

Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
 Data File : BM013799.D  
 Acq On : 25 Jan 2018 09:43  
 Operator : SJ/JU  
 Sample : J1222-13  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6A95

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 1 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011018.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         | 8.810       | 985           | 990         | 999          | rVB      | 2081790        | 3249436       | 1.12%           | 0.264%        |
| 2         | 10.351      | 1244          | 1252        | 1260         | rBV2     | 4359627        | 8296567       | 2.85%           | 0.673%        |
| 3         | 11.398      | 1424          | 1430        | 1434         | rBV      | 2930392        | 4868154       | 1.67%           | 0.395%        |
| 4         | 11.804      | 1491          | 1499        | 1504         | rBV      | 6940830        | 10758390      | 3.70%           | 0.873%        |
| 5         | 12.774      | 1660          | 1664        | 1669         | rVB      | 3543550        | 4650814       | 1.60%           | 0.377%        |
| 6         | 13.074      | 1710          | 1715        | 1722         | rVB      | 12230346       | 17252703      | 5.93%           | 1.400%        |
| 7         | 13.657      | 1810          | 1814        | 1819         | rVB3     | 2871070        | 4585741       | 1.58%           | 0.372%        |
| 8         | 13.739      | 1820          | 1828        | 1832         | rBV2     | 6921030        | 10719259      | 3.68%           | 0.870%        |
| 9         | 13.780      | 1832          | 1835        | 1843         | rVB2     | 2360971        | 3993749       | 1.37%           | 0.324%        |
| 10        | 14.168      | 1894          | 1901        | 1905         | rBV      | 18033188       | 23173244      | 7.96%           | 1.881%        |
| 11        | 14.239      | 1905          | 1913        | 1921         | rVV2     | 6419709        | 12072539      | 4.15%           | 0.980%        |
| 12        | 14.321      | 1921          | 1927        | 1935         | rVV      | 7342892        | 12142296      | 4.17%           | 0.985%        |
| 13        | 14.398      | 1936          | 1940        | 1947         | rVB      | 2948521        | 4113548       | 1.41%           | 0.334%        |
| 14        | 14.657      | 1980          | 1984        | 1990         | rVB2     | 5502863        | 8299465       | 2.85%           | 0.674%        |
| 15        | 14.786      | 2001          | 2006        | 2008         | rBV2     | 5022851        | 8021696       | 2.76%           | 0.651%        |
| 16        | 14.851      | 2013          | 2017        | 2019         | rVV      | 2610935        | 3353326       | 1.15%           | 0.272%        |
| 17        | 14.892      | 2019          | 2024        | 2030         | rVB      | 6595096        | 9988817       | 3.43%           | 0.811%        |
| 18        | 15.021      | 2042          | 2046        | 2054         | rVB4     | 1581742        | 3577443       | 1.23%           | 0.290%        |
| 19        | 15.127      | 2058          | 2064        | 2068         | rVB      | 24099899       | 29733408      | 10.21%          | 2.413%        |
| 20        | 15.310      | 2090          | 2095        | 2099         | rBV      | 5374453        | 7927319       | 2.72%           | 0.643%        |
| 21        | 15.527      | 2125          | 2132        | 2137         | rBV      | 9958583        | 16035910      | 5.51%           | 1.301%        |
| 22        | 15.621      | 2143          | 2148        | 2153         | rVB4     | 2241131        | 4212922       | 1.45%           | 0.342%        |
| 23        | 15.674      | 2154          | 2157        | 2161         | rBV2     | 2675585        | 4022738       | 1.38%           | 0.326%        |
| 24        | 15.727      | 2161          | 2166        | 2177         | rVB4     | 9764915        | 16761930      | 5.76%           | 1.360%        |
| 25        | 15.998      | 2205          | 2212        | 2214         | rBV      | 28582956       | 47428311      | 16.29%          | 3.849%        |
| 26        | 16.021      | 2214          | 2216        | 2219         | rVV      | 30136602       | 33435695      | 11.49%          | 2.713%        |
| 27        | 16.051      | 2219          | 2221        | 2226         | rVB      | 5379363        | 6110128       | 2.10%           | 0.496%        |
| 28        | 16.504      | 2294          | 2298        | 2301         | rBV      | 2648808        | 3286310       | 1.13%           | 0.267%        |
| 29        | 16.721      | 2321          | 2335        | 2340         | rVV2     | 111078217      | 291094809     | 100.00%         | 23.623%       |
| 30        | 16.792      | 2340          | 2347        | 2351         | rVV      | 38310268       | 52639158      | 18.08%          | 4.272%        |
| 31        | 16.839      | 2351          | 2355        | 2365         | rVB2     | 13544991       | 20543883      | 7.06%           | 1.667%        |
| 32        | 16.921      | 2366          | 2369        | 2373         | rBV2     | 2376879        | 3360937       | 1.15%           | 0.273%        |
| 33        | 17.027      | 2382          | 2387        | 2390         | rBV3     | 3917853        | 7055733       | 2.42%           | 0.573%        |
| 34        | 17.074      | 2390          | 2395        | 2399         | rVV      | 22371101       | 34201778      | 11.75%          | 2.776%        |

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ALS Vial : 33 Sample Multiplier: 1

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Max Peaks: 100  
Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
Peak separation: 5

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011018.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |        |        |
|----|--------|------|------|------|------|----------|----------|--------|--------|
| 35 | 17.121 | 2399 | 2403 | 2406 | rVV3 | 5908606  | 9688134  | 3.33%  | 0.786% |
| 36 | 17.156 | 2406 | 2409 | 2414 | rVB  | 8623416  | 11235918 | 3.86%  | 0.912% |
| 37 | 17.245 | 2420 | 2424 | 2430 | rVB3 | 5007953  | 7894581  | 2.71%  | 0.641% |
| 38 | 17.321 | 2434 | 2437 | 2443 | rVB4 | 2649196  | 3810665  | 1.31%  | 0.309% |
| 39 | 17.533 | 2467 | 2473 | 2476 | rBV  | 35289737 | 46166798 | 15.86% | 3.747% |
| 40 | 17.803 | 2515 | 2519 | 2527 | rBV4 | 3833381  | 7395561  | 2.54%  | 0.600% |
| 41 | 18.227 | 2585 | 2591 | 2595 | rBV  | 35951151 | 44678403 | 15.35% | 3.626% |
| 42 | 18.445 | 2625 | 2628 | 2631 | rBV2 | 2870146  | 3460283  | 1.19%  | 0.281% |
| 43 | 18.862 | 2694 | 2699 | 2706 | rVB  | 35967532 | 44977526 | 15.45% | 3.650% |
| 44 | 18.962 | 2713 | 2716 | 2722 | rBV  | 3718803  | 5542638  | 1.90%  | 0.450% |
| 45 | 19.015 | 2723 | 2725 | 2732 | rVV5 | 2416064  | 3749828  | 1.29%  | 0.304% |
| 46 | 19.098 | 2732 | 2739 | 2745 | rVB  | 28442565 | 43691337 | 15.01% | 3.546% |
| 47 | 19.445 | 2791 | 2798 | 2807 | rVB2 | 47599957 | 79993964 | 27.48% | 6.492% |
| 48 | 19.756 | 2849 | 2851 | 2861 | rVB3 | 5075951  | 10666903 | 3.66%  | 0.866% |
| 49 | 19.980 | 2885 | 2889 | 2893 | rVB  | 33610281 | 37641174 | 12.93% | 3.055% |
| 50 | 20.480 | 2969 | 2974 | 2978 | rBV  | 28247687 | 31326320 | 10.76% | 2.542% |
| 51 | 20.939 | 3049 | 3052 | 3058 | rVB  | 26214655 | 30099953 | 10.34% | 2.443% |
| 52 | 21.209 | 3095 | 3098 | 3101 | rBV2 | 4478038  | 6887126  | 2.37%  | 0.559% |
| 53 | 21.374 | 3123 | 3126 | 3132 | rVB  | 18894244 | 20496211 | 7.04%  | 1.663% |
| 54 | 21.803 | 3195 | 3199 | 3209 | rVB  | 14970596 | 19273209 | 6.62%  | 1.564% |
| 55 | 22.262 | 3273 | 3277 | 3283 | rVB  | 9033479  | 11690194 | 4.02%  | 0.949% |
| 56 | 22.756 | 3357 | 3361 | 3363 | rBV  | 8479470  | 11931664 | 4.10%  | 0.968% |
| 57 | 23.303 | 3450 | 3454 | 3458 | rBV3 | 4866593  | 8964668  | 3.08%  | 0.728% |

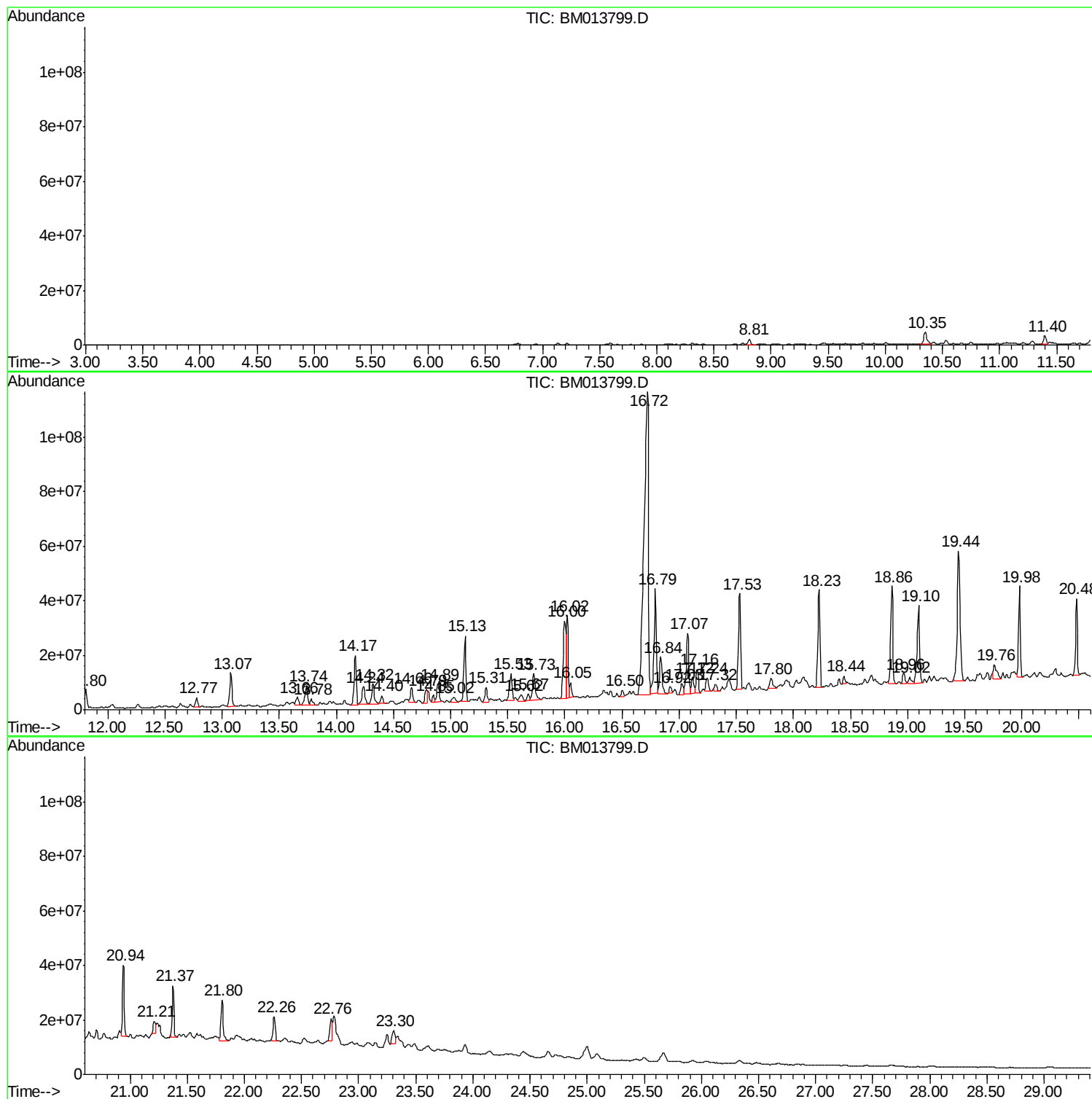
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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



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Quant Title : SVOA CALIBRATION

