

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
 Data File : BM024586.D  
 Acq On : 22 Jan 2020 17:46  
 Operator : JU  
 Sample : L1118-11DL2 20X  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GASK2DL2

## 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-BM011520MA.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         | 3.757       | 144           | 148         | 152          | rBV      | 49484          | 63137         | 1.33%           | 0.075%        |
| 2         | 5.028       | 359           | 364         | 370          | rVB      | 33369          | 53773         | 1.13%           | 0.064%        |
| 3         | 5.151       | 377           | 385         | 396          | rBV      | 160181         | 332638        | 7.02%           | 0.395%        |
| 4         | 5.510       | 435           | 446         | 456          | rVV3     | 112002         | 288076        | 6.08%           | 0.342%        |
| 5         | 6.569       | 619           | 626         | 631          | rBV      | 62866          | 99495         | 2.10%           | 0.118%        |
| 6         | 6.628       | 631           | 636         | 643          | rVV2     | 36816          | 88243         | 1.86%           | 0.105%        |
| 7         | 6.698       | 644           | 648         | 654          | rVB2     | 72538          | 100717        | 2.13%           | 0.120%        |
| 8         | 6.992       | 689           | 698         | 704          | rVB6     | 27757          | 62163         | 1.31%           | 0.074%        |
| 9         | 7.116       | 712           | 719         | 726          | rBV2     | 365533         | 589656        | 12.44%          | 0.700%        |
| 10        | 7.198       | 731           | 733         | 739          | rVB2     | 46398          | 68936         | 1.45%           | 0.082%        |
| 11        | 7.451       | 770           | 776         | 784          | rBV      | 776162         | 1204560       | 25.42%          | 1.430%        |
| 12        | 7.569       | 790           | 796         | 803          | rVV      | 69968          | 111874        | 2.36%           | 0.133%        |
| 13        | 7.822       | 834           | 839         | 847          | rVB      | 35328          | 54166         | 1.14%           | 0.064%        |
| 14        | 7.981       | 858           | 866         | 878          | rBV      | 430641         | 720647        | 15.21%          | 0.856%        |
| 15        | 8.098       | 882           | 886         | 891          | rVB2     | 40405          | 54393         | 1.15%           | 0.065%        |
| 16        | 8.163       | 891           | 897         | 903          | rVB6     | 28825          | 59978         | 1.27%           | 0.071%        |
| 17        | 8.557       | 958           | 964         | 968          | rBV3     | 51688          | 108650        | 2.29%           | 0.129%        |
| 18        | 8.704       | 983           | 989         | 999          | rBV2     | 116507         | 221609        | 4.68%           | 0.263%        |
| 19        | 8.916       | 1014          | 1025        | 1029         | rBV2     | 36199          | 91791         | 1.94%           | 0.109%        |
| 20        | 9.128       | 1056          | 1061        | 1070         | rVB2     | 45441          | 95601         | 2.02%           | 0.114%        |
| 21        | 9.651       | 1137          | 1150        | 1156         | rBV3     | 156851         | 408764        | 8.63%           | 0.485%        |
| 22        | 9.722       | 1156          | 1162        | 1166         | rBV      | 105328         | 175244        | 3.70%           | 0.208%        |
| 23        | 9.839       | 1178          | 1182        | 1192         | rVB5     | 39921          | 78757         | 1.66%           | 0.094%        |
| 24        | 10.210      | 1237          | 1245        | 1249         | rVV      | 1140375        | 2006874       | 42.35%          | 2.383%        |
| 25        | 10.263      | 1249          | 1254        | 1261         | rVV      | 2353932        | 3968155       | 83.74%          | 4.712%        |
| 26        | 10.375      | 1265          | 1273        | 1277         | rVV2     | 61488          | 119196        | 2.52%           | 0.142%        |
| 27        | 11.280      | 1421          | 1427        | 1438         | rVB6     | 41769          | 135485        | 2.86%           | 0.161%        |
| 28        | 11.410      | 1444          | 1449        | 1461         | rVB6     | 34387          | 78440         | 1.66%           | 0.093%        |
| 29        | 11.686      | 1491          | 1496        | 1501         | rVV      | 165607         | 257205        | 5.43%           | 0.305%        |
| 30        | 11.739      | 1501          | 1505        | 1516         | rVV10    | 29278          | 99391         | 2.10%           | 0.118%        |
| 31        | 11.886      | 1523          | 1530        | 1542         | rVB      | 1154553        | 1882870       | 39.73%          | 2.236%        |
| 32        | 12.110      | 1557          | 1568        | 1573         | rBV      | 549313         | 896815        | 18.92%          | 1.065%        |
| 33        | 12.669      | 1659          | 1663        | 1667         | rVB2     | 71599          | 96046         | 2.03%           | 0.114%        |
| 34        | 12.922      | 1701          | 1706        | 1709         | rVV      | 113876         | 171023        | 3.61%           | 0.203%        |

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 ClientSampleId :  
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Integrator: RTE  
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 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-BM011520MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 12.963 | 1709 | 1713 | 1722 | rVV  | 179387  | 306233  | 6.46%  | 0.364% |
| 36 | 13.057 | 1722 | 1729 | 1733 | rVV9 | 31146   | 76504   | 1.61%  | 0.091% |
| 37 | 13.121 | 1734 | 1740 | 1749 | rVB  | 98419   | 222554  | 4.70%  | 0.264% |
| 38 | 13.263 | 1758 | 1764 | 1765 | rBV  | 190095  | 316367  | 6.68%  | 0.376% |
| 39 | 13.416 | 1785 | 1790 | 1795 | rBV  | 245443  | 408896  | 8.63%  | 0.486% |
| 40 | 13.469 | 1795 | 1799 | 1804 | rBV  | 163386  | 238751  | 5.04%  | 0.284% |
| 41 | 13.545 | 1805 | 1812 | 1819 | rVB4 | 188821  | 394230  | 8.32%  | 0.468% |
| 42 | 13.633 | 1820 | 1827 | 1842 | rVB2 | 390284  | 909005  | 19.18% | 1.079% |
| 43 | 13.816 | 1849 | 1858 | 1867 | rBV  | 682630  | 1289278 | 27.21% | 1.531% |
| 44 | 13.945 | 1876 | 1880 | 1883 | rBV3 | 67516   | 97074   | 2.05%  | 0.115% |
| 45 | 13.992 | 1885 | 1888 | 1892 | rVV6 | 65629   | 92214   | 1.95%  | 0.110% |
| 46 | 14.051 | 1893 | 1898 | 1901 | rVV  | 484342  | 630948  | 13.31% | 0.749% |
| 47 | 14.098 | 1901 | 1906 | 1913 | rVV  | 1874586 | 2866975 | 60.50% | 3.404% |
| 48 | 14.157 | 1913 | 1916 | 1921 | rVV4 | 96733   | 157263  | 3.32%  | 0.187% |
| 49 | 14.321 | 1940 | 1944 | 1950 | rBV7 | 78690   | 169701  | 3.58%  | 0.202% |
| 50 | 14.498 | 1970 | 1974 | 1981 | rVB4 | 174800  | 353779  | 7.47%  | 0.420% |
| 51 | 14.563 | 1981 | 1985 | 1992 | rBV4 | 214619  | 420304  | 8.87%  | 0.499% |
| 52 | 14.674 | 2000 | 2004 | 2008 | rBV  | 380496  | 486503  | 10.27% | 0.578% |
| 53 | 14.716 | 2008 | 2011 | 2013 | rBV2 | 74564   | 103204  | 2.18%  | 0.123% |
| 54 | 14.780 | 2020 | 2022 | 2030 | rVB4 | 95218   | 136675  | 2.88%  | 0.162% |
| 55 | 14.910 | 2040 | 2044 | 2051 | rVB  | 242012  | 342909  | 7.24%  | 0.407% |
| 56 | 14.980 | 2052 | 2056 | 2057 | rBV2 | 165434  | 213372  | 4.50%  | 0.253% |
| 57 | 15.010 | 2057 | 2061 | 2065 | rVB  | 1266860 | 1515005 | 31.97% | 1.799% |
| 58 | 15.098 | 2073 | 2076 | 2081 | rVB3 | 241101  | 321780  | 6.79%  | 0.382% |
| 59 | 15.151 | 2082 | 2085 | 2093 | rVV  | 196768  | 344891  | 7.28%  | 0.410% |
| 60 | 15.421 | 2123 | 2131 | 2135 | rBV  | 1659088 | 2643709 | 55.79% | 3.139% |
| 61 | 15.568 | 2152 | 2156 | 2160 | rBV2 | 323805  | 450333  | 9.50%  | 0.535% |
| 62 | 15.615 | 2161 | 2164 | 2165 | rBV2 | 175674  | 185322  | 3.91%  | 0.220% |
| 63 | 15.745 | 2182 | 2186 | 2190 | rBV3 | 186846  | 284666  | 6.01%  | 0.338% |
| 64 | 15.874 | 2204 | 2208 | 2210 | rBV  | 2305768 | 2831404 | 59.75% | 3.362% |
| 65 | 15.904 | 2210 | 2213 | 2221 | rVB  | 2559014 | 3515878 | 74.19% | 4.175% |
| 66 | 16.221 | 2263 | 2267 | 2274 | rBV5 | 327474  | 842787  | 17.78% | 1.001% |
| 67 | 16.280 | 2274 | 2277 | 2283 | rVB  | 457812  | 615651  | 12.99% | 0.731% |
| 68 | 16.333 | 2284 | 2286 | 2289 | rBV2 | 124593  | 122257  | 2.58%  | 0.145% |
| 69 | 16.668 | 2338 | 2343 | 2347 | rBV  | 2417635 | 2994183 | 63.18% | 3.556% |
| 70 | 16.721 | 2347 | 2352 | 2362 | rVB  | 1872111 | 2895582 | 61.10% | 3.438% |
| 71 | 16.851 | 2369 | 2374 | 2378 | rBV  | 2162258 | 2896910 | 61.13% | 3.440% |

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |        |
|-----|--------|------|------|------|------|---------|---------|---------|--------|
| 72  | 16.892 | 2378 | 2381 | 2388 | rVV2 | 835745  | 1318617 | 27.83%  | 1.566% |
| 73  | 16.980 | 2392 | 2396 | 2404 | rVB2 | 413364  | 756863  | 15.97%  | 0.899% |
| 74  | 17.133 | 2419 | 2422 | 2427 | rBV2 | 250413  | 373277  | 7.88%   | 0.443% |
| 75  | 17.204 | 2431 | 2434 | 2439 | rVB  | 323773  | 357980  | 7.55%   | 0.425% |
| 76  | 17.257 | 2440 | 2443 | 2448 | rBV  | 193484  | 279988  | 5.91%   | 0.332% |
| 77  | 17.327 | 2451 | 2455 | 2460 | rVV  | 626780  | 901528  | 19.02%  | 1.071% |
| 78  | 17.404 | 2464 | 2468 | 2474 | rVB  | 2558621 | 2927812 | 61.78%  | 3.477% |
| 79  | 17.680 | 2512 | 2515 | 2519 | rBV2 | 229844  | 320243  | 6.76%   | 0.380% |
| 80  | 17.851 | 2541 | 2544 | 2550 | rVB3 | 316773  | 509482  | 10.75%  | 0.605% |
| 81  | 17.915 | 2551 | 2555 | 2559 | rBV5 | 297462  | 666991  | 14.07%  | 0.792% |
| 82  | 18.098 | 2582 | 2586 | 2592 | rBV  | 2241541 | 2577840 | 54.40%  | 3.061% |
| 83  | 18.568 | 2662 | 2666 | 2671 | rVB  | 514550  | 642164  | 13.55%  | 0.763% |
| 84  | 18.739 | 2691 | 2695 | 2700 | rVB  | 2033107 | 2318213 | 48.92%  | 2.753% |
| 85  | 18.915 | 2720 | 2725 | 2730 | rBV2 | 578492  | 933446  | 19.70%  | 1.108% |
| 86  | 19.286 | 2784 | 2788 | 2791 | rVB2 | 517574  | 663717  | 14.01%  | 0.788% |
| 87  | 19.327 | 2791 | 2795 | 2801 | rVB  | 1683502 | 1905001 | 40.20%  | 2.262% |
| 88  | 19.862 | 2883 | 2886 | 2893 | rBV  | 1380769 | 1646602 | 34.75%  | 1.955% |
| 89  | 20.086 | 2921 | 2924 | 2928 | rVB  | 435208  | 513407  | 10.83%  | 0.610% |
| 90  | 20.362 | 2968 | 2971 | 2982 | rVB2 | 1094319 | 1347411 | 28.43%  | 1.600% |
| 91  | 20.827 | 3047 | 3050 | 3059 | rVB2 | 1088870 | 1780645 | 37.58%  | 2.114% |
| 92  | 21.062 | 3085 | 3090 | 3103 | rVB2 | 3090421 | 4738840 | 100.00% | 5.627% |
| 93  | 21.268 | 3122 | 3125 | 3129 | rVB  | 719250  | 750473  | 15.84%  | 0.891% |
| 94  | 21.686 | 3193 | 3196 | 3201 | rVB  | 542195  | 615736  | 12.99%  | 0.731% |
| 95  | 22.127 | 3268 | 3271 | 3275 | rVB  | 388885  | 444253  | 9.37%   | 0.528% |
| 96  | 22.715 | 3367 | 3371 | 3376 | rBV  | 436789  | 617890  | 13.04%  | 0.734% |
| 97  | 22.997 | 3415 | 3419 | 3424 | rVB  | 925026  | 1457768 | 30.76%  | 1.731% |
| 98  | 23.086 | 3430 | 3434 | 3439 | rVB  | 585732  | 965798  | 20.38%  | 1.147% |
| 99  | 23.180 | 3445 | 3450 | 3455 | rVB  | 1761419 | 2778071 | 58.62%  | 3.299% |
| 100 | 25.244 | 3796 | 3801 | 3813 | rVB  | 611030  | 1565131 | 33.03%  | 1.859% |

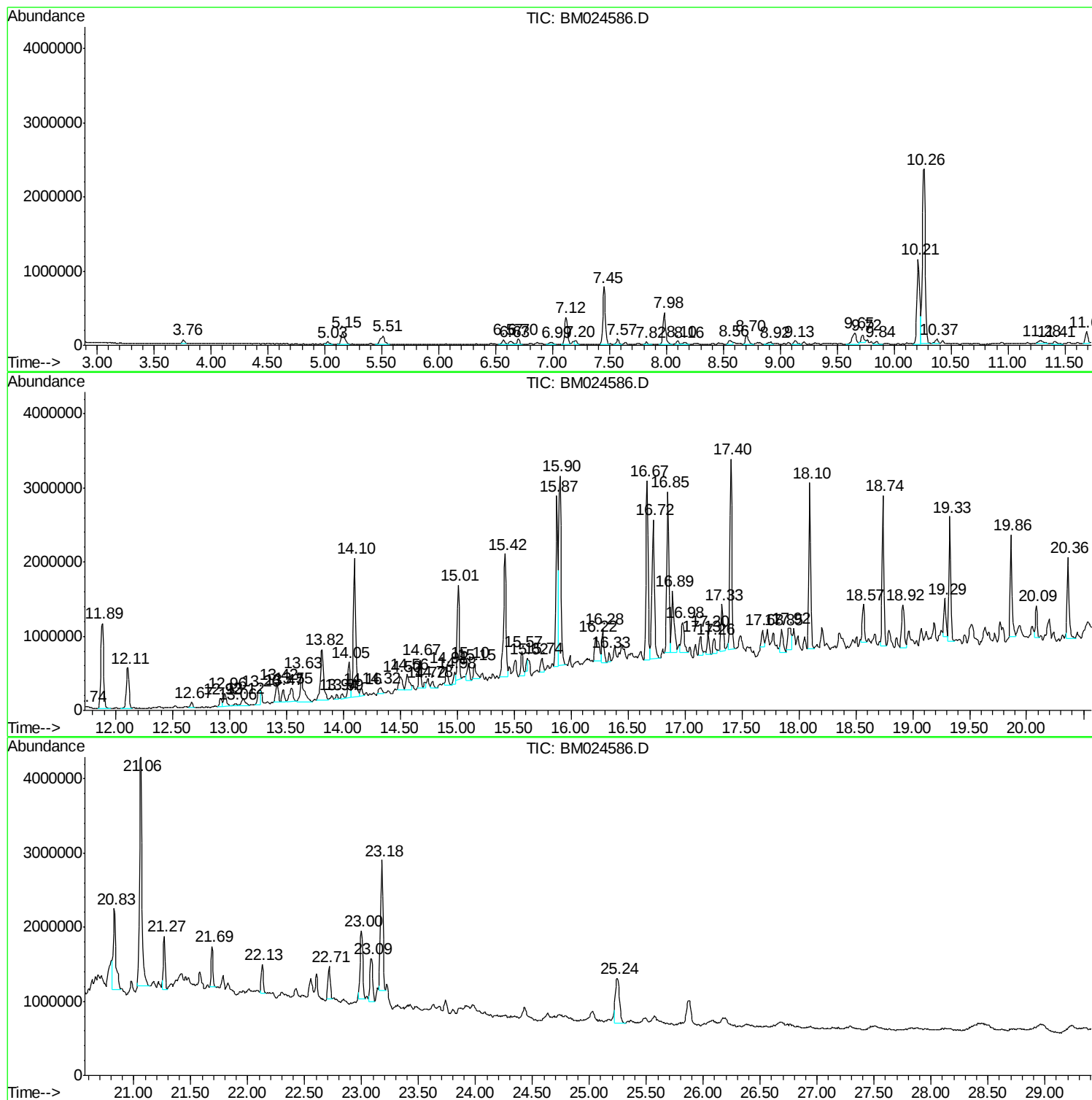
Sum of corrected areas: 84212652

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

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



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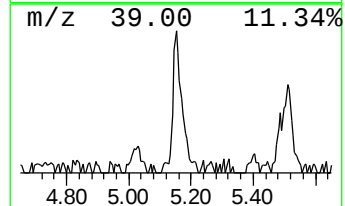
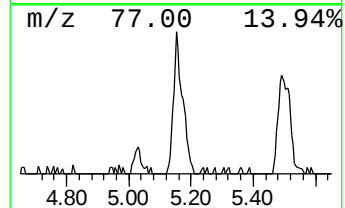
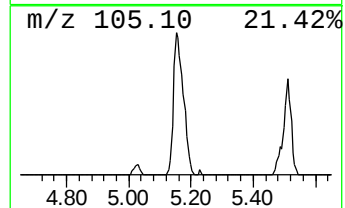
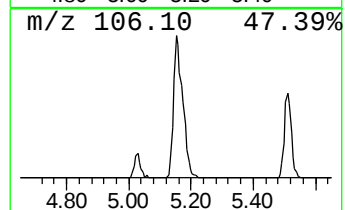
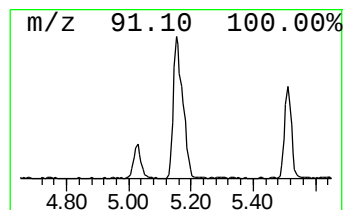
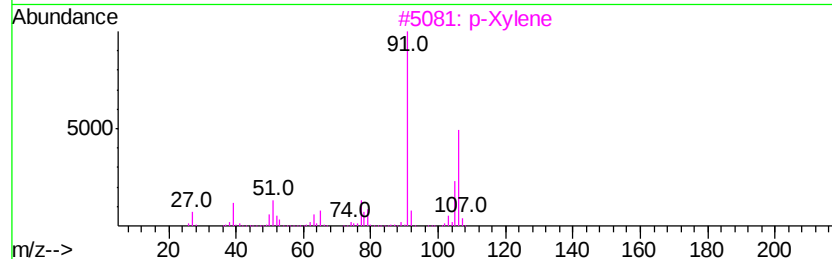
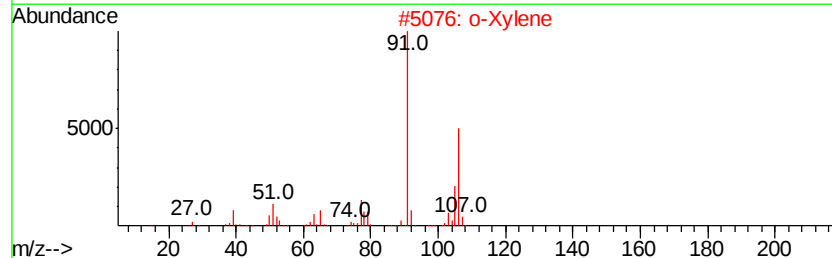
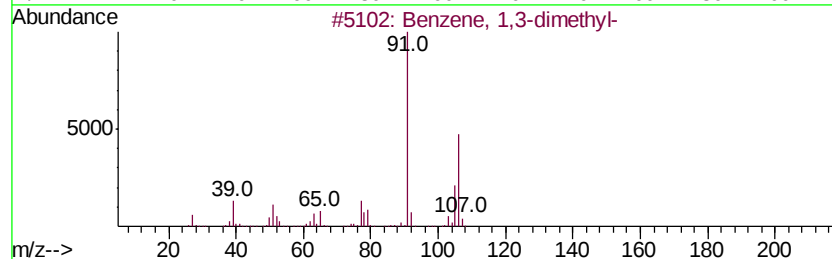
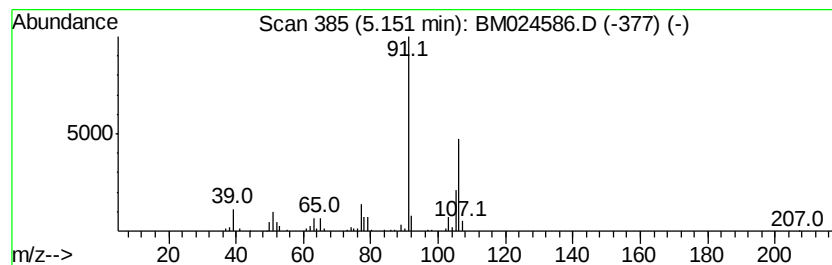
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 5.15 | 5.52 ng/ul | 332638 | 1,4-Dichlorobenzene-d4 | 7.45 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 2    |    |   | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |
| 3    |    |   | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |
| 4    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 5    |    |   | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 95   |



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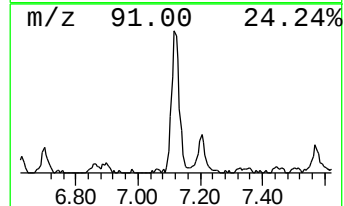
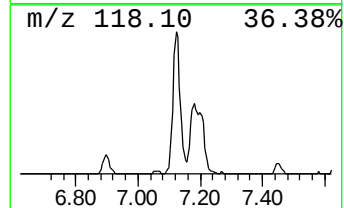
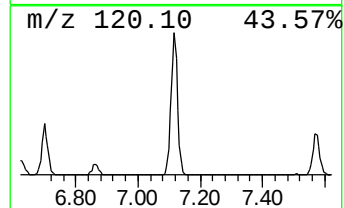
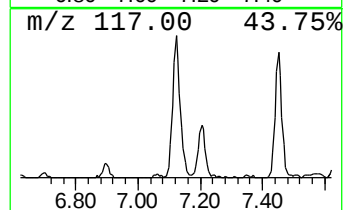
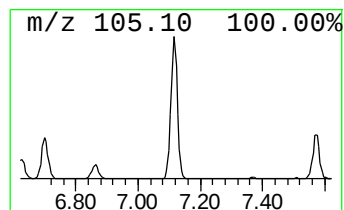
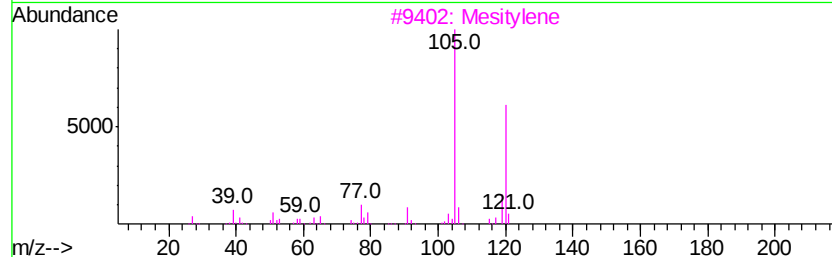
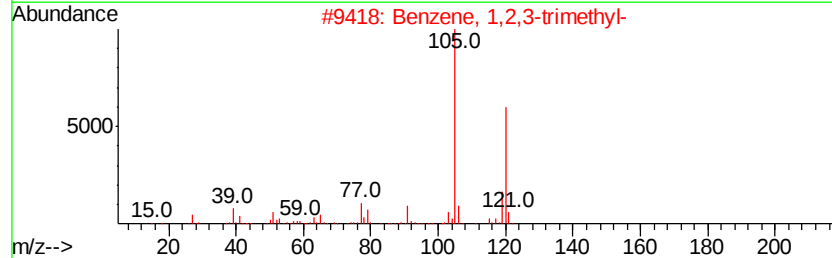
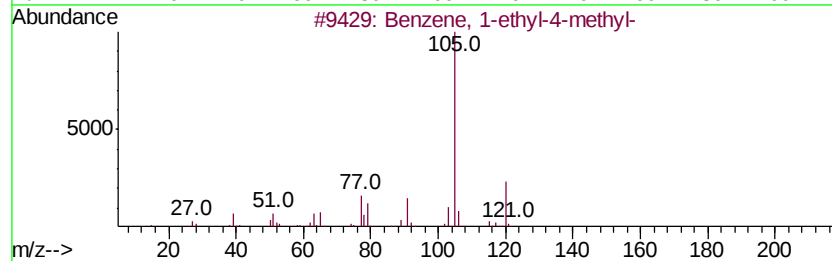
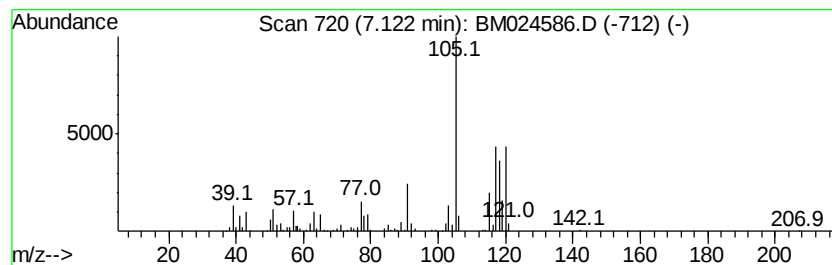
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 12

