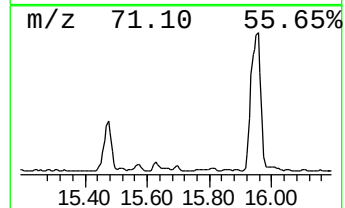
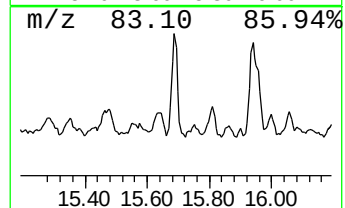
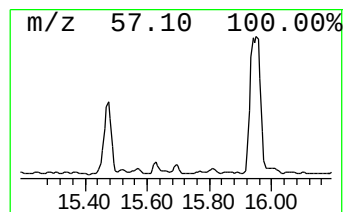
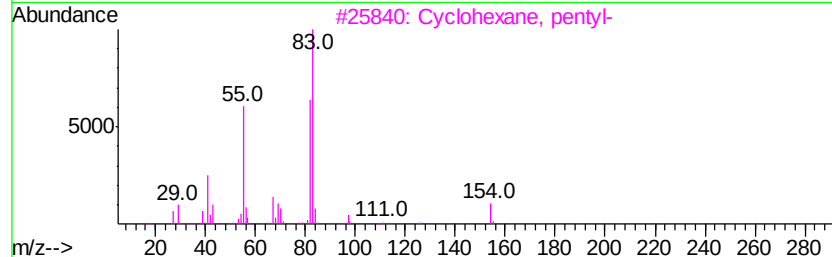
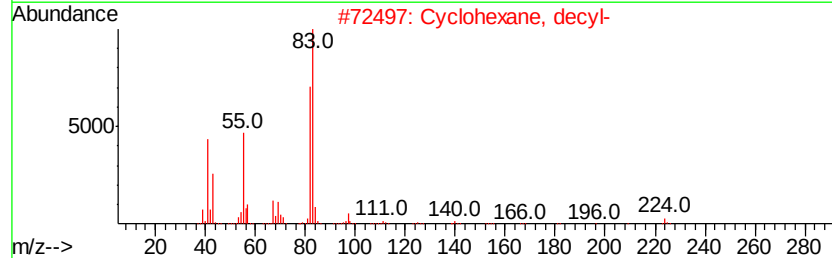
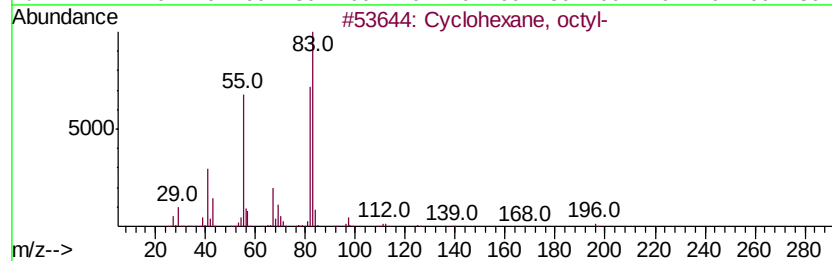
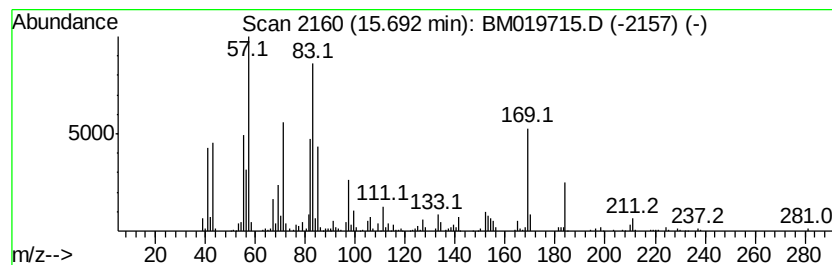
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 (DEL) Alkane: Cyclic15.69 Concentration Rank 60

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.69 | 3.06 ng/ul | 410221 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, octyl-   | 196 | C14H28  | 001795-15-9 | 42   |
| 2    |    |   | Cyclohexane, decyl-   | 224 | C16H32  | 001795-16-0 | 30   |
| 3    |    |   | Cyclohexane, pentyl-  | 154 | C11H22  | 004292-92-6 | 30   |
| 4    |    |   | Cyclohexane, undecyl- | 238 | C17H34  | 054105-66-7 | 30   |
| 5    |    |   | Cyclohexane, pentyl-  | 154 | C11H22  | 004292-92-6 | 30   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

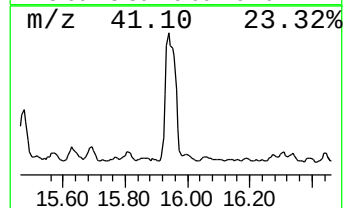
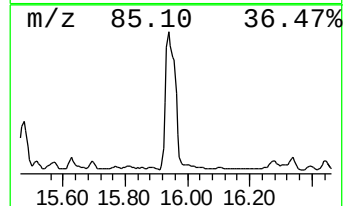
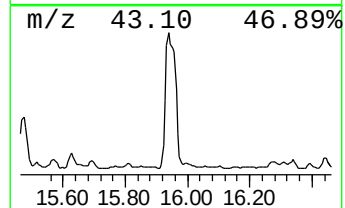
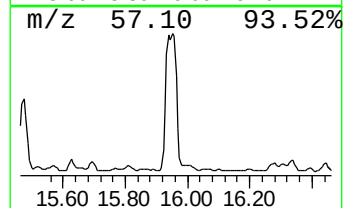
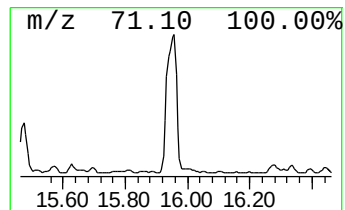
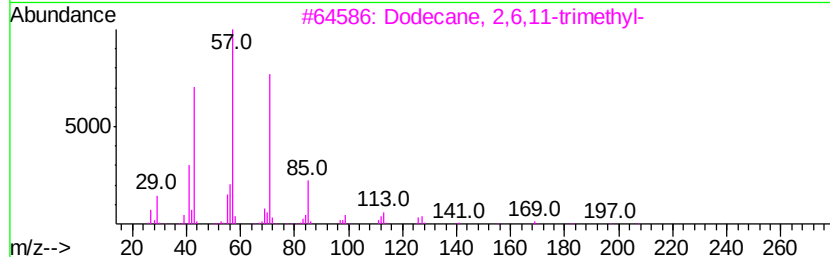
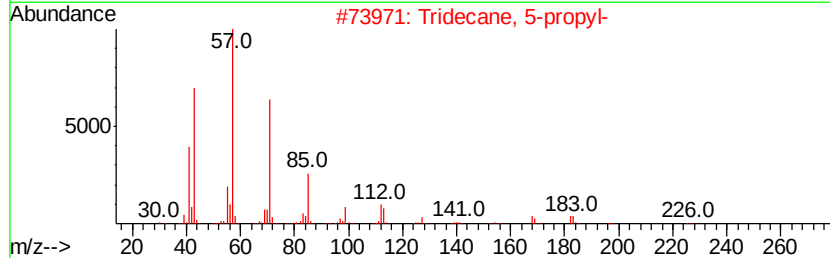
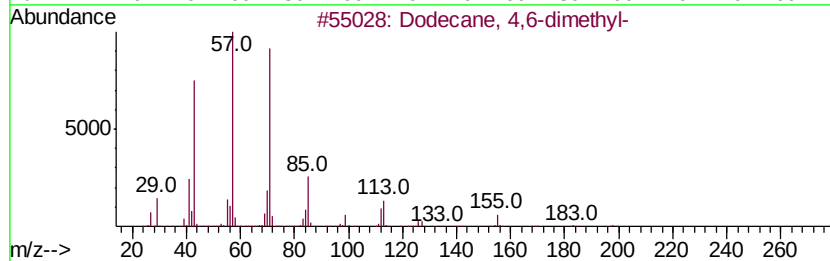
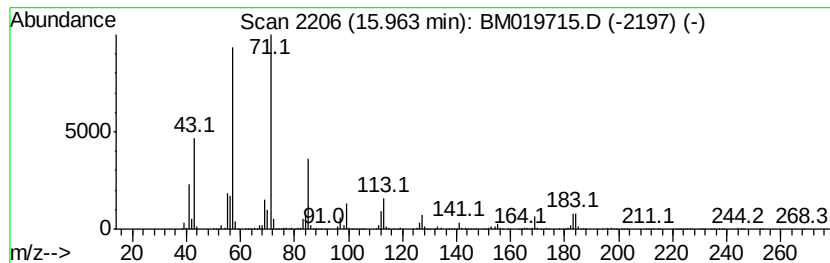
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.96 | 55.38 ng/ul | 7423740 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 4,6-dimethyl-             | 198 | C14H30  | 061141-72-8 | 86   |
| 2    |    | Tridecane, 5-propyl-                | 226 | C16H34  | 055045-11-9 | 72   |
| 3    |    | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32  | 031295-56-4 | 64   |
| 4    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 58   |
| 5    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

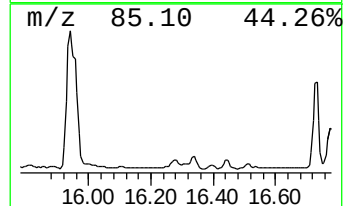
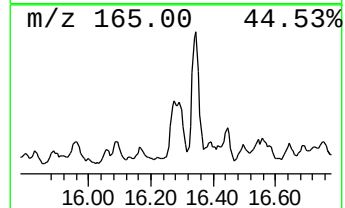
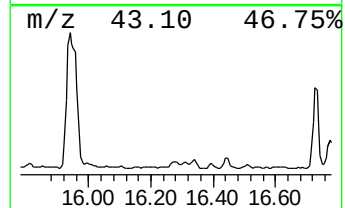
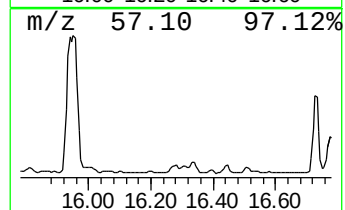
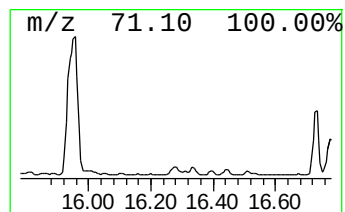
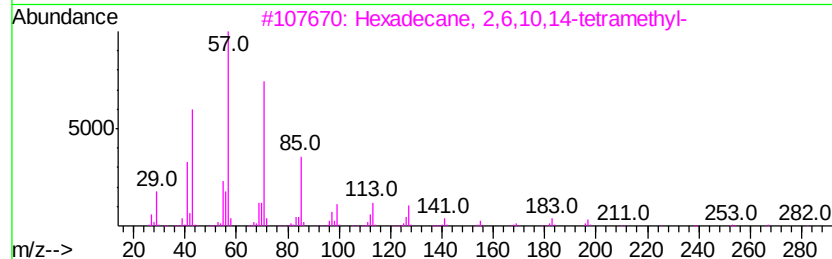
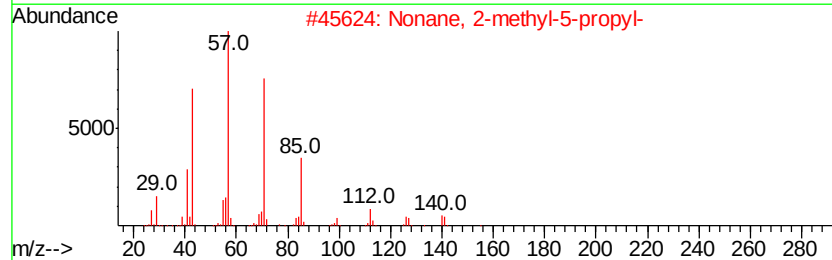
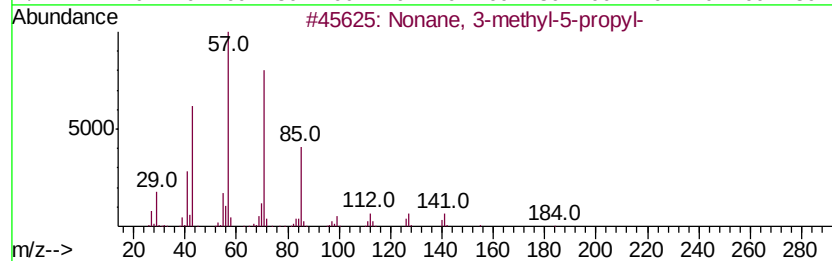
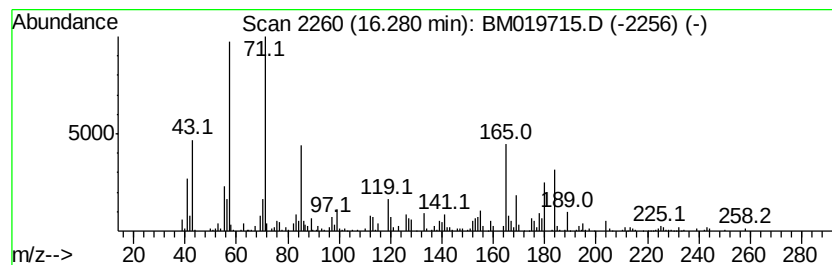
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 66