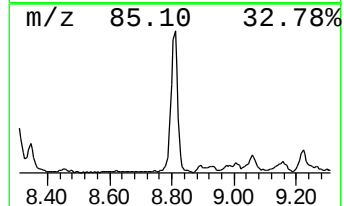
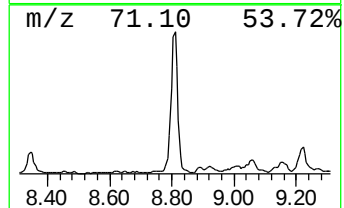
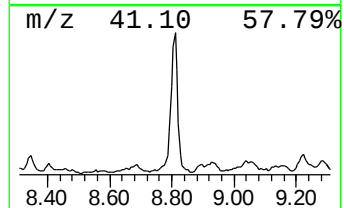
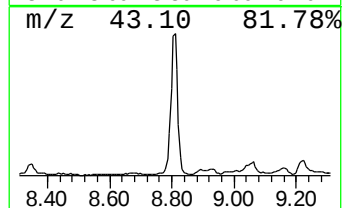
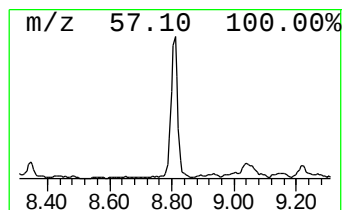
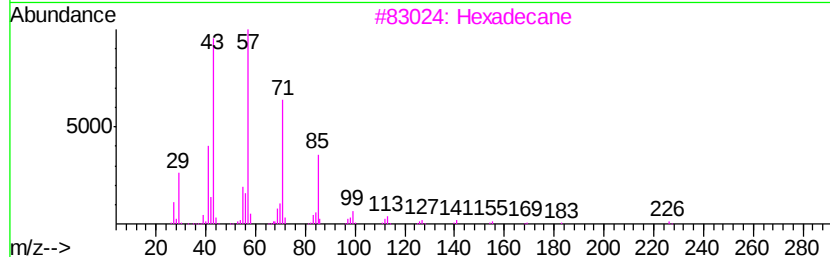
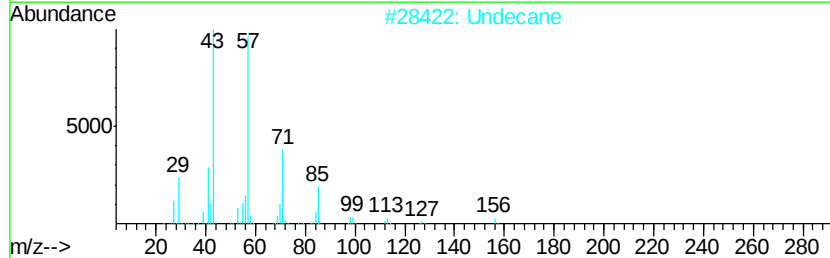
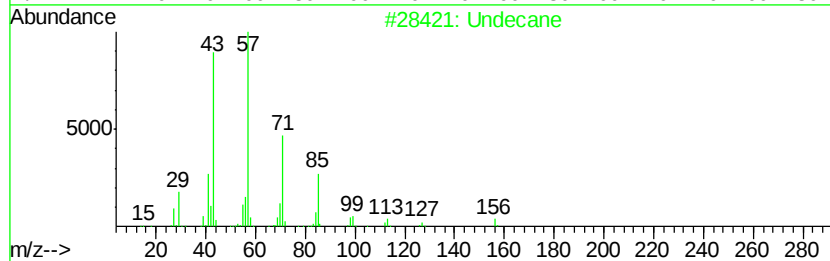
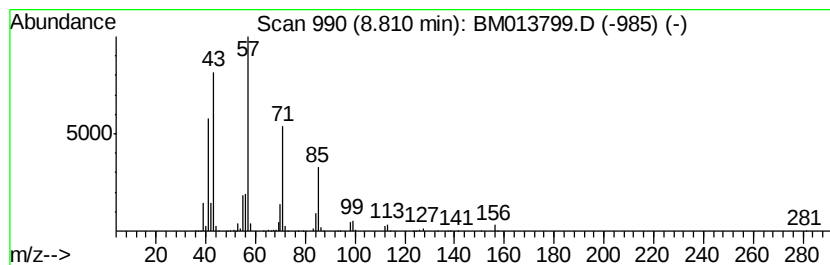
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.81 | 20.00 ng/ul | 3249440 | 1,4-Dichlorobenzene-d4 | 7.59 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Undecane     | 156 | C11H24  | 001120-21-4 | 91   |
| 2    |    | Undecane     | 156 | C11H24  | 001120-21-4 | 91   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 83   |
| 5    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 78   |



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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

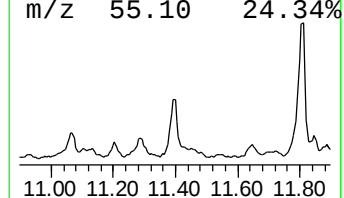
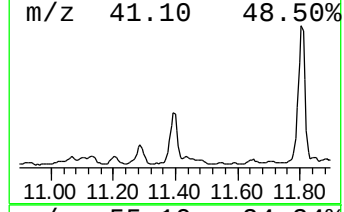
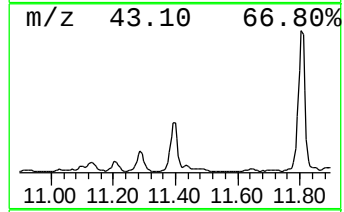
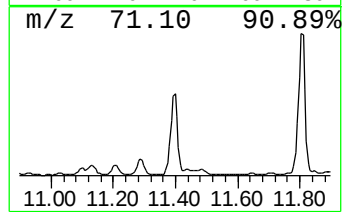
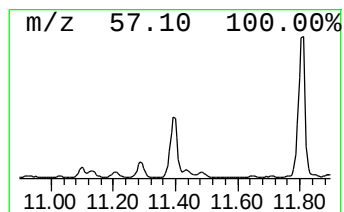
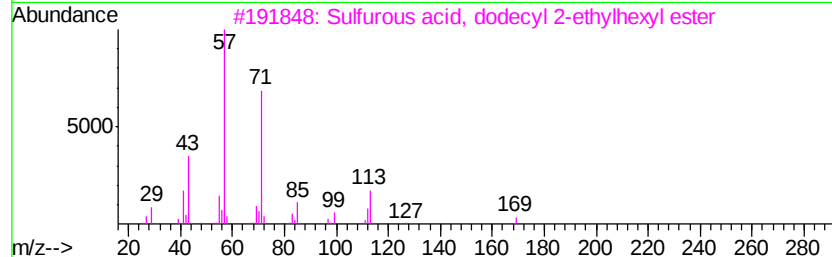
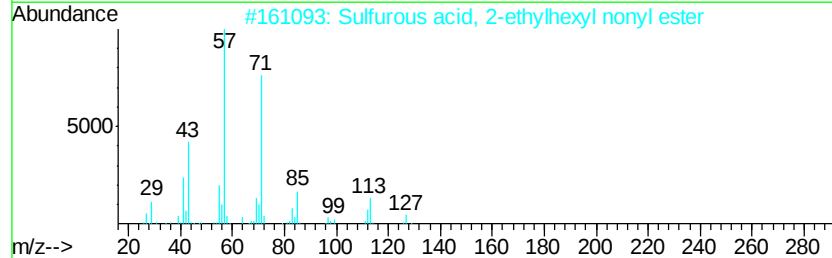
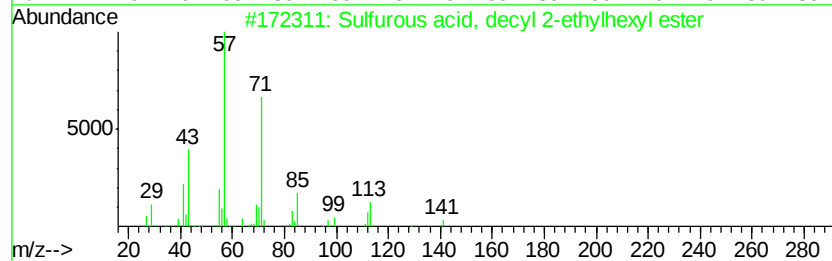
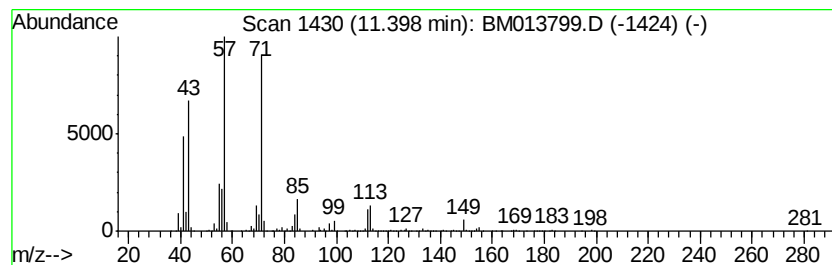
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Sulfurous acid, decyl 2-eth... Concentration Rank 26

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 11.40 | 11.74 ng/ul | 4868150 | Napthalene-d8    | 10.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 80   |
| 2    |    | Sulfurous acid, 2-ethylhexyl non... | 320 | C17H36O3S | 1000309-19-2 | 72   |
| 3    |    | Sulfurous acid, dodecyl 2-ethylh... | 362 | C20H42O3S | 1000309-19-5 | 72   |
| 4    |    | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 72   |
| 5    |    | Nonane, 4-methyl-5-propyl-          | 184 | C13H28    | 062185-55-1  | 64   |



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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

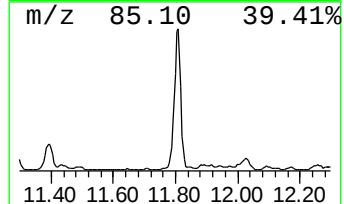
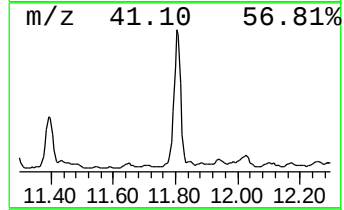
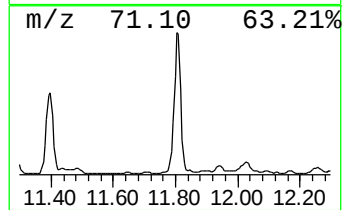
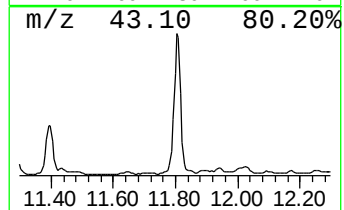
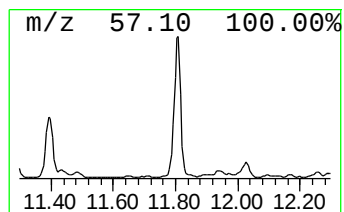
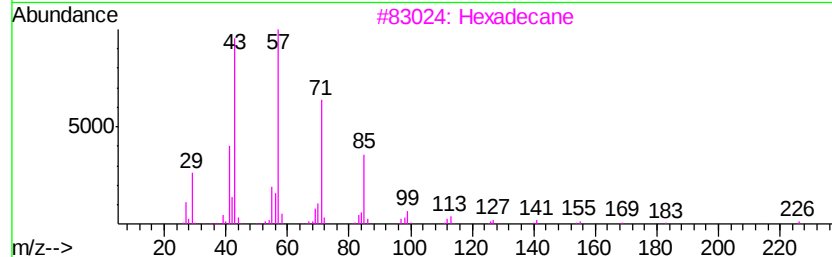
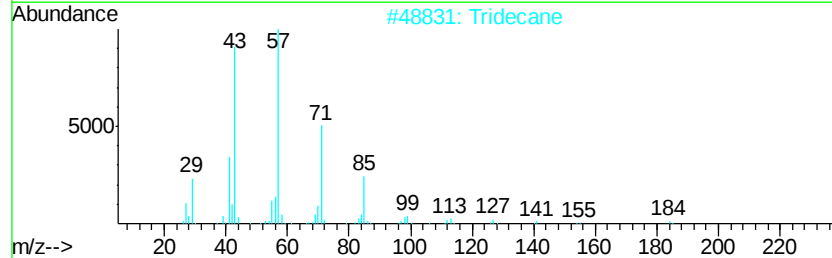
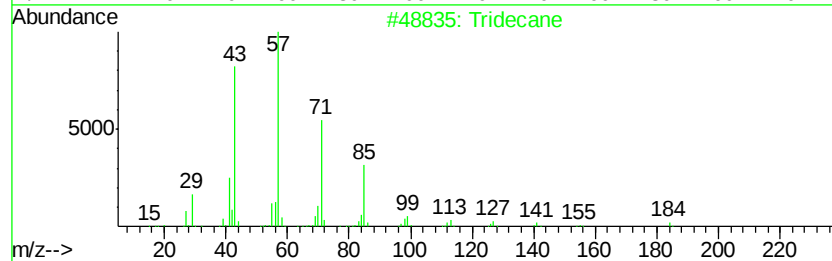
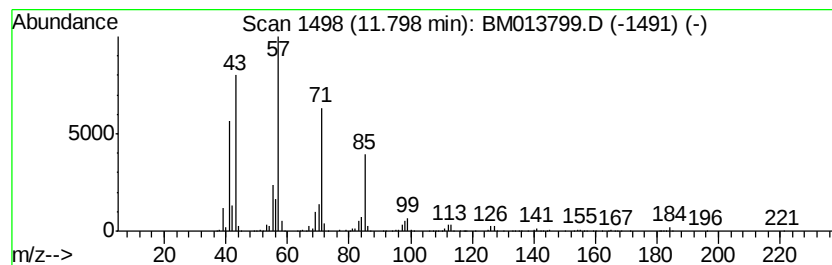
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 11.80 | 25.93 ng/ul | 10758400 | Napthalene-d8    | 10.37 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane    | 184 | C13H28  | 000629-50-5 | 94   |
| 2    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 93   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 91   |
| 4    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 91   |
| 5    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 90   |