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.12 | 9.79 ng/ul | 589656 | 1,4-Dichlorobenzene-d4 | 7.45 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 87   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 64   |
| 3    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 64   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 60   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

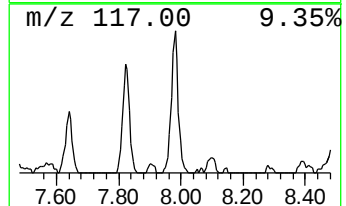
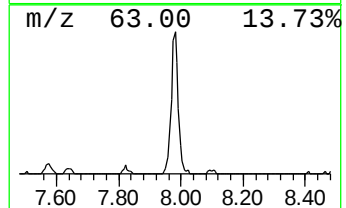
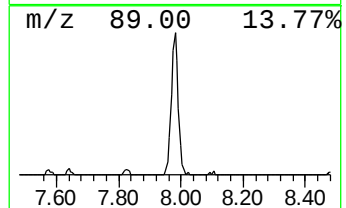
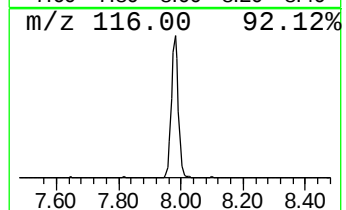
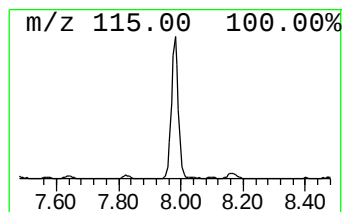
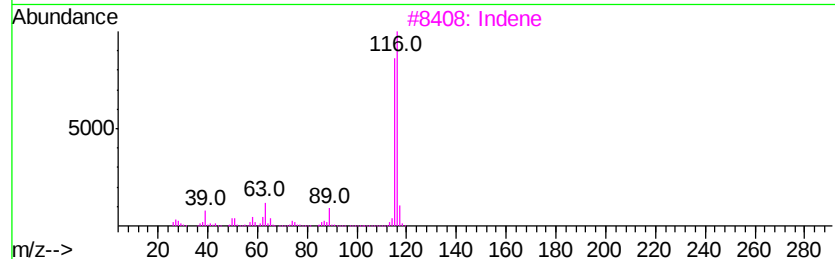
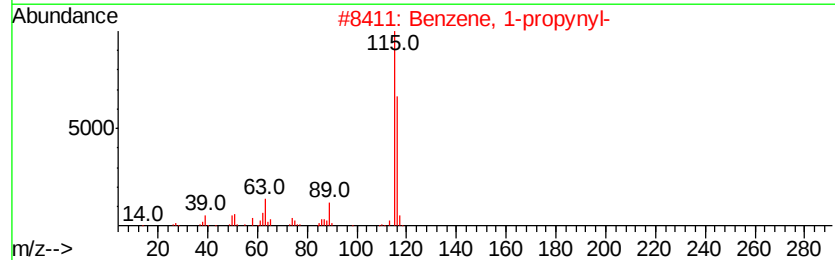
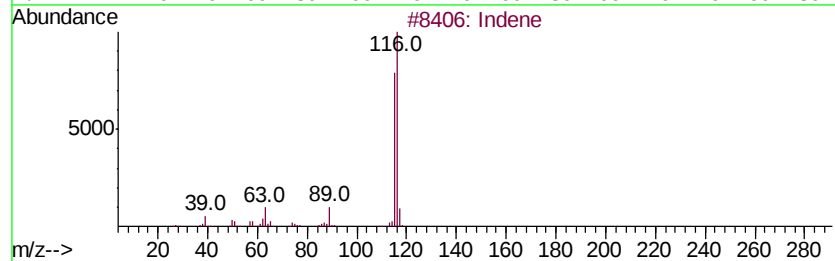
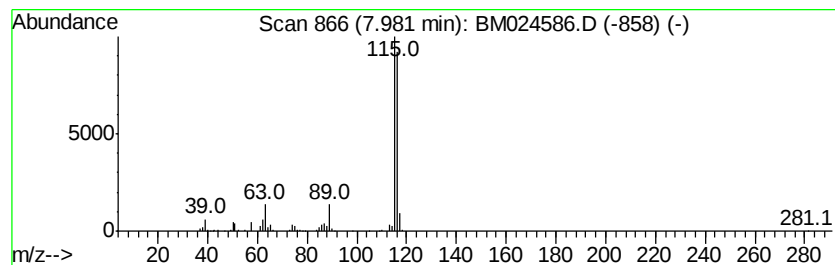
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Indene Concentration Rank 9

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 7.98 | 11.97 ng/ul | 720647 | 1,4-Dichlorobenzene-d4 | 7.45 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Indene                  | 116 | C9H8    | 000095-13-6 | 95   |
| 2    |    | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 94   |
| 3    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 94   |
| 4    |    | 3-Methylphenylacetylene | 116 | C9H8    | 000766-82-5 | 94   |
| 5    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

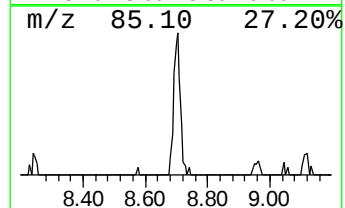
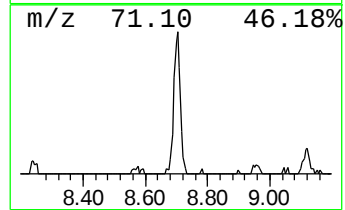
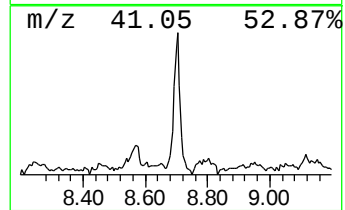
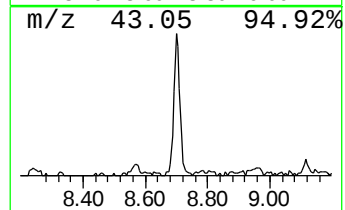
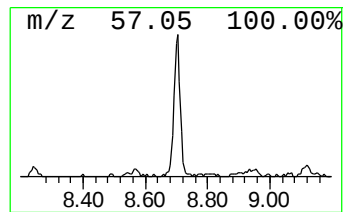
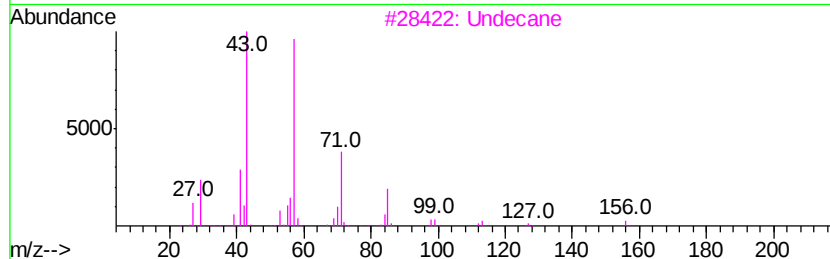
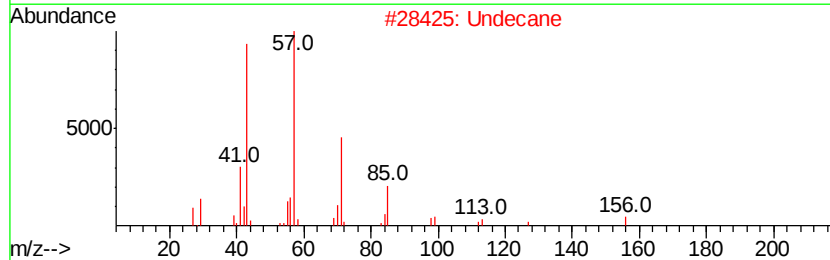
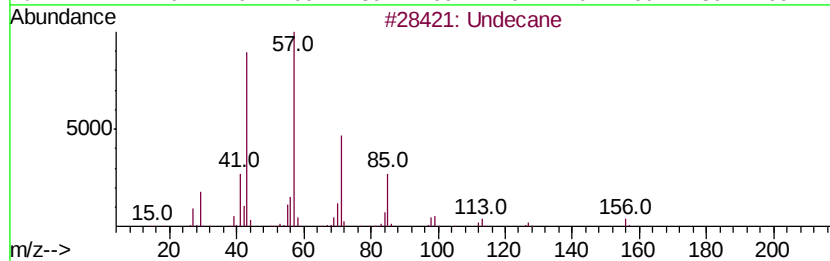
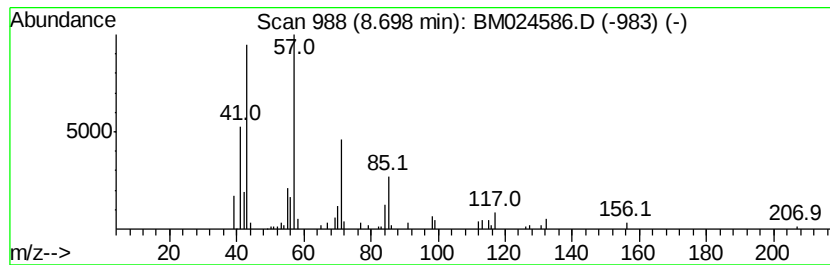
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.70 | 3.68 ng/ul | 221609 | 1,4-Dichlorobenzene-d4 | 7.45 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Undecane     | 156 | C11H24  | 001120-21-4 | 76   |
| 2    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 70   |
| 3    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 62   |
| 4    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 58   |
| 5    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
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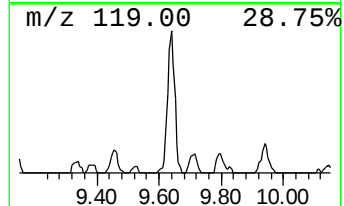
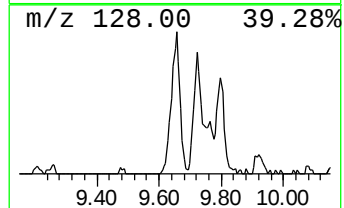
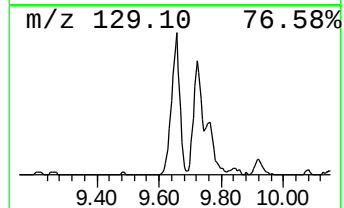
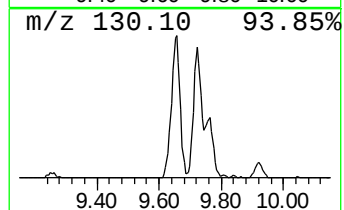
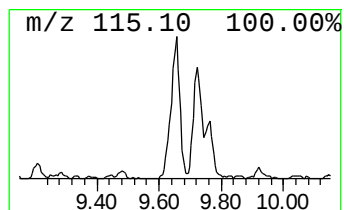
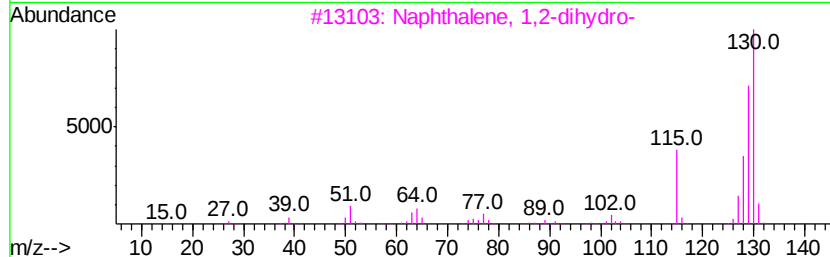
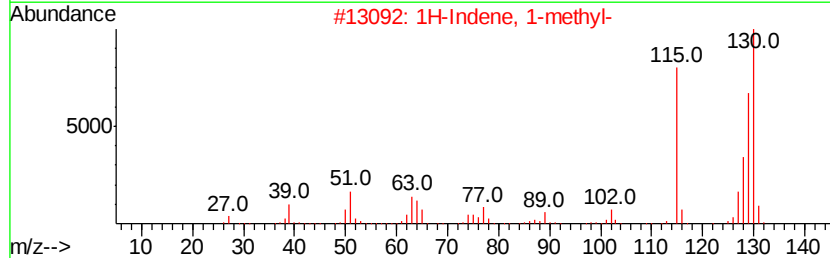
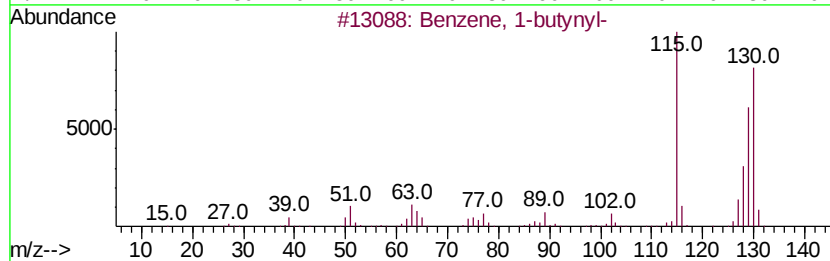
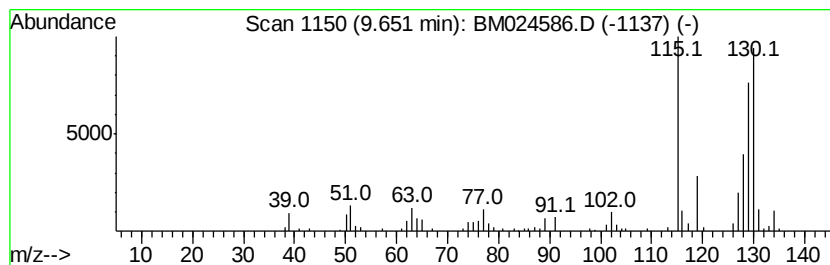
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1-butynyl- Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.65 | 4.07 ng/ul | 408764 | Naphthalene-d8   | 10.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-butynyl-                 | 130 | C10H10  | 000622-76-4 | 95   |
| 2    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 90   |
| 3    |    | Naphthalene, 1,2-dihydro-           | 130 | C10H10  | 000447-53-0 | 90   |
| 4    |    | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 90   |
| 5    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

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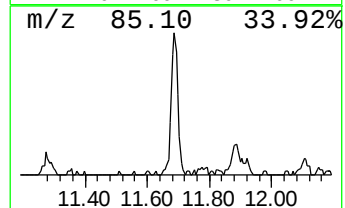
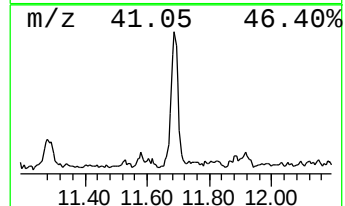
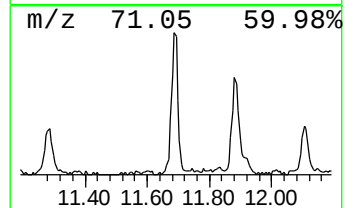
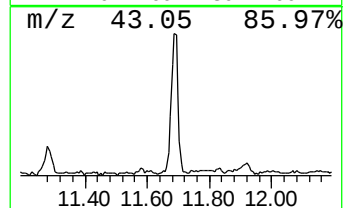
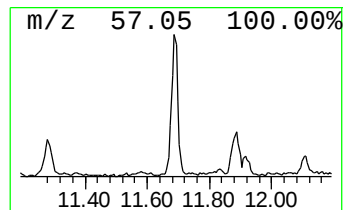
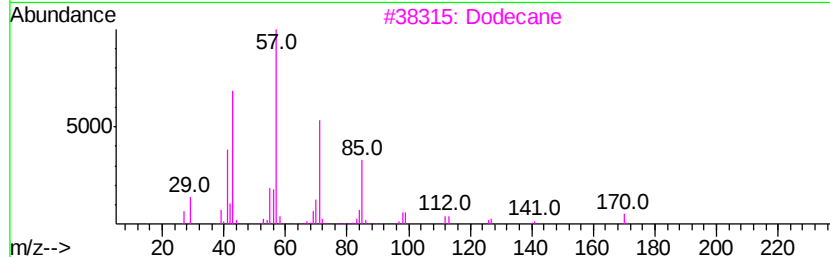
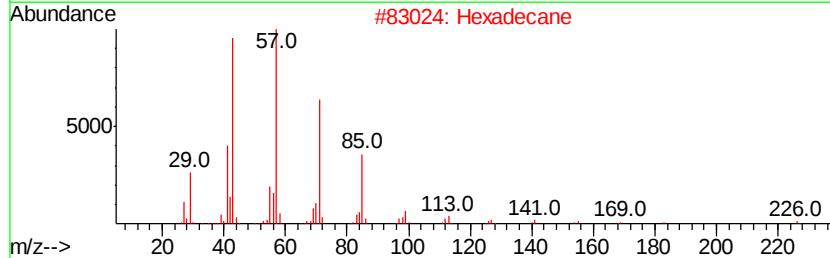
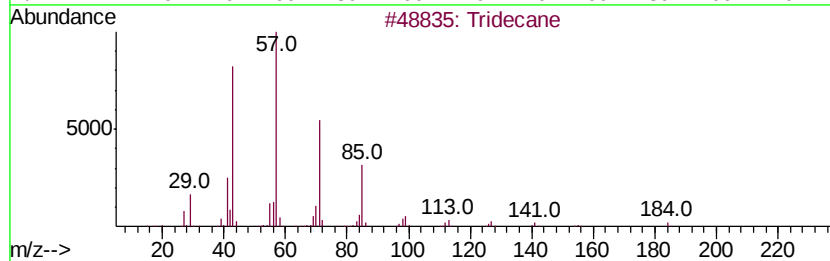
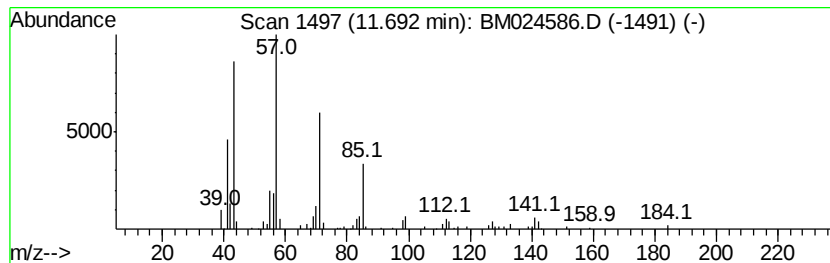
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.69 | 2.56 ng/ul | 257205 | Naphthalene-d8   | 10.21 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane            | 184 | C13H28  | 000629-50-5 | 93   |
| 2    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 86   |
| 3    |    | Dodecane             | 170 | C12H26  | 000112-40-3 | 86   |
| 4    |    | Hexadecane           | 226 | C16H34  | 000544-76-3 | 80   |
| 5    |    | Tetradecane, 1-iodo- | 324 | C14H29I | 019218-94-1 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

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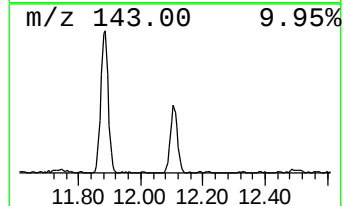
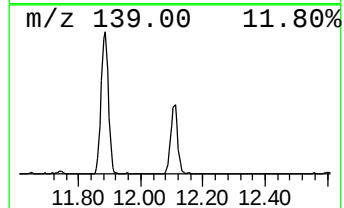
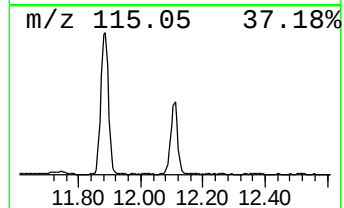
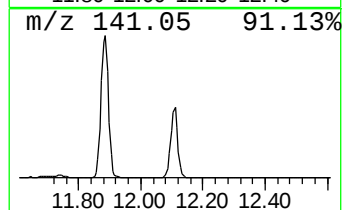
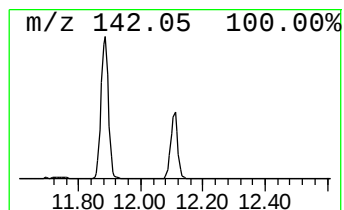
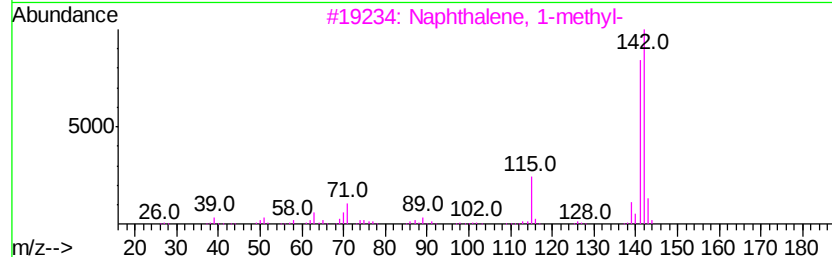
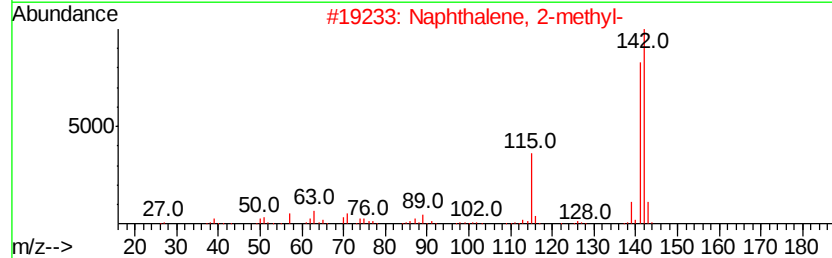
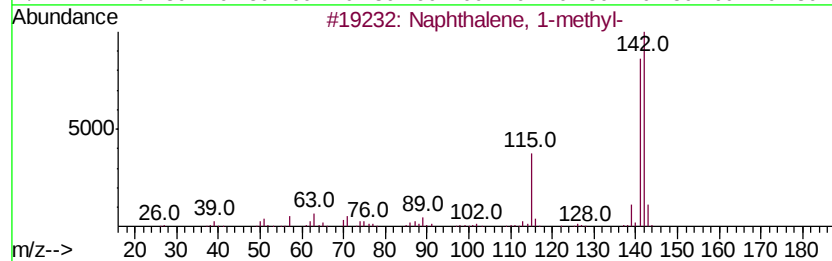
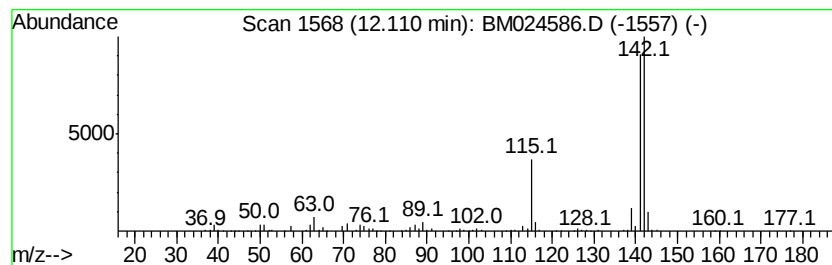
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.11 | 8.94 ng/ul | 896815 | Naphthalene-d8   | 10.21 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 2    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 3    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 94   |
| 4    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 93   |
| 5    |    | Benzocycloheptatriene  | 142 | C11H10  | 000264-09-5 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