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.28 | 2.67 ng/ul | 358274 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Nonane, 3-methyl-5-propyl-          | 184 | C13H28  | 031081-18-2  | 38   |
| 2    |    |   | Nonane, 2-methyl-5-propyl-          | 184 | C13H28  | 031081-17-1  | 38   |
| 3    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8  | 35   |
| 4    |    |   | 3-Methyl-4-(methoxycarbonyl)hexa... | 184 | C9H12O4 | 1000104-10-8 | 25   |
| 5    |    |   | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32  | 031295-56-4  | 25   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

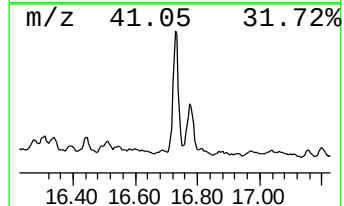
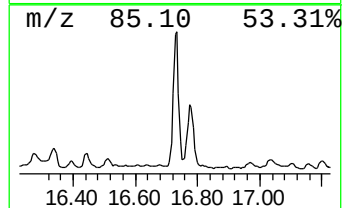
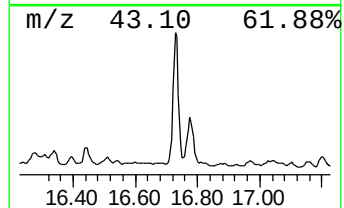
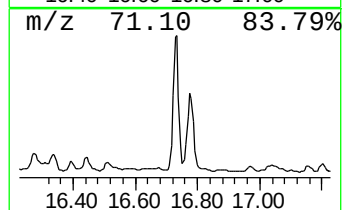
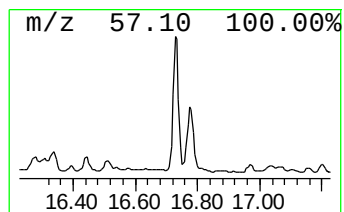
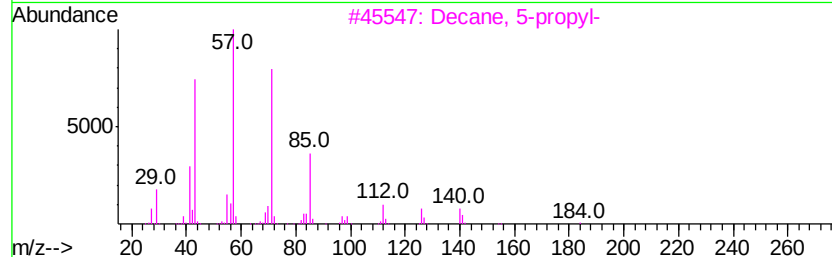
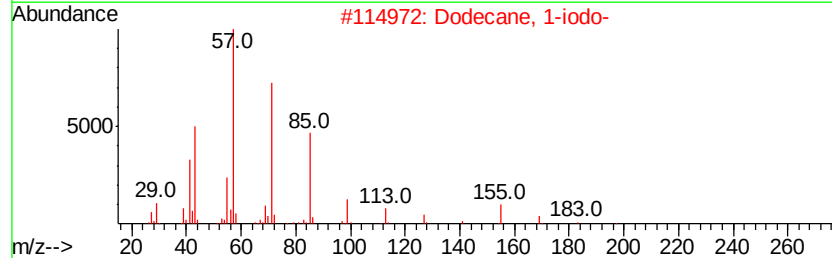
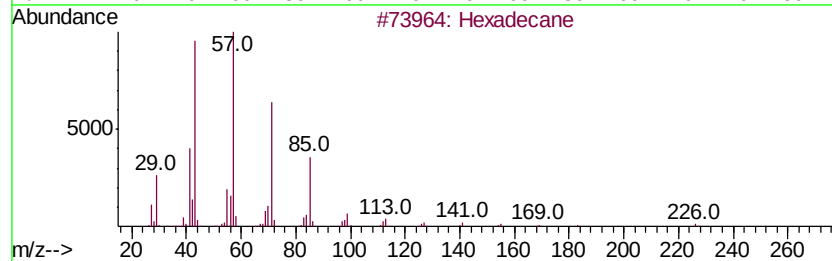
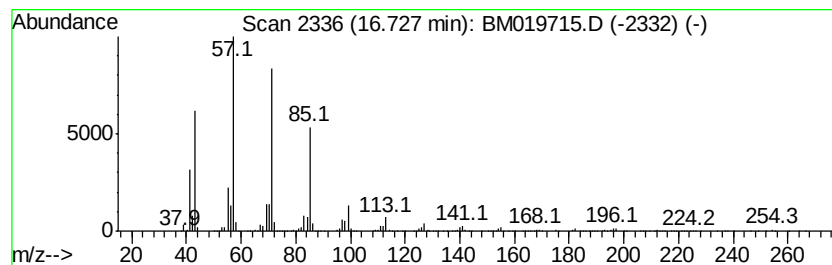
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.73 | 15.95 ng/ul | 2137650 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane         | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    | Dodecane, 1-iodo-  | 296 | C12H25I | 004292-19-7 | 86   |
| 3    |    | Decane, 5-propyl-  | 184 | C13H28  | 017312-62-8 | 83   |
| 4    |    | Undecane, 6-ethyl- | 184 | C13H28  | 017312-60-6 | 81   |
| 5    |    | Dodecane           | 170 | C12H26  | 000112-40-3 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

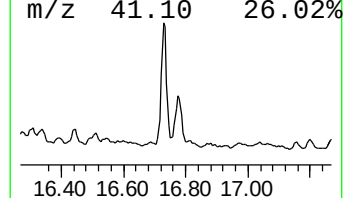
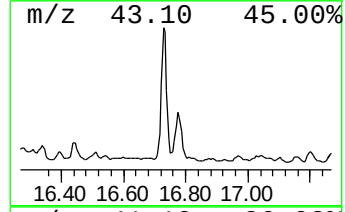
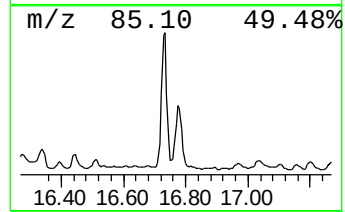
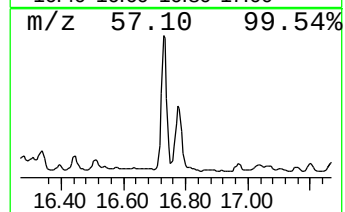
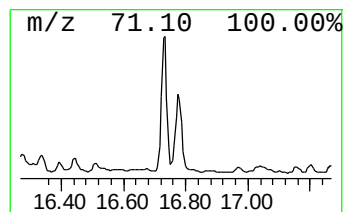
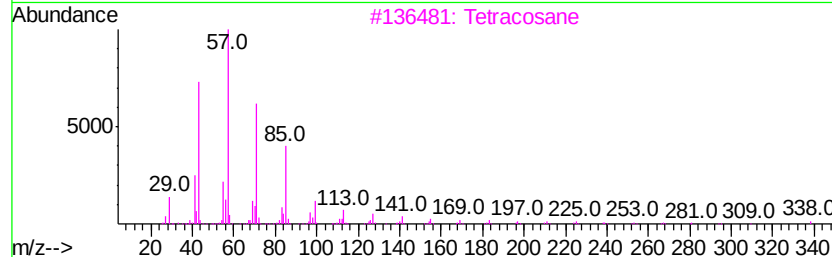
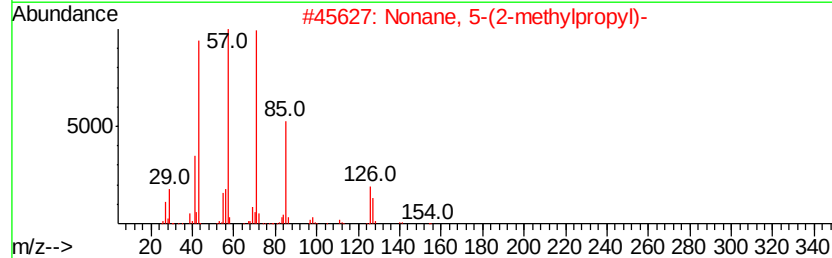
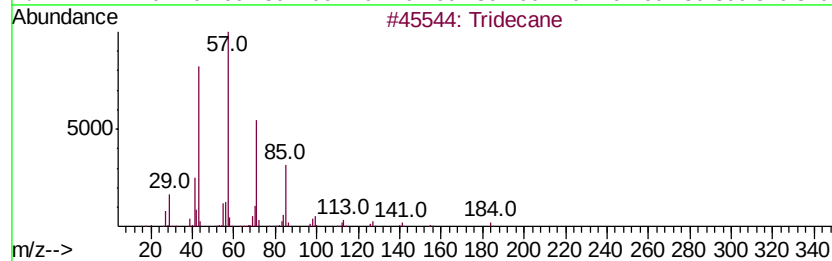
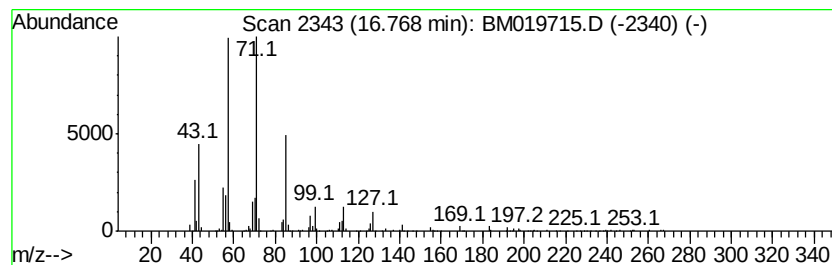
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.77 | 8.26 ng/ul | 1106910 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane                           | 184 | C13H28  | 000629-50-5 | 91   |
| 2    |    |   | Nonane, 5-(2-methylpropyl)-         | 184 | C13H28  | 062185-53-9 | 64   |
| 3    |    |   | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 64   |
| 4    |    |   | Nonane, 3,7-dimethyl-               | 156 | C11H24  | 017302-32-8 | 58   |
| 5    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

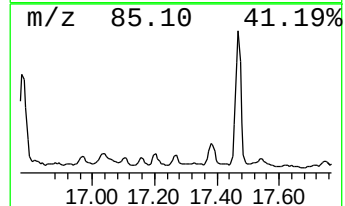
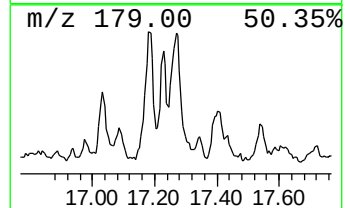
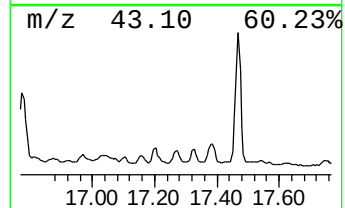
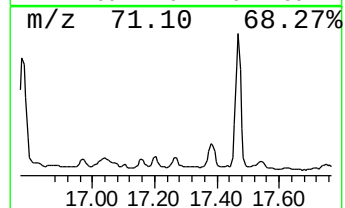
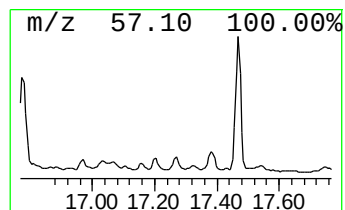
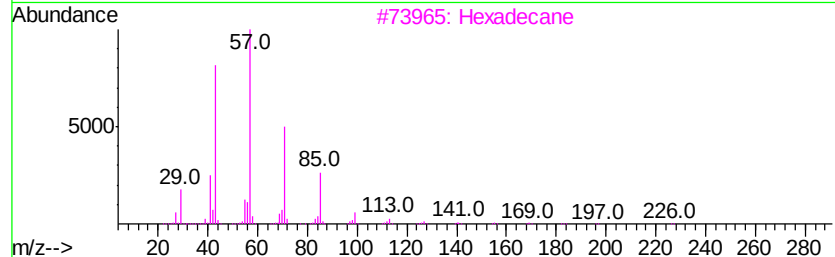
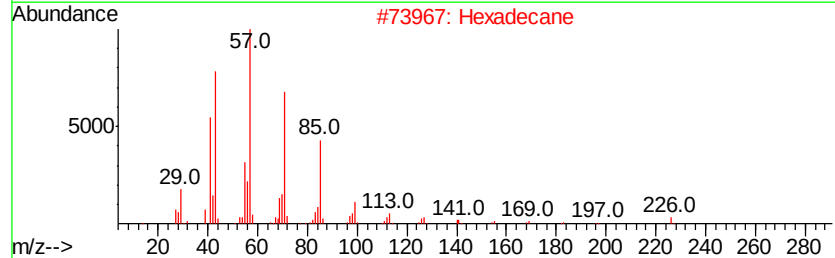
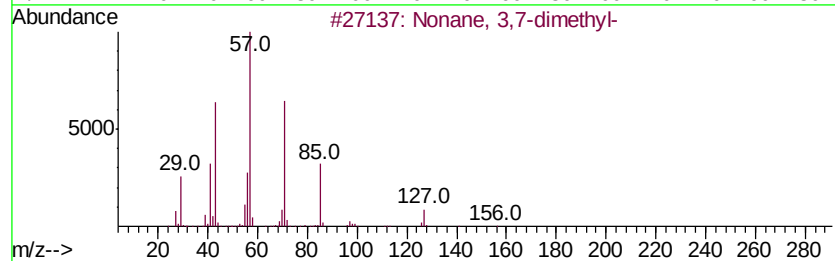
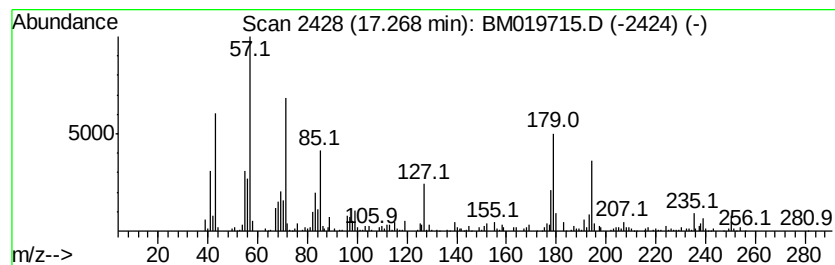
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 67