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Quant Title : SVOA CALIBRATION

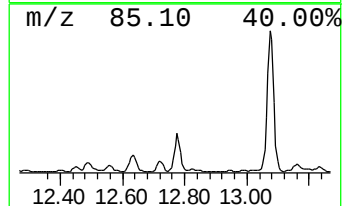
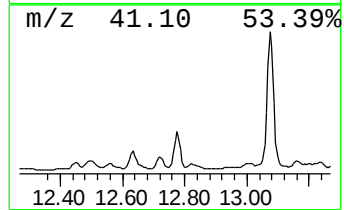
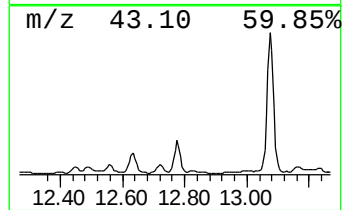
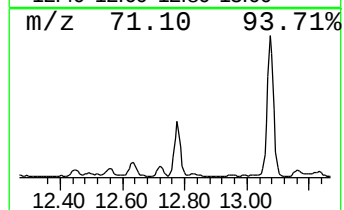
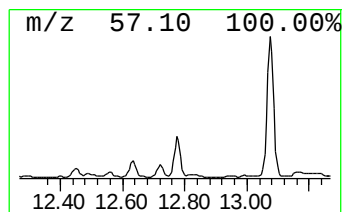
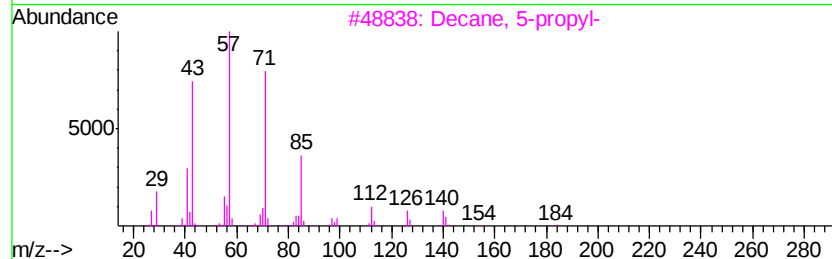
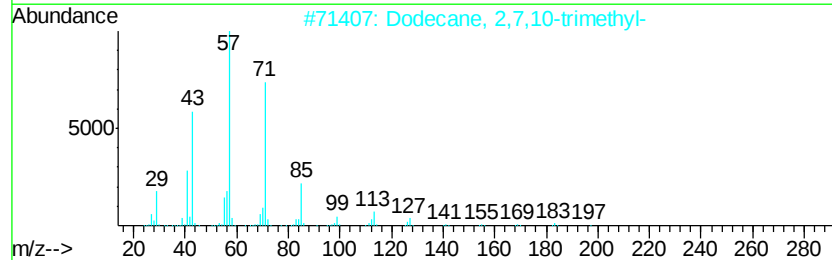
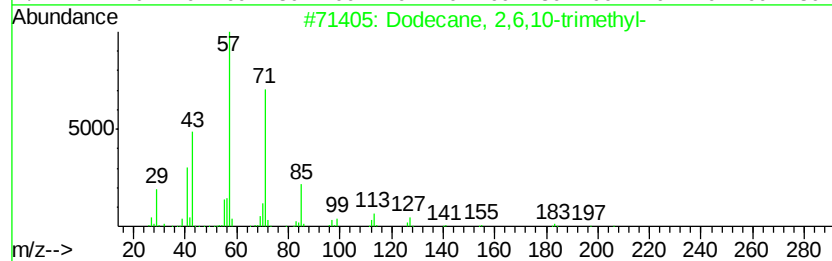
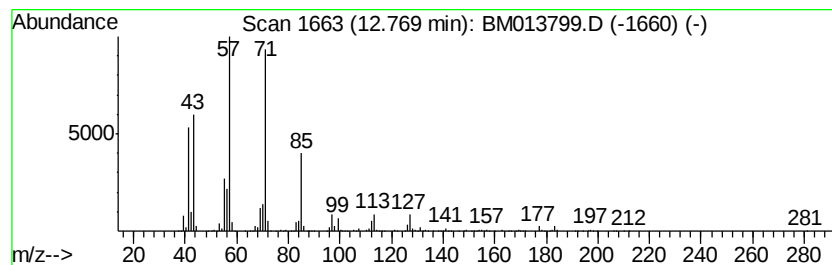
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 12.77 | 7.70 ng/ul | 4650810 | Acenaphthene-d10 | 14.25 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 80   |
| 2    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 80   |
| 3    |    | Decane, 5-propyl-           | 184 | C13H28  | 017312-62-8 | 74   |
| 4    |    | Nonane, 3,7-dimethyl-       | 156 | C11H24  | 017302-32-8 | 68   |
| 5    |    | Octane, 6-ethyl-2-methyl-   | 156 | C11H24  | 062016-19-7 | 64   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

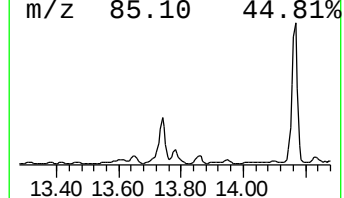
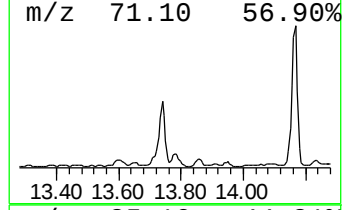
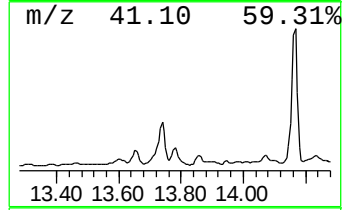
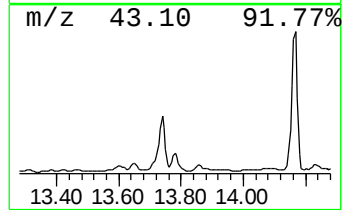
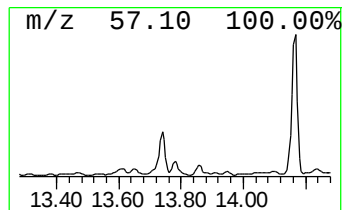
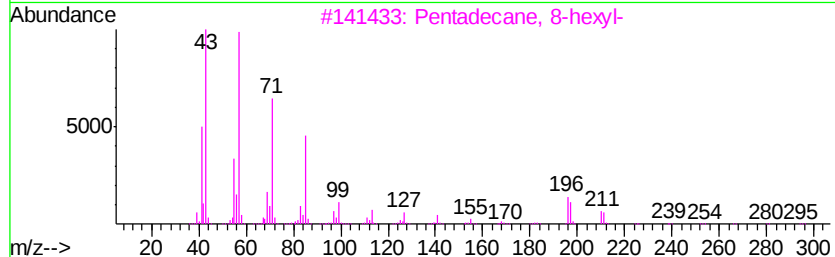
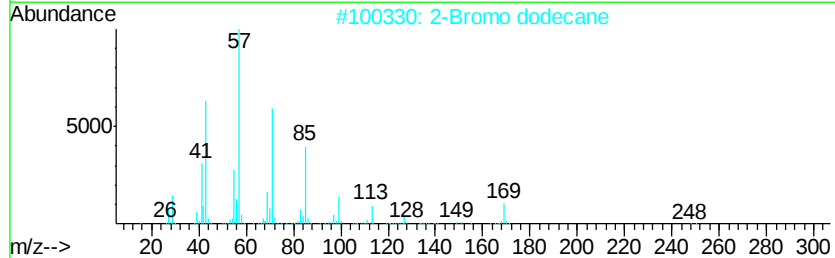
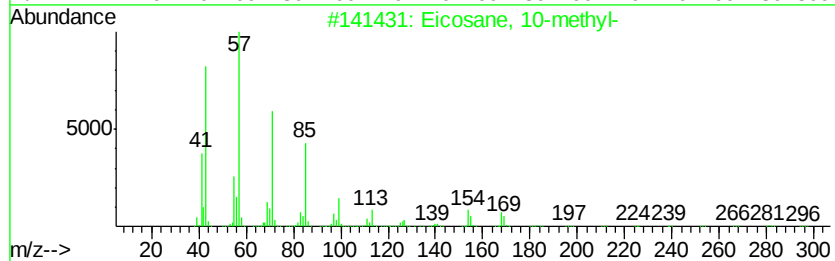
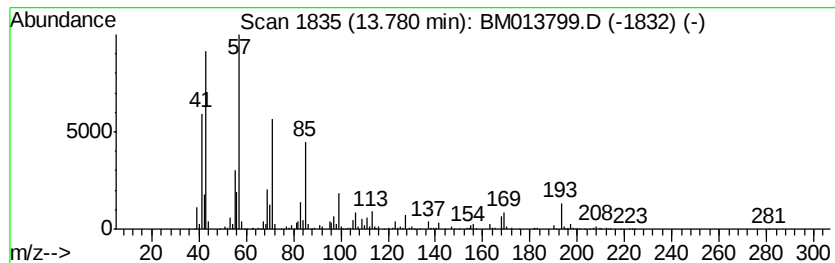
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.      | EstConc               | Area    | Relative to ISTD | R.T.        |      |
|-----------|-----------------------|---------|------------------|-------------|------|
| 13.78     | 6.62 ng/ul            | 3993750 | Acenaphthene-d10 | 14.25       |      |
| Hit# of 5 | Tentative ID          | MW      | MolForm          | CAS#        | Qual |
| 1         | Eicosane, 10-methyl-  | 296     | C21H44           | 054833-23-7 | 80   |
| 2         | 2-Bromo dodecane      | 248     | C12H25Br         | 013187-99-0 | 80   |
| 3         | Pentadecane, 8-hexyl- | 296     | C21H44           | 013475-75-7 | 72   |
| 4         | Heneicosane           | 296     | C21H44           | 000629-94-7 | 72   |
| 5         | Tetracosane           | 338     | C24H50           | 000646-31-1 | 72   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

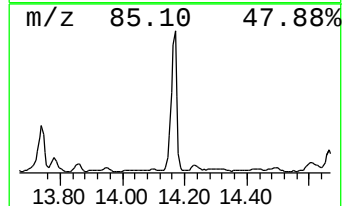
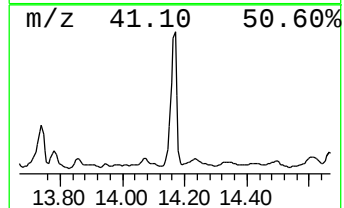
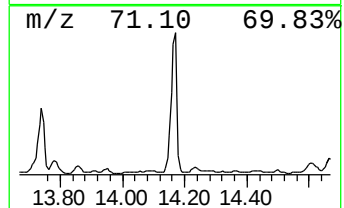
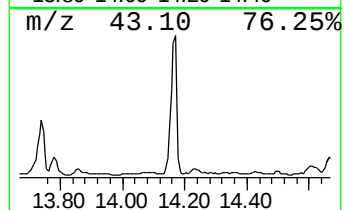
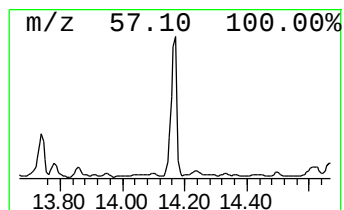
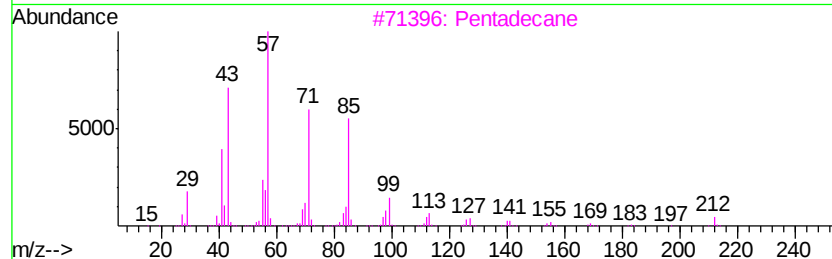
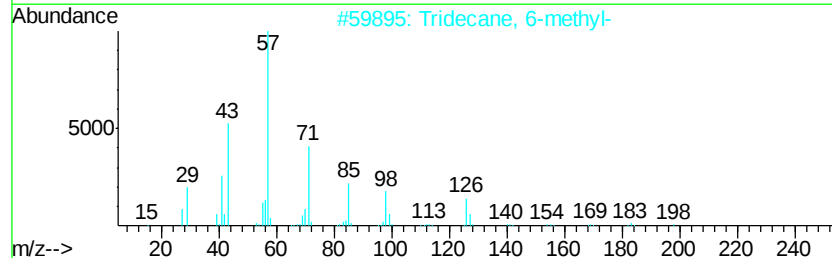
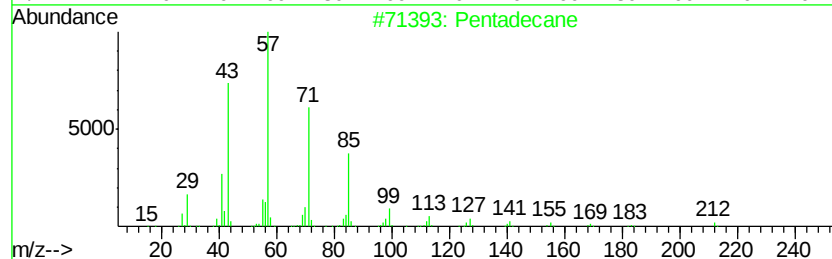
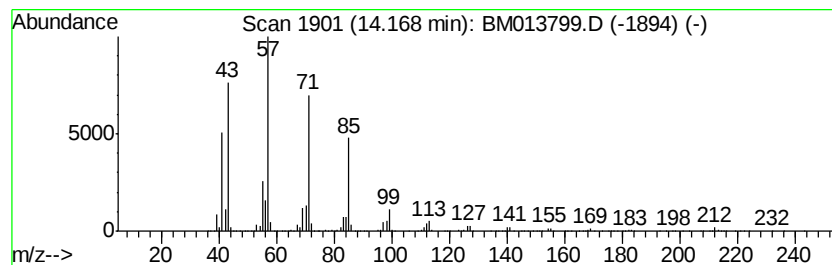
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 15