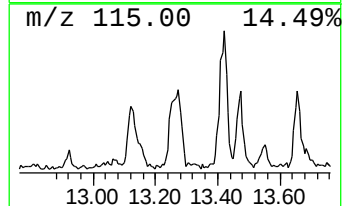
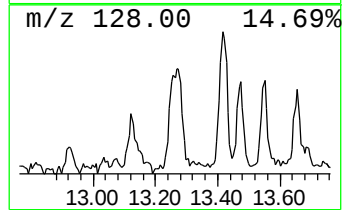
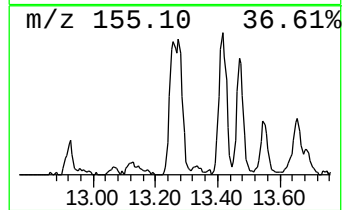
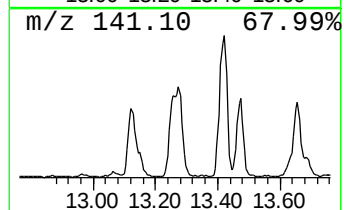
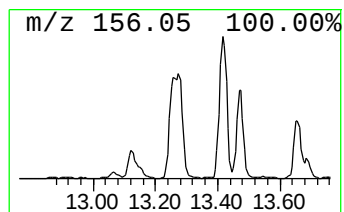
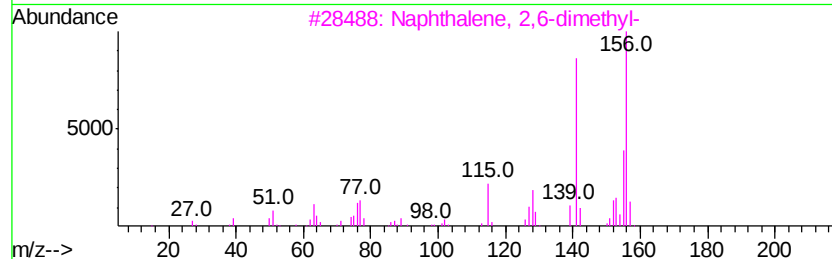
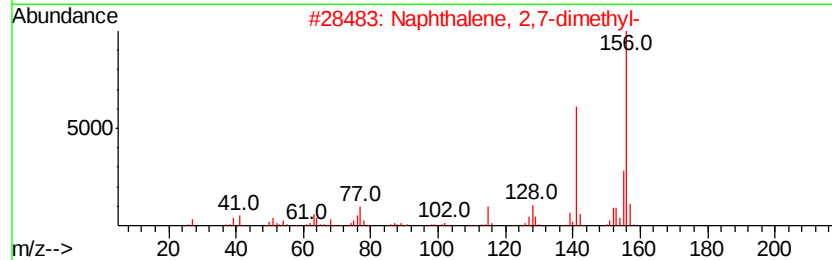
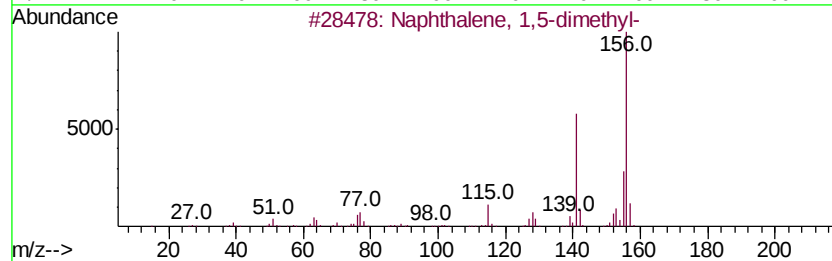
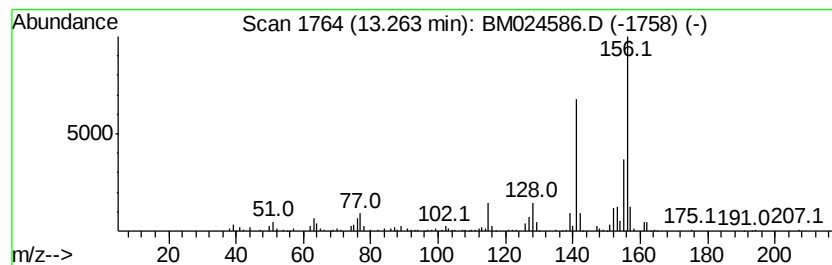
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 1,5-dimethyl- Concentration Rank 43

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.26 | 2.21 ng/ul | 316367 | Acenaphthene-d10 | 14.10 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 2    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |
| 3    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 4    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 5    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

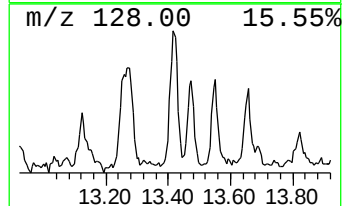
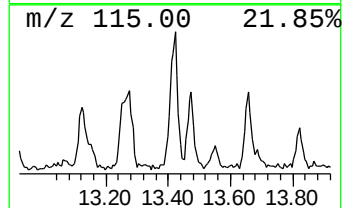
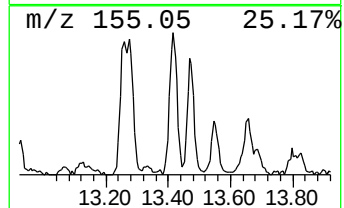
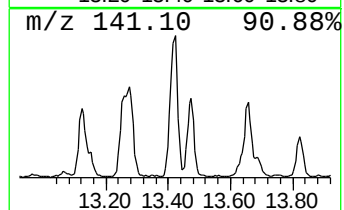
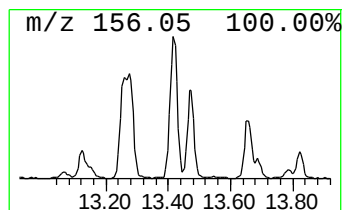
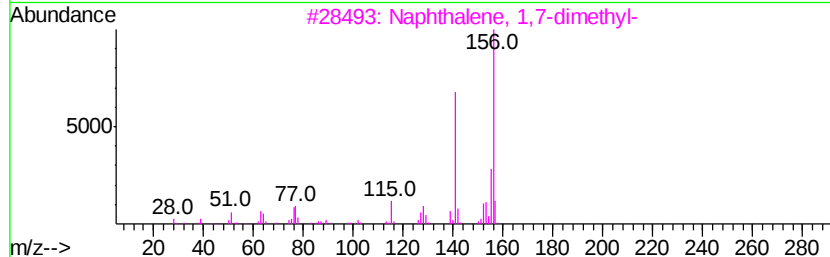
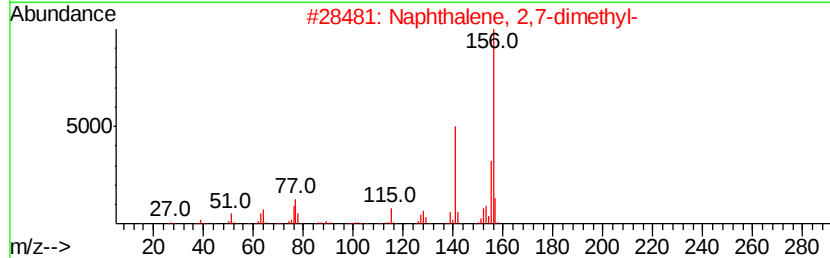
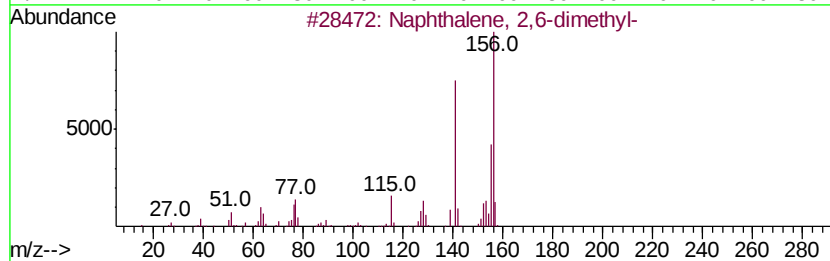
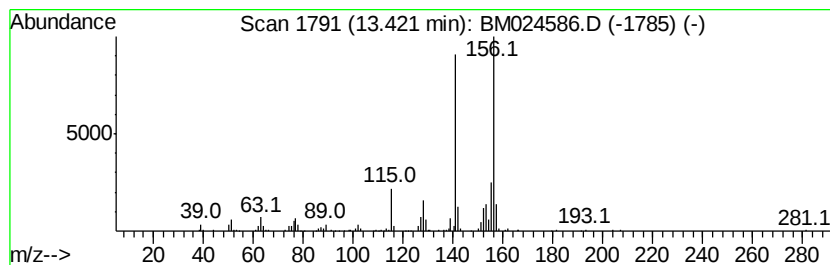
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Naphthalene, 2,6-dimethyl- Concentration Rank 34

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.42 | 2.85 ng/ul | 408896 | Acenaphthene-d10 | 14.10 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 2    |    | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 3    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 97   |
| 5    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 97   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

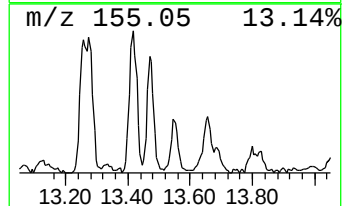
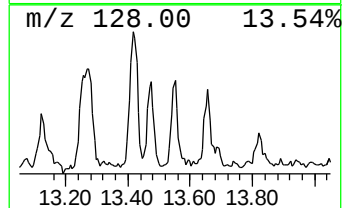
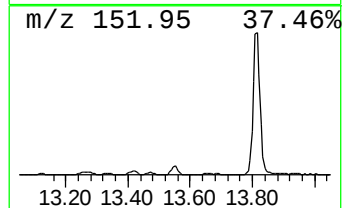
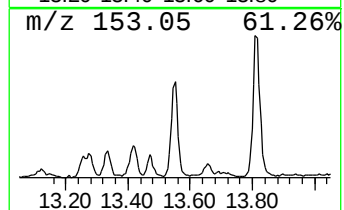
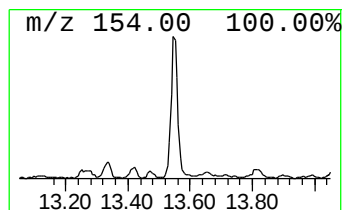
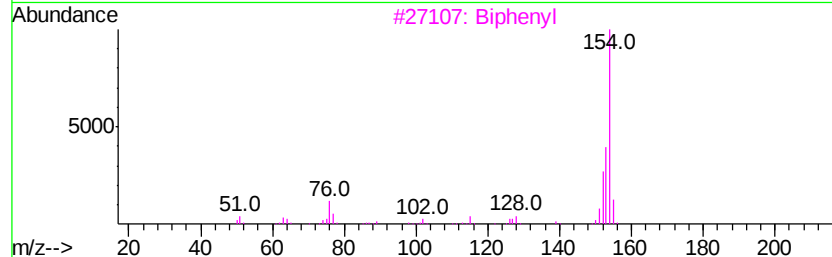
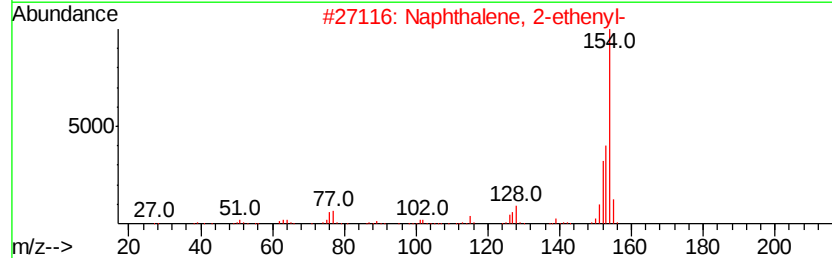
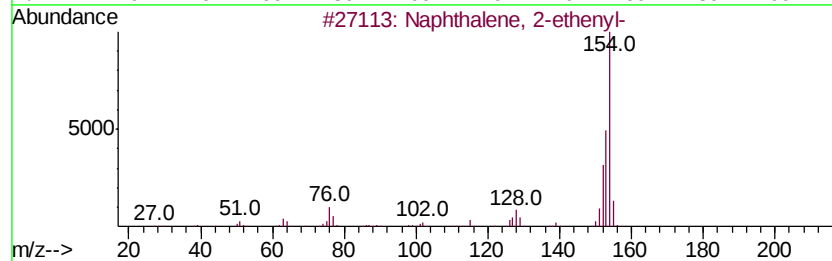
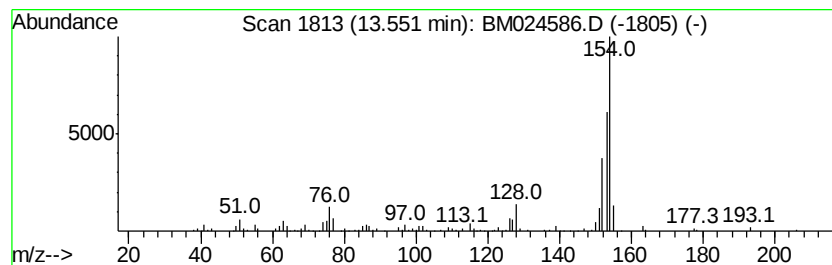
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Naphthalene, 2-ethenyl- Concentration Rank 35

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.55 | 2.75 ng/ul | 394230 | Acenaphthene-d10 | 14.10 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 96   |
| 2    |    | Naphthalene, 2-ethenyl- | 154 | C12H10  | 000827-54-3 | 93   |
| 3    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 87   |
| 4    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 87   |
| 5    |    | Biphenyl                | 154 | C12H10  | 000092-52-4 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

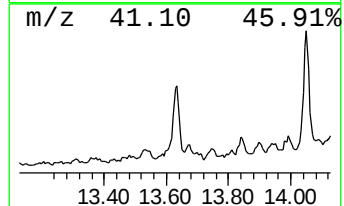
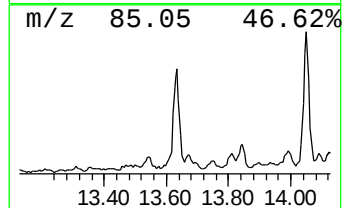
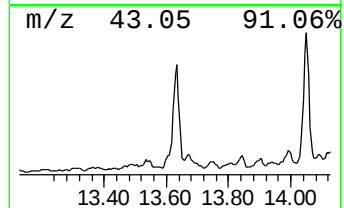
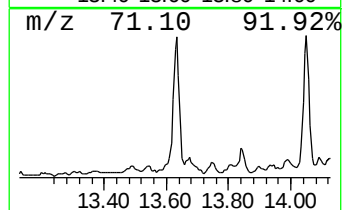
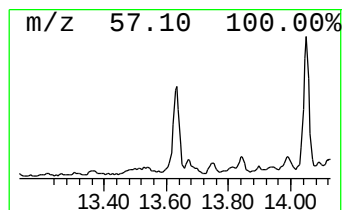
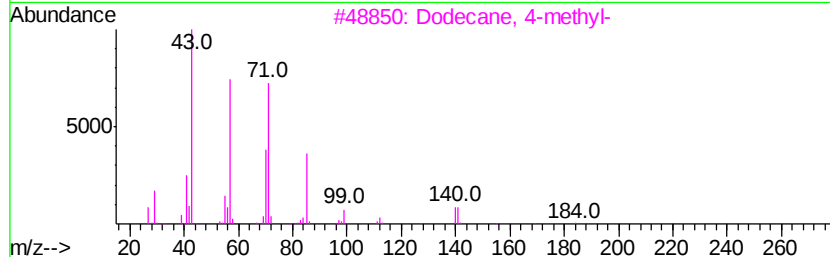
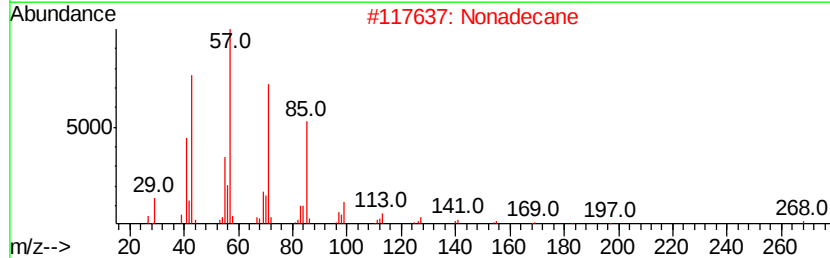
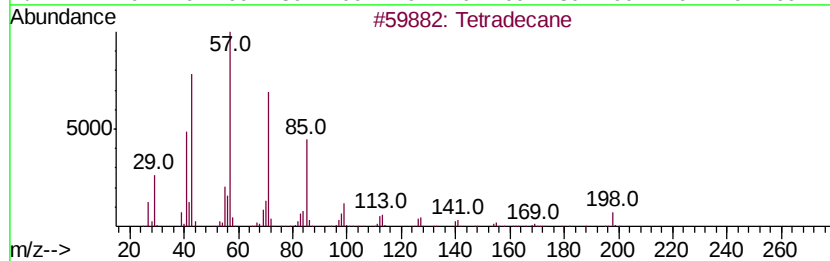
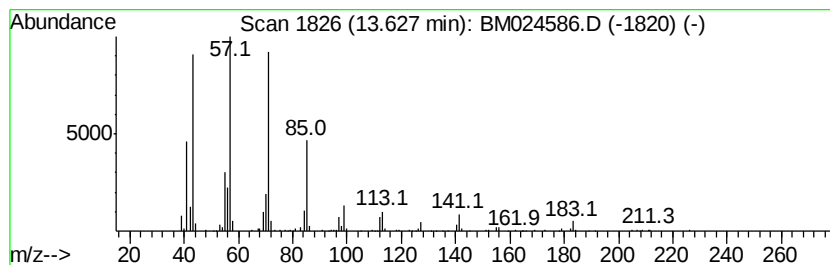
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.63 | 6.34 ng/ul | 909005 | Acenaphthene-d10 | 14.10 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane         | 198 | C14H30  | 000629-59-4 | 90   |
| 2    |    |   | Nonadecane          | 268 | C19H40  | 000629-92-5 | 86   |
| 3    |    |   | Dodecane, 4-methyl- | 184 | C13H28  | 006117-97-1 | 70   |
| 4    |    |   | Dodecane, 4-methyl- | 184 | C13H28  | 006117-97-1 | 68   |
| 5    |    |   | Decane, 5-propyl-   | 184 | C13H28  | 017312-62-8 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

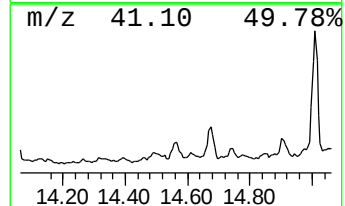
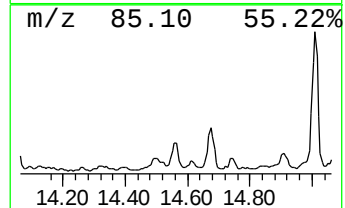
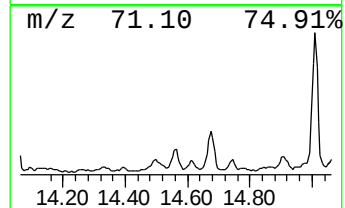
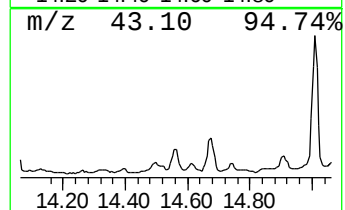
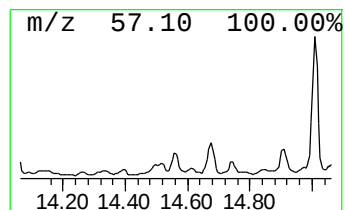
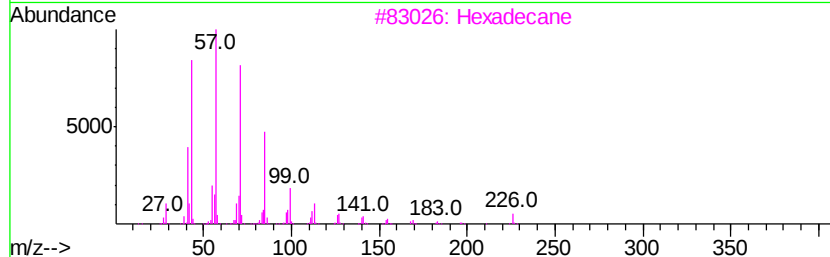
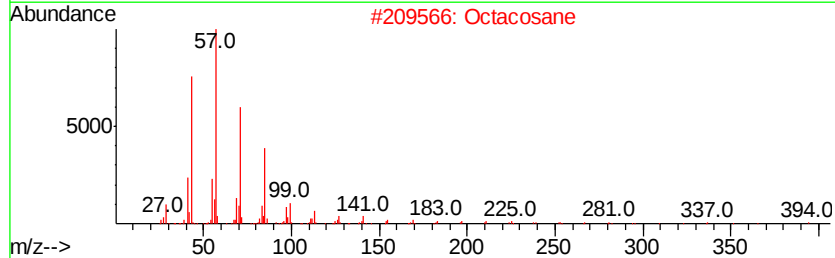
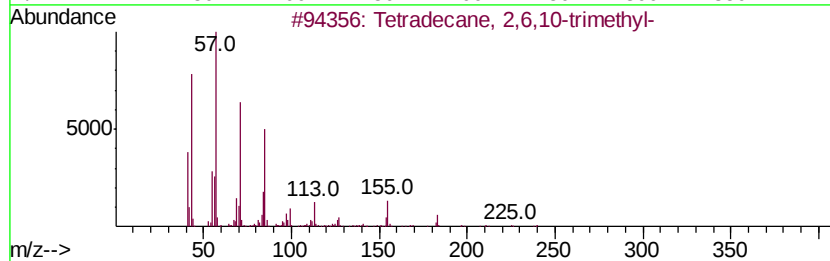
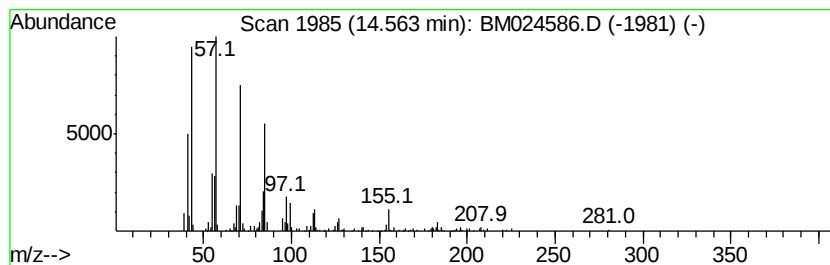
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.56 | 2.93 ng/ul | 420304 | Acenaphthene-d10 | 14.10 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7 | 86   |
| 2    |    |   | Octacosane                     | 394 | C28H58  | 000630-02-4 | 74   |
| 3    |    |   | Hexadecane                     | 226 | C16H34  | 000544-76-3 | 72   |
| 4    |    |   | Octadecane, 1-iodo-            | 380 | C18H37I | 000629-93-6 | 72   |
| 5    |    |   | Hexacosane                     | 366 | C26H54  | 000630-01-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

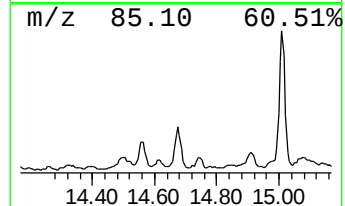
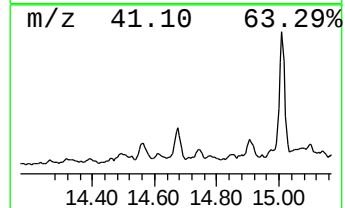
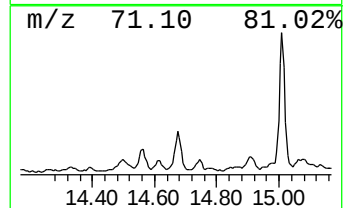
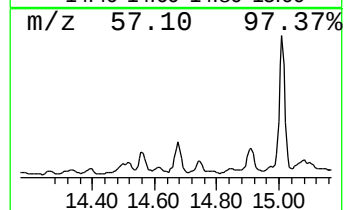
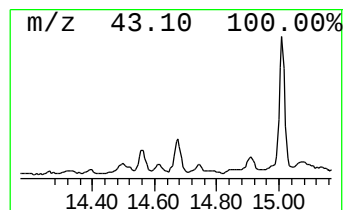
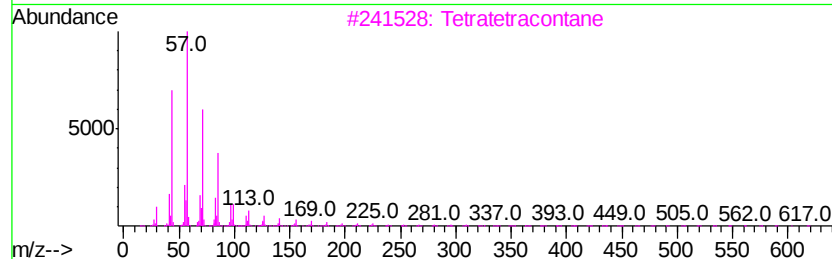
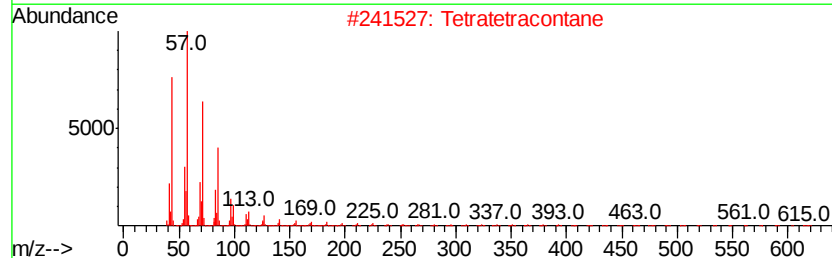
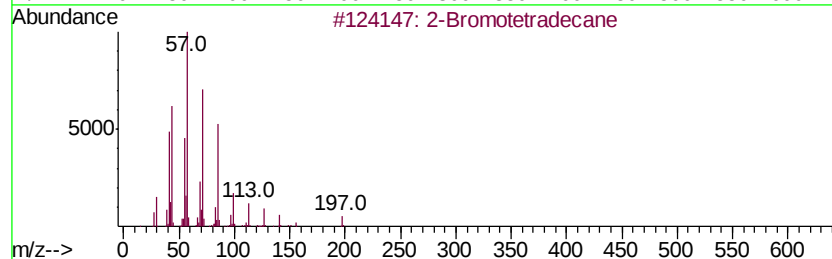
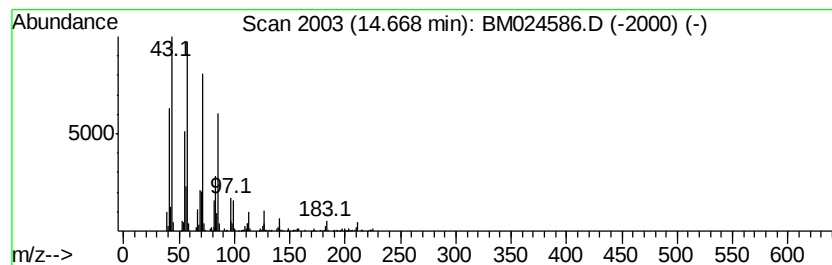
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 2-Bromotetradecane Concentration Rank 29