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.27 | 2.65 ng/ul | 355258 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 3,7-dimethyl- | 156 | C11H24  | 017302-32-8 | 46   |
| 2    |    | Hexadecane            | 226 | C16H34  | 000544-76-3 | 42   |
| 3    |    | Hexadecane            | 226 | C16H34  | 000544-76-3 | 42   |
| 4    |    | Octadecane            | 254 | C18H38  | 000593-45-3 | 38   |
| 5    |    | Octadecane            | 254 | C18H38  | 000593-45-3 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
 Data File : BM019715.D  
 Acq On : 03 Apr 2019 05:41  
 Operator : JU/SJ  
 Sample : K2168-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION

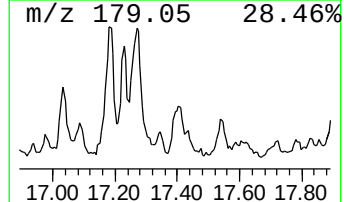
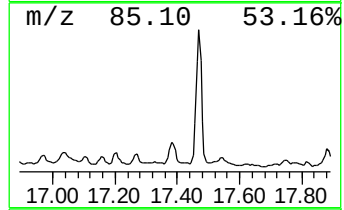
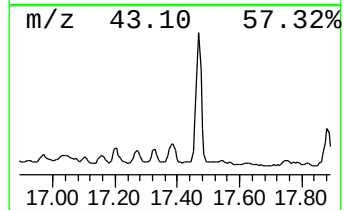
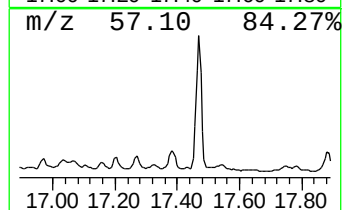
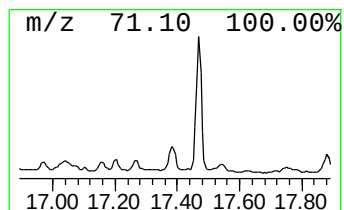
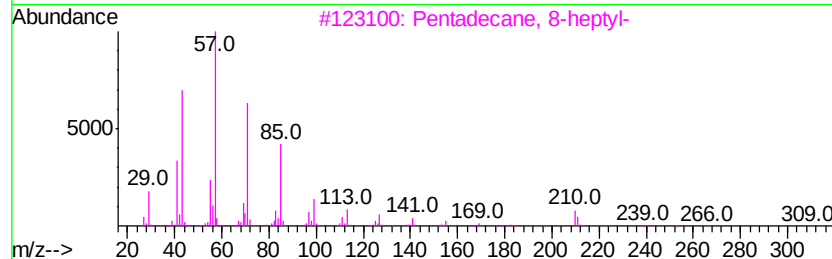
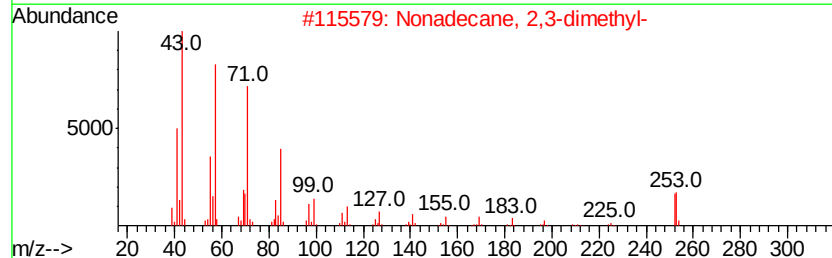
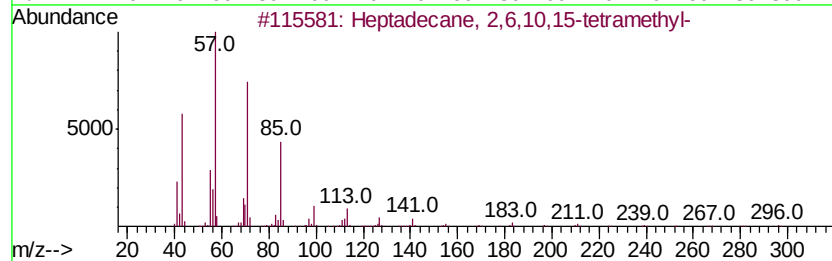
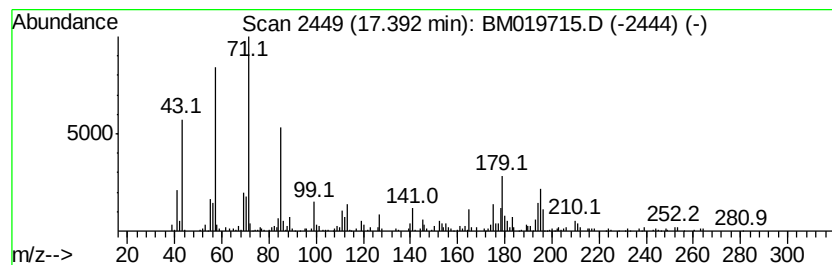
TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 68

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.39 | 2.63 ng/ul | 352696 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 64   |
| 2    |    |   | Nonadecane, 2,3-dimethyl-           | 296 | C21H44  | 075163-99-4 | 64   |
| 3    |    |   | Pentadecane, 8-heptyl-              | 310 | C22H46  | 071005-15-7 | 64   |
| 4    |    |   | Tetradecane, 4-methyl-              | 212 | C15H32  | 025117-24-2 | 62   |
| 5    |    |   | Eicosane                            | 282 | C20H42  | 000112-95-8 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

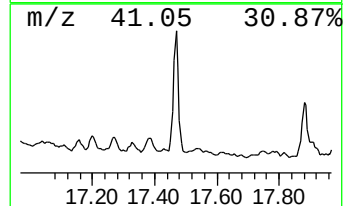
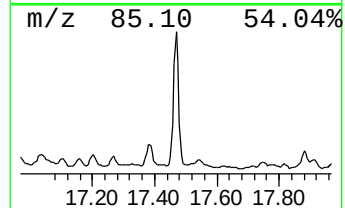
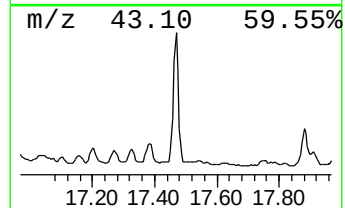
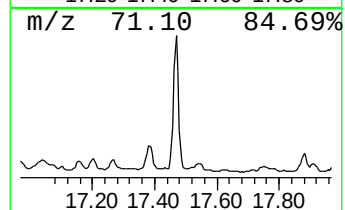
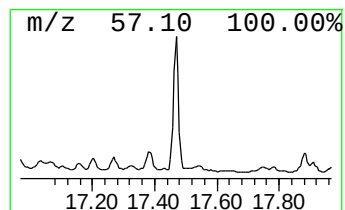
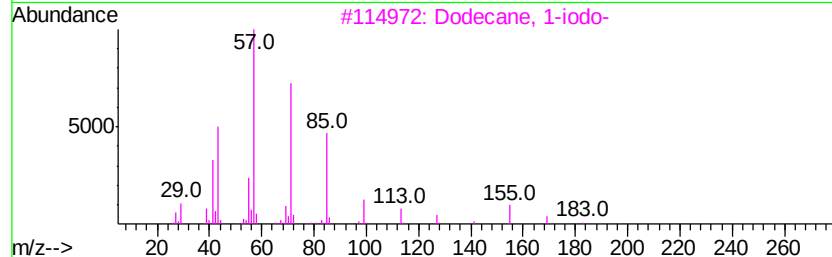
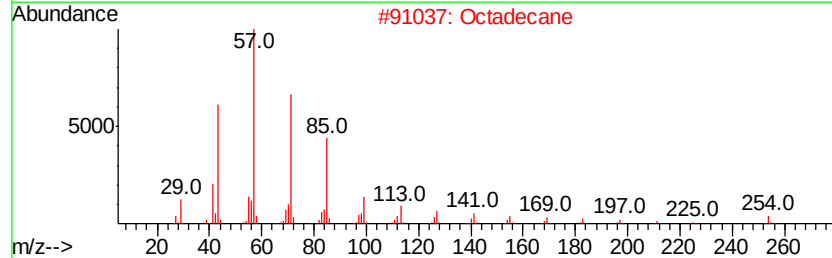
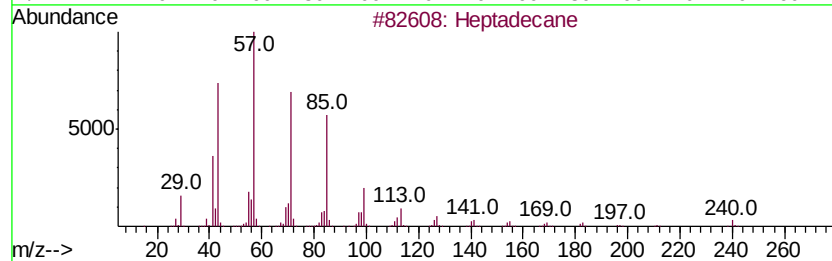
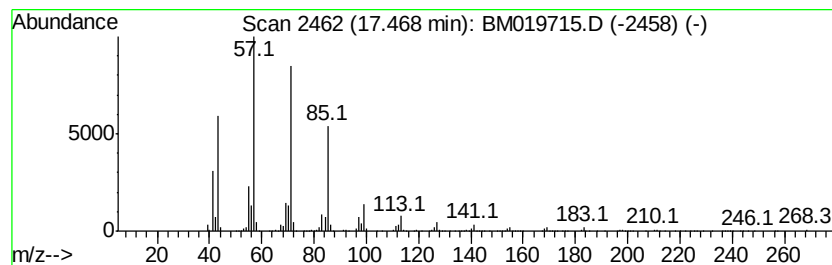
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.47 | 11.08 ng/ul | 1485280 | Phenanthrene-d10 | 16.99 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane       | 240 | C17H36  | 000629-78-7 | 95   |
| 2    |    | Octadecane        | 254 | C18H38  | 000593-45-3 | 90   |
| 3    |    | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 86   |
| 4    |    | Tetradecane       | 198 | C14H30  | 000629-59-4 | 83   |
| 5    |    | Heptacosane       | 380 | C27H56  | 000593-49-7 | 83   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

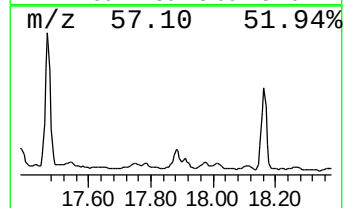
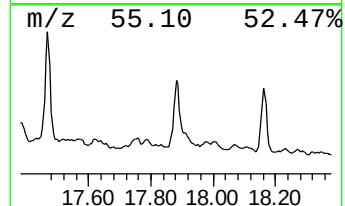
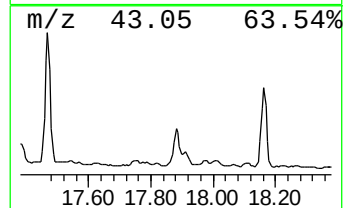
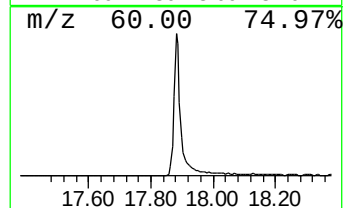
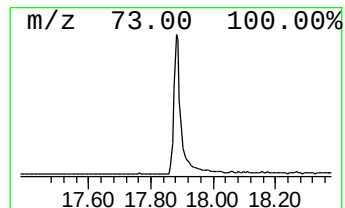
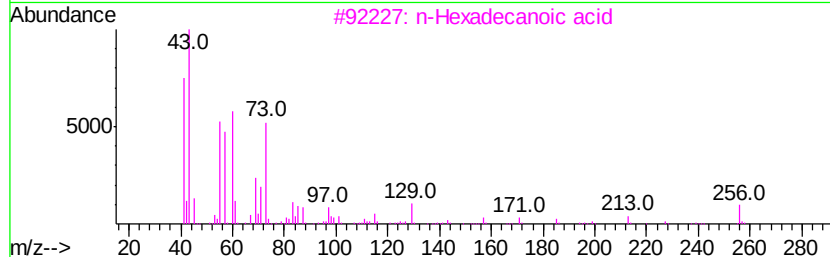
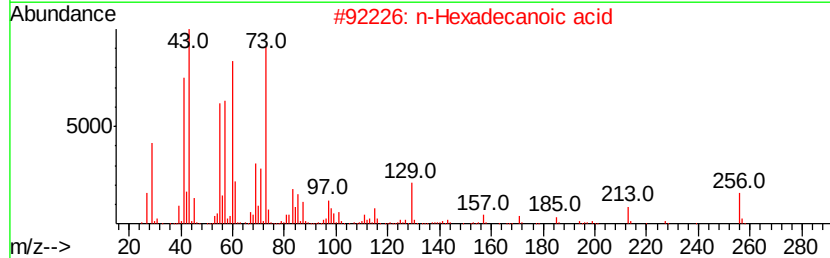
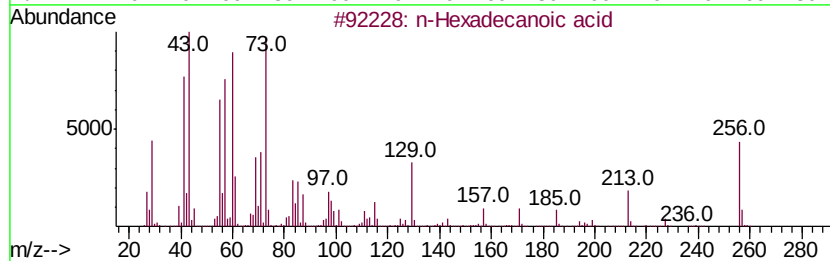
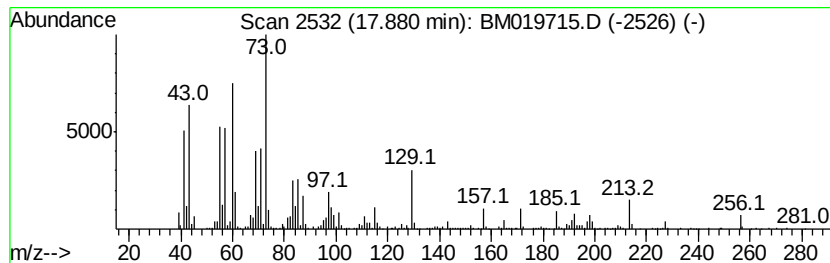
Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 54 n-Hexadecanoic acid Concentration Rank 27