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 14.17 | 38.39 ng/ul | 23173200 | Acenaphthene-d10 | 14.25 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane          | 212 | C15H32  | 000629-62-9 | 93   |
| 2    |    | Tridecane, 6-methyl- | 198 | C14H30  | 013287-21-3 | 91   |
| 3    |    | Pentadecane          | 212 | C15H32  | 000629-62-9 | 91   |
| 4    |    | Nonadecane           | 268 | C19H40  | 000629-92-5 | 91   |
| 5    |    | Tetradecane          | 198 | C14H30  | 000629-59-4 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

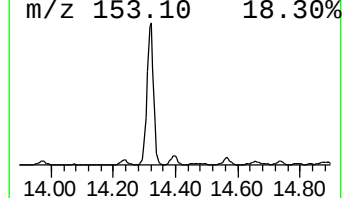
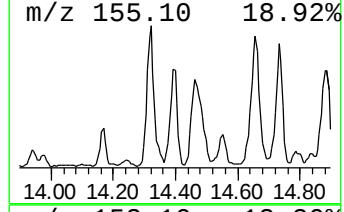
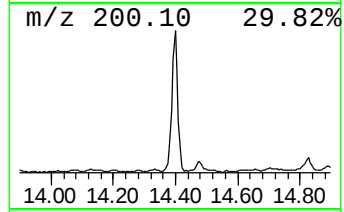
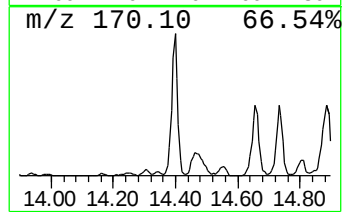
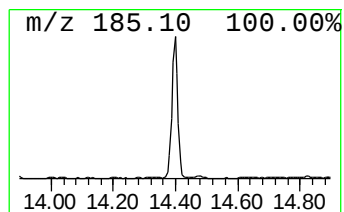
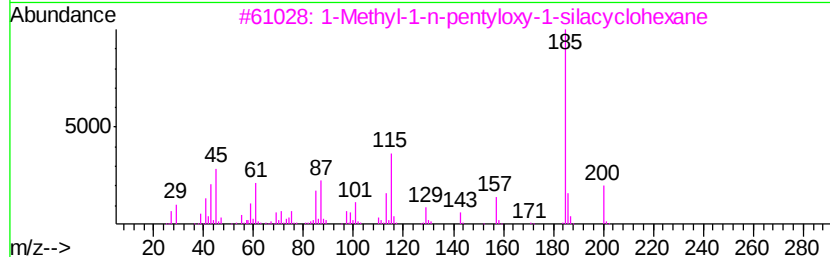
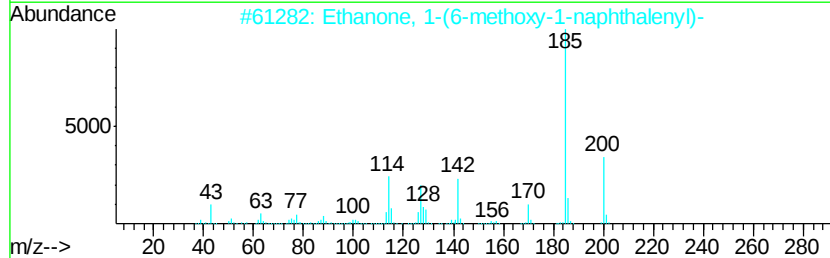
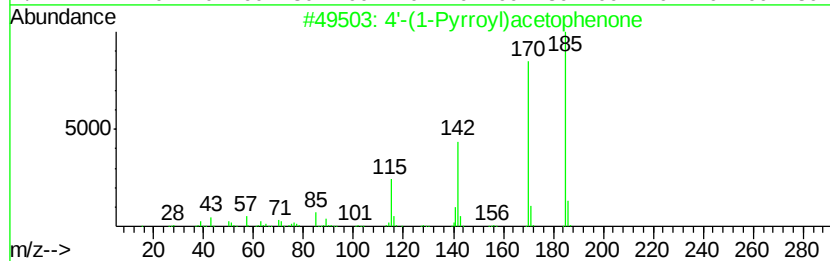
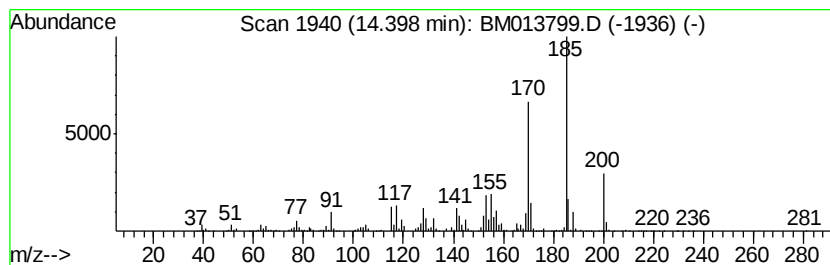
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ClientSampleId :  
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Quant Title : SVOA CALIBRATION

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

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Peak Number 7 4'-(1-Pyrrolyl)acetophenone Concentration Rank 35

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 14.40     | 6.81 ng/ul                          | 4113550 | Acenaphthene-d10 | 14.25        |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 4'-(1-Pyrrovl)acetophenone          | 185     | C12H11NO         | 022106-37-2  | 50   |
| 2         | Ethanone, 1-(6-methoxy-1-naphtha... | 200     | C13H12O2         | 058149-89-6  | 49   |
| 3         | 1-Methyl-1-n-pentvloxy-1-silacvc... | 200     | C11H24OSi        | 232270-61-0  | 46   |
| 4         | 1H-Pyrazole, 5-(t-butyl)-3-phenyl-  | 200     | C13H16N2         | 1000261-34-8 | 46   |
| 5         | 9-Methyl-S-octahydroanthracene      | 200     | C15H20           | 004948-51-0  | 38   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

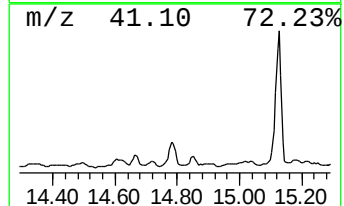
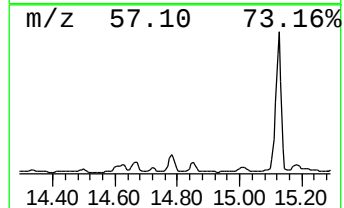
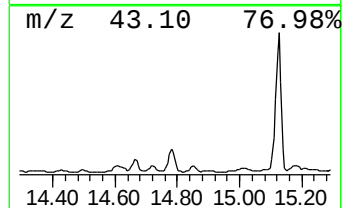
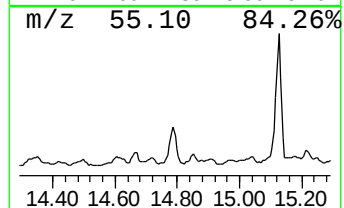
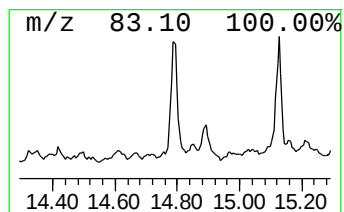
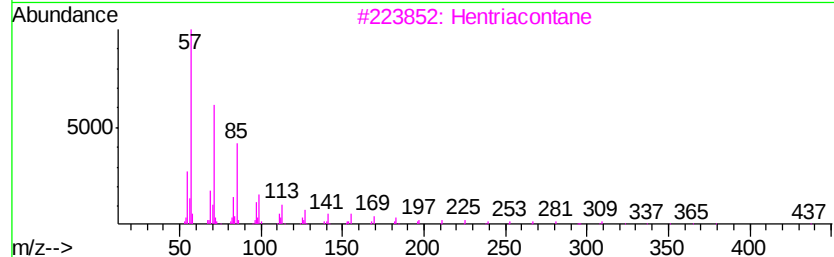
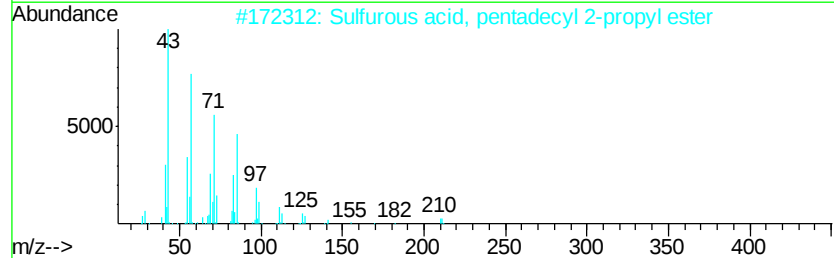
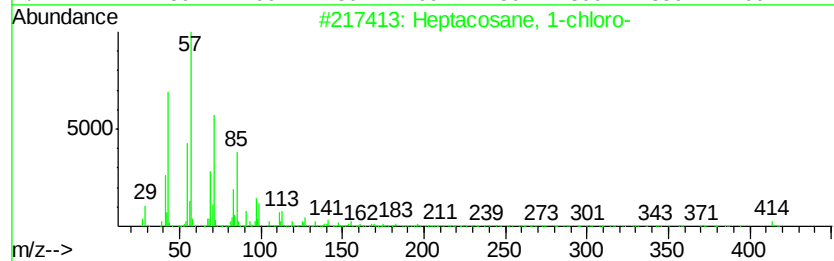
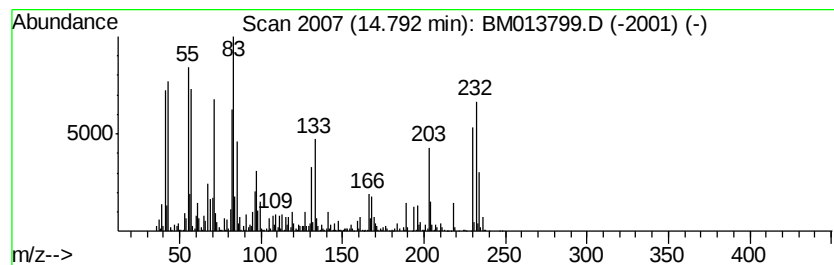
Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Heptacosane, 1-chloro- Concentration Rank 25

| R.T.    | EstConc                             | Area          | Relative to ISTD | R.T.      |
|---------|-------------------------------------|---------------|------------------|-----------|
| 14.79   | 13.29 ng/ul                         | 8021700       | Acenaphthene-d10 | 14.25     |
| Hit# of | 5                                   | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | Heptacosane, 1-chloro-              | 414 C27H55Cl  | 062016-79-9      | 52        |
| 2       | Sulfurous acid, pentadecyl 2-pro... | 334 C18H38O3S | 1000309-12-6     | 52        |
| 3       | Hentriacontane                      | 437 C31H64    | 000630-04-6      | 49        |
| 4       | Hexacosane                          | 366 C26H54    | 000630-01-3      | 49        |
| 5       | 2-Bromotetradecane                  | 276 C14H29Br  | 074036-95-6      | 49        |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