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.67 | 3.39 ng/ul | 486503 | Acenaphthene-d10 | 14.10 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Bromotetradecane | 276 | C14H29Br | 074036-95-6 | 86   |
| 2    |    |   | Tetratetracontane  | 619 | C44H90   | 007098-22-8 | 86   |
| 3    |    |   | Tetratetracontane  | 619 | C44H90   | 007098-22-8 | 86   |
| 4    |    |   | Heptacosane        | 380 | C27H56   | 000593-49-7 | 86   |
| 5    |    |   | Heneicosane        | 296 | C21H44   | 000629-94-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

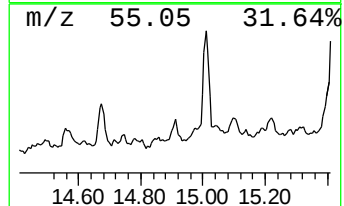
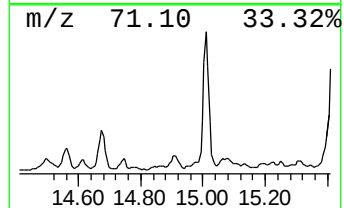
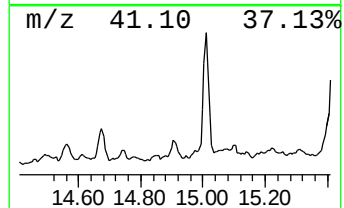
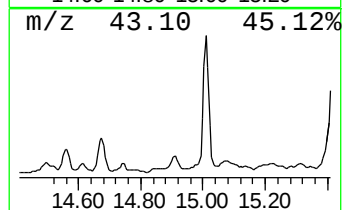
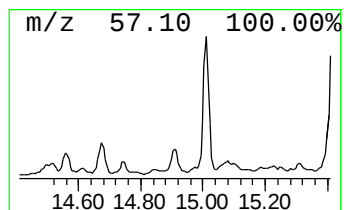
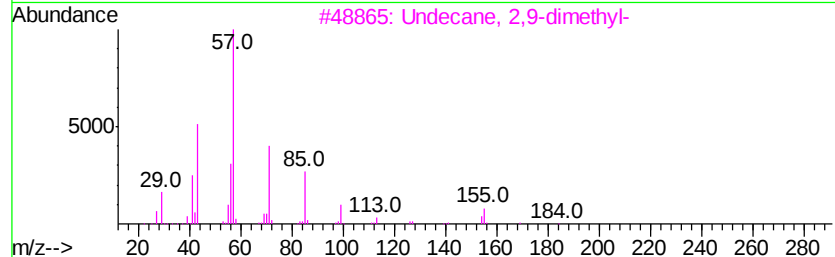
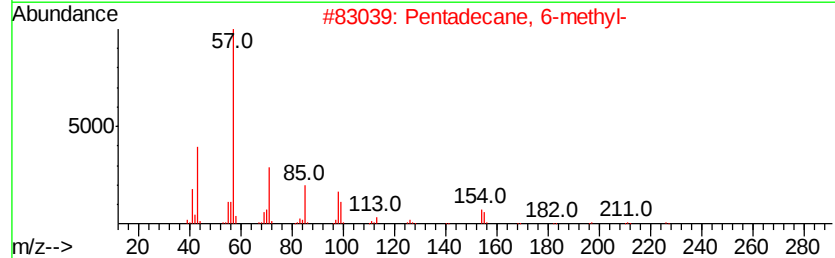
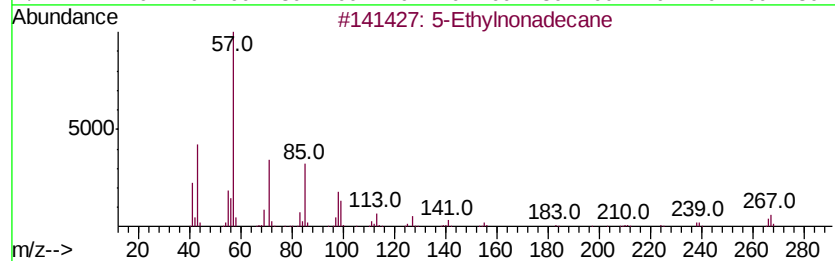
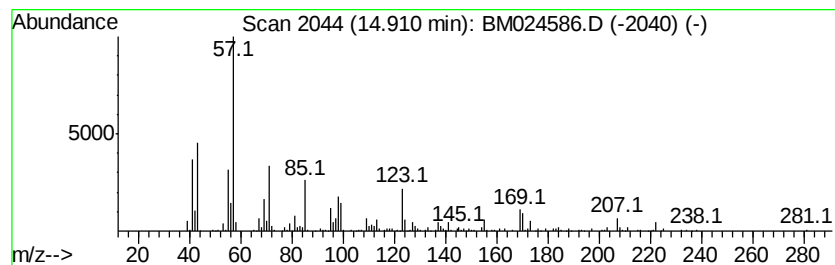
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.91 | 2.39 ng/ul | 342909 | Acenaphthene-d10 | 14.10 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------|-----|----------|--------------|------|
| 1    |    |   | 5-Ethylnonadecane       | 296 | C21H44   | 1000360-43-1 | 49   |
| 2    |    |   | Pentadecane, 6-methyl-  | 226 | C16H34   | 010105-38-1  | 47   |
| 3    |    |   | Undecane, 2,9-dimethyl- | 184 | C13H28   | 017301-26-7  | 43   |
| 4    |    |   | 2-Bromo dodecane        | 248 | C12H25Br | 013187-99-0  | 38   |
| 5    |    |   | Tetratetracontane       | 619 | C44H90   | 007098-22-8  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

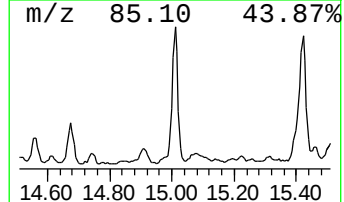
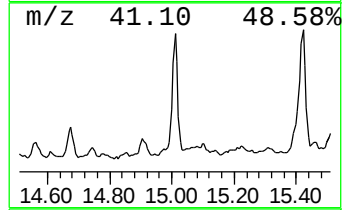
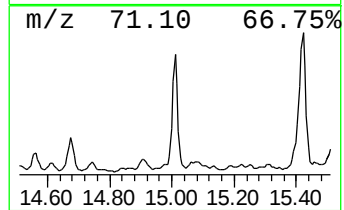
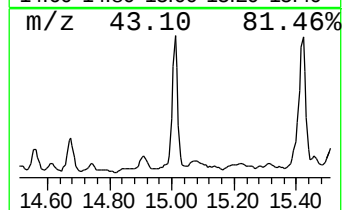
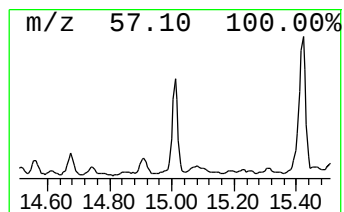
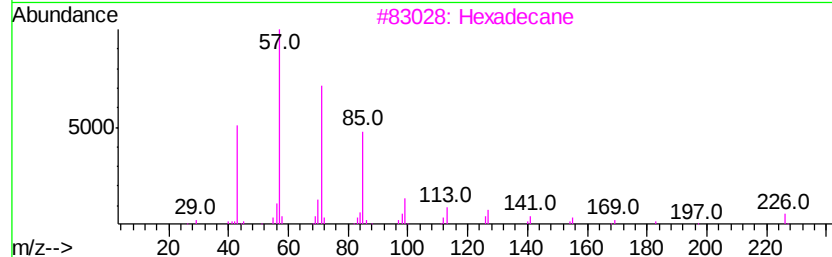
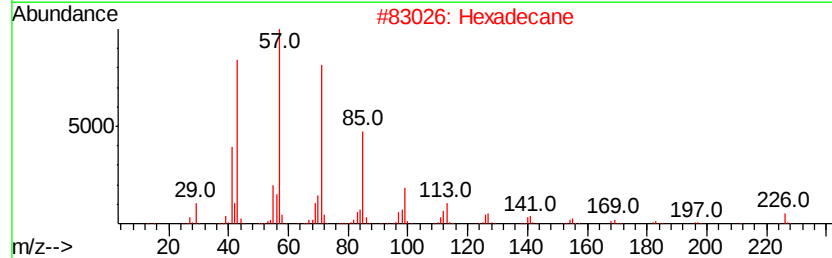
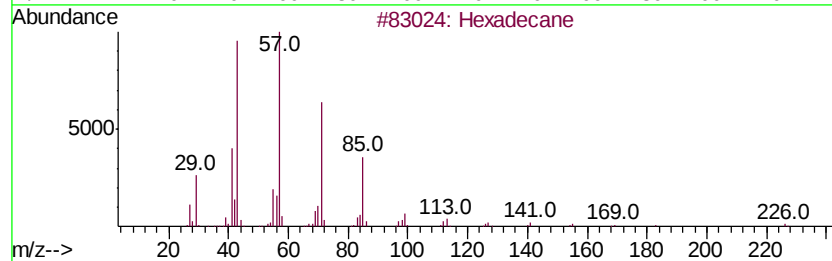
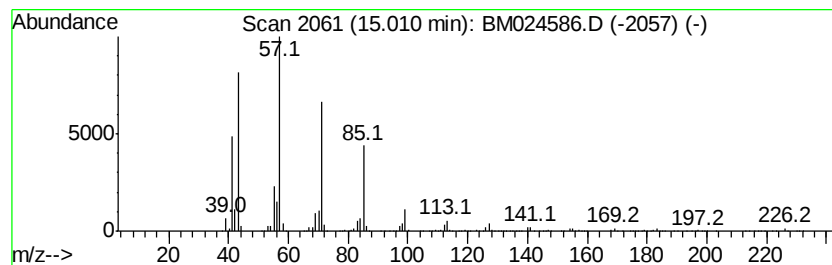
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.01 | 10.57 ng/ul | 1515010 | Acenaphthene-d10 | 14.10 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 2    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 89   |
| 5    |    | Pentadecane  | 212 | C15H32  | 000629-62-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

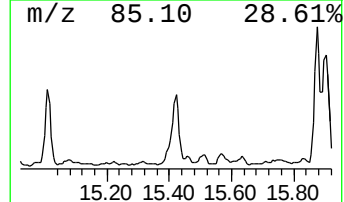
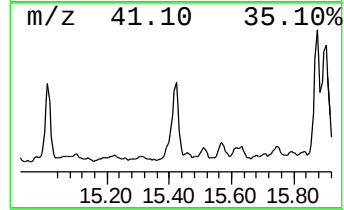
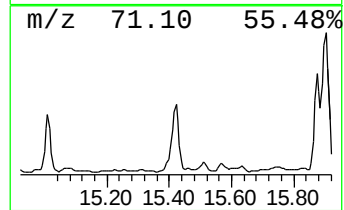
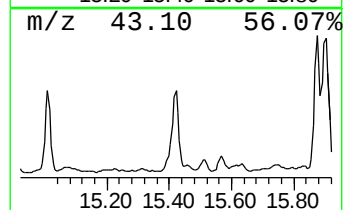
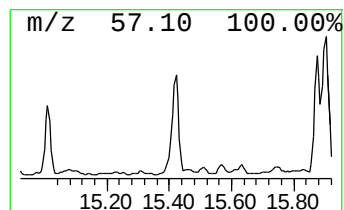
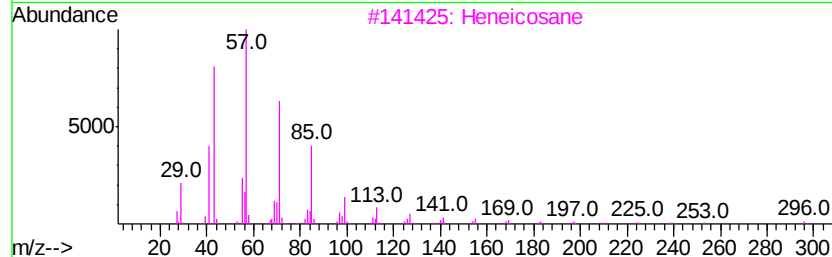
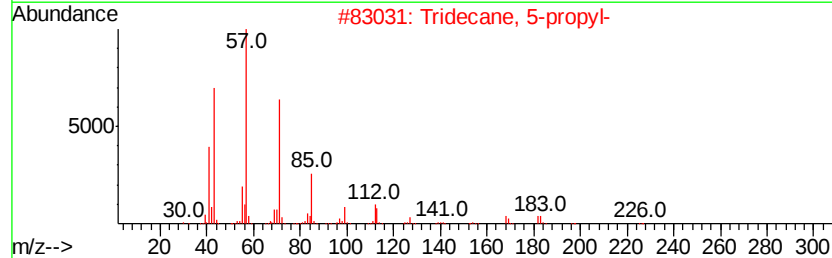
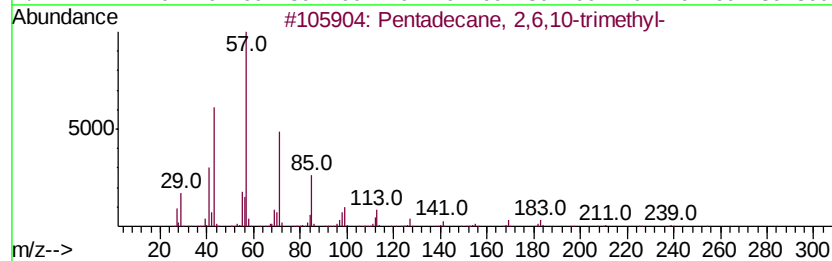
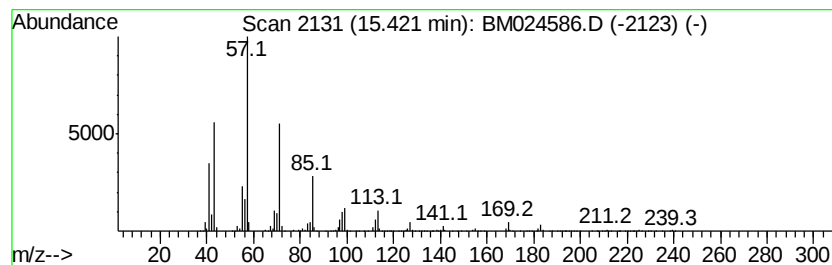
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.42 | 18.44 ng/ul | 2643710 | Acenaphthene-d10 | 14.10 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl-     | 254 | C18H38  | 003892-00-0 | 91   |
| 2    |    | Tridecane, 5-propyl-               | 226 | C16H34  | 055045-11-9 | 90   |
| 3    |    | Heneicosane                        | 296 | C21H44  | 000629-94-7 | 83   |
| 4    |    | Octacosane                         | 394 | C28H58  | 000630-02-4 | 80   |
| 5    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

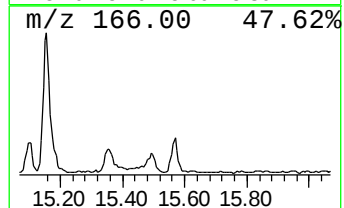
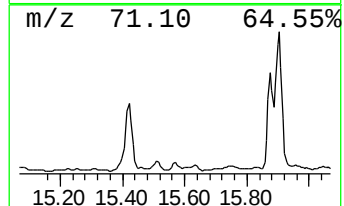
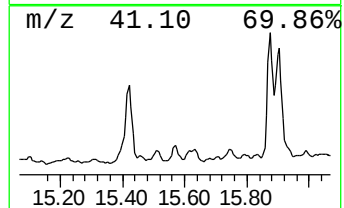
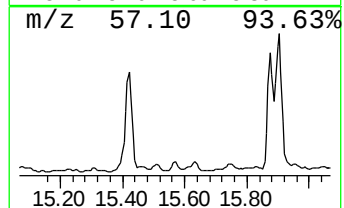
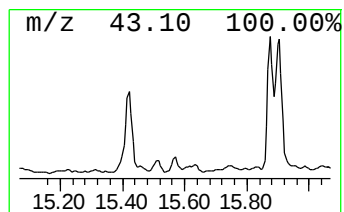
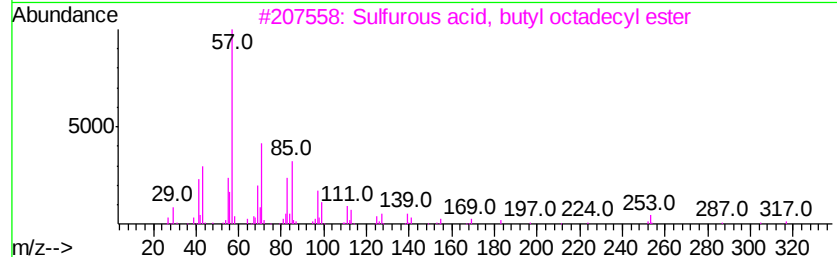
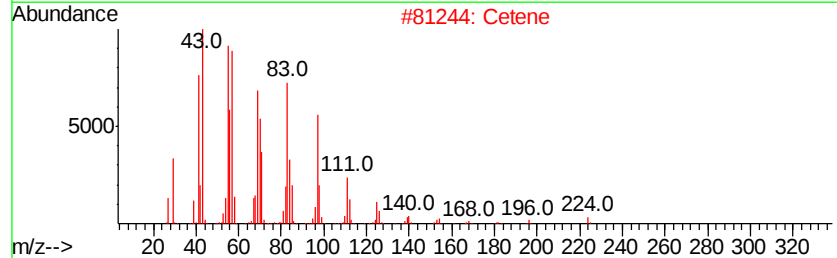
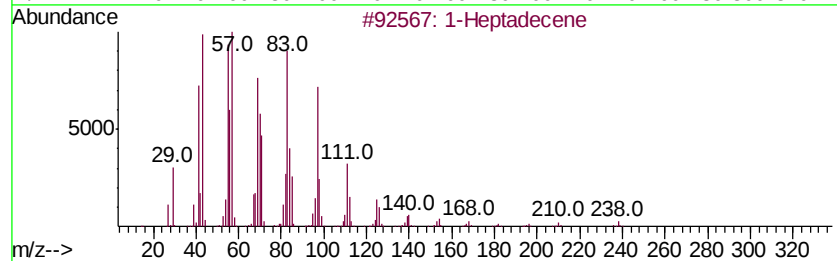
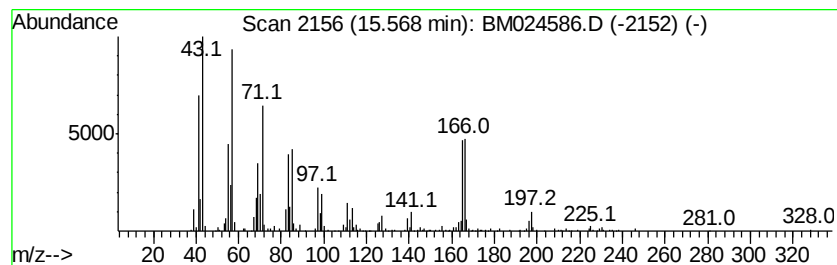
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Cyclic15.57 Concentration Rank 32

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.57 | 3.11 ng/ul | 450333 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1-Heptadecene                       | 238 | C17H34    | 006765-39-5  | 55   |
| 2    |    |   | Cetene                              | 224 | C16H32    | 000629-73-2  | 53   |
| 3    |    |   | Sulfurous acid, butyl octadecyl ... | 390 | C22H46O3S | 1000309-18-5 | 52   |
| 4    |    |   | Oxalic acid, isobutyl tetradecyl... | 342 | C20H38O4  | 1000309-37-9 | 46   |
| 5    |    |   | 1-Chloroeicosane                    | 316 | C20H41Cl  | 042217-02-7  | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

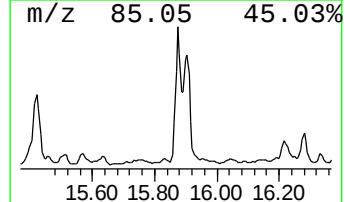
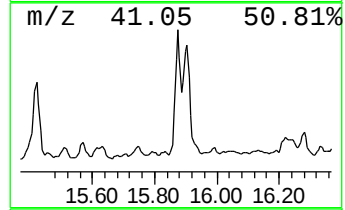
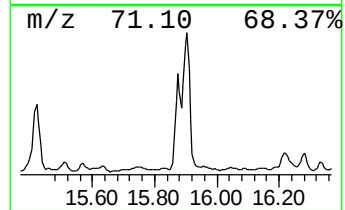
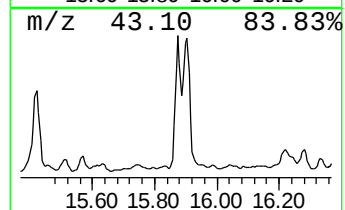
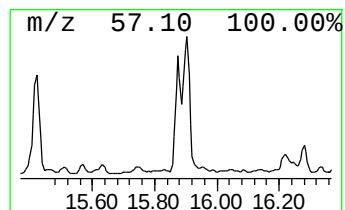
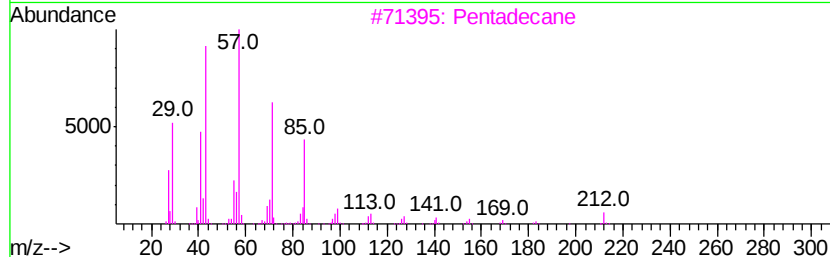
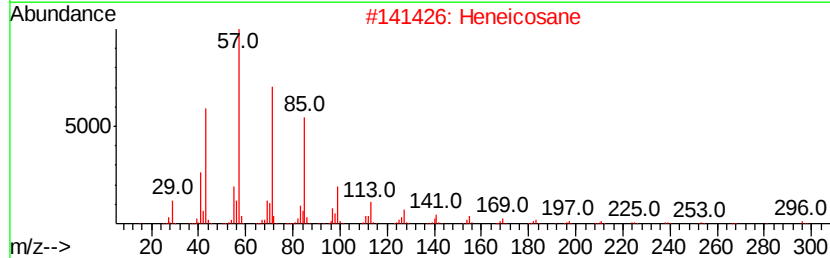
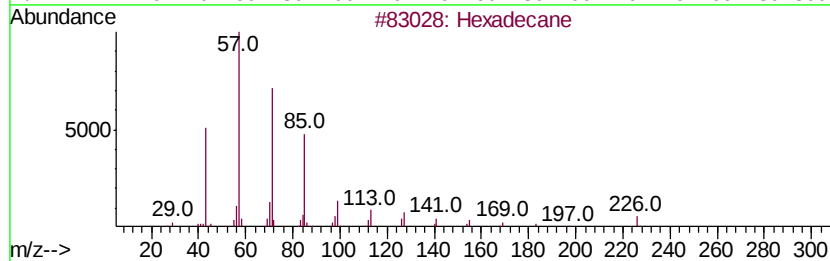
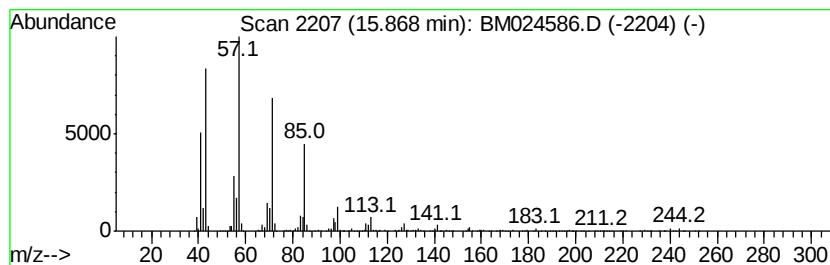
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.87 | 19.55 ng/ul | 2831400 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane           | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |    | Heneicosane          | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Pentadecane          | 212 | C15H32  | 000629-62-9 | 87   |
| 4    |    | Eicosane, 10-methyl- | 296 | C21H44  | 054833-23-7 | 87   |
| 5    |    | Tetradecane          | 198 | C14H30  | 000629-59-4 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