| R.T.    | EstConc             | Area         | Relative to ISTD | R.T.      |
|---------|---------------------|--------------|------------------|-----------|
| 17.88   | 8.03 ng/ul          | 1076290      | Phenanthrene-d10 | 16.99     |
| Hit# of | 5                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | n-Hexadecanoic acid | 256 C16H32O2 | 000057-10-3      | 98        |
| 2       | n-Hexadecanoic acid | 256 C16H32O2 | 000057-10-3      | 96        |
| 3       | n-Hexadecanoic acid | 256 C16H32O2 | 000057-10-3      | 95        |
| 4       | Tetradecanoic acid  | 228 C14H28O2 | 000544-63-8      | 80        |
| 5       | Pentadecanoic acid  | 242 C15H30O2 | 001002-84-2      | 78        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

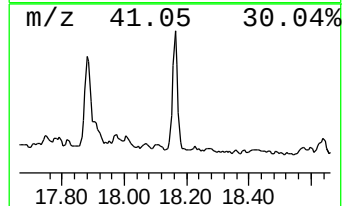
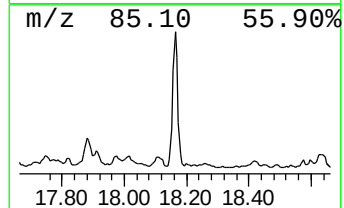
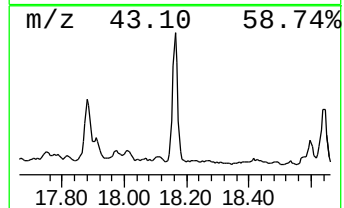
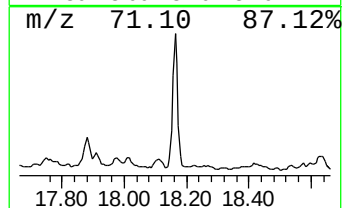
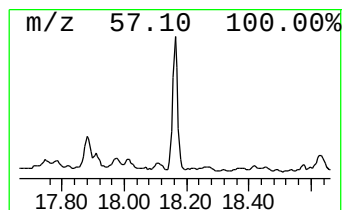
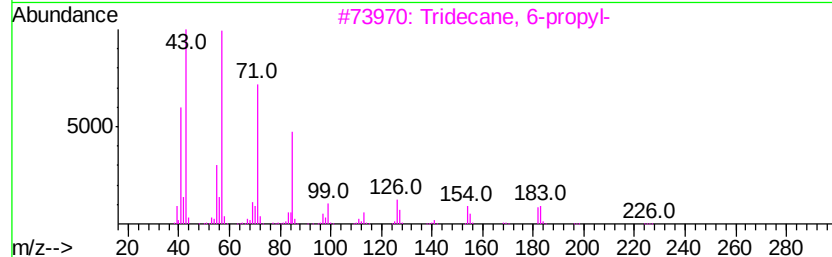
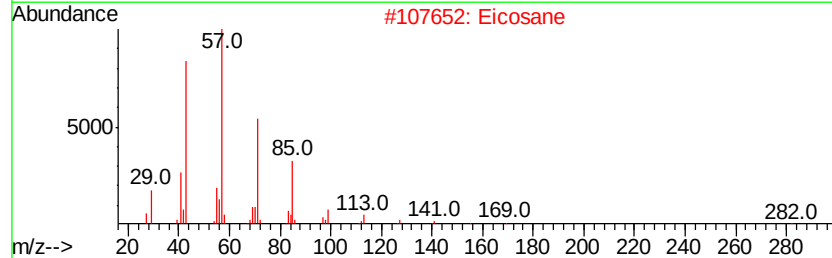
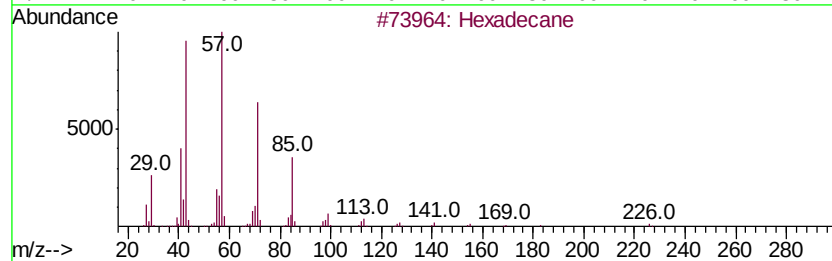
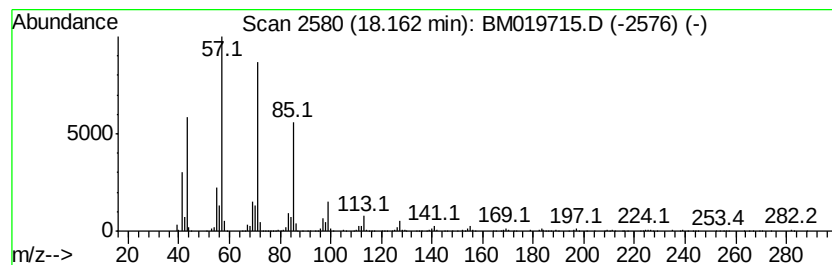
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.16 | 6.69 ng/ul | 896556 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane           | 226 | C16H34  | 000544-76-3 | 96   |
| 2    |    |   | Eicosane             | 282 | C20H42  | 000112-95-8 | 91   |
| 3    |    |   | Tridecane, 6-propyl- | 226 | C16H34  | 055045-10-8 | 90   |
| 4    |    |   | Dodecane, 1-iodo-    | 296 | C12H25I | 004292-19-7 | 87   |
| 5    |    |   | Heptacosane          | 380 | C27H56  | 000593-49-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

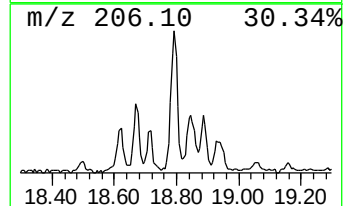
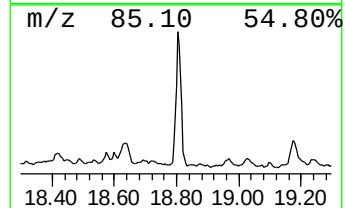
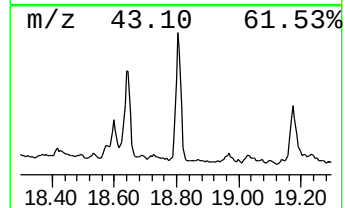
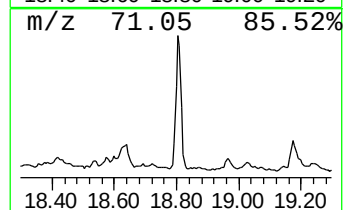
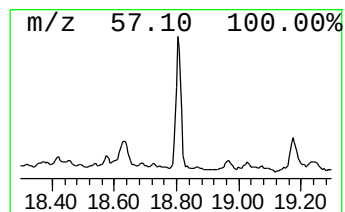
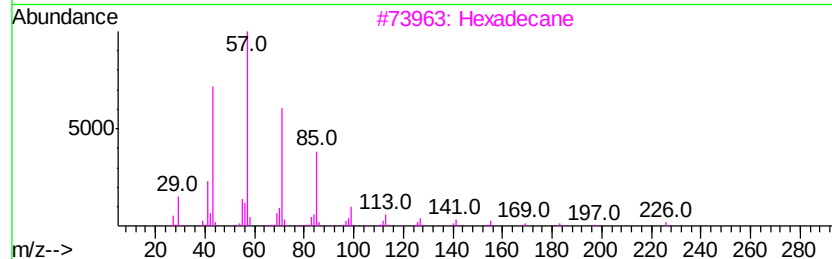
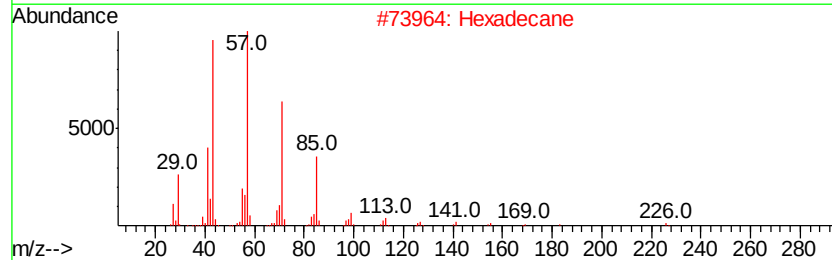
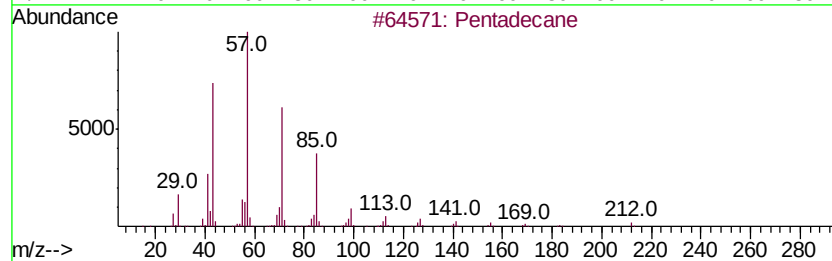
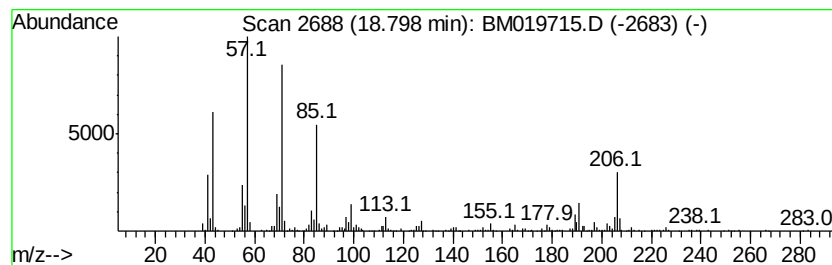
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.80 | 5.44 ng/ul | 728971 | Phenanthrene-d10 | 16.99 |

| Hit# | of | 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane       | 212 | C15H32  | 000629-62-9 | 95   |
| 2    |    |   | Hexadecane        | 226 | C16H34  | 000544-76-3 | 95   |
| 3    |    |   | Hexadecane        | 226 | C16H34  | 000544-76-3 | 83   |
| 4    |    |   | Nonadecane        | 268 | C19H40  | 000629-92-5 | 83   |
| 5    |    |   | Dodecane, 1-iodo- | 296 | C12H25I | 004292-19-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