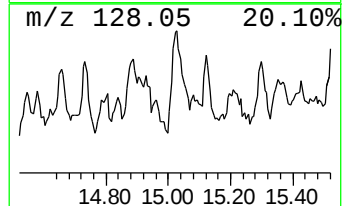
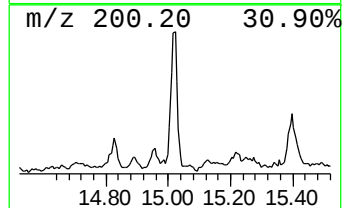
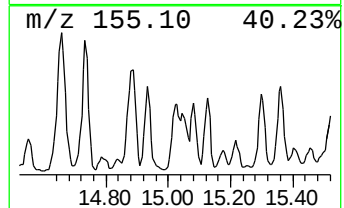
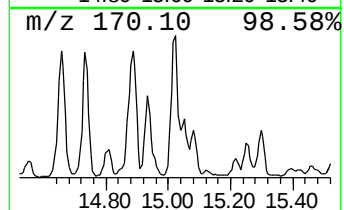
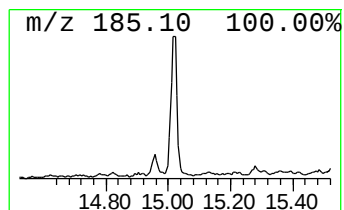
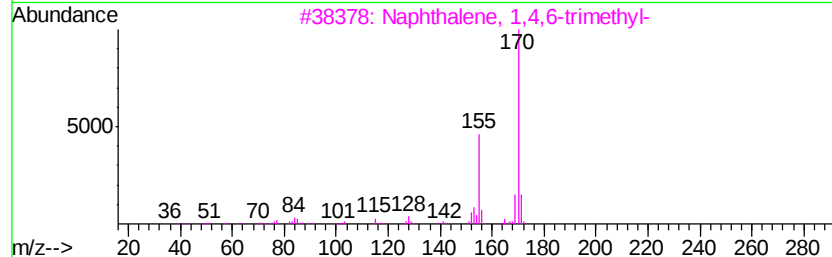
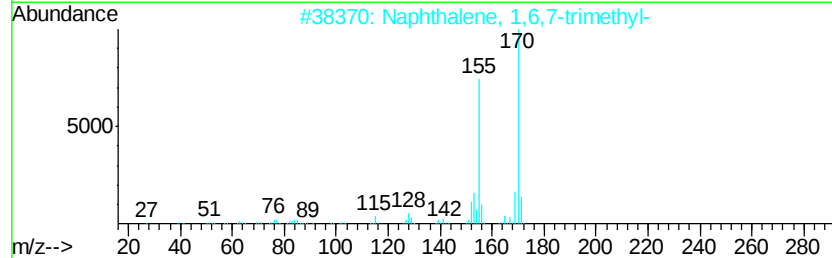
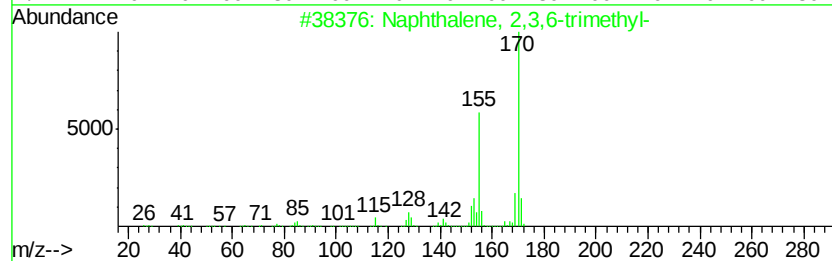
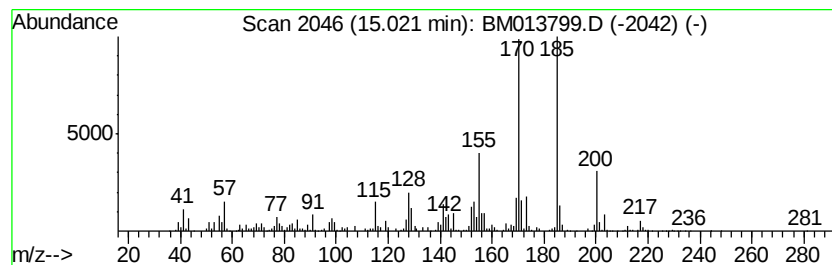
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 37

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.02 | 5.93 ng/ul | 3577440 | Acenaphthene-d10 | 14.25 |

| Hit# | of | Tentative ID                  | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------|-----|----------|-------------|------|
| 1    | 5  | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14   | 000829-26-5 | 90   |
| 2    |    | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14   | 002245-38-7 | 70   |
| 3    |    | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14   | 002131-42-2 | 70   |
| 4    |    | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14   | 000829-26-5 | 70   |
| 5    |    | 4'-(1-Pyrrolyl)acetophenone   | 185 | C12H11NO | 022106-37-2 | 53   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

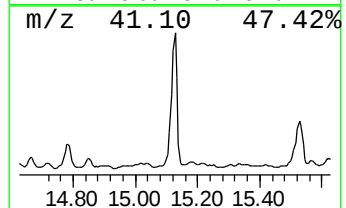
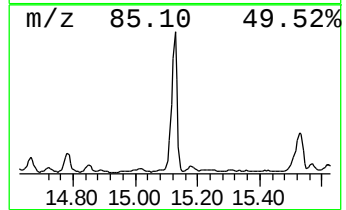
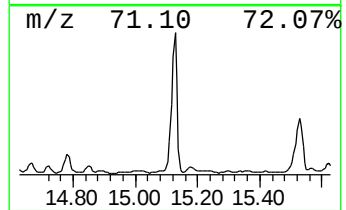
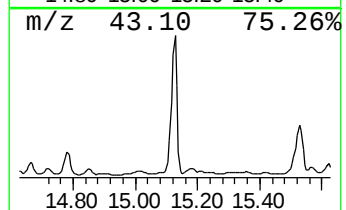
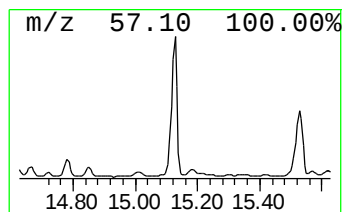
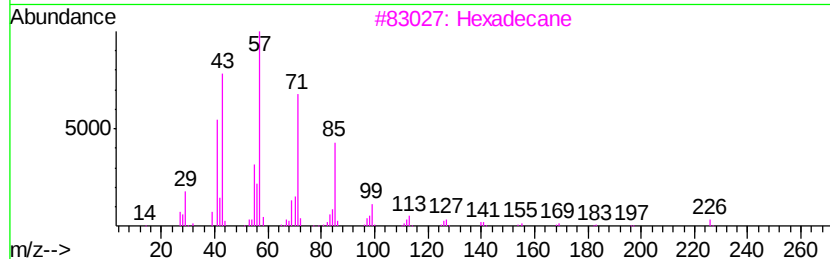
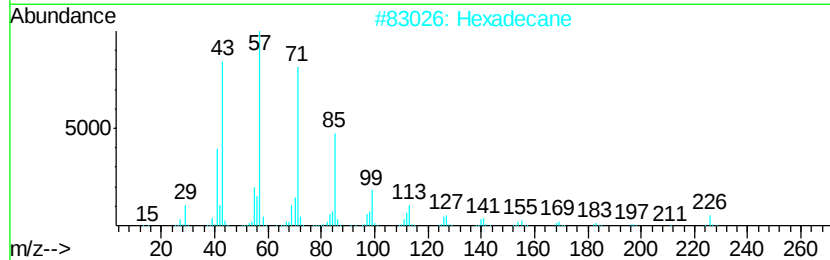
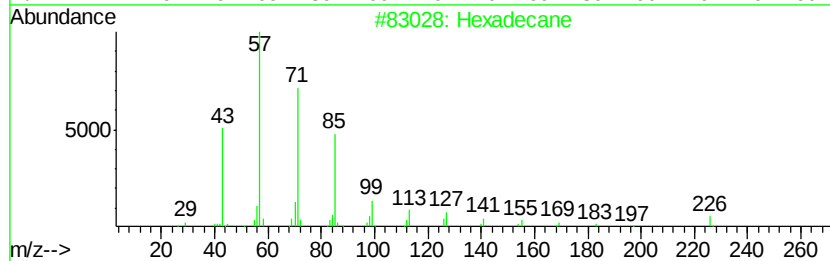
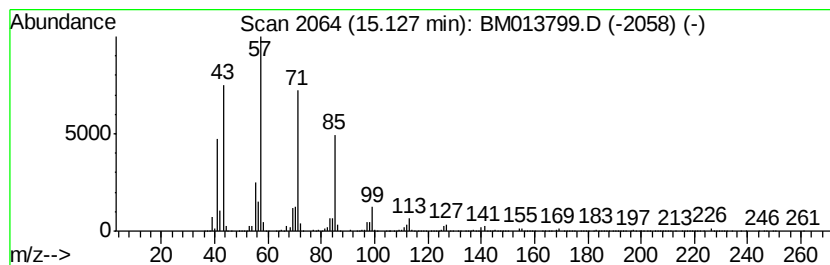
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 15.13 | 49.26 ng/ul | 29733400 | Acenaphthene-d10 | 14.25 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane   | 226 | C16H34  | 000544-76-3 | 97   |
| 2    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 95   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 5    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

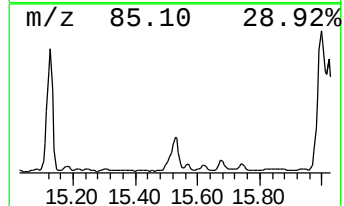
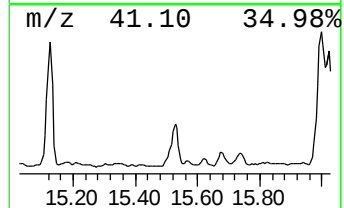
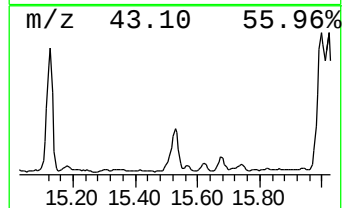
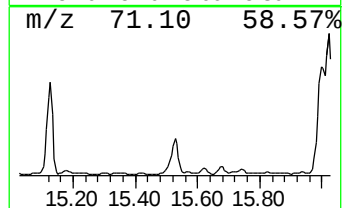
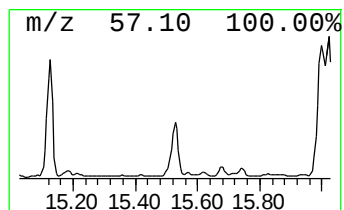
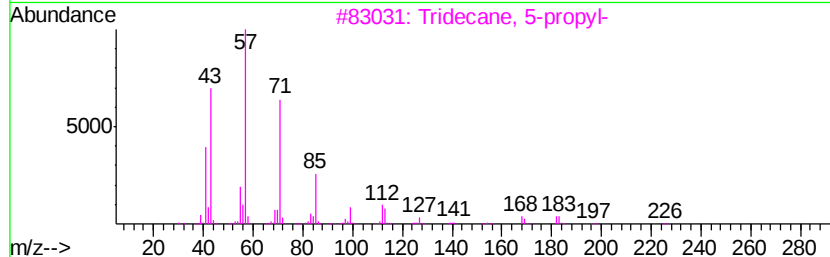
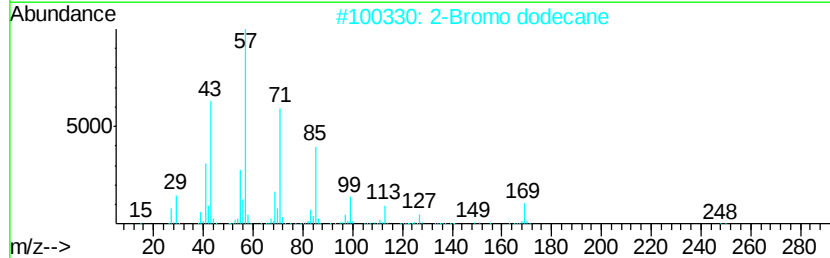
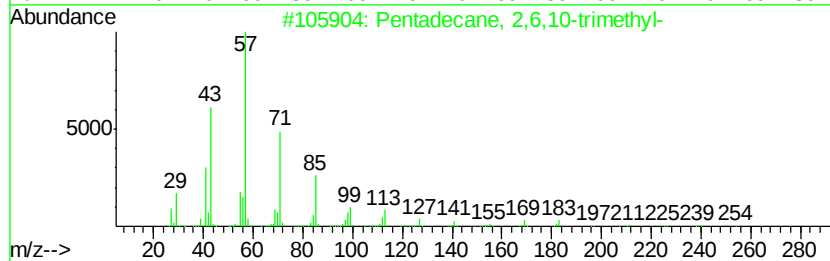
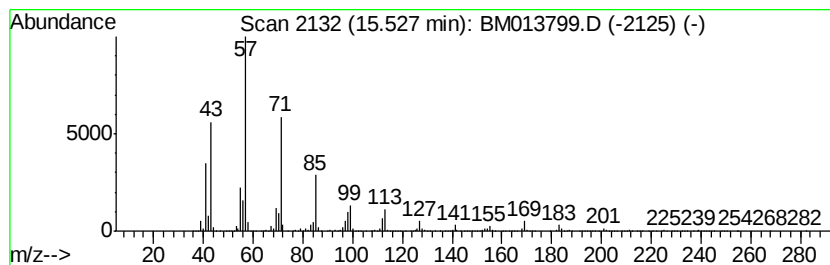
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 18