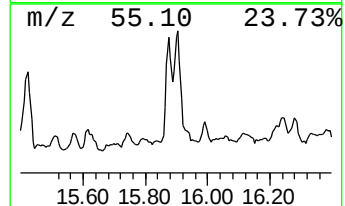
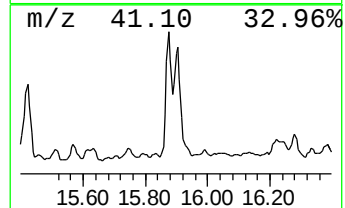
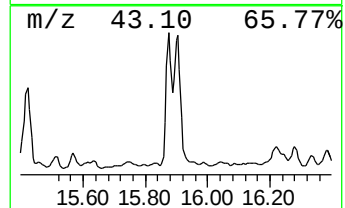
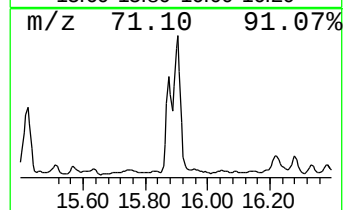
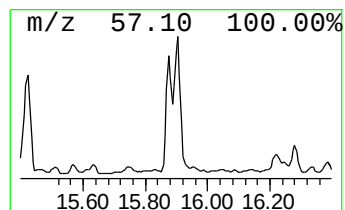
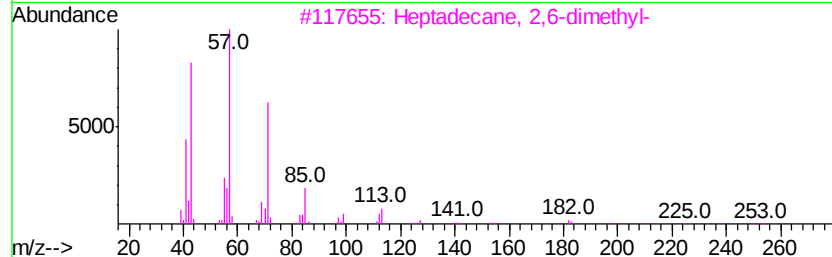
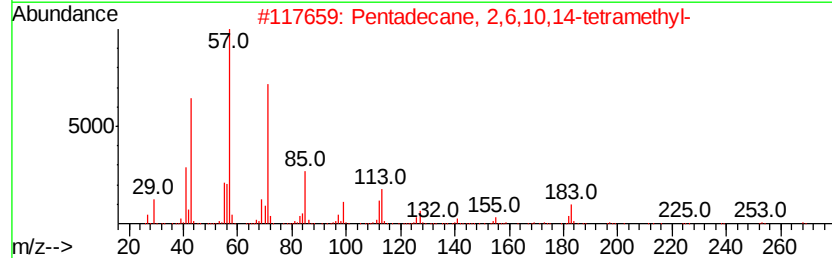
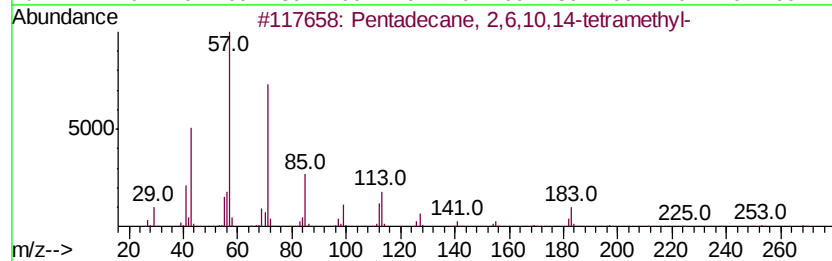
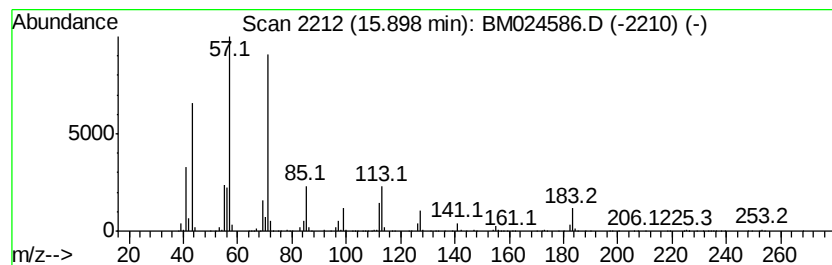
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.90 | 24.27 ng/ul | 3515880 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40    | 001921-70-6  | 87   |
| 2    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40    | 001921-70-6  | 87   |
| 3    |    | Heptadecane, 2,6-dimethyl-          | 268 | C19H40    | 054105-67-8  | 81   |
| 4    |    | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1000309-19-2 | 72   |
| 5    |    | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

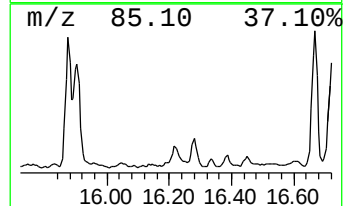
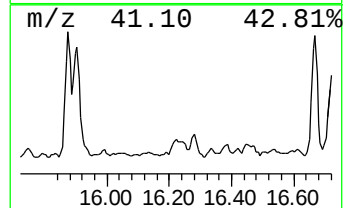
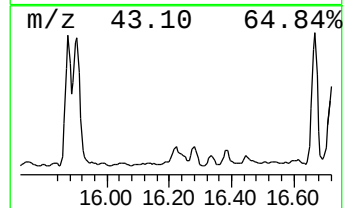
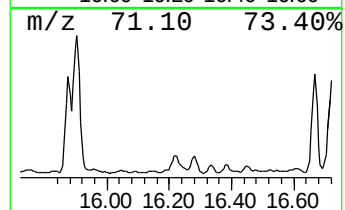
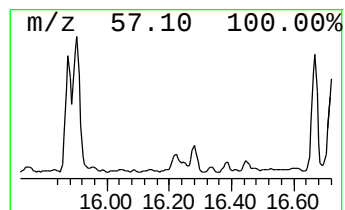
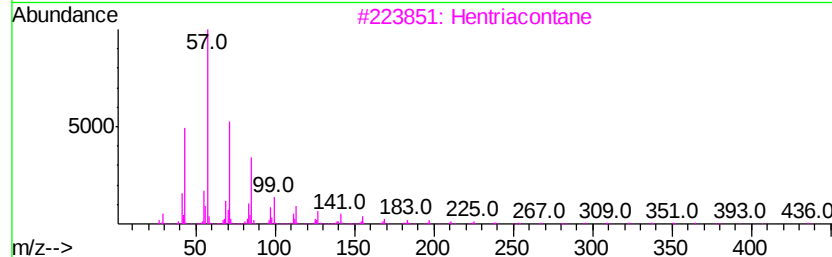
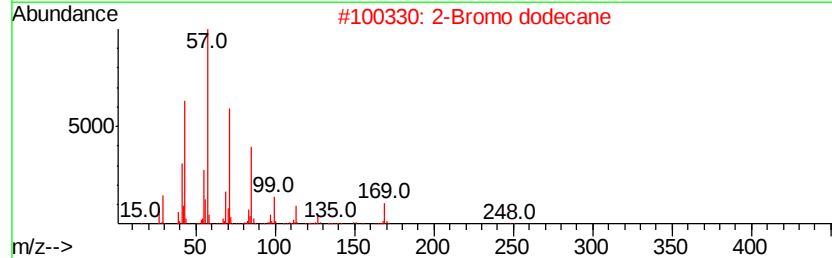
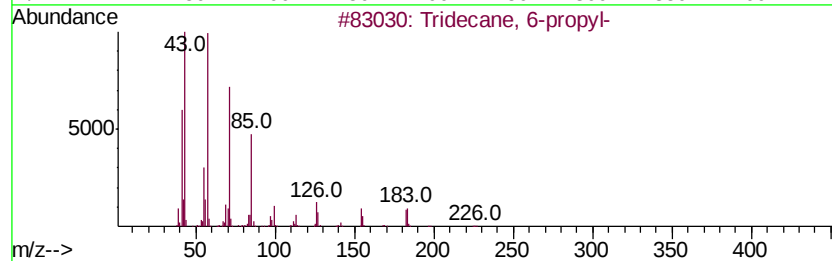
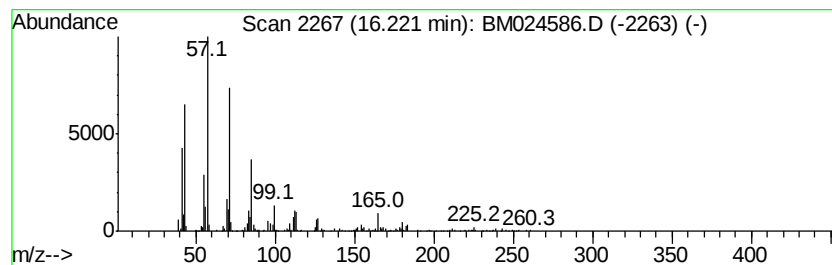
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.22 | 5.82 ng/ul | 842787 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------|-----|----------|-------------|------|
| 1    |    |   | Tridecane, 6-propyl- | 226 | C16H34   | 055045-10-8 | 87   |
| 2    |    |   | 2-Bromo dodecane     | 248 | C12H25Br | 013187-99-0 | 83   |
| 3    |    |   | Hentriacontane       | 437 | C31H64   | 000630-04-6 | 83   |
| 4    |    |   | Hexadecane           | 226 | C16H34   | 000544-76-3 | 83   |
| 5    |    |   | Tetracosane          | 338 | C24H50   | 000646-31-1 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

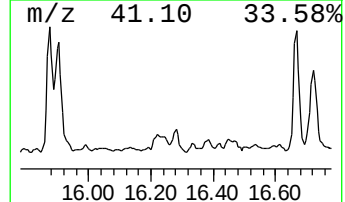
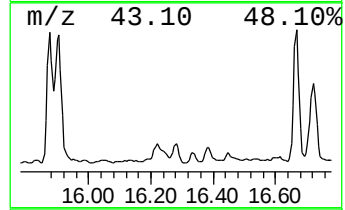
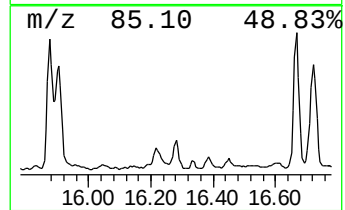
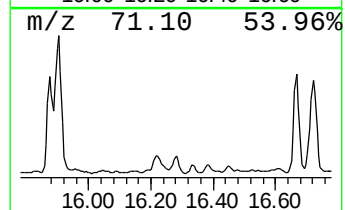
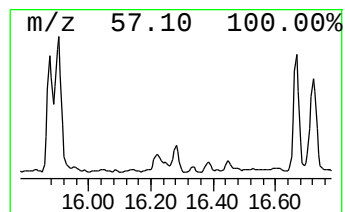
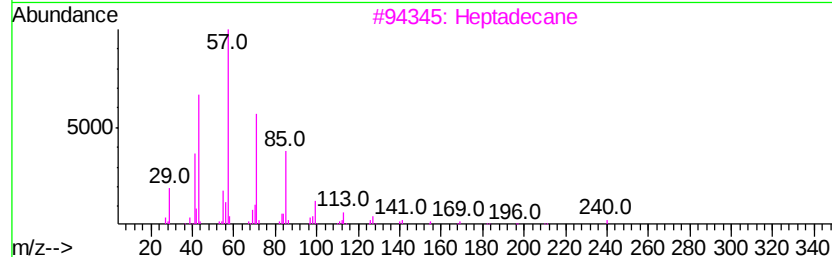
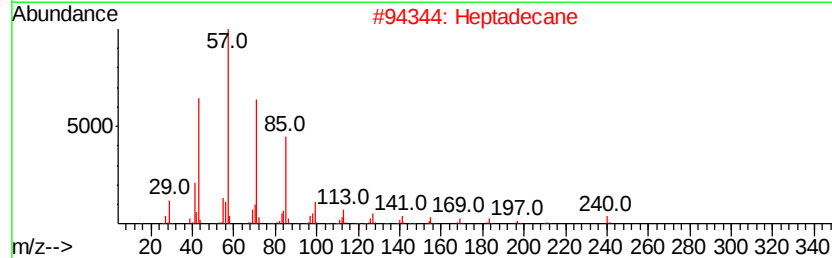
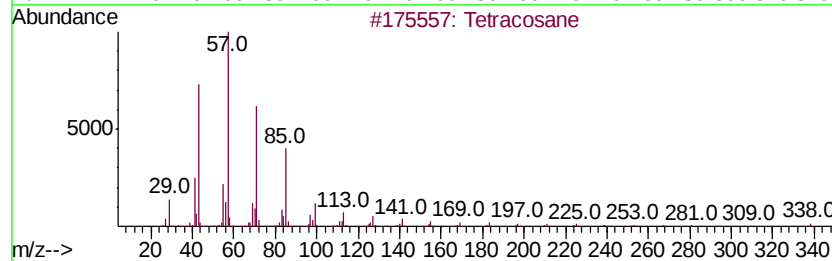
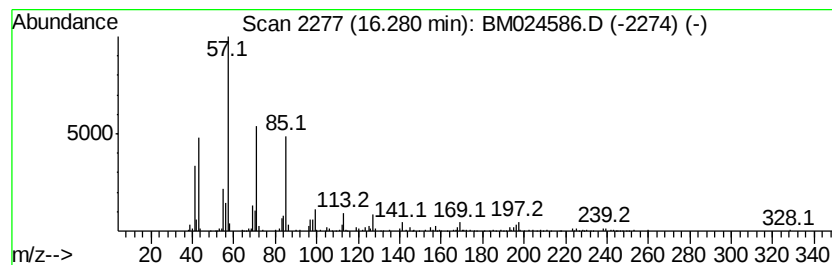
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.28 | 4.25 ng/ul | 615651 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane  | 338 | C24H50  | 000646-31-1 | 86   |
| 2    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 86   |
| 3    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 86   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 5    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

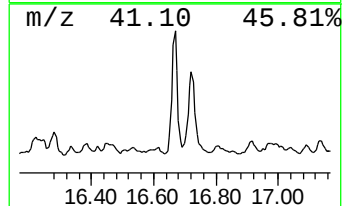
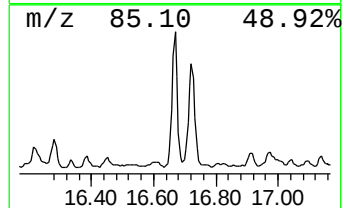
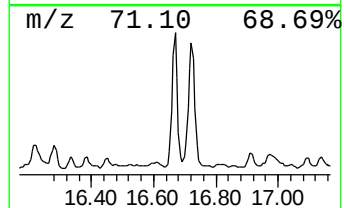
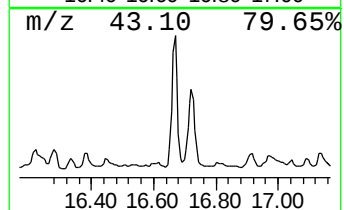
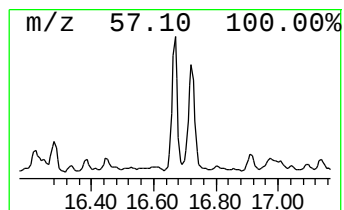
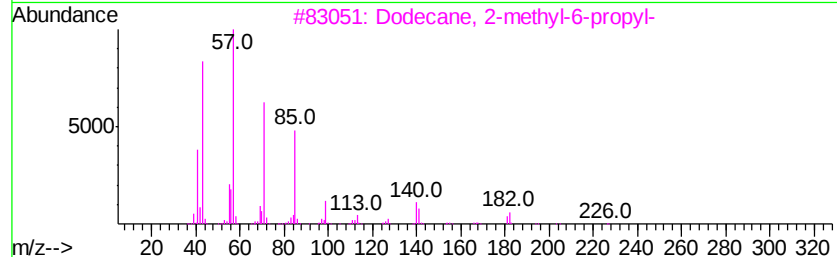
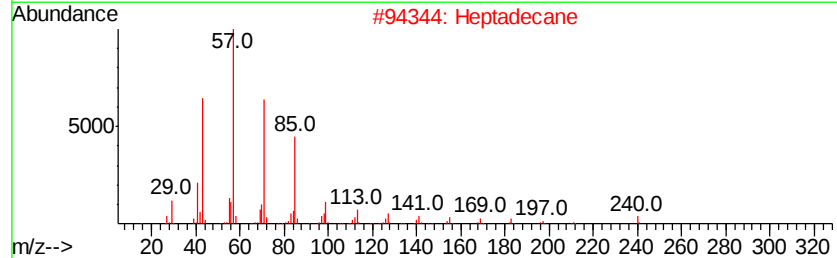
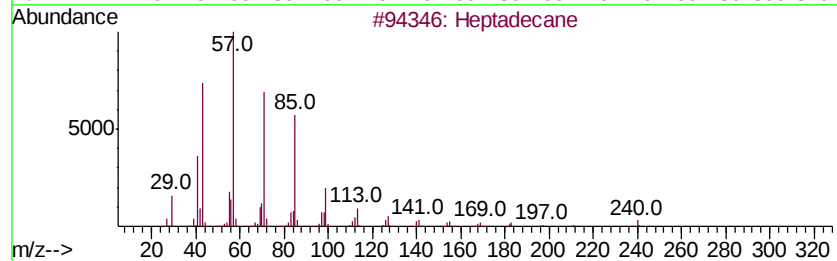
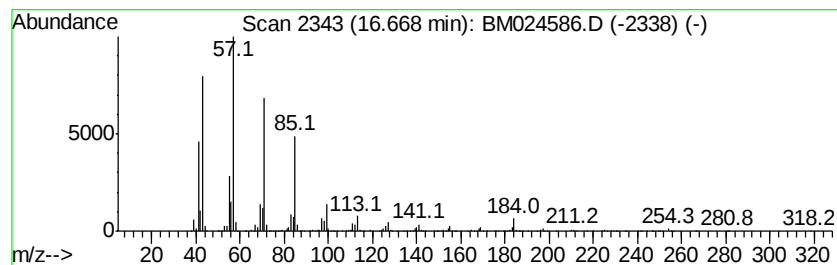
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.67 | 20.67 ng/ul | 2994180 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane                  | 240 | C17H36  | 000629-78-7 | 97   |
| 2    |    | Heptadecane                  | 240 | C17H36  | 000629-78-7 | 94   |
| 3    |    | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 94   |
| 4    |    | Tridecane                    | 184 | C13H28  | 000629-50-5 | 93   |
| 5    |    | Nonadecane                   | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

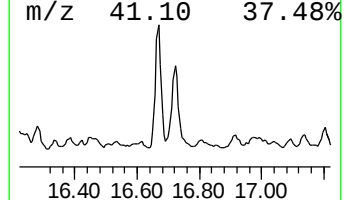
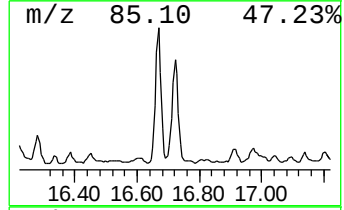
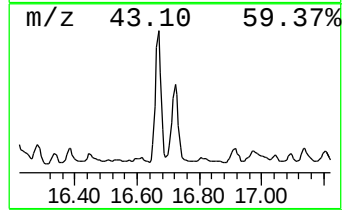
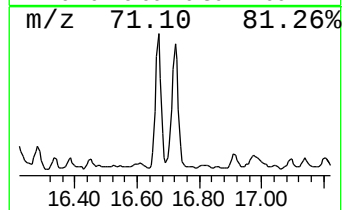
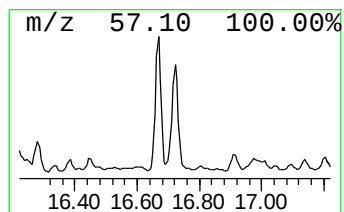
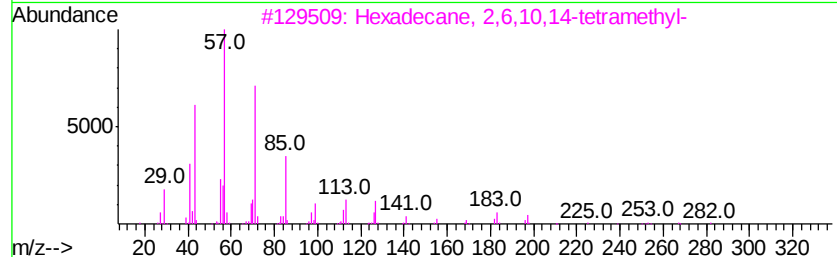
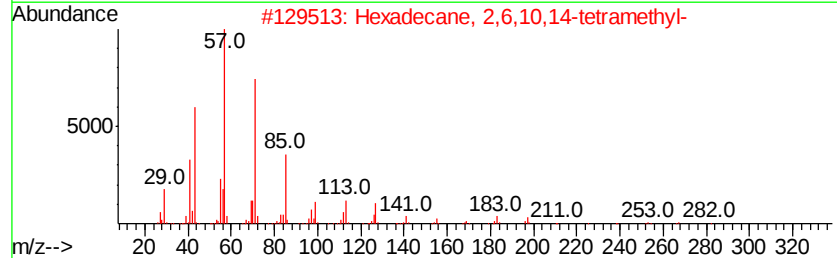
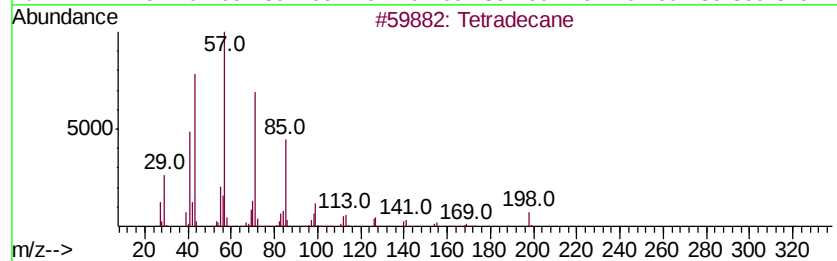
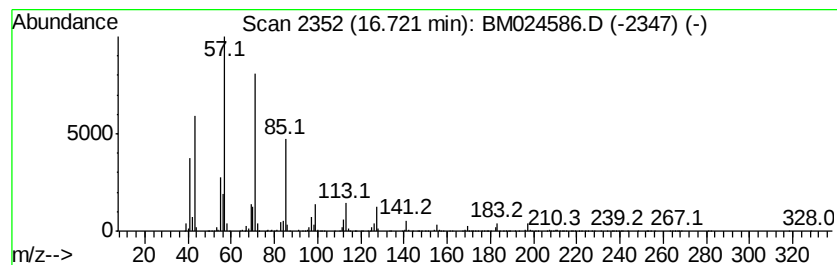
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.72 | 19.99 ng/ul | 2895580 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#         | Qual |
|------|----|------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Tetradecane                        | 198 | C14H30  | 000629-59-4  | 93   |
| 2    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8  | 90   |
| 3    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8  | 90   |
| 4    |    | 1-Iodo-2-methylnonane              | 268 | C10H21I | 1000101-47-9 | 87   |
| 5    |    | Tridecane, 7-hexyl-                | 268 | C19H40  | 007225-66-3  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