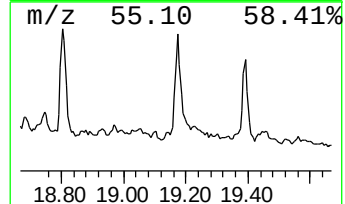
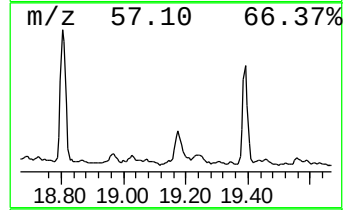
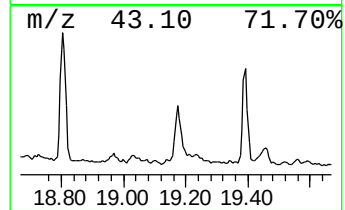
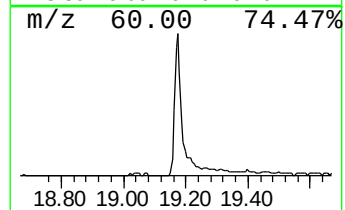
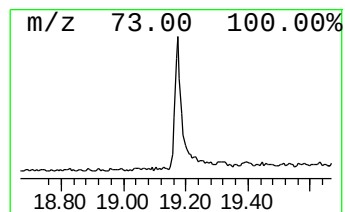
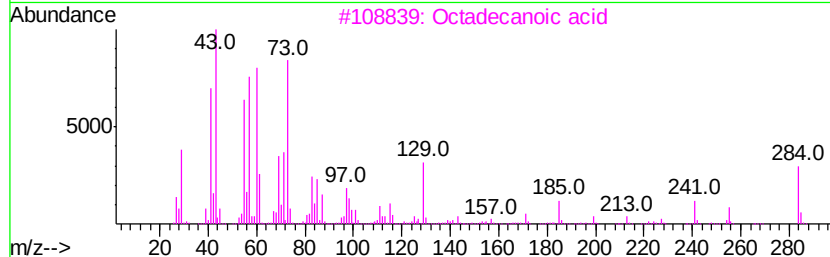
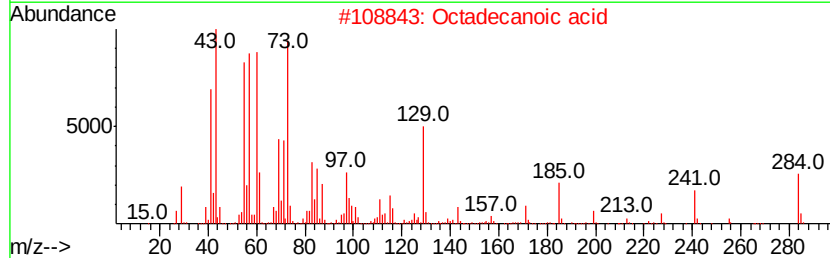
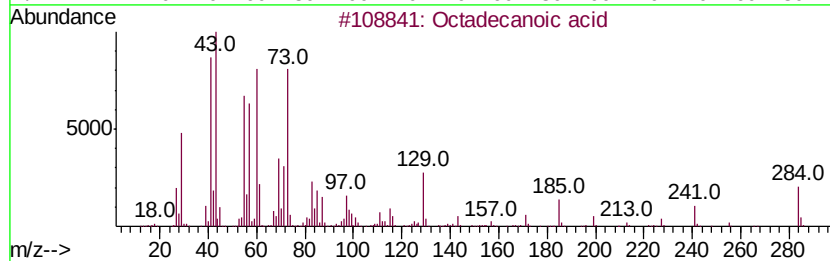
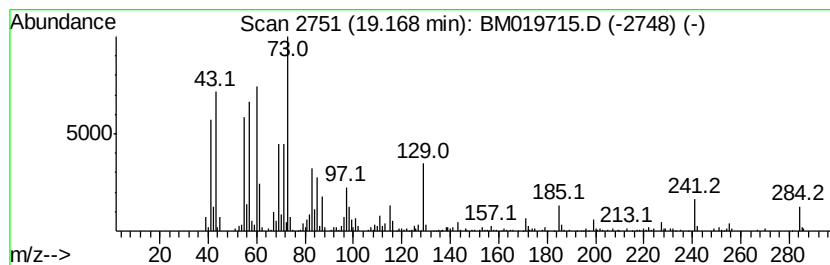
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 57 Octadecanoic acid Concentration Rank 57

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.17 | 3.36 ng/ul | 536441 | Chrysene-d12     | 21.20 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 96   |
| 2    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 95   |
| 3    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 91   |
| 4    |    |   | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 74   |
| 5    |    |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

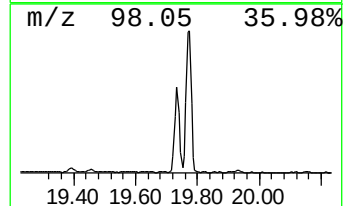
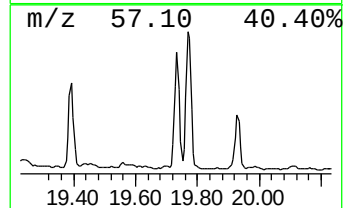
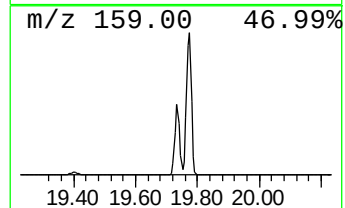
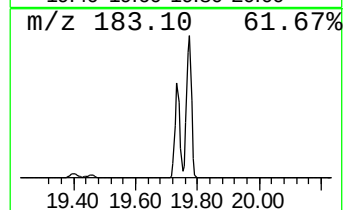
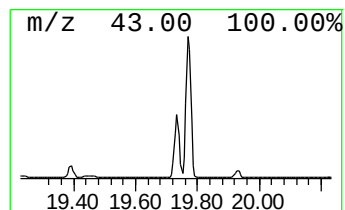
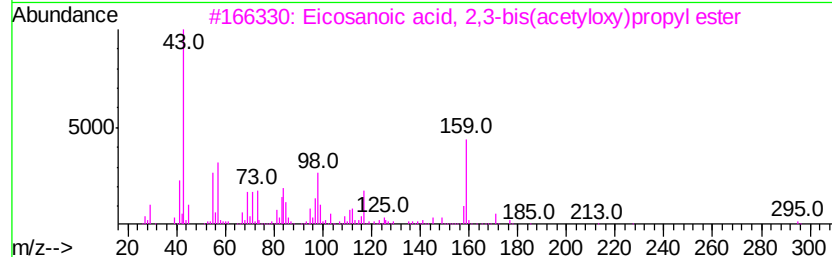
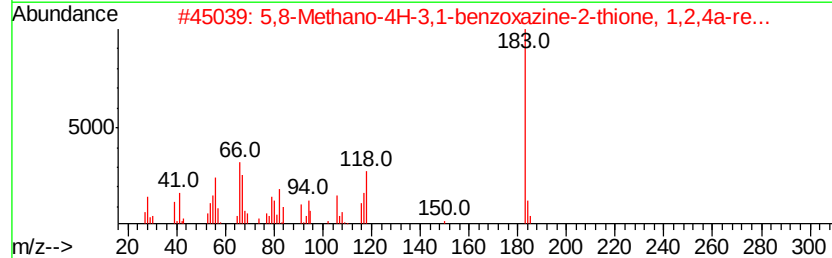
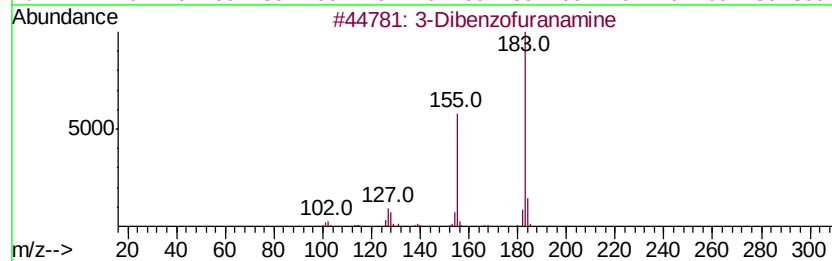
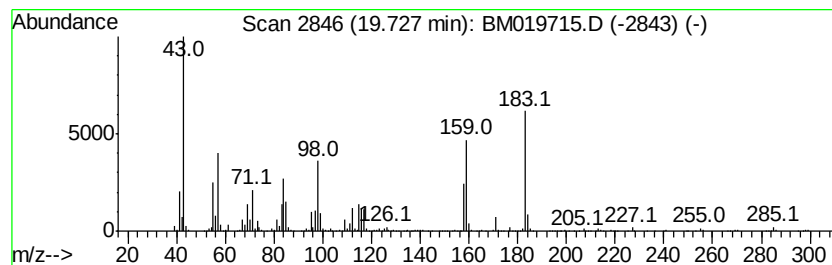
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 unknown-04 Concentration Rank 17

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.73 | 10.61 ng/ul | 1695590 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Dibenzofuranamine                 | 183 | C12H9NO   | 004106-66-5  | 38   |
| 2    |    | 5,8-Methano-4H-3,1-benzoxazine-2... | 183 | C9H13NOS  | 1000142-76-7 | 27   |
| 3    |    | Eicosanoic acid, 2,3-bis(acetylo... | 470 | C27H50O6  | 055429-67-9  | 25   |
| 4    |    | Eicosanoic acid, 2-(acetyloxy)-1... | 470 | C27H50O6  | 055429-68-0  | 22   |
| 5    |    | 4-Methyl-5-trifluoromethyl-4H-1,... | 183 | C4H4F3N3S | 030682-81-6  | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

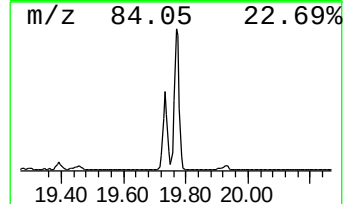
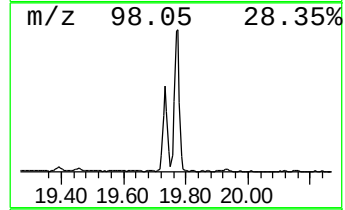
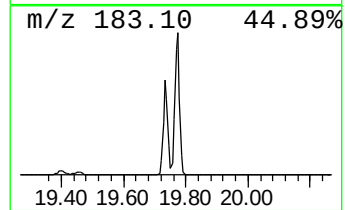
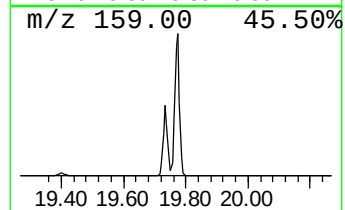
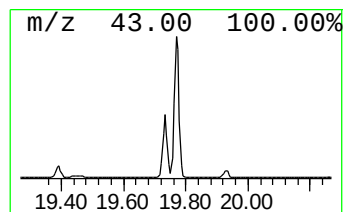
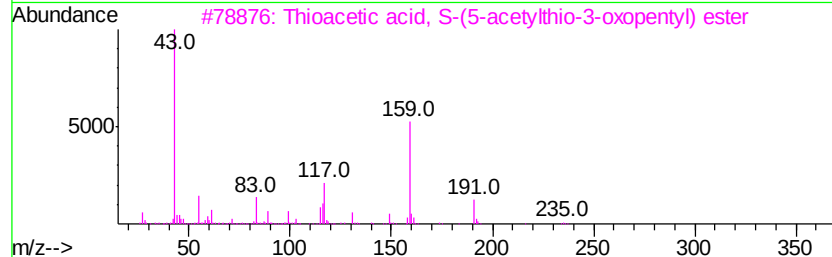
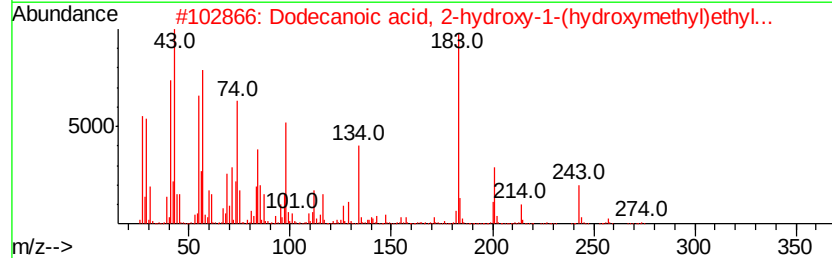
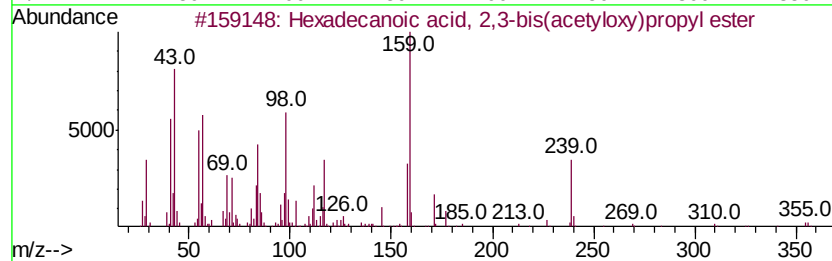
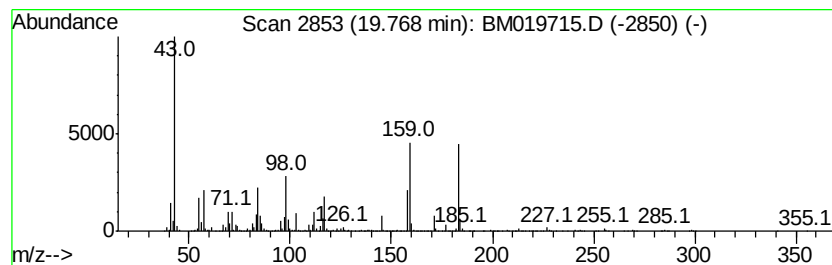
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 59 unknown-05 Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.77 | 20.37 ng/ul | 3256780 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Hexadecanoic acid, 2,3-bis(acety... | 414 | C23H42O6   | 055268-70-7  | 38   |
| 2    |    | Dodecanoic acid, 2-hydroxy-1-(hy... | 274 | C15H30O4   | 001678-45-1  | 35   |
| 3    |    | Thioacetic acid, S-(5-acetylthio... | 234 | C9H14O3S2  | 1000185-15-4 | 18   |
| 4    |    | 2,3,4-Trifluorobenzoic acid, 4-m... | 260 | C13H15F3O2 | 1000282-29-8 | 14   |
| 5    |    | Lauric anhydride                    | 382 | C24H46O3   | 000645-66-9  | 14   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