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 15.53 | 26.57 ng/ul | 16035900 | Acenaphthene-d10 | 14.25 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38   | 003892-00-0 | 95   |
| 2    |    |   | 2-Bromo dodecane               | 248 | C12H25Br | 013187-99-0 | 94   |
| 3    |    |   | Tridecane, 5-propyl-           | 226 | C16H34   | 055045-11-9 | 87   |
| 4    |    |   | Heneicosane                    | 296 | C21H44   | 000629-94-7 | 83   |
| 5    |    |   | Tetracosane                    | 338 | C24H50   | 000646-31-1 | 80   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

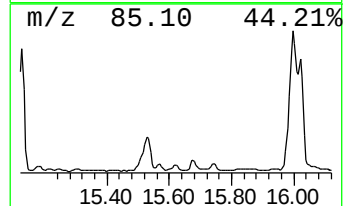
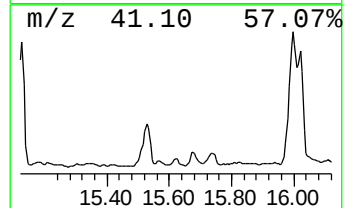
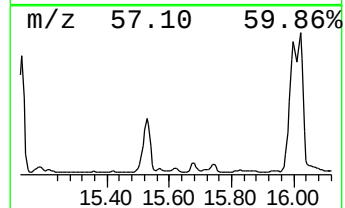
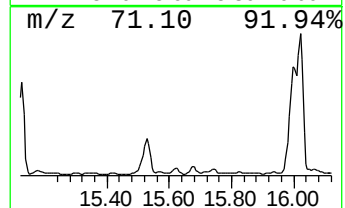
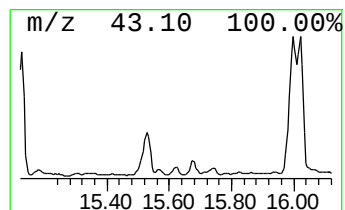
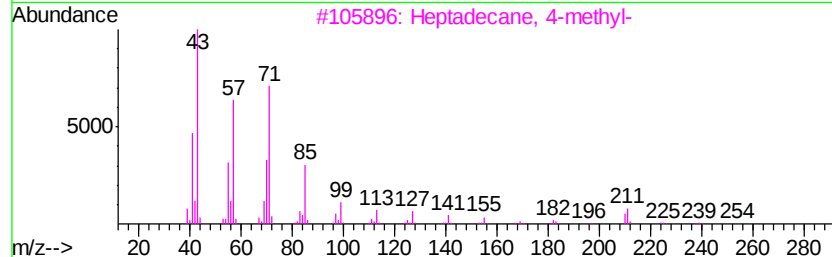
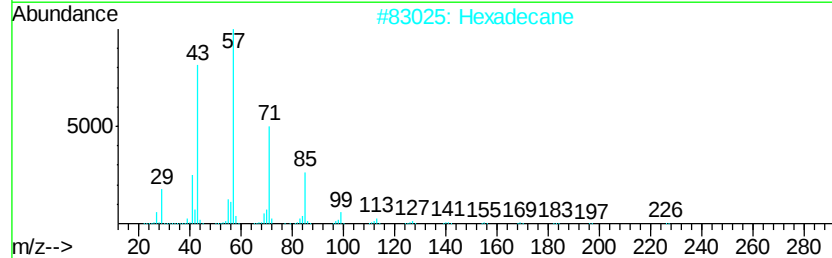
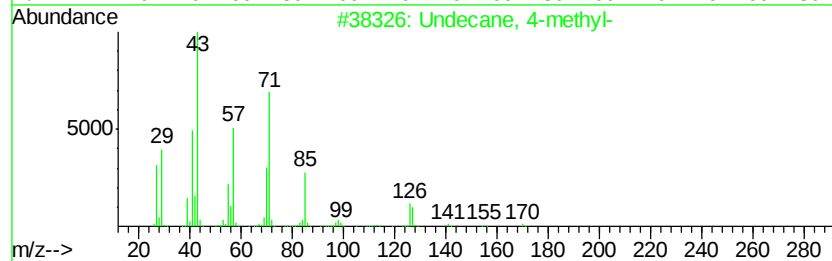
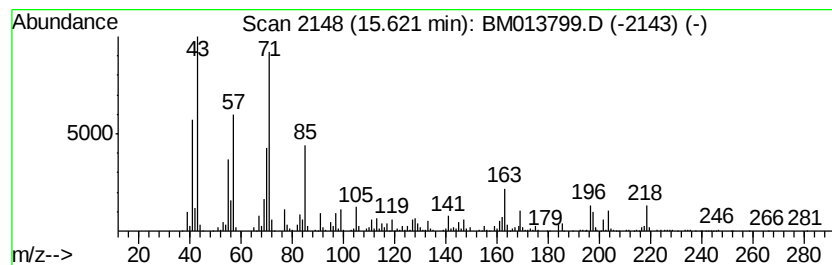
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 34

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 15.62 | 6.98 ng/ul | 4212920 | Acenaphthene-d10 | 14.25 |

| Hit# | of                         | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Undecane, 4-methyl-        |   |              | 170 | C12H26  | 002980-69-0 | 58   |
| 2    | Hexadecane                 |   |              | 226 | C16H34  | 000544-76-3 | 55   |
| 3    | Heptadecane, 4-methyl-     |   |              | 254 | C18H38  | 026429-11-8 | 53   |
| 4    | Tetradecane, 4-methyl-     |   |              | 212 | C15H32  | 025117-24-2 | 52   |
| 5    | 1-Hexene, 3,4,5-trimethyl- |   |              | 126 | C9H18   | 056728-10-0 | 50   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

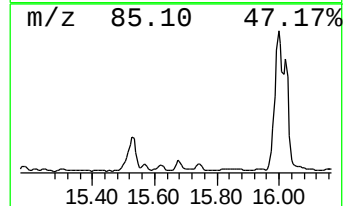
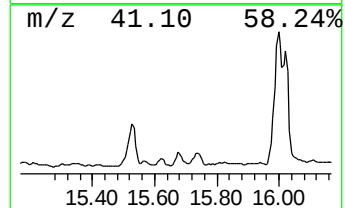
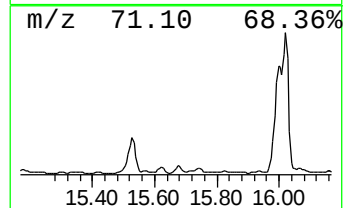
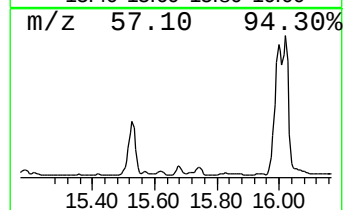
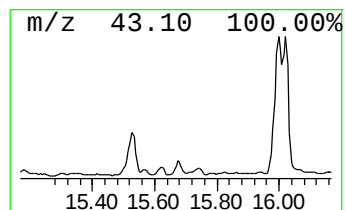
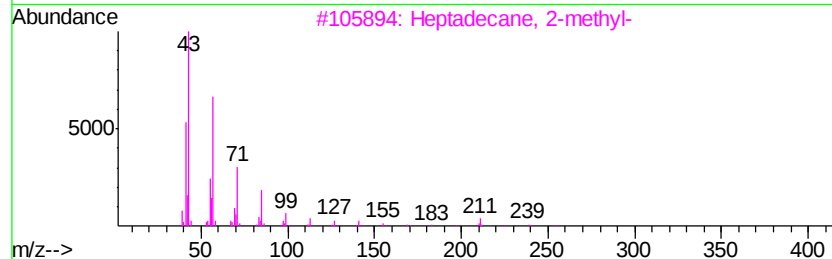
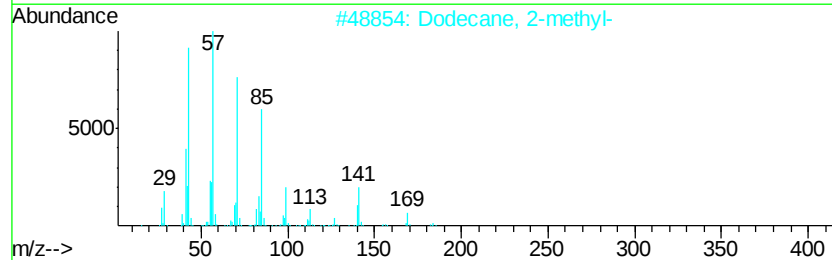
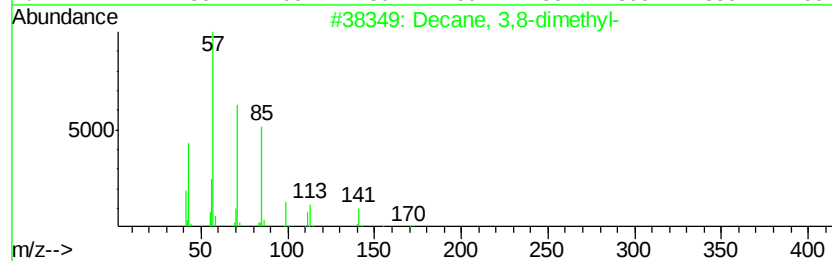
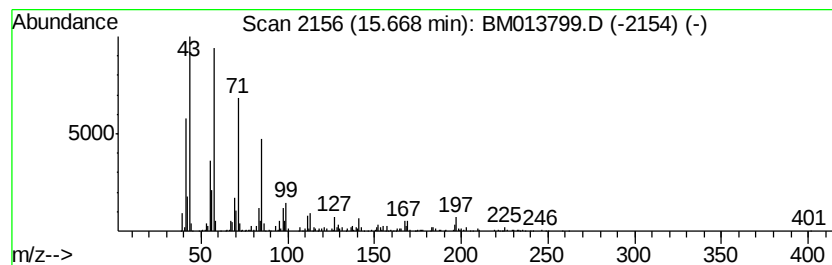
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 15.67 | 11.40 ng/ul | 4022740 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 3,8-dimethyl-  | 170 | C12H26  | 017312-55-9 | 93   |
| 2    |    | Dodecane, 2-methyl-    | 184 | C13H28  | 001560-97-0 | 93   |
| 3    |    | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 91   |
| 4    |    | 10-Methylnonadecane    | 282 | C20H42  | 056862-62-5 | 87   |
| 5    |    | Nonadecane, 2-methyl-  | 282 | C20H42  | 001560-86-7 | 87   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

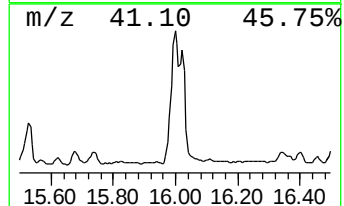
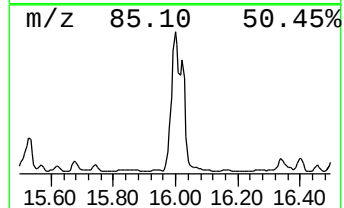
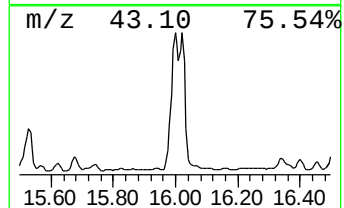
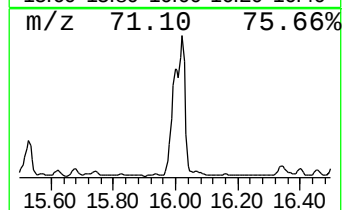
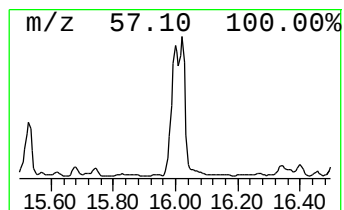
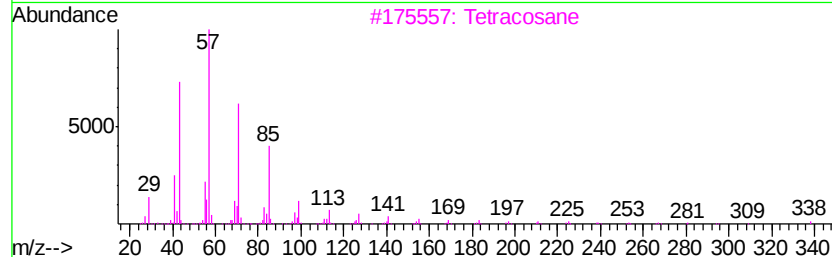
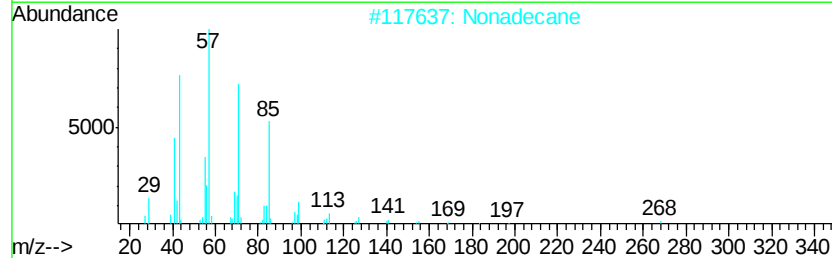
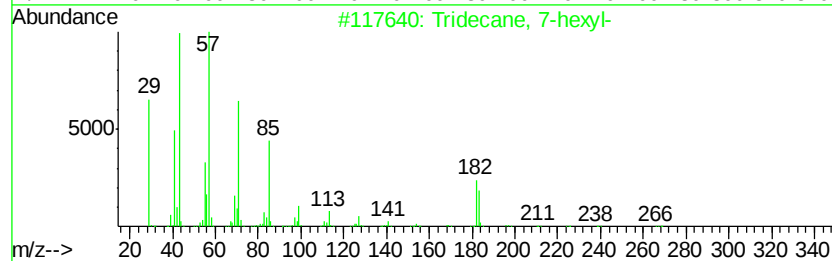
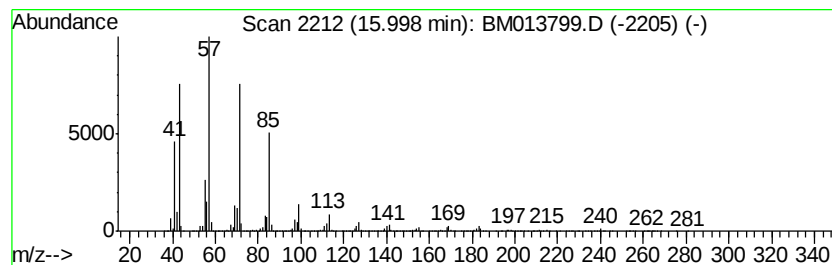
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 16.00 | 134.44 ng/ul | 47428300 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 7-hexyl- | 268 | C19H40  | 007225-66-3 | 95   |
| 2    |    |   | Nonadecane          | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |    |   | Tetracosane         | 338 | C24H50  | 000646-31-1 | 91   |
| 4    |    |   | Heneicosane         | 296 | C21H44  | 000629-94-7 | 91   |
| 5    |    |   | Octacosane          | 394 | C28H58  | 000630-02-4 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