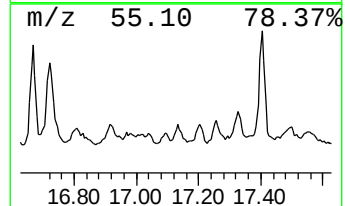
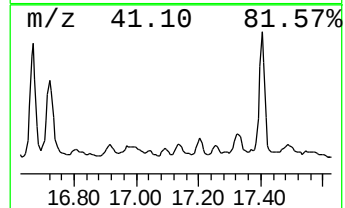
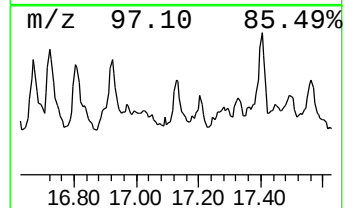
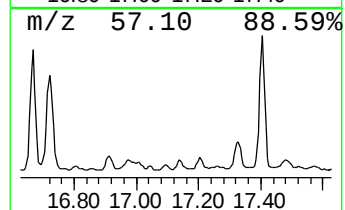
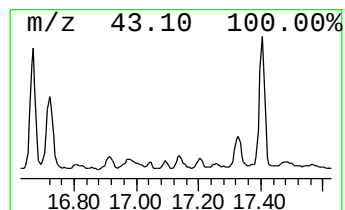
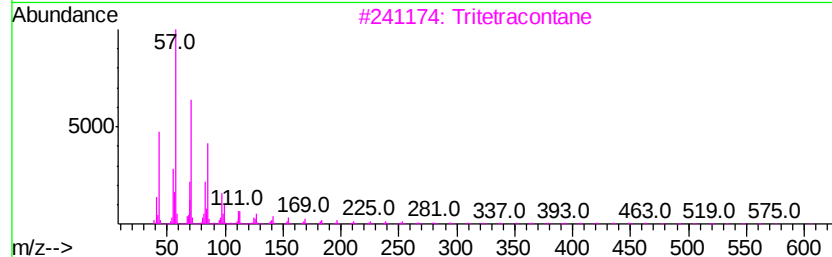
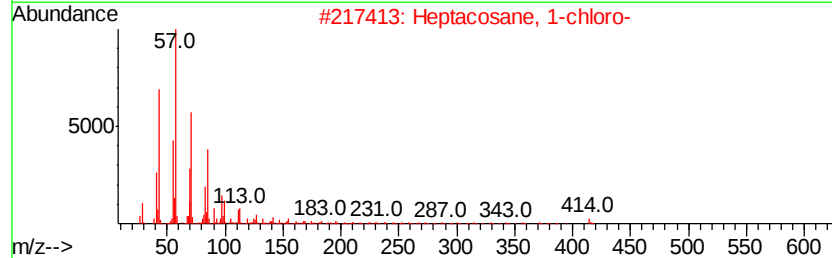
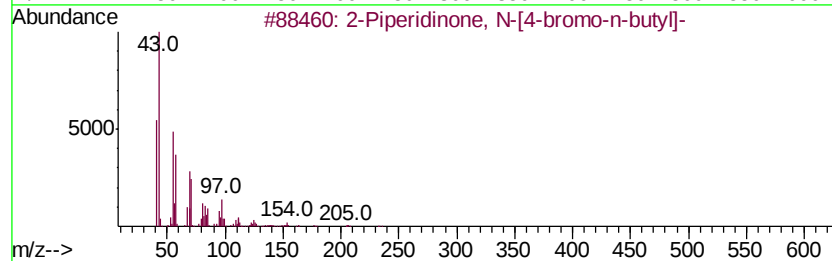
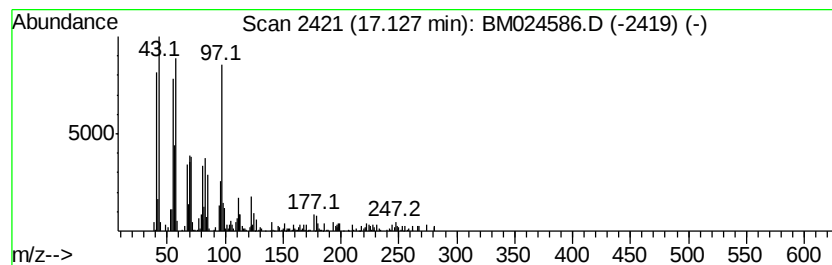
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 26 2-Piperidinone, N-[4-bromo-... Concentration Rank 37

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.13 | 2.58 ng/ul | 373277 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | 2-Piperidinone, N-[4-bromo-n-but... | 233 | C9H16BrNO | 195194-80-0 | 64   |
| 2    |    |   | Heptacosane, 1-chloro-              | 414 | C27H55Cl  | 062016-79-9 | 58   |
| 3    |    |   | Tritetracontane                     | 605 | C43H88    | 007098-21-7 | 52   |
| 4    |    |   | Heptadecane, 4-methyl-              | 254 | C18H38    | 026429-11-8 | 50   |
| 5    |    |   | Hexadecane, 2-methyl-               | 240 | C17H36    | 001560-92-5 | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

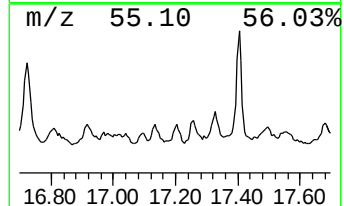
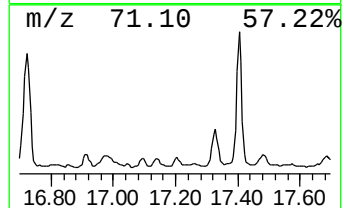
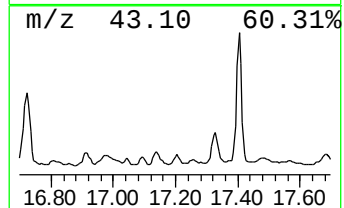
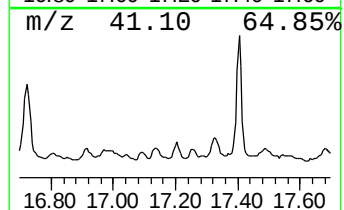
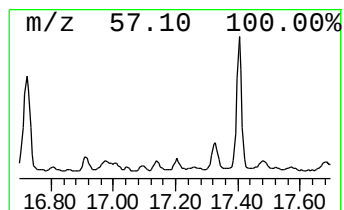
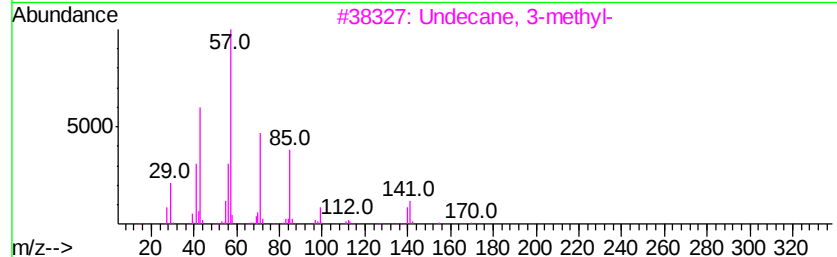
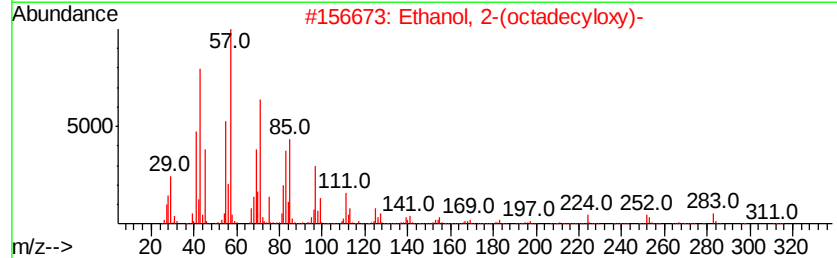
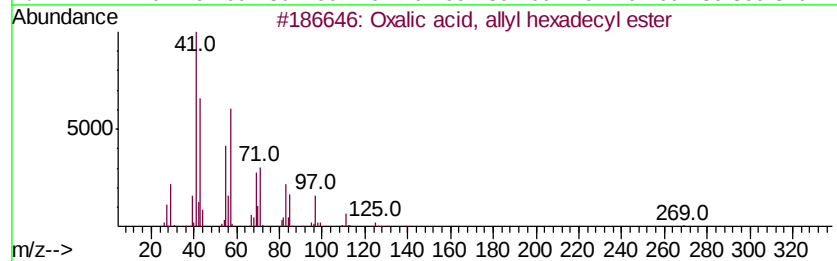
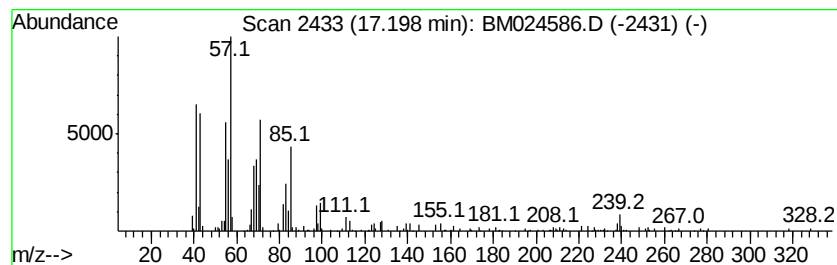
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 Oxalic acid, allyl hexadecy... Concentration Rank 39

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.20 | 2.47 ng/ul | 357980 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Oxalic acid, allyl hexadecyl ester  | 354 | C21H38O4   | 1000309-24-4 | 62   |
| 2    |    |   | Ethanol, 2-(octadecyloxy)-          | 314 | C20H42O2   | 002136-72-3  | 52   |
| 3    |    |   | Undecane, 3-methyl-                 | 170 | C12H26     | 001002-43-3  | 52   |
| 4    |    |   | 2-Chloropropionic acid, hexadec...  | 332 | C19H37ClO2 | 086711-81-1  | 52   |
| 5    |    |   | Propanal, 2-methyl-, 2-propenylh... | 126 | C7H14N2    | 066075-08-9  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

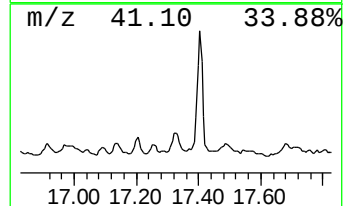
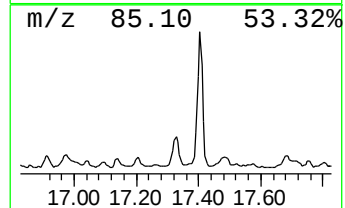
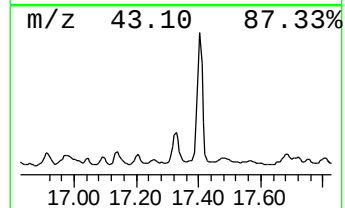
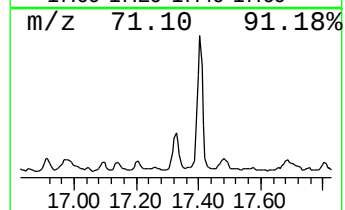
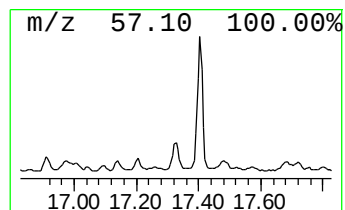
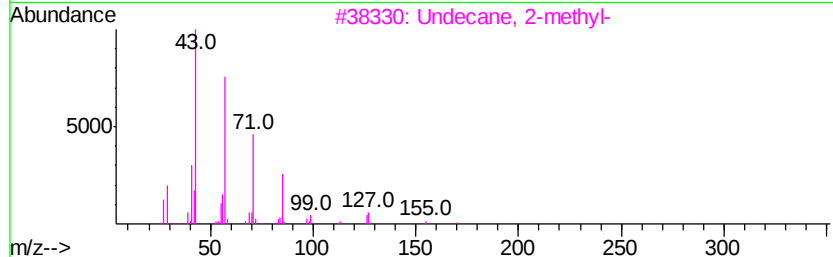
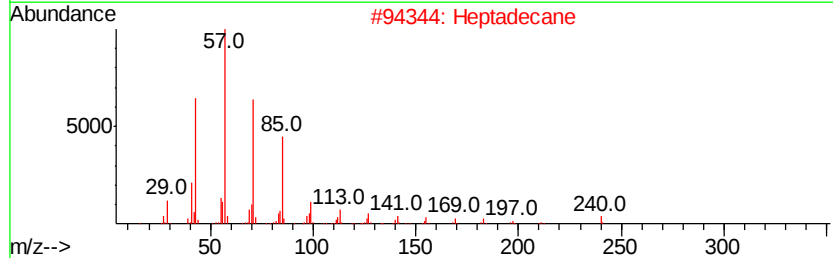
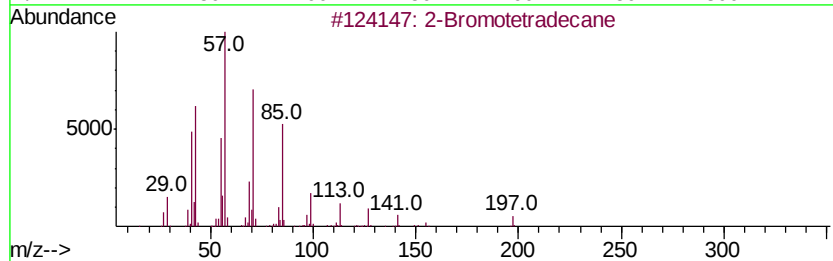
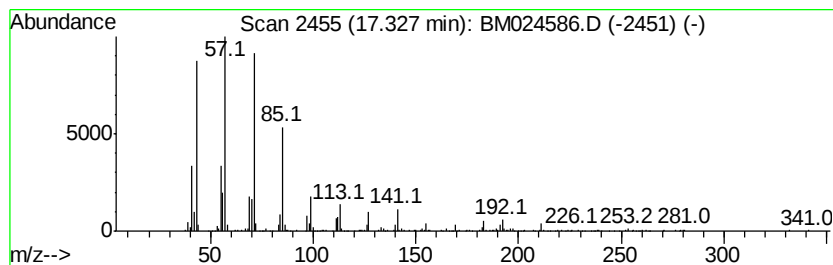
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.33 | 6.22 ng/ul | 901528 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Bromotetradecane                 | 276 | C14H29Br | 074036-95-6 | 90   |
| 2    |    |   | Heptadecane                        | 240 | C17H36   | 000629-78-7 | 87   |
| 3    |    |   | Undecane, 2-methyl-                | 170 | C12H26   | 007045-71-8 | 83   |
| 4    |    |   | Octadecane                         | 254 | C18H38   | 000593-45-3 | 80   |
| 5    |    |   | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42   | 000638-36-8 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

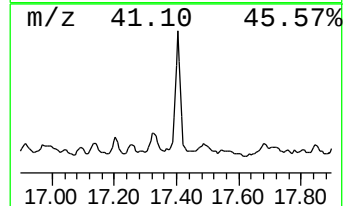
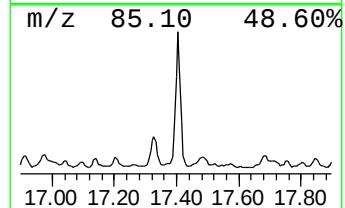
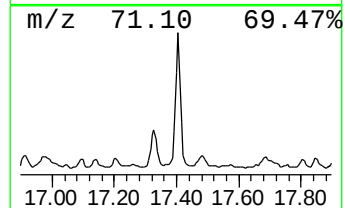
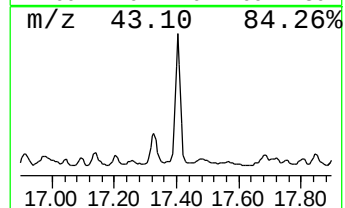
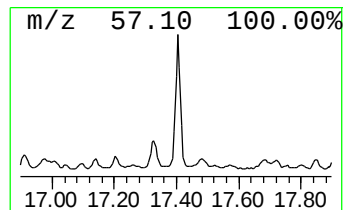
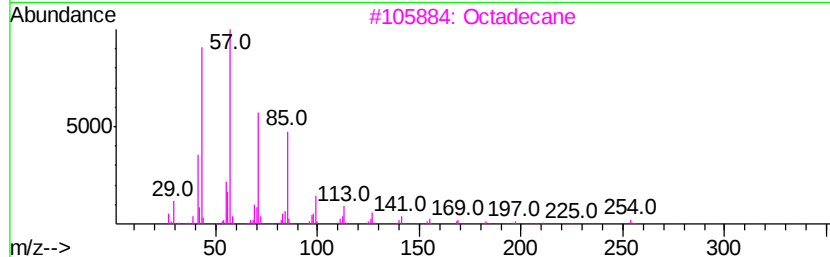
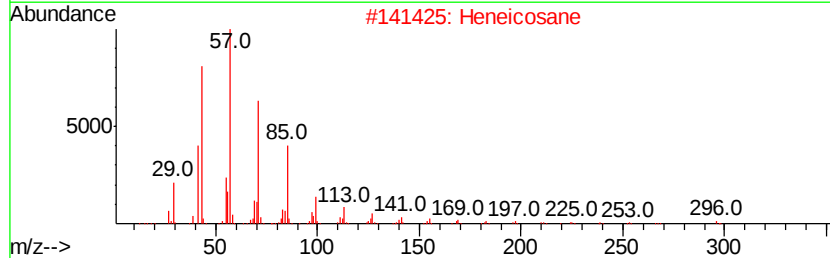
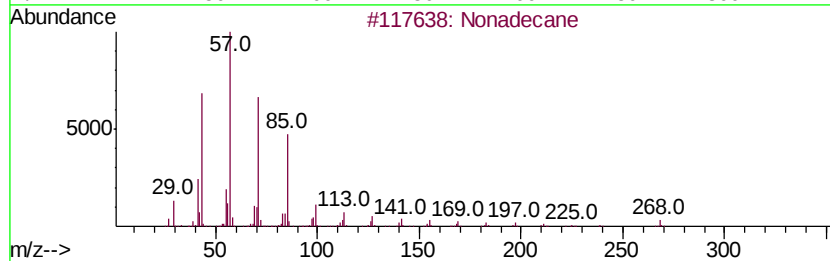
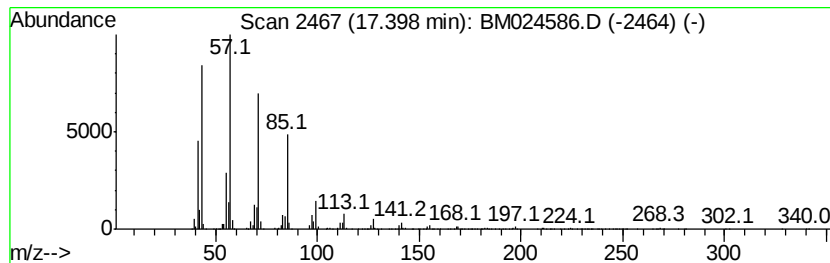
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.40 | 20.21 ng/ul | 2927810 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane   | 268 | C19H40  | 000629-92-5 | 96   |
| 2    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 91   |
| 4    |    | Eicosane     | 282 | C20H42  | 000112-95-8 | 91   |
| 5    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

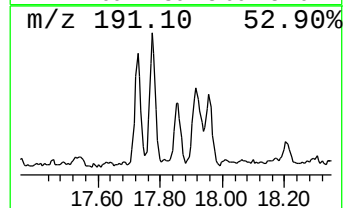
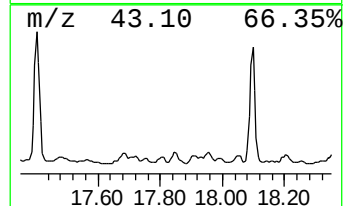
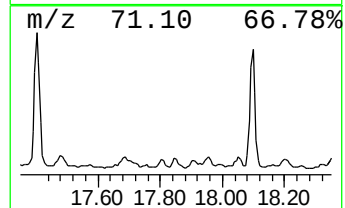
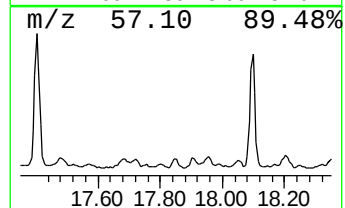
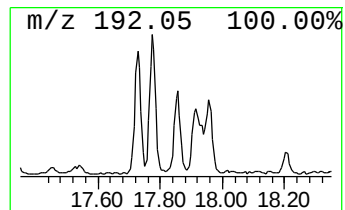
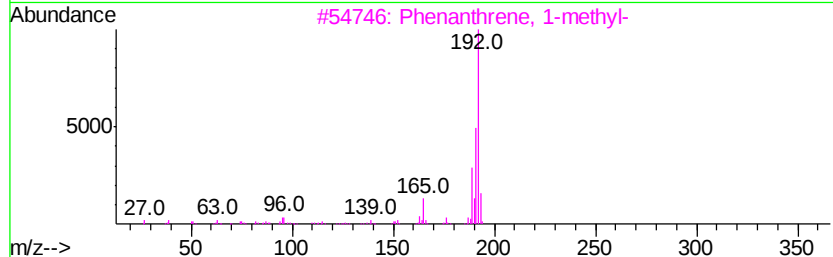
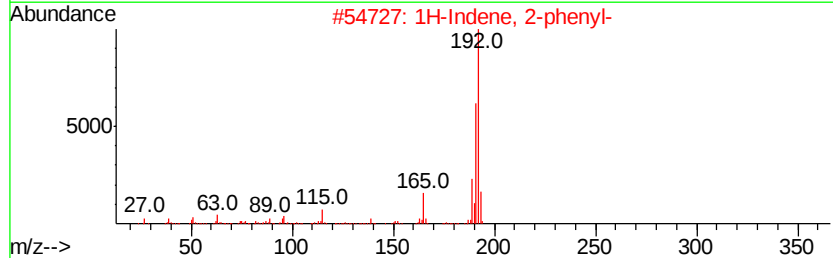
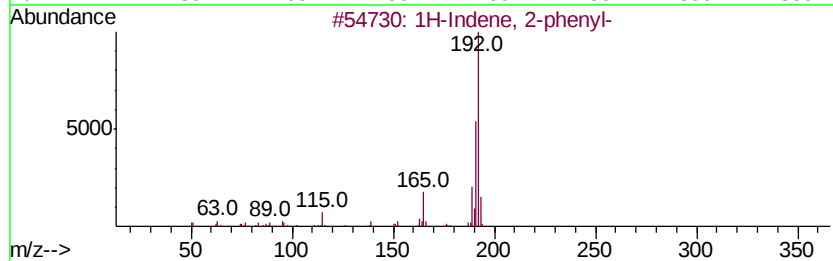
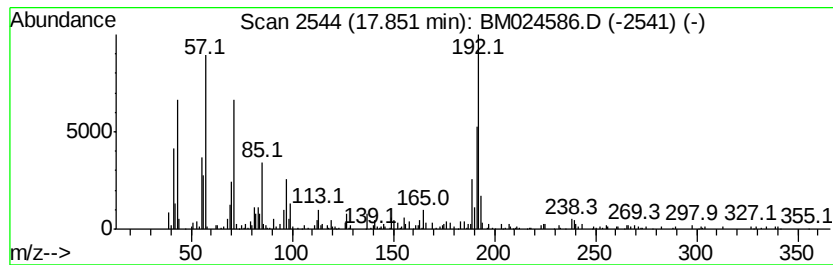
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 31 1H-Indene, 2-phenyl- Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.85 | 3.52 ng/ul | 509482 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID                            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2-phenyl-                    | 192 | C15H12  | 004505-48-0 | 70   |
| 2    |    |   | 1H-Indene, 2-phenyl-                    | 192 | C15H12  | 004505-48-0 | 55   |
| 3    |    |   | Phenanthrene, 1-methyl-                 | 192 | C15H12  | 000832-69-9 | 55   |
| 4    |    |   | 1H-Cyclopropa[1,1]phenanthrene, 1a, ... | 192 | C15H12  | 000949-41-7 | 55   |
| 5    |    |   | 1H-Indene, 2-phenyl-                    | 192 | C15H12  | 004505-48-0 | 55   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