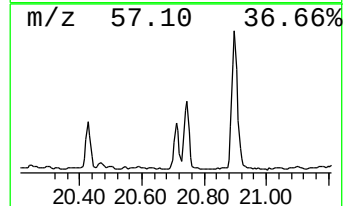
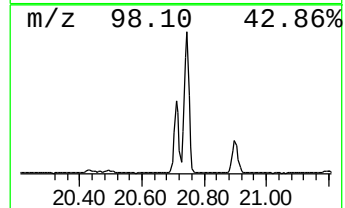
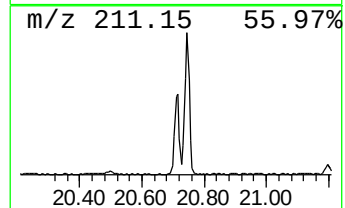
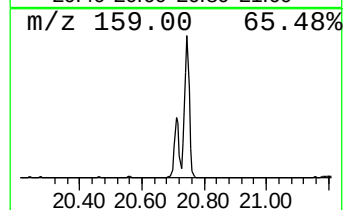
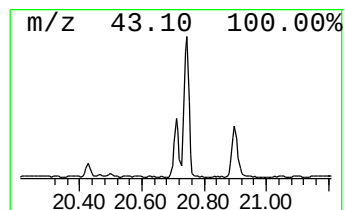
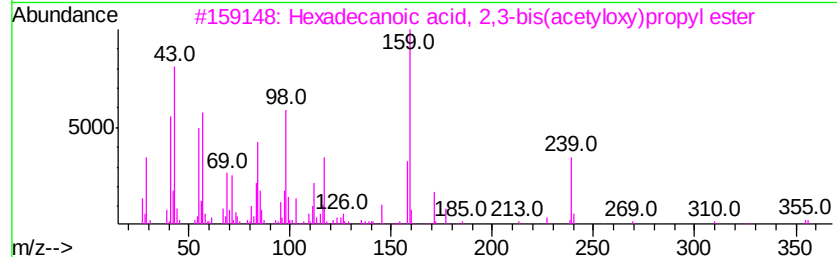
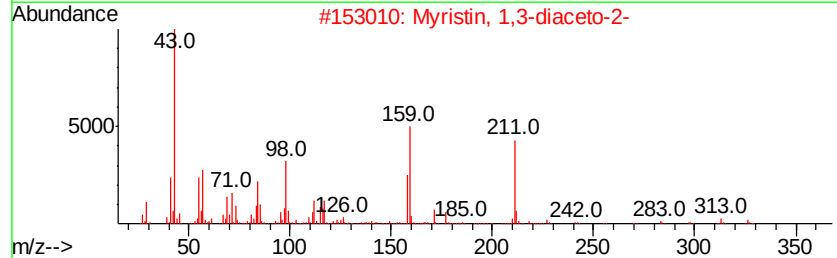
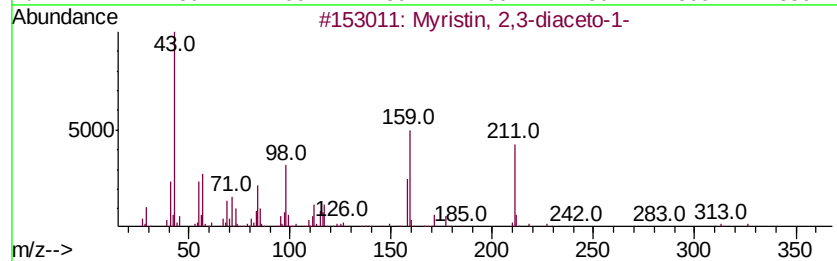
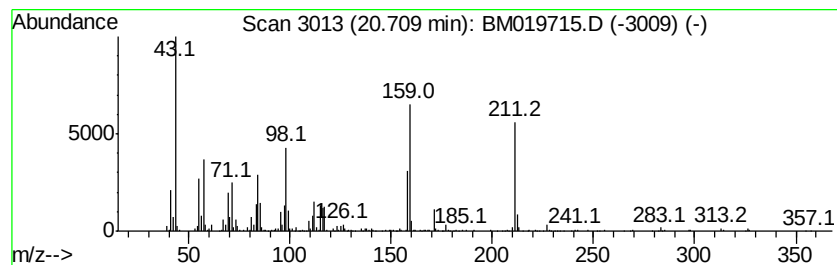
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 60 Myristin, 2,3-diaceto-1- Concentration Rank 48

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.71 | 5.07 ng/ul | 809892 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID                                      | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Myristin, 2,3-diaceto-1-                          | 386 | C21H38O6 | 014473-55-3 | 91   |
| 2    |    | Myristin, 1,3-diaceto-2-                          | 386 | C21H38O6 | 014290-23-4 | 91   |
| 3    |    | Hexadecanoic acid, 2,3-bis(acetyloxy)propyl ester | 414 | C23H42O6 | 055268-70-7 | 70   |
| 4    |    | Myristin, 2,3-diaceto-1-                          | 386 | C21H38O6 | 014473-55-3 | 38   |
| 5    |    | Eicosanoic acid, 2-(acetyloxy)-1-                 | 470 | C27H50O6 | 055429-68-0 | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

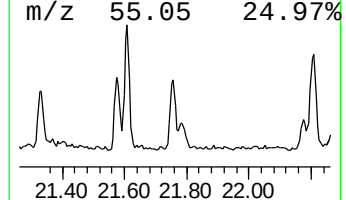
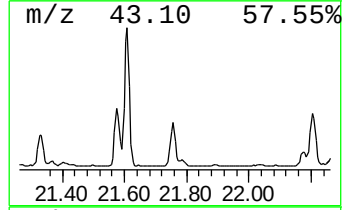
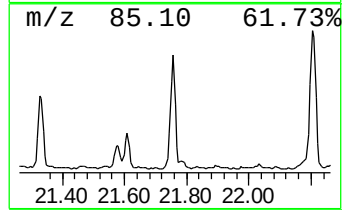
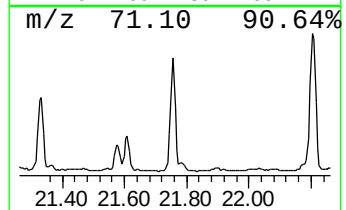
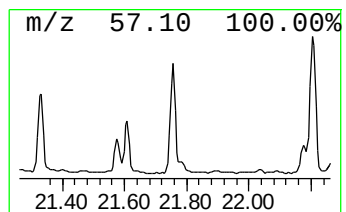
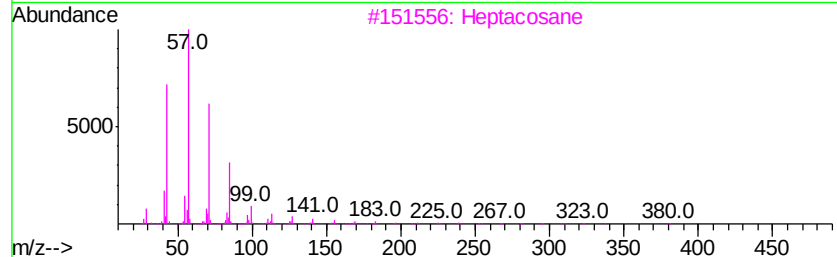
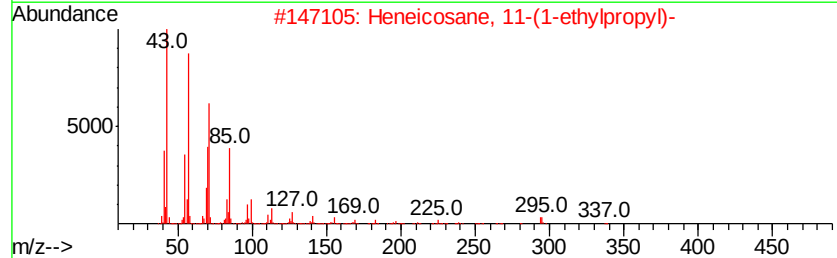
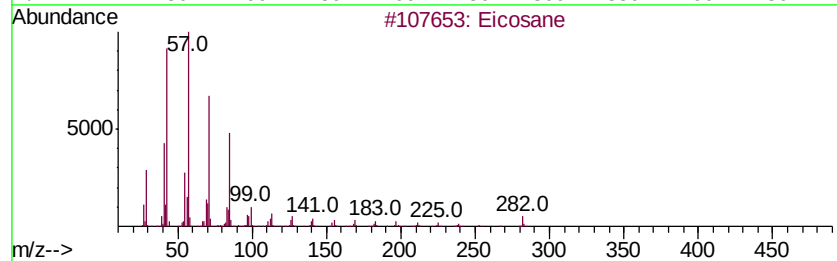
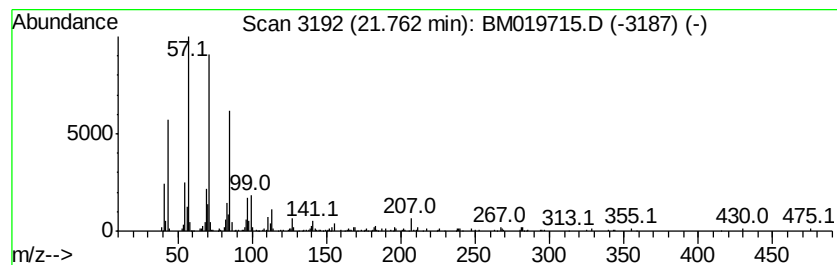
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 70

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.76 | 2.42 ng/ul | 387104 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane                         | 282 | C20H42  | 000112-95-8 | 90   |
| 2    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 87   |
| 3    |    | Heptacosane                      | 380 | C27H56  | 000593-49-7 | 87   |
| 4    |    | Tetradecane, 2,6,10-trimethyl-   | 240 | C17H36  | 014905-56-7 | 83   |
| 5    |    | Hexadecane, 7-methyl-            | 240 | C17H36  | 026730-20-1 | 81   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

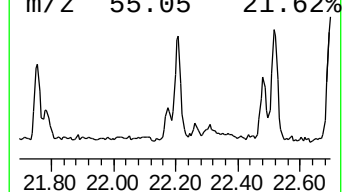
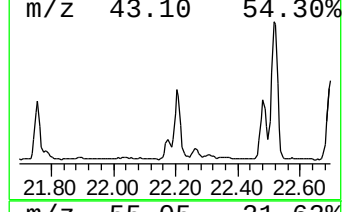
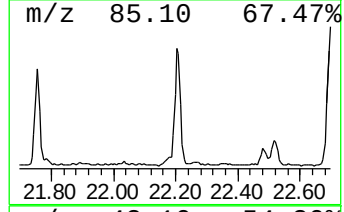
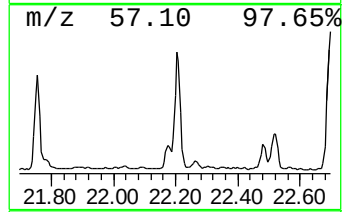
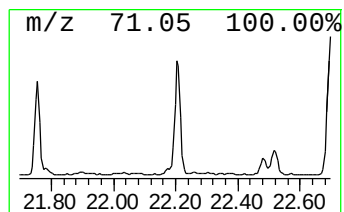
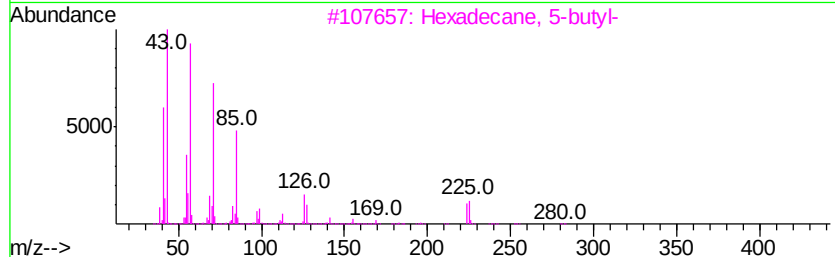
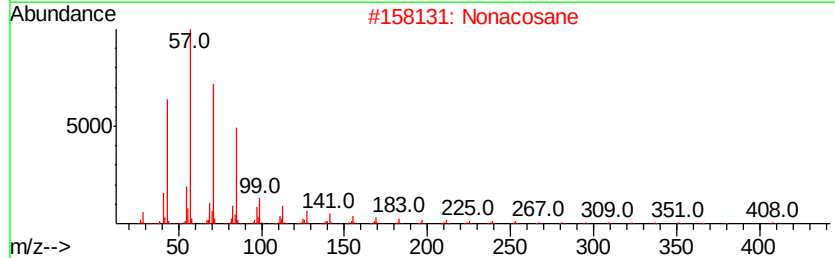
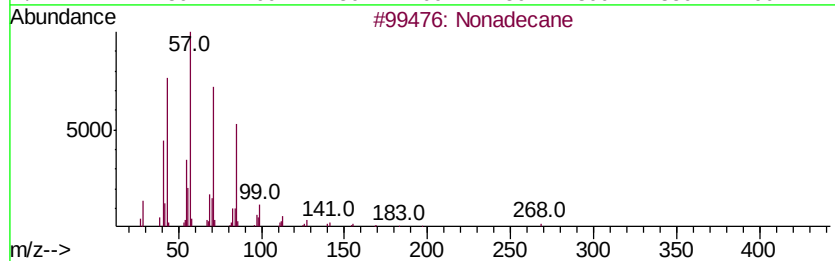
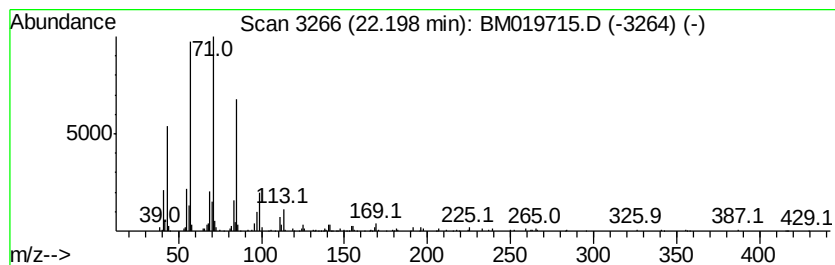
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 56