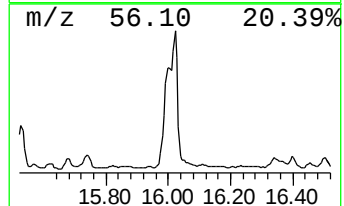
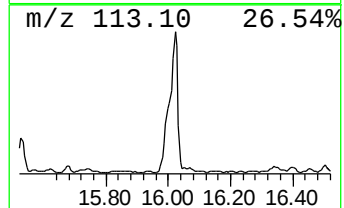
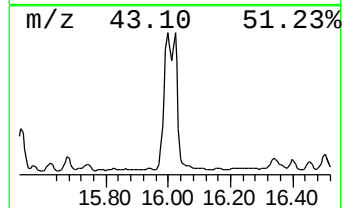
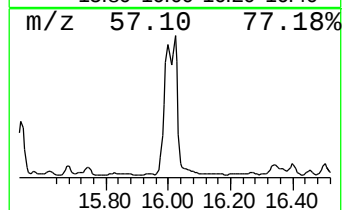
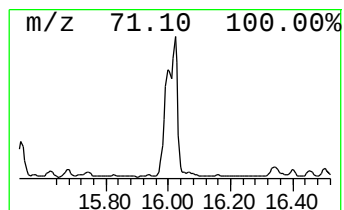
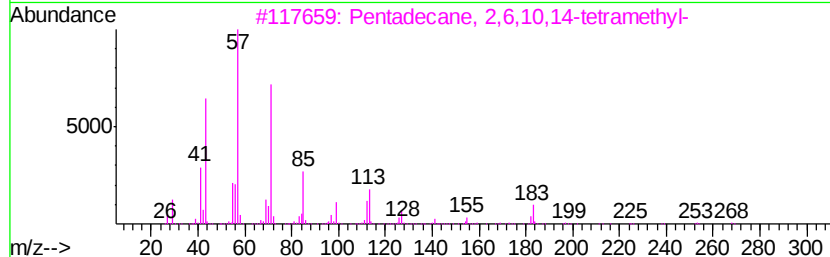
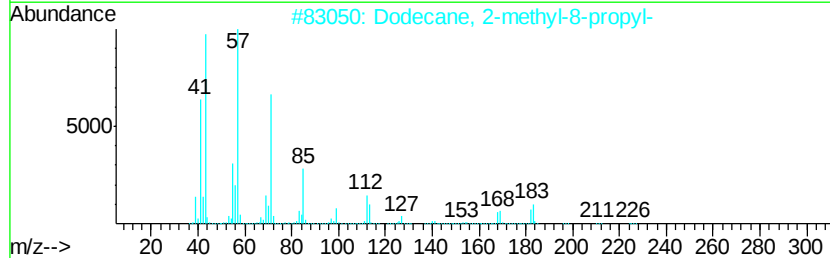
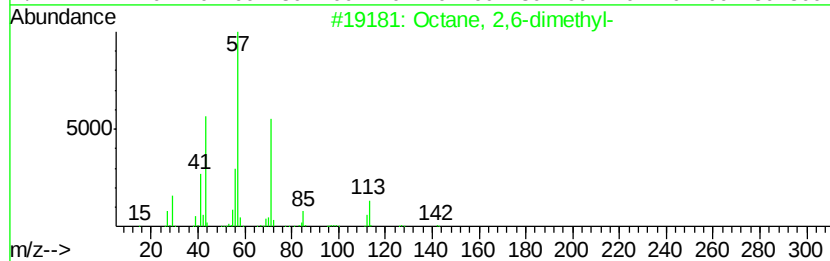
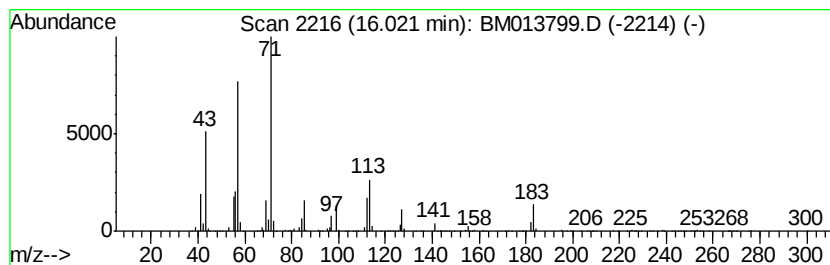
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 16.02 | 94.78 ng/ul | 33435700 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 72   |
| 2    |    | Dodecane, 2-methyl-8-propyl-        | 226 | C16H34    | 055045-07-3  | 68   |
| 3    |    | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40    | 001921-70-6  | 58   |
| 4    |    | Sulfurous acid, 2-ethylhexyl tet... | 390 | C22H46O3S | 1000309-19-7 | 53   |
| 5    |    | Dodecane, 4-methyl-                 | 184 | C13H28    | 006117-97-1  | 50   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

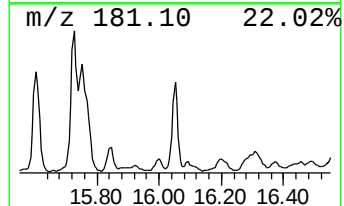
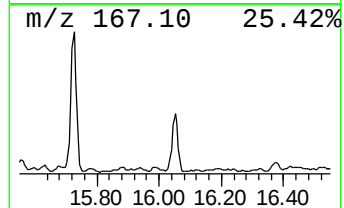
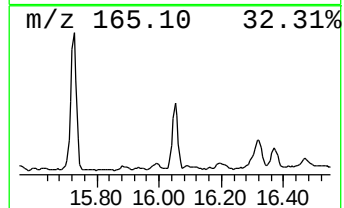
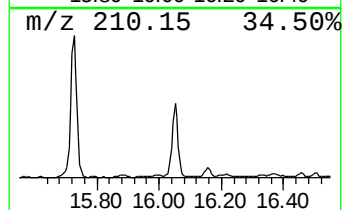
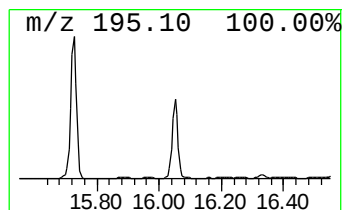
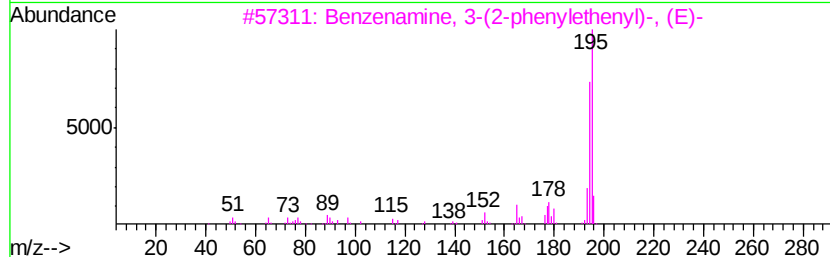
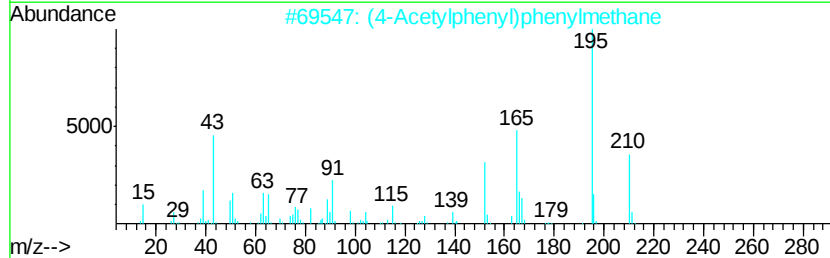
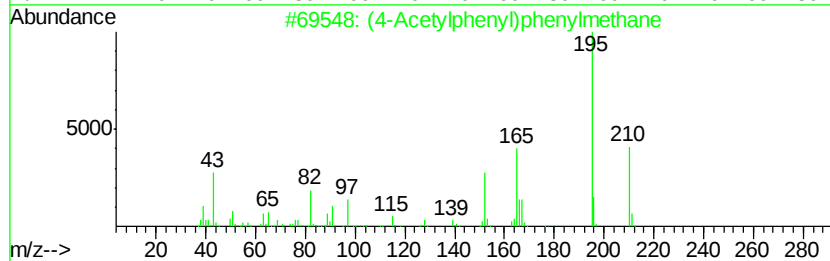
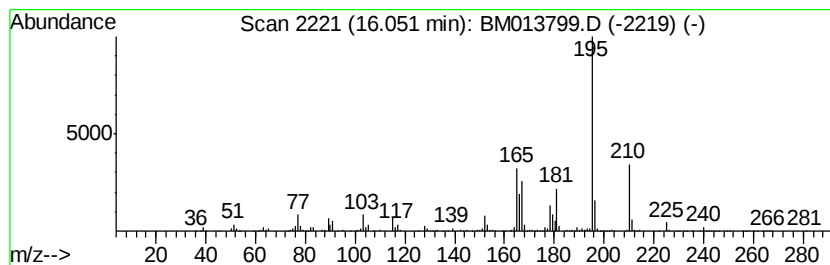
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (4-Acetylphenyl)phenylmethane Concentration Rank 23

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.05 | 17.32 ng/ul | 6110130 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | (4-Acetylphenyl)phenylmethane       | 210 | C15H14O | 000782-92-3 | 62   |
| 2    |    |   | (4-Acetylphenyl)phenylmethane       | 210 | C15H14O | 000782-92-3 | 58   |
| 3    |    |   | Benzenamine, 3-(2-phenylethenyl)... | 195 | C14H13N | 014064-82-5 | 47   |
| 4    |    |   | 2',3',4',5',6'-Pentafluoroacetop... | 210 | C8H3F5O | 000652-29-9 | 47   |
| 5    |    |   | 4-Amino-9-fluorenone                | 195 | C13H9NO | 004269-15-2 | 47   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

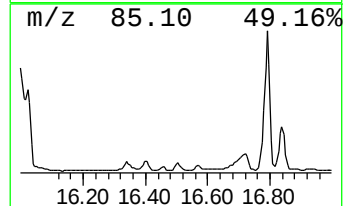
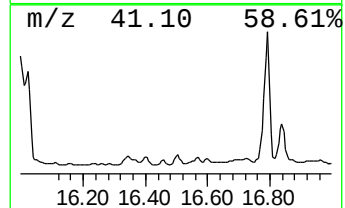
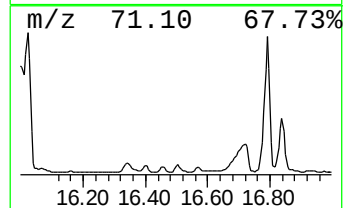
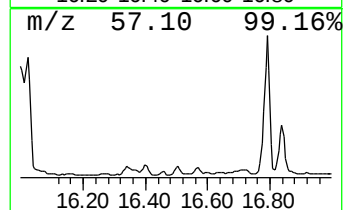
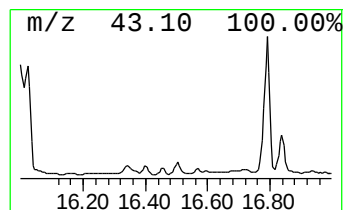
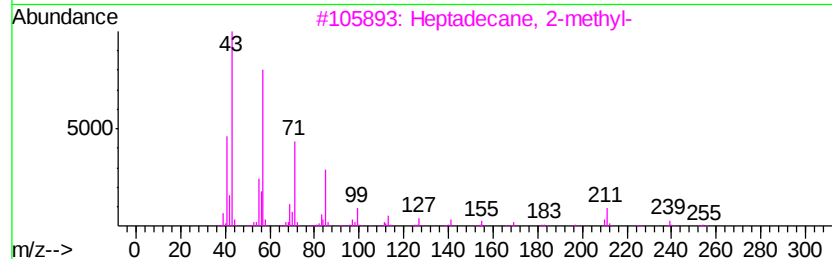
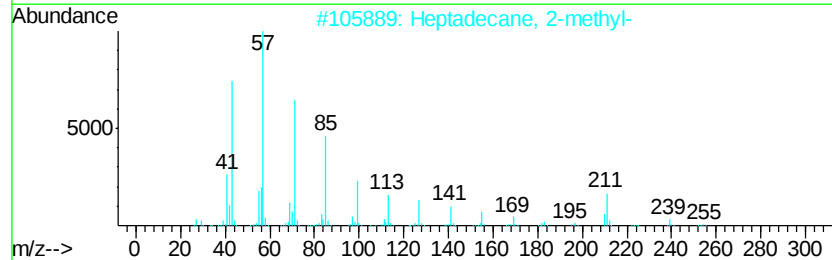
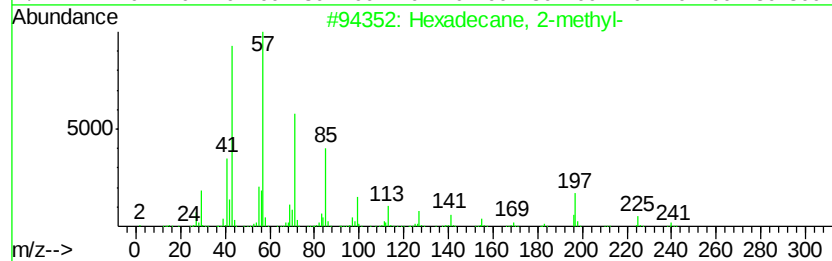
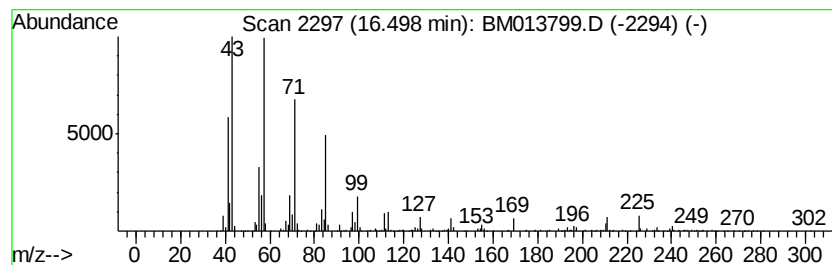
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.50 | 9.32 ng/ul | 3286310 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2-methyl-  | 240 | C17H36  | 001560-92-5 | 97   |
| 2    |    |   | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 94   |
| 3    |    |   | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0 | 90   |
| 4    |    |   | Heptadecane            | 240 | C17H36  | 000629-78-7 | 87   |
| 5    |    |   | Tetradecane, 2-methyl- | 212 | C15H32  | 001560-95-8 | 87   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

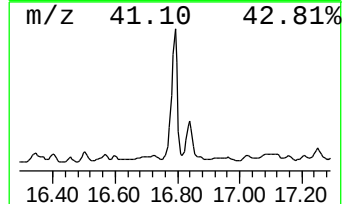
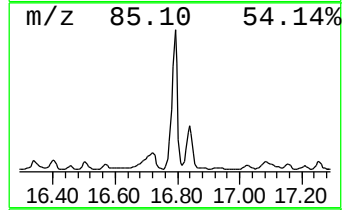
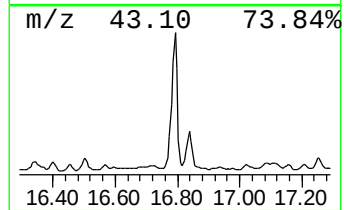
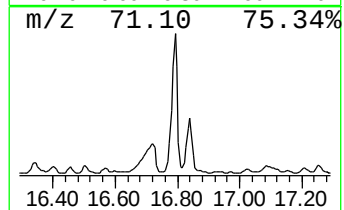
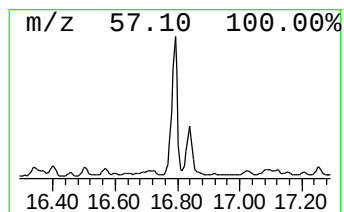
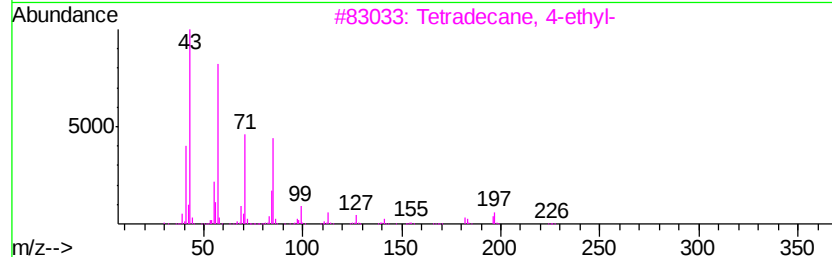
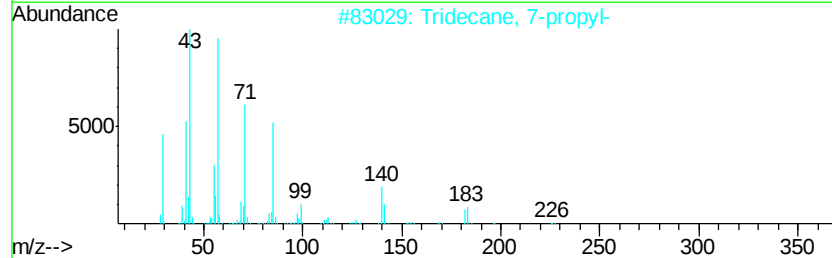
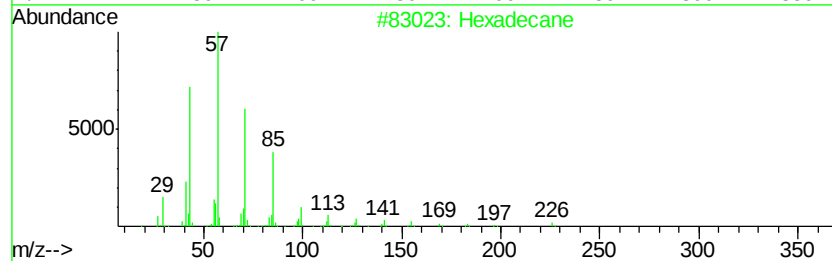
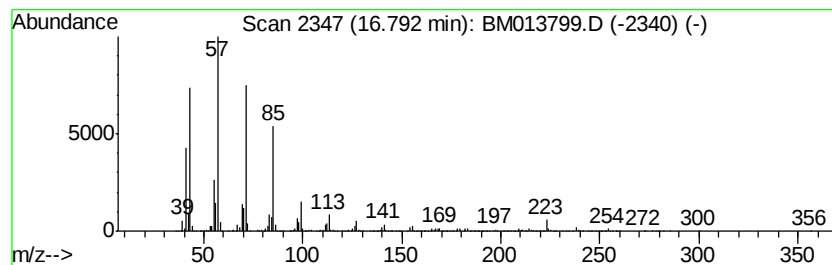
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 16.79 | 149.21 ng/ul | 52639200 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane            | 226 | C16H34  | 000544-76-3 | 94   |
| 2    |    | Tridecane, 7-propyl-  | 226 | C16H34  | 055045-09-5 | 91   |
| 3    |    | Tetradecane, 4-ethyl- | 226 | C16H34  | 055045-14-2 | 90   |
| 4    |    | Tridecane, 6-propyl-  | 226 | C16H34  | 055045-10-8 | 87   |
| 5    |    | Tetradecane           | 198 | C14H30  | 000629-59-4 | 87   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

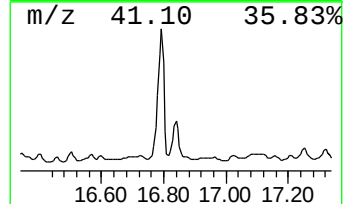
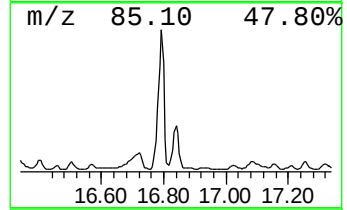
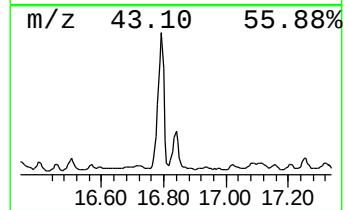
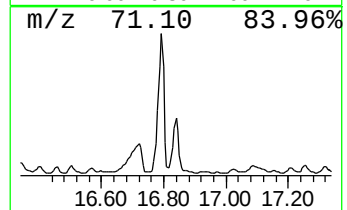
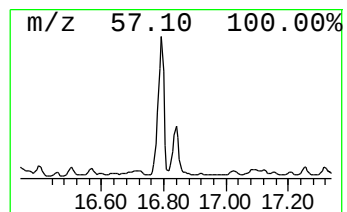
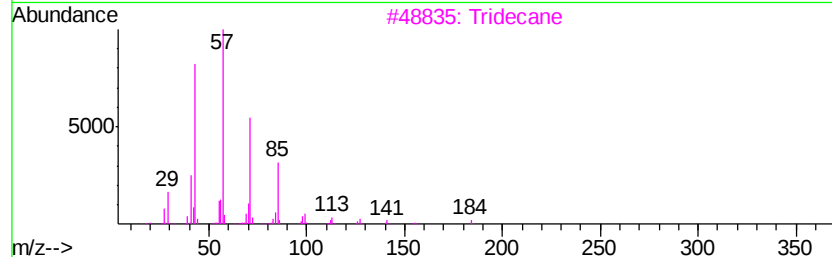
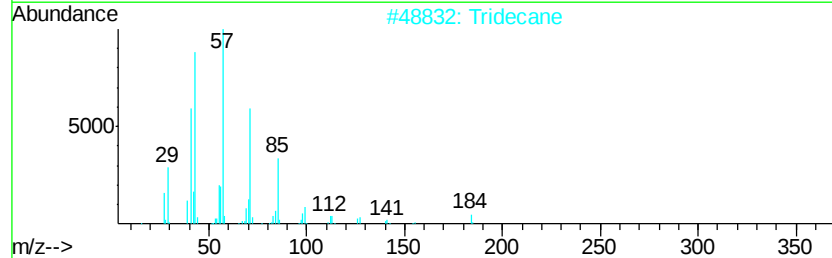
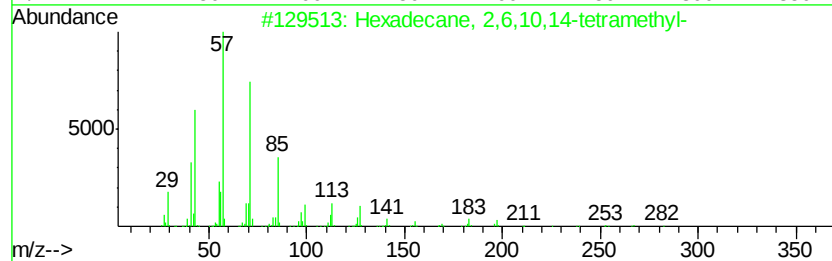
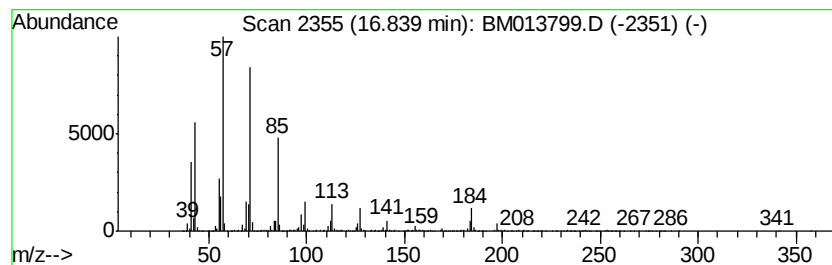
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 16.84 | 58.23 ng/ul | 20543900 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 91   |
| 2    |    | Tridecane                          | 184 | C13H28  | 000629-50-5 | 91   |
| 3    |    | Tridecane                          | 184 | C13H28  | 000