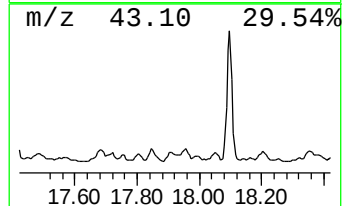
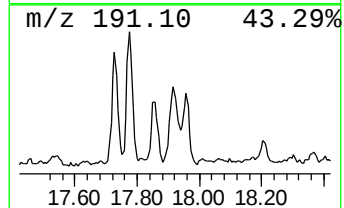
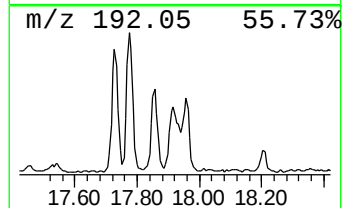
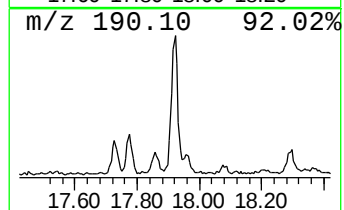
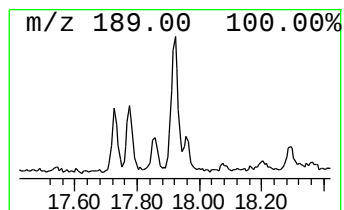
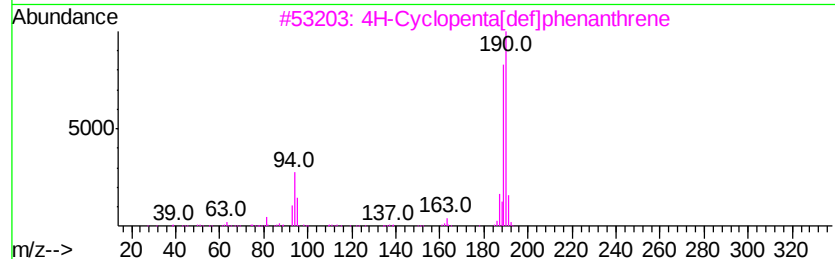
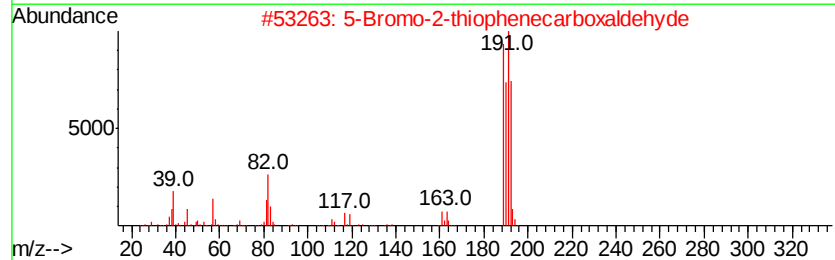
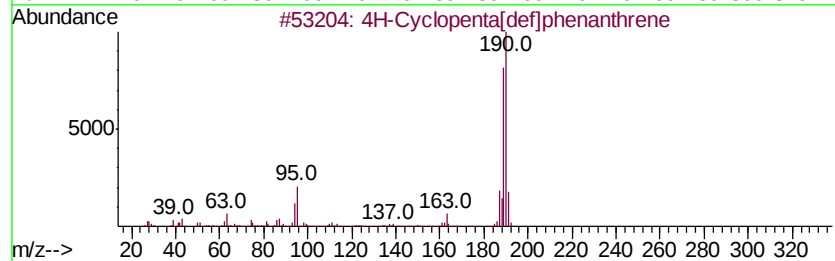
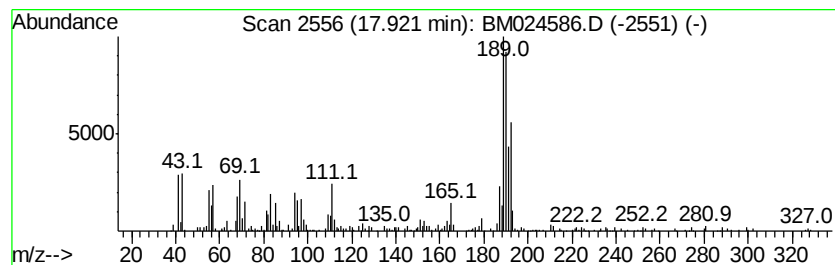
Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 32 4H-Cyclopenta[def]phenanthrene Concentration Rank 22

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.     |             |      |
|---------|------------|-------------------------------------|------------------|----------|-------------|------|
| 17.92   | 4.60 ng/ul | 666991                              | Phenanthrene-d10 | 16.85    |             |      |
| Hit# of | 5          | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1       |            | 4H-Cvclopenta[def]phenanthrene      | 190              | C15H10   | 000203-64-5 | 55   |
| 2       |            | 5-Bromo-2-thiophenecarboxaldehyde   | 190              | C5H3BrOS | 004701-17-1 | 50   |
| 3       |            | 4H-Cvclopenta[def]phenanthrene      | 190              | C15H10   | 000203-64-5 | 46   |
| 4       |            | Phenanthrene, 4-methyl-             | 192              | C15H12   | 000832-64-4 | 42   |
| 5       |            | 8,9-Dihydrocyclopenta[def]phenan... | 192              | C15H12   | 027410-55-5 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

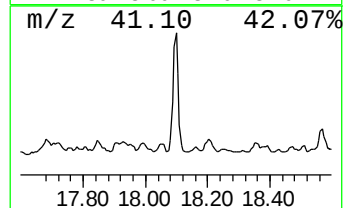
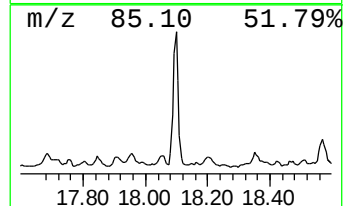
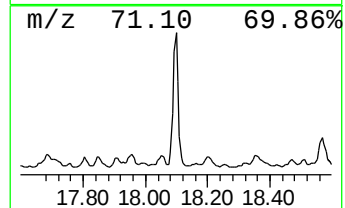
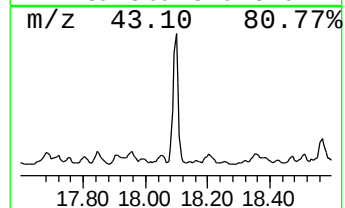
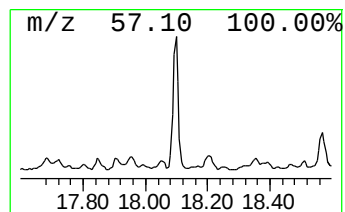
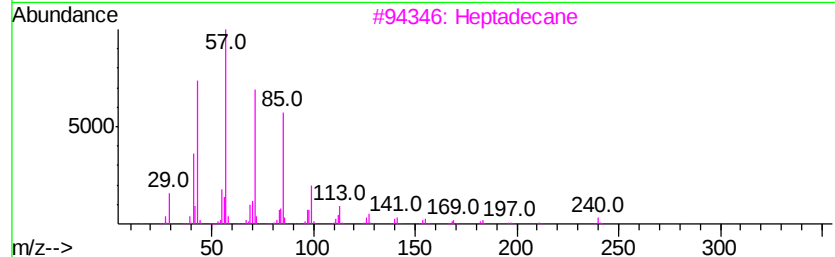
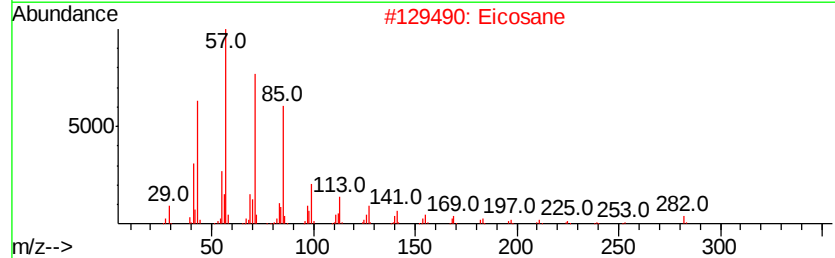
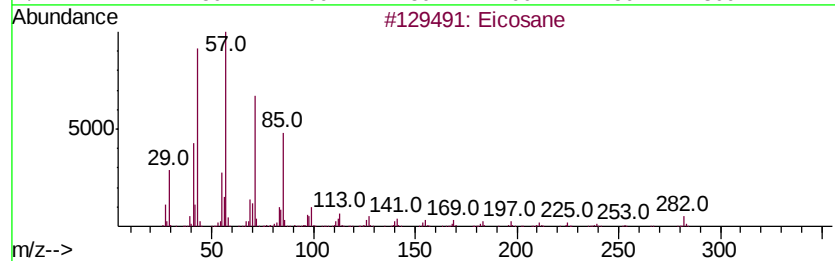
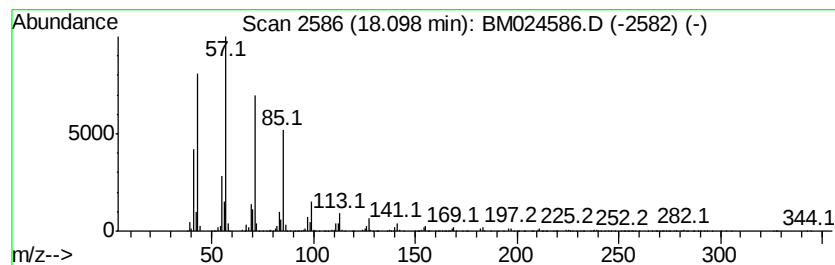
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.10 | 17.80 ng/ul | 2577840 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane              | 282 | C20H42  | 000112-95-8 | 97   |
| 2    |    | Eicosane              | 282 | C20H42  | 000112-95-8 | 96   |
| 3    |    | Heptadecane           | 240 | C17H36  | 000629-78-7 | 95   |
| 4    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 94   |
| 5    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

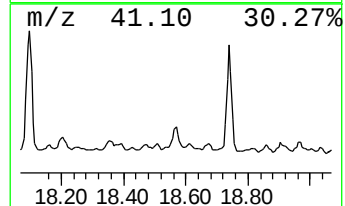
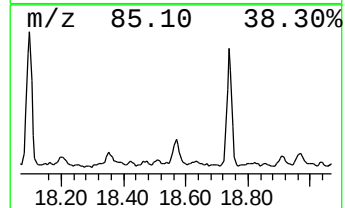
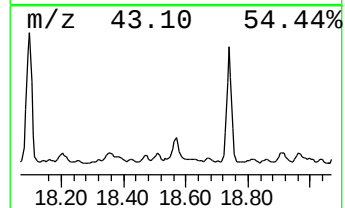
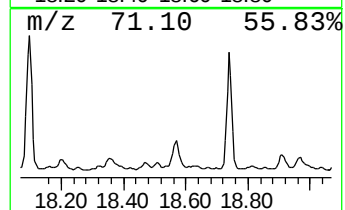
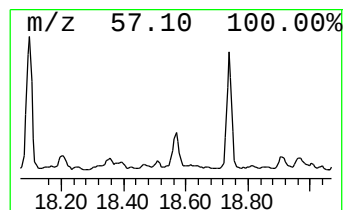
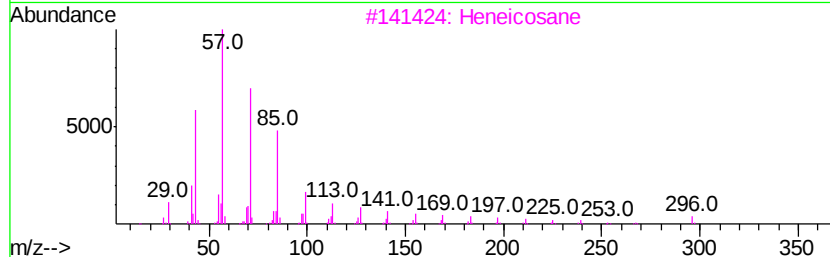
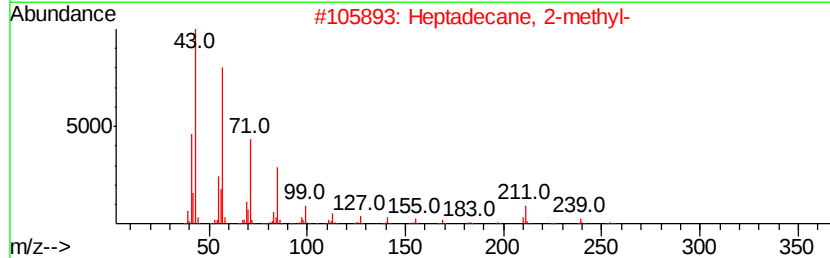
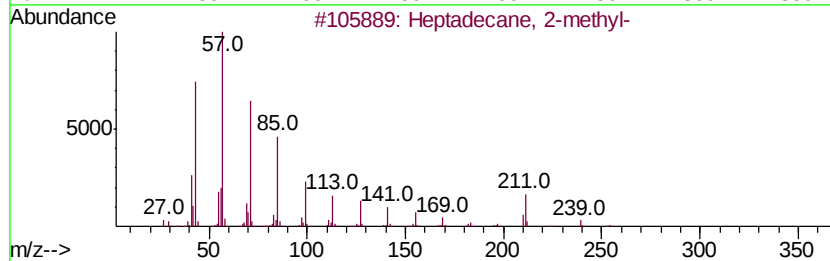
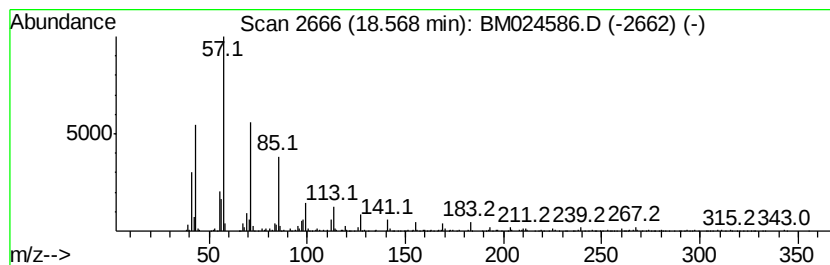
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.57 | 4.43 ng/ul | 642164 | Phenanthrene-d10 | 16.85 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 95   |
| 2    |    |   | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 91   |
| 3    |    |   | Heneicosane            | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    |   | Heptadecane, 3-methyl- | 254 | C18H38  | 006418-44-6 | 90   |
| 5    |    |   | Heptadecane, 3-methyl- | 254 | C18H38  | 006418-44-6 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

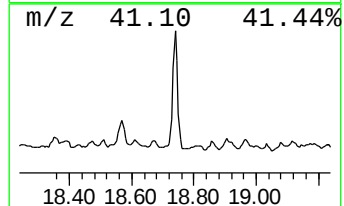
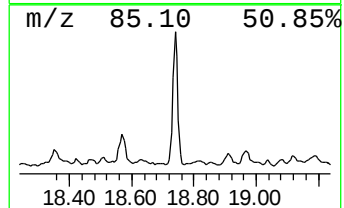
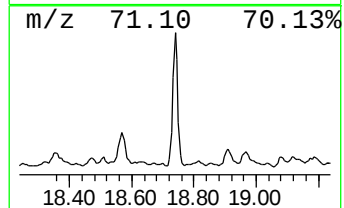
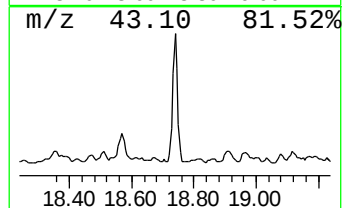
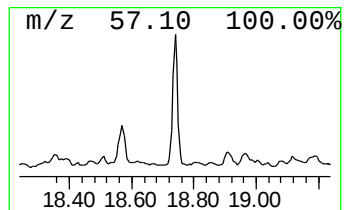
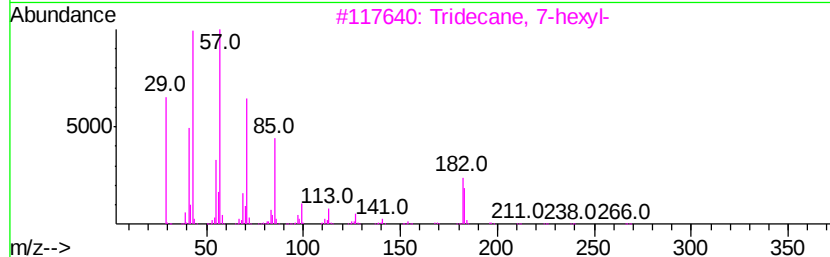
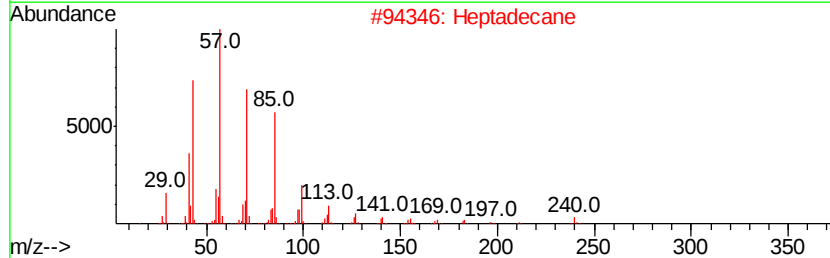
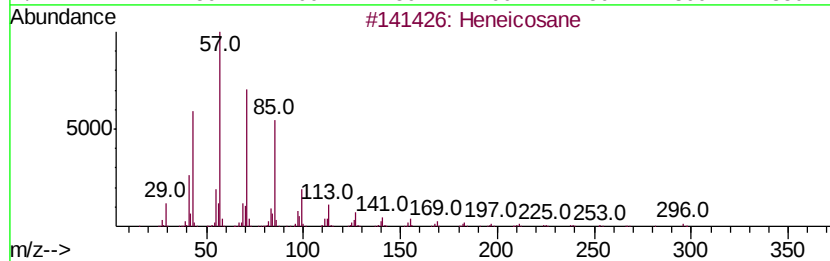
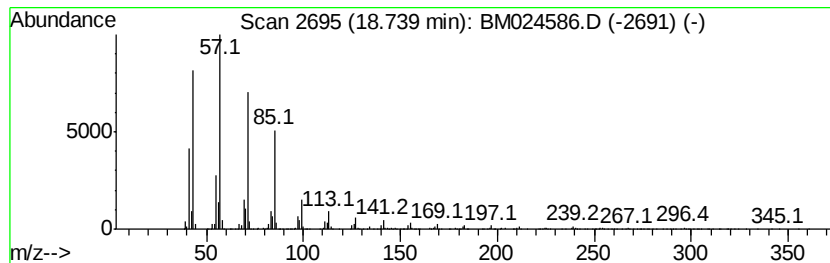
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.74 | 16.00 ng/ul | 2318210 | Phenanthrene-d10 | 16.85 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 98   |
| 2    |    | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 97   |
| 3    |    | Tridecane, 7-hexyl-                 | 268 | C19H40  | 007225-66-3 | 96   |
| 4    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 96   |
| 5    |    | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

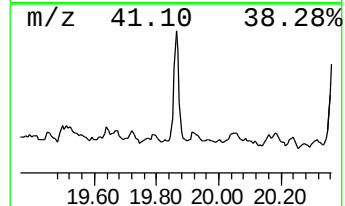
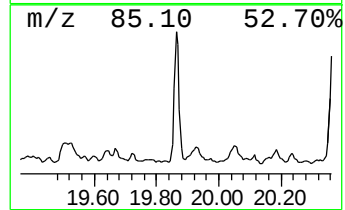
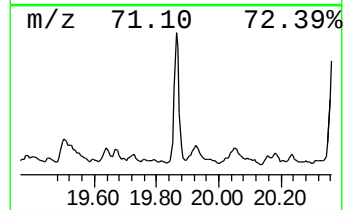
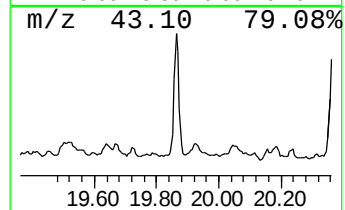
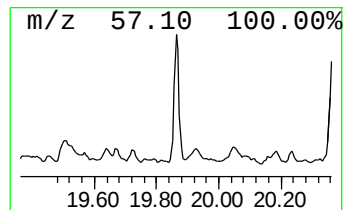
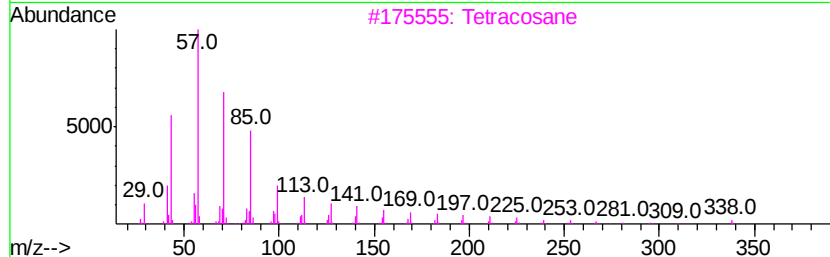
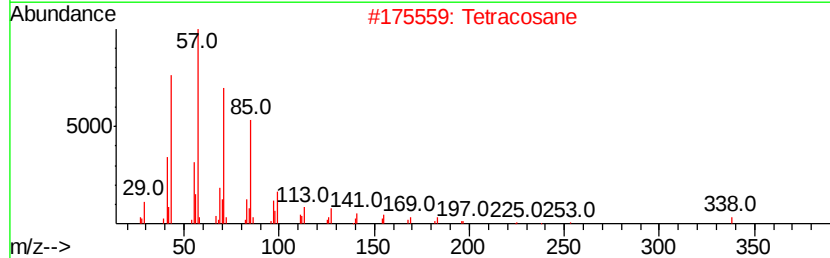
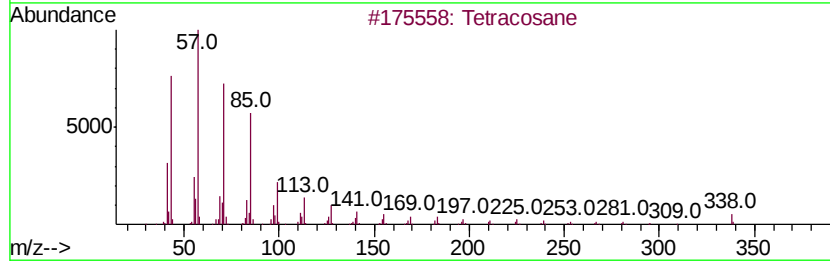
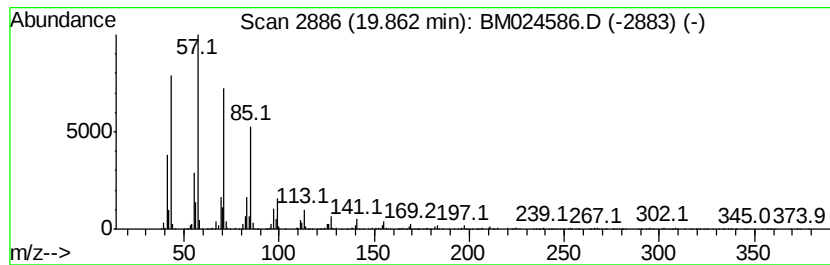
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 19.86 | 6.95 ng/ul | 1646600 | Chrysene-d12     | 21.06 |

| Hit# | of | 5 | Tentative ID     | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------|-----|----------|-------------|------|
| 1    |    |   | Tetracosane      | 338 | C24H50   | 000646-31-1 | 98   |
| 2    |    |   | Tetracosane      | 338 | C24H50   | 000646-31-1 | 98   |
| 3    |    |   | Tetracosane      | 338 | C24H50   | 000646-31-1 | 95   |
| 4    |    |   | 2-Bromo dodecane | 248 | C12H25Br | 013187-99-0 | 93   |
| 5    |    |   | Tetracosane      | 338 | C24H50   | 000646-31-1 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

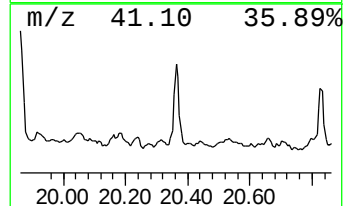
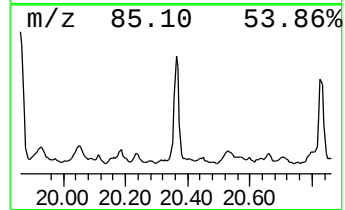
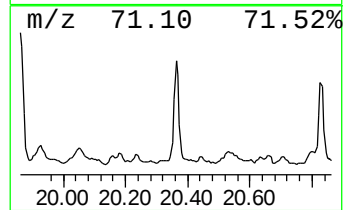
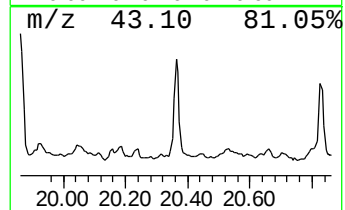
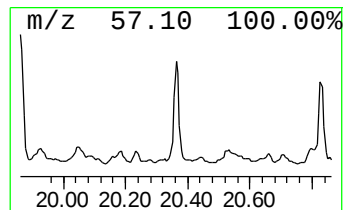
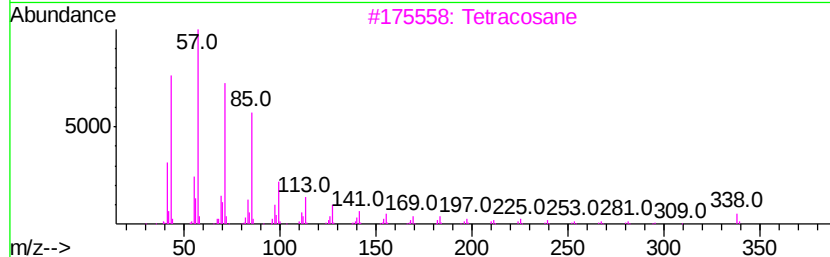
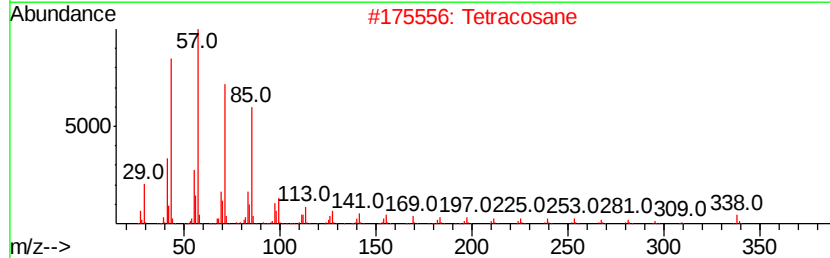
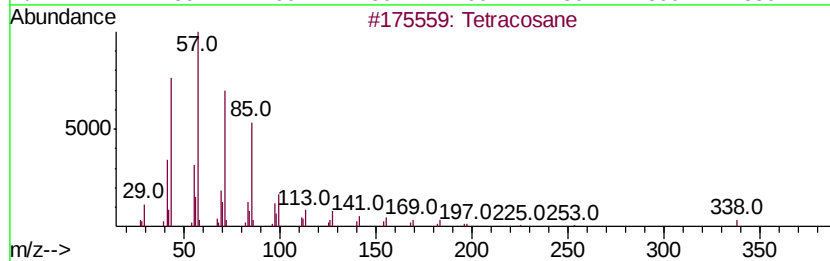
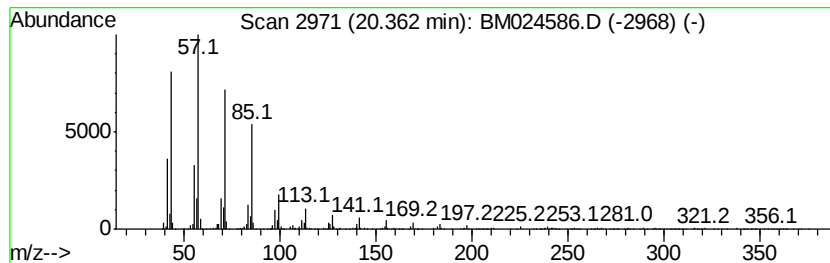
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 20.36 | 5.69 ng/ul | 1347410 | Chrysene-d12     | 21.06 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 98   |
| 2    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 98   |
| 3    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1 | 98   |
| 4    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 95   |
| 5    |    |   | Heptacosane           | 380 | C27H56  | 000593-49-7 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