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.20 | 3.37 ng/ul | 538063 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 91   |
| 2    |    | Nonacosane                          | 408 | C29H60  | 000630-03-5 | 90   |
| 3    |    | Hexadecane, 5-butyl-                | 282 | C20H42  | 006912-07-8 | 90   |
| 4    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 78   |
| 5    |    | Tridecane, 6-propyl-                | 226 | C16H34  | 055045-10-8 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

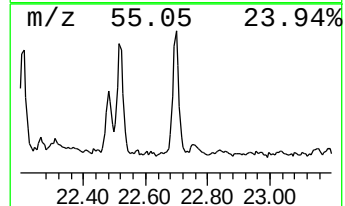
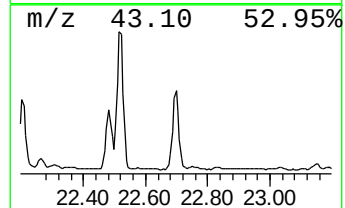
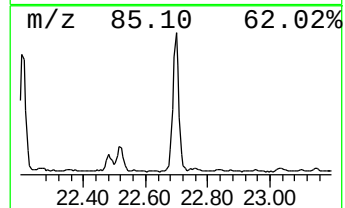
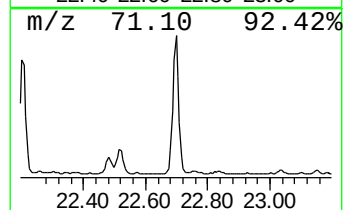
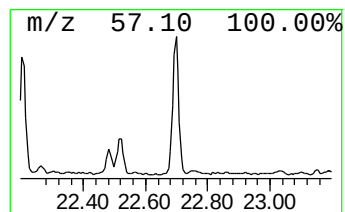
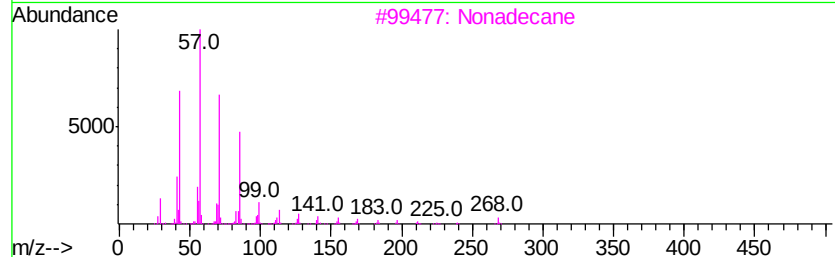
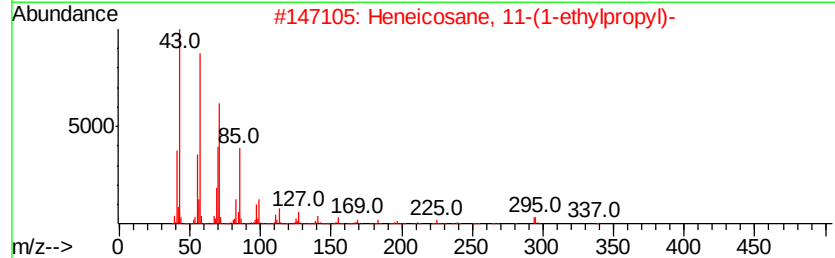
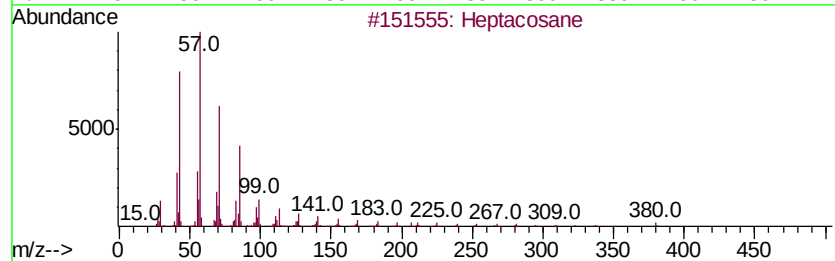
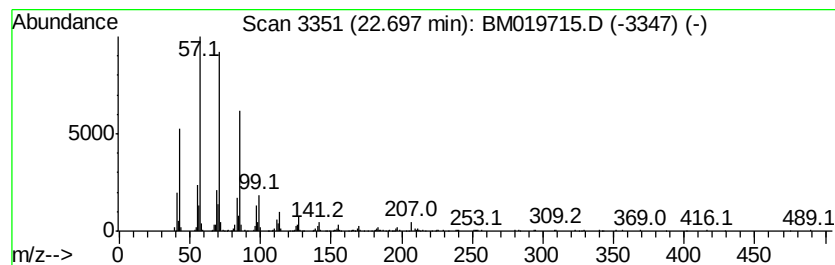
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.70 | 4.61 ng/ul | 701970 | Perylene-d12     | 23.40 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptacosane                      | 380 | C27H56  | 000593-49-7 | 87   |
| 2    |    |   | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 87   |
| 3    |    |   | Nonadecane                       | 268 | C19H40  | 000629-92-5 | 86   |
| 4    |    |   | Docosane, 5-butyl-               | 366 | C26H54  | 055282-16-1 | 86   |
| 5    |    |   | Octadecane                       | 254 | C18H38  | 000593-45-3 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

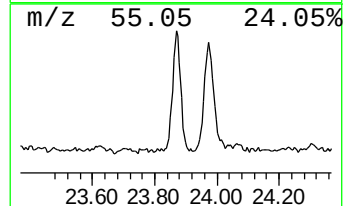
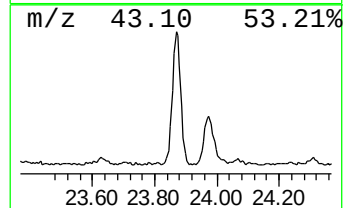
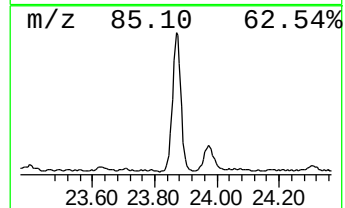
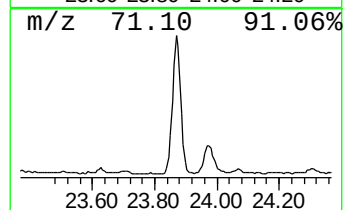
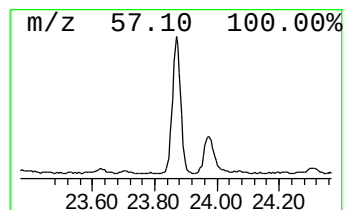
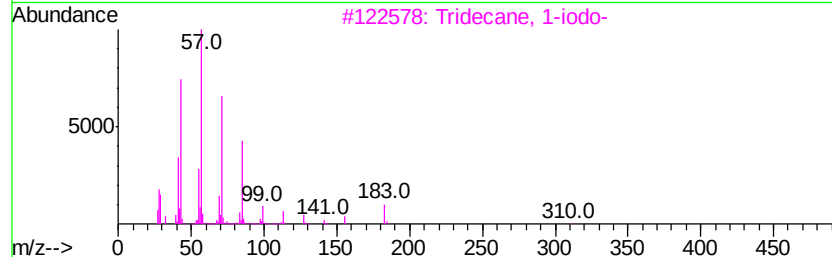
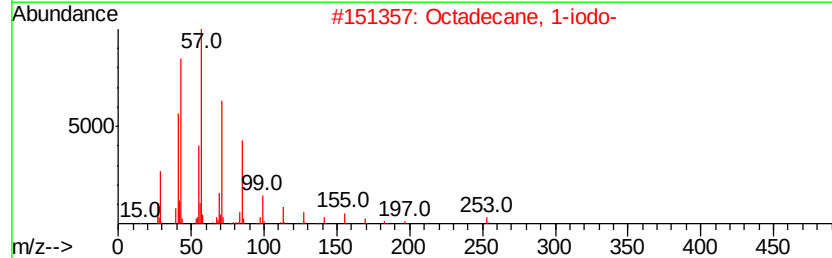
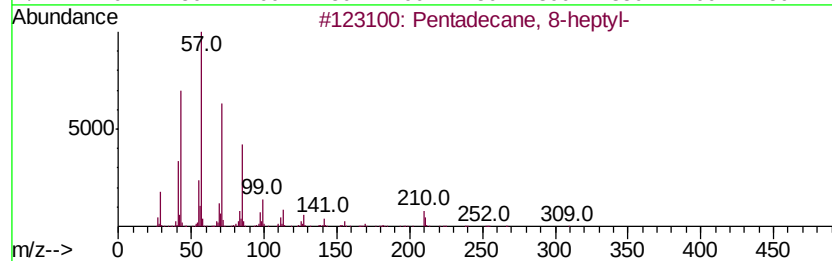
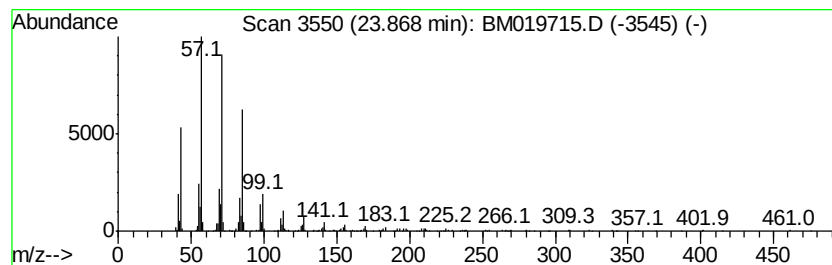
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.87 | 4.81 ng/ul | 733067 | Perylene-d12     | 23.40 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 90   |
| 2    |    |   | Octadecane, 1-iodo-    | 380 | C18H37I | 000629-93-6 | 87   |
| 3    |    |   | Tridecane, 1-iodo-     | 310 | C13H27I | 035599-77-0 | 87   |
| 4    |    |   | Tridecane, 6-propyl-   | 226 | C16H34  | 055045-10-8 | 87   |
| 5    |    |   | Dodecane               | 170 | C12H26  | 000112-40-3 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

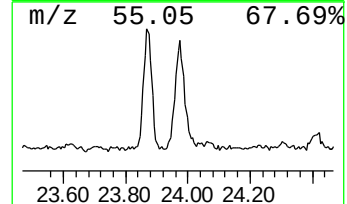
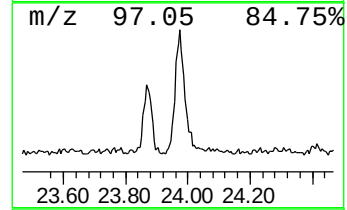
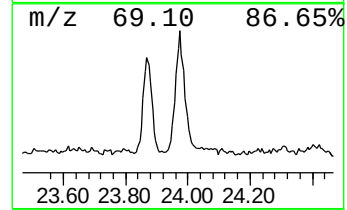
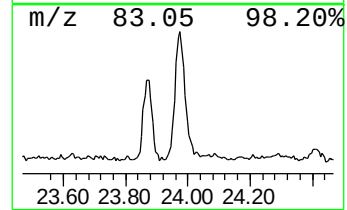
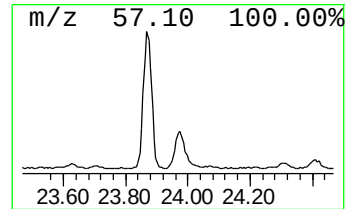
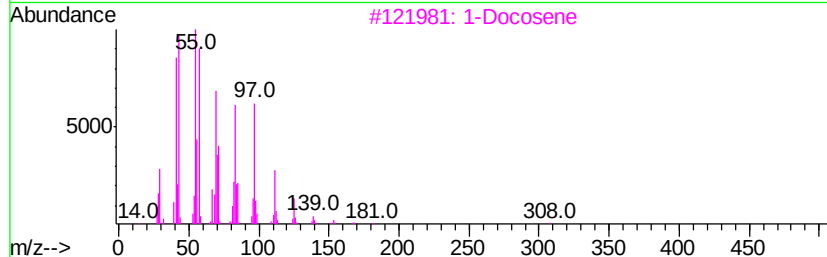
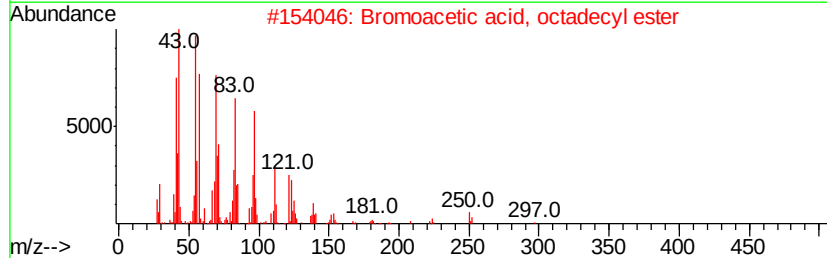
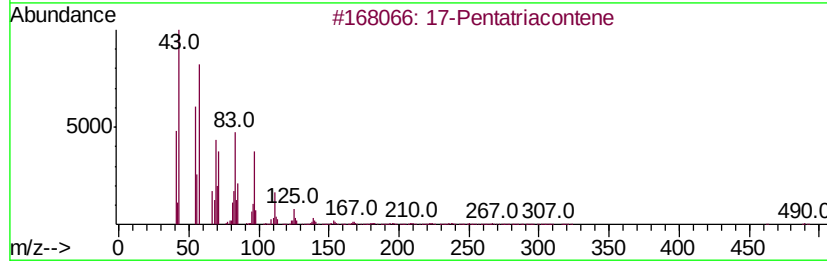
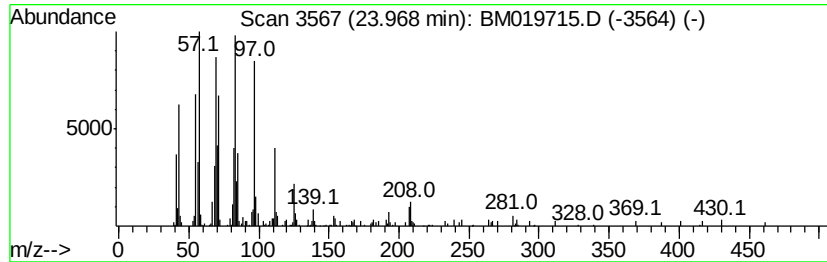
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 71 (DEL) Alkane: Cyclic23.97 Concentration Rank 61

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 23.97 | 2.94 ng/ul | 447346 | Perylene-d12     | 23.40 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0  | 91   |
| 2    |    | Bromoacetic acid, octadecyl ester   | 390 | C20H39BrO2  | 018992-03-5  | 91   |
| 3    |    | 1-Docosene                          | 308 | C22H44      | 001599-67-3  | 86   |
| 4    |    | Trichloroacetic acid, hexadecyl ... | 386 | C18H33Cl3O2 | 074339-54-1  | 83   |
| 5    |    | Dichloroacetic acid, tridecyl ester | 310 | C15H28Cl2O2 | 1000280-48-3 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040219\  
Data File : BM019715.D  
Acq On : 03 Apr 2019 05:41  
Operator : JU/SJ  
Sample : K2168-13  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
Quant Title : SVOA CALIBRATION

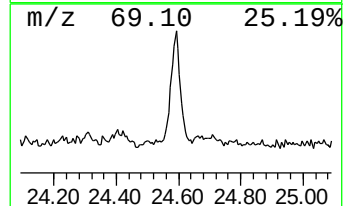
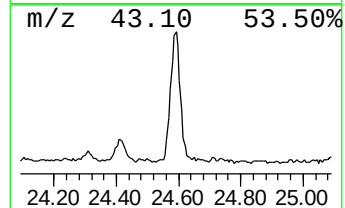
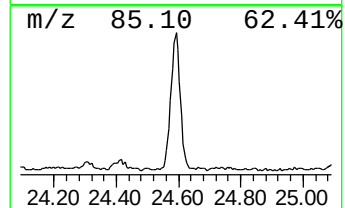
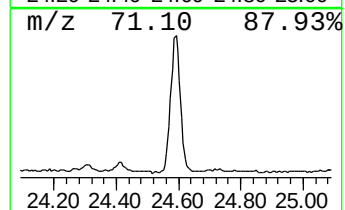
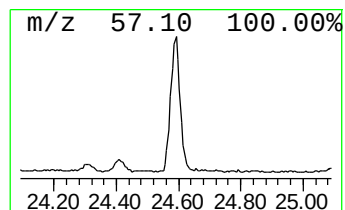
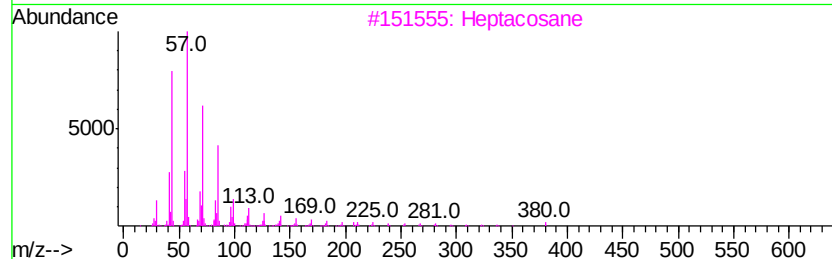
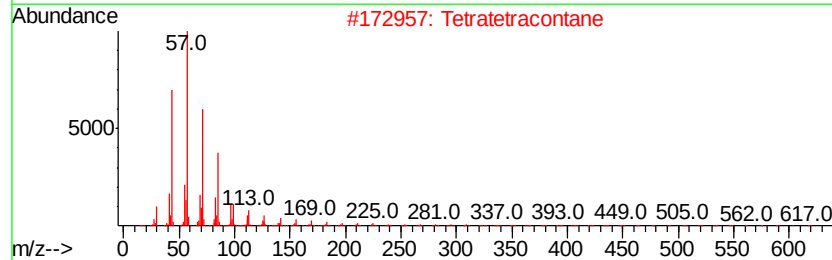
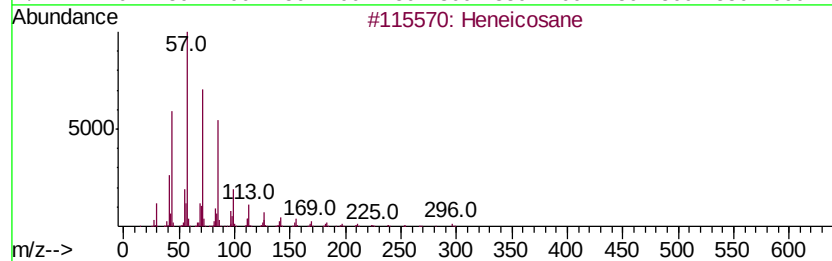
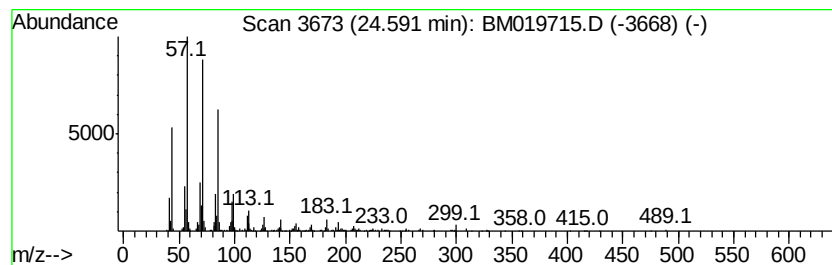
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 72 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 24.59 | 3.74 ng/ul | 570469 | Perylene-d12     | 23.40 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Heneicosane            | 296 | C21H44  | 000629-94-7 | 87   |
| 2    |    |   | Tetratetracontane      | 619 | C44H90  | 007098-22-8 | 83   |
| 3    |    |   | Heptacosane            | 380 | C27H56  | 000593-49-7 | 83   |
| 4    |    |   | Octacosane             | 394 | C28H58  | 000630-02-4 | 83   |
| 5    |    |   | Pentadecane, 8-heptyl- | 310 | C22H46  | 071005-15-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM040219\  
 Data File : BM019715.D  
 Acq On : 03 Apr 2019 05:41  
 Operator : JU/SJ  
 Sample : K2168-13  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF638

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM032219MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| (DEL) Alkane: Str... | 5.58  | 8.1     | ng/ul | 419571   | 1                     | 7.57  | 1032860 | 20.0 |
| (DEL) Alkane: Str... | 7.16  | 12.2    | ng/ul | 628030   | 1                     | 7.57  | 1032860 | 20.0 |
| (DEL) Alkane: Str... | 7.50  | 6.6     | ng/ul | 339015   | 1                     | 7.57  | 1032860 | 20.0 |
| 3a,6-Methano-3aH-... | 8.19  | 9.8     | ng/ul | 507717   | 1                     | 7.57  | 1032860 | 20.0 |
| Benzene, 1-methyl... | 8.65  | 10.6    | ng/ul | 546754   | 1                     | 7.57  | 1032860 | 20.0 |
| Benzene, 1-ethyl...  | 8.98  | 6.3     | ng/ul | 348786   | 2                     | 10.35 | 1113660 | 20.0 |
| Benzene, 1-ethyl...  | 9.16  | 6.4     | ng/ul | 357562   | 2                     | 10.35 | 1113660 | 20.0 |
| Benzene, 1-methyl... | 9.74  | 14.3    | ng/ul | 796024   | 2                     | 10.35 | 1113660 | 20.0 |
| Benzene, 1-ethyl...  | 9.92  | 5.3     | ng/ul | 296811   | 2                     | 10.35 | 1113660 | 20.0 |
| Benzene, 1-methyl... | 10.03 | 5.1     | ng/ul | 283508   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 10.29 | 37.7    | ng/ul | 2100840  | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Cyc... | 10.80 | 6.3     | ng/ul | 349946   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Cyc... | 11.02 | 6.6     | ng/ul | 367296   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Cyc... | 11.15 | 5.0     | ng/ul | 279268   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 11.23 | 10.4    | ng/ul | 577041   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 11.33 | 27.9    | ng/ul | 1554410  | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Cyc... | 11.60 | 6.3     | ng/ul | 348574   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 11.75 | 49.2    | ng/ul | 2739840  | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 11.97 | 7.8     | ng/ul | 432755   | 2                     | 10.35 | 1113660 | 20.0 |
| Naphthalene, 1-me... | 12.25 | 8.6     | ng/ul | 480429   | 2                     | 10.35 | 1113660 | 20.0 |
| (DEL) Alkane: Str... | 12.39 | 4.6     | ng/ul | 522198   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Cyc... | 12.46 | 4.7     | ng/ul | 530255   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 12.58 | 2.7     | ng/ul | 308269   | 3                     | 14.23 | 2274550 | 20.0 |
| Dodecane, 1-iodo-    | 12.67 | 2.9     | ng/ul | 330916   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 12.72 | 14.7    | ng/ul | 1675200  | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 13.02 | 37.2    | ng/ul | 4233160  | 3                     | 14.23 | 2274550 | 20.0 |
| unknown-01           | 13.25 | 6.4     | ng/ul | 729088   | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 2,6-... | 13.38 | 5.1     | ng/ul | 578559   | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 2,3-... | 13.54 | 12.2    | ng/ul | 1392560  | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 1,6-... | 13.60 | 11.0    | ng/ul | 1248850  | 3                     | 14.23 | 2274550 | 20.0 |
| unknown-02           | 14.03 | 6.0     | ng/ul | 687488   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 14.11 | 38.7    | ng/ul | 4402870  | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 14.17 | 5.1     | ng/ul | 577910   | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 1,4,... | 14.62 | 8.1     | ng/ul | 915618   | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 1,6,... | 14.70 | 5.4     | ng/ul | 612146   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 14.73 | 7.9     | ng/ul | 899171   | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 2,3,... | 14.85 | 6.3     | ng/ul | 722645   | 3                     | 14.23 | 2274550 | 20.0 |
| unknown-03           | 14.90 | 4.5     | ng/ul | 510660   | 3                     | 14.23 | 2274550 | 20.0 |
| Azulene, 4,6,8-tr... | 15.00 | 3.2     | ng/ul | 366263   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 15.07 | 31.5    | ng/ul | 3584150  | 3                     | 14.23 | 2274550 | 20.0 |
| Naphthalene, 2-me... | 15.33 | 2.9     | ng/ul | 333546   | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 15.47 | 18.5    | ng/ul | 2103450  | 3                     | 14.23 | 2274550 | 20.0 |
| (DEL) Alkane: Str... | 15.63 | 3.1     | ng/ul | 416572   | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Cyc... | 15.69 | 3.1     | ng/ul | 410221   | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 15.96 | 55.4    | ng/ul | 7423740  | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 16.28 | 2.7     | ng/ul | 358274   | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 16.73 | 15.9    | ng/ul | 2137650  | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 16.77 | 8.3     | ng/ul | 1106910  | 4                     | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 17.27 | 2.6     | ng/ul | 355258   | 4                     | 16.99 | 2681030 | 20.0 |

|                      |       |            |         |   |       |         |      |
|----------------------|-------|------------|---------|---|-------|---------|------|
| (DEL) Alkane: Str... | 17.39 | 2.6 ng/ul  | 352696  | 4 | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 17.47 | 11.1 ng/ul | 1485280 | 4 | 16.99 | 2681030 | 20.0 |
| n-Hexadecanoic acid  | 17.88 | 8.0 ng/ul  | 1076290 | 4 | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 18.16 | 6.7 ng/ul  | 896556  | 4 | 16.99 | 2681030 | 20.0 |
| (DEL) Alkane: Str... | 18.80 | 5.4 ng/ul  | 728971  | 4 | 16.99 | 2681030 | 20.0 |
| Octadecanoic acid    | 19.17 | 3.4 ng/ul  | 536441  | 5 | 21.20 | 3197390 | 20.0 |
| unknown-04           | 19.73 | 10.6 ng/ul | 1695590 | 5 | 21.20 | 3197390 | 20.0 |
| unknown-05           | 19.77 | 20.4 ng/ul | 3256780 | 5 | 21.20 | 3197390 | 20.0 |
| Myristin, 2,3-dia... | 20.71 | 5.1 ng/ul  | 809892  | 5 | 21.20 | 3197390 | 20.0 |
| (DEL) Alkane: Str... | 21.76 | 2.4 ng/ul  | 387104  | 5 | 21.20 | 3197390 | 20.0 |
| (DEL) Alkane: Str... | 22.20 | 3.4 ng/ul  | 538063  | 5 | 21.20 | 3197390 | 20.0 |
| (DEL) Alkane: Str... | 22.70 | 4.6 ng/ul  | 701970  | 6 | 23.40 | 3047070 | 20.0 |
| (DEL) Alkane: Str... | 23.87 | 4.8 ng/ul  | 733067  | 6 | 23.40 | 3047070 | 20.0 |
| (DEL) Alkane: Cyc... | 23.97 | 2.9 ng/ul  | 447346  | 6 | 23.40 | 3047070 | 20.0 |
| (DEL) Alkane: Str... | 24.59 | 3.7 ng/ul  | 570469  | 6 | 23.40 | 3047070 | 20.0 |

Instrument :

BNA\_M

ClientSampleId :

BF638