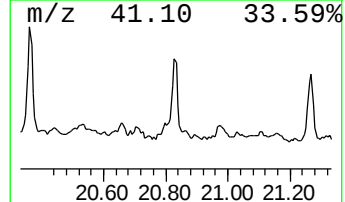
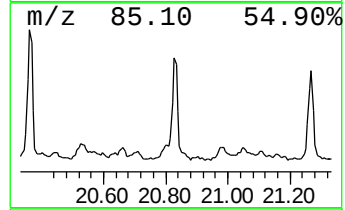
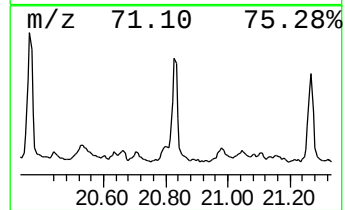
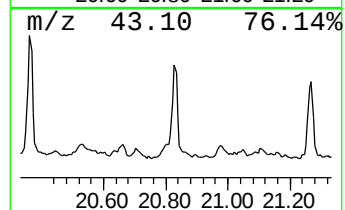
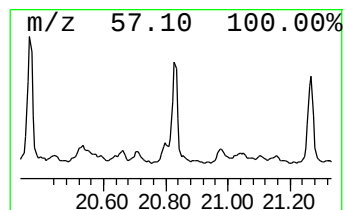
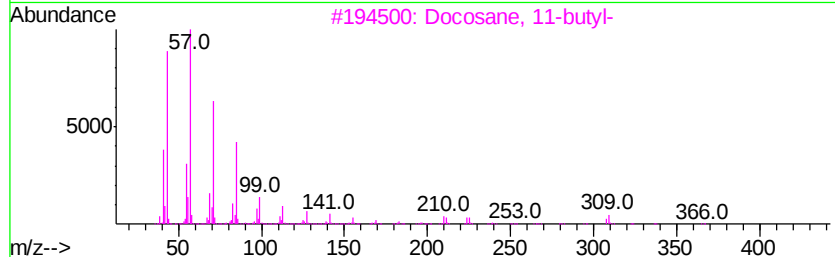
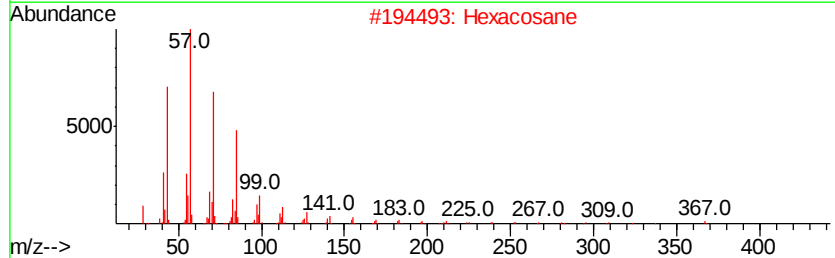
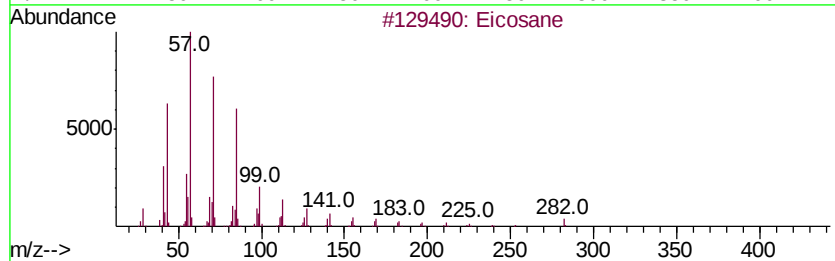
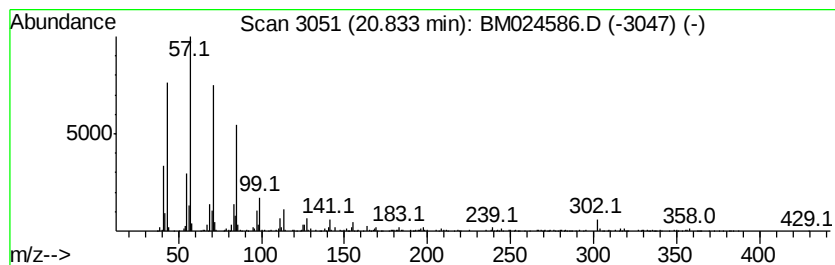
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 20.83 | 7.52 ng/ul | 1780650 | Chrysene-d12     | 21.06 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane            | 282 | C20H42  | 000112-95-8 | 93   |
| 2    |    |   | Hexacosane          | 366 | C26H54  | 000630-01-3 | 90   |
| 3    |    |   | Docosane, 11-butyl- | 366 | C26H54  | 013475-76-8 | 90   |
| 4    |    |   | Heneicosane         | 296 | C21H44  | 000629-94-7 | 87   |
| 5    |    |   | Tetracosane         | 338 | C24H50  | 000646-31-1 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

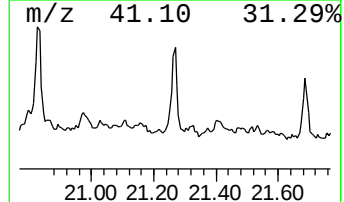
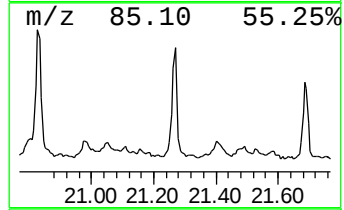
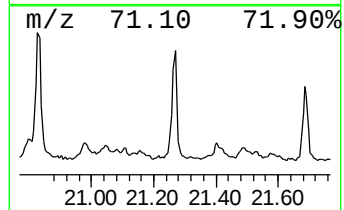
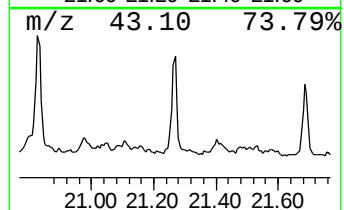
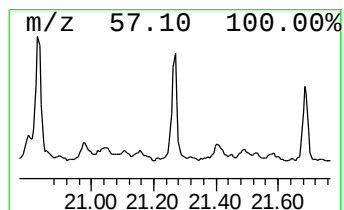
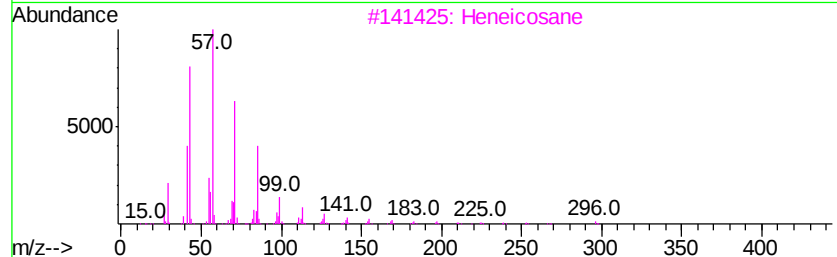
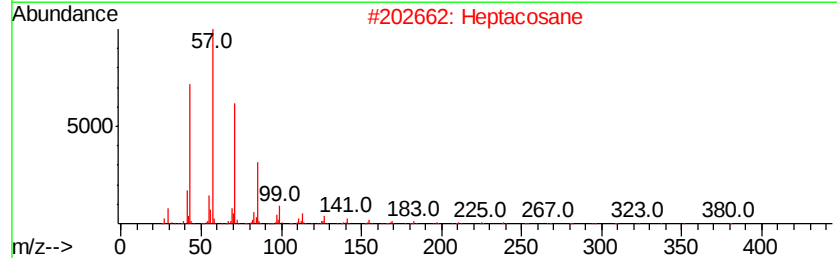
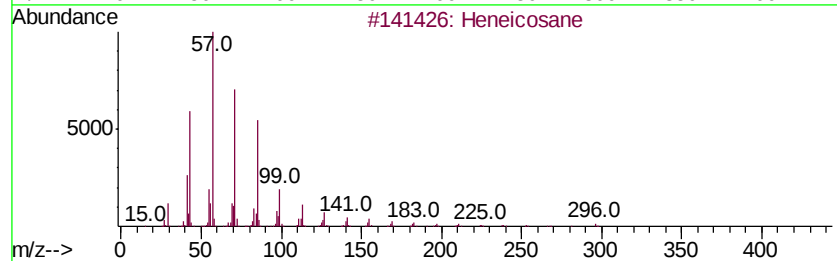
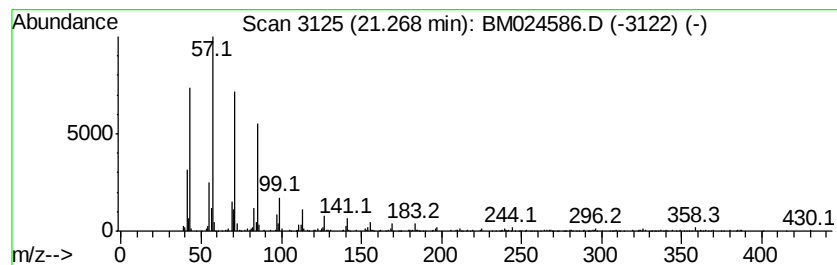
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.27 | 3.17 ng/ul | 750473 | Chrysene-d12     | 21.06 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 98   |
| 2    |    |   | Heptacosane                         | 380 | C27H56  | 000593-49-7 | 95   |
| 3    |    |   | Heneicosane                         | 296 | C21H44  | 000629-94-7 | 93   |
| 4    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 93   |
| 5    |    |   | Eicosane                            | 282 | C20H42  | 000112-95-8 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

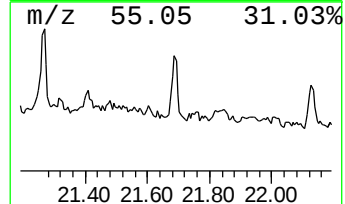
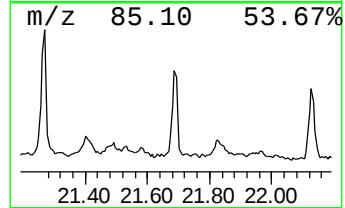
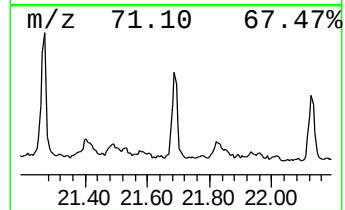
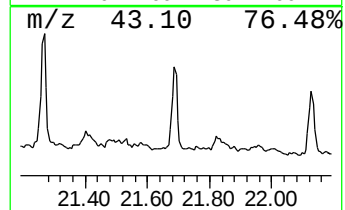
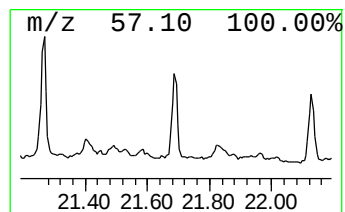
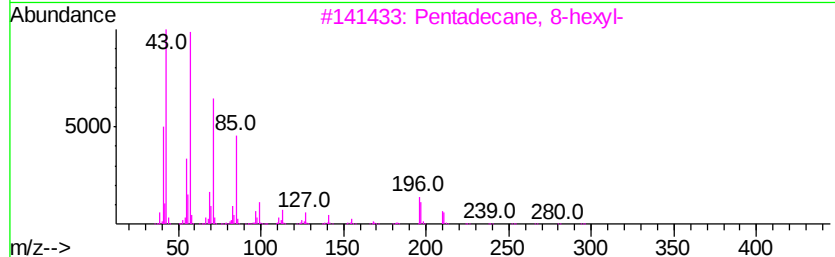
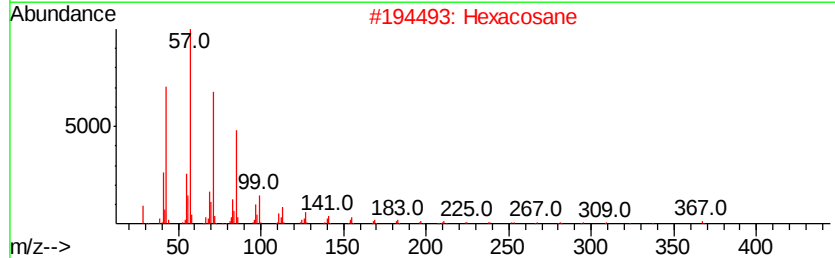
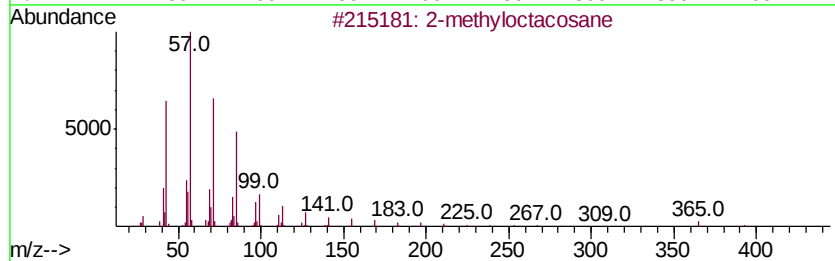
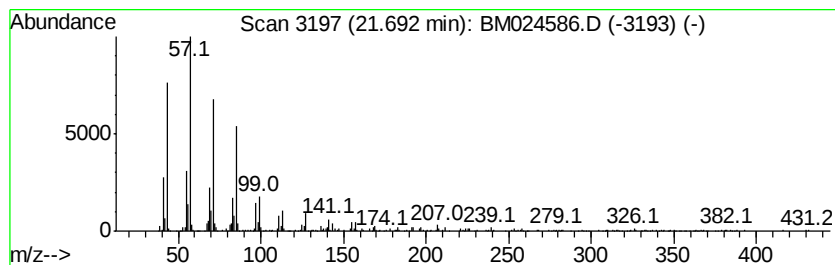
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.69 | 2.60 ng/ul | 615736 | Chrysene-d12     | 21.06 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|----|-----------------------|-----|---------|--------------|------|
| 1    | 5  | 2-methyloctacosane    | 408 | C29H60  | 1000376-72-8 | 90   |
| 2    |    | Hexacosane            | 366 | C26H54  | 000630-01-3  | 87   |
| 3    |    | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7  | 87   |
| 4    |    | Tricosane             | 324 | C23H48  | 000638-67-5  | 87   |
| 5    |    | Eicosane, 10-methyl-  | 296 | C21H44  | 054833-23-7  | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM012220\  
Data File : BM024586.D  
Acq On : 22 Jan 2020 17:46  
Operator : JU  
Sample : L1118-11DL2 20X  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION

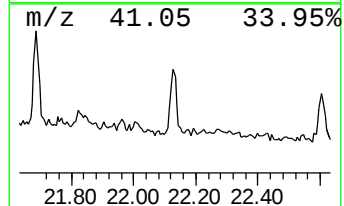
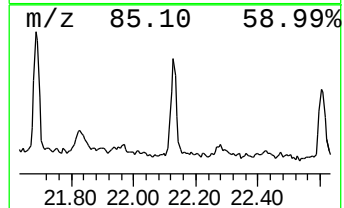
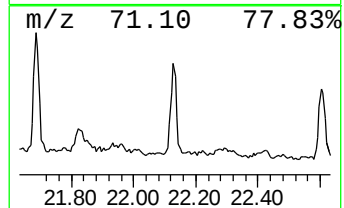
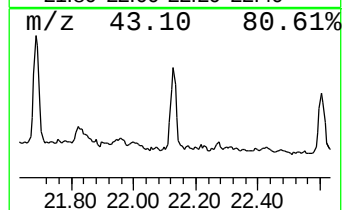
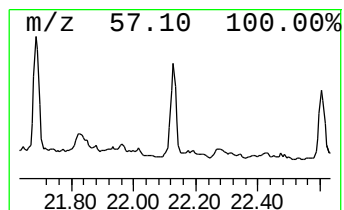
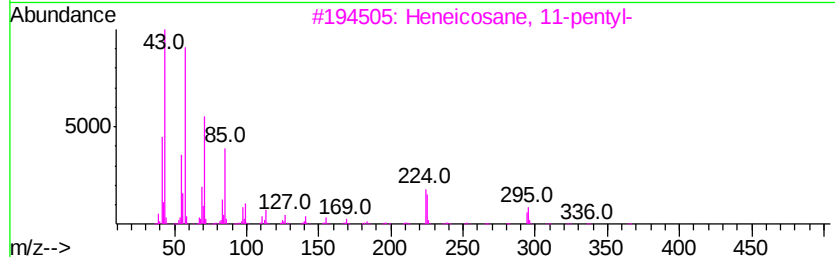
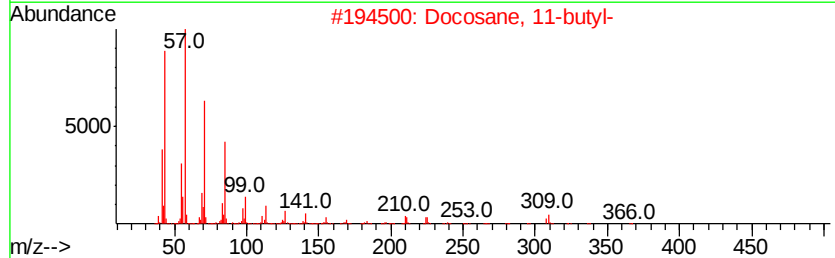
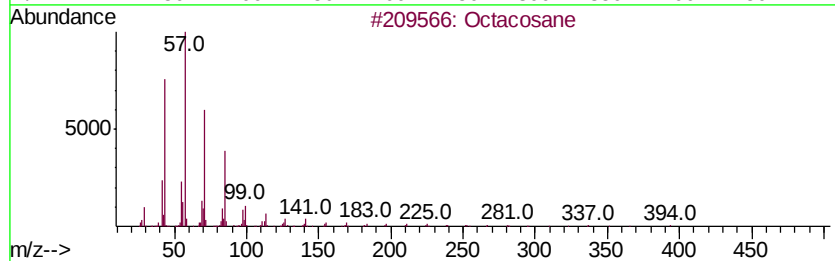
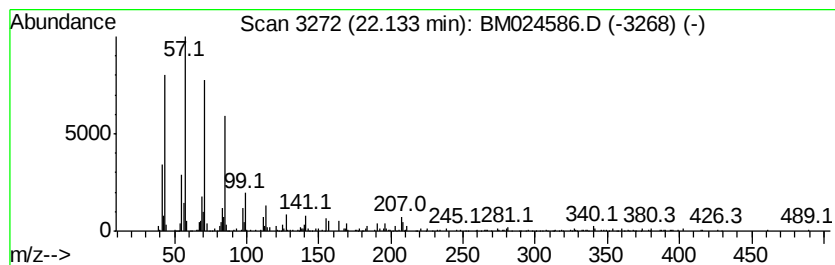
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 30

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.13 | 3.20 ng/ul | 444253 | Perylene-d12     | 23.18 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane              | 394 | C28H58  | 000630-02-4 | 97   |
| 2    |    |   | Docosane, 11-butyl-     | 366 | C26H54  | 013475-76-8 | 93   |
| 3    |    |   | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1 | 93   |
| 4    |    |   | Tricosane               | 324 | C23H48  | 000638-67-5 | 91   |
| 5    |    |   | Heneicosane             | 296 | C21H44  | 000629-94-7 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM012220\  
 Data File : BM024586.D  
 Acq On : 22 Jan 2020 17:46  
 Operator : JU  
 Sample : L1118-11DL2 20X  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 GASK2DL2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Quant Title : SVOA CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Benzene, 1,3-dime... | 5.15  | 5.5     | ng/ul | 332638   | 1                     | 7.45  | 1204560 | 20.0 |
| Benzene, 1-ethyl-... | 7.12  | 9.8     | ng/ul | 589656   | 1                     | 7.45  | 1204560 | 20.0 |
| Indene               | 7.98  | 12.0    | ng/ul | 720647   | 1                     | 7.45  | 1204560 | 20.0 |
| (DEL) Alkane: Str... | 8.70  | 3.7     | ng/ul | 221609   | 1                     | 7.45  | 1204560 | 20.0 |
| Benzene, 1-butyryl-  | 9.65  | 4.1     | ng/ul | 408764   | 2                     | 10.21 | 2006870 | 20.0 |
| (DEL) Alkane: Str... | 11.69 | 2.6     | ng/ul | 257205   | 2                     | 10.21 | 2006870 | 20.0 |
| Naphthalene, 1-me... | 12.11 | 8.9     | ng/ul | 896815   | 2                     | 10.21 | 2006870 | 20.0 |
| Naphthalene, 1,5-... | 13.26 | 2.2     | ng/ul | 316367   | 3                     | 14.10 | 2866980 | 20.0 |
| Naphthalene, 2,6-... | 13.42 | 2.9     | ng/ul | 408896   | 3                     | 14.10 | 2866980 | 20.0 |
| Naphthalene, 2-et... | 13.55 | 2.8     | ng/ul | 394230   | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Str... | 13.63 | 6.3     | ng/ul | 909005   | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Str... | 14.56 | 2.9     | ng/ul | 420304   | 3                     | 14.10 | 2866980 | 20.0 |
| 2-Bromotetradecane   | 14.67 | 3.4     | ng/ul | 486503   | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Str... | 14.91 | 2.4     | ng/ul | 342909   | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Str... | 15.01 | 10.6    | ng/ul | 1515010  | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Str... | 15.42 | 18.4    | ng/ul | 2643710  | 3                     | 14.10 | 2866980 | 20.0 |
| (DEL) Alkane: Cyc... | 15.57 | 3.1     | ng/ul | 450333   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 15.87 | 19.6    | ng/ul | 2831400  | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 15.90 | 24.3    | ng/ul | 3515880  | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 16.22 | 5.8     | ng/ul | 842787   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 16.28 | 4.3     | ng/ul | 615651   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 16.67 | 20.7    | ng/ul | 2994180  | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 16.72 | 20.0    | ng/ul | 2895580  | 4                     | 16.85 | 2896910 | 20.0 |
| 2-Piperidinone, N... | 17.13 | 2.6     | ng/ul | 373277   | 4                     | 16.85 | 2896910 | 20.0 |
| Oxalic acid, ally... | 17.20 | 2.5     | ng/ul | 357980   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 17.33 | 6.2     | ng/ul | 901528   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 17.40 | 20.2    | ng/ul | 2927810  | 4                     | 16.85 | 2896910 | 20.0 |
| 1H-Indene, 2-phenyl- | 17.85 | 3.5     | ng/ul | 509482   | 4                     | 16.85 | 2896910 | 20.0 |
| 4H-Cyclopenta[def... | 17.92 | 4.6     | ng/ul | 666991   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 18.10 | 17.8    | ng/ul | 2577840  | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 18.57 | 4.4     | ng/ul | 642164   | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 18.74 | 16.0    | ng/ul | 2318210  | 4                     | 16.85 | 2896910 | 20.0 |
| (DEL) Alkane: Str... | 19.86 | 7.0     | ng/ul | 1646600  | 5                     | 21.06 | 4738840 | 20.0 |
| (DEL) Alkane: Str... | 20.36 | 5.7     | ng/ul | 1347410  | 5                     | 21.06 | 4738840 | 20.0 |
| (DEL) Alkane: Str... | 20.83 | 7.5     | ng/ul | 1780650  | 5                     | 21.06 | 4738840 | 20.0 |
| (DEL) Alkane: Str... | 21.27 | 3.2     | ng/ul | 750473   | 5                     | 21.06 | 4738840 | 20.0 |
| (DEL) Alkane: Str... | 21.69 | 2.6     | ng/ul | 615736   | 5                     | 21.06 | 4738840 | 20.0 |
| (DEL) Alkane: Str... | 22.13 | 3.2     | ng/ul | 444253   | 6                     | 23.18 | 2778070 | 20.0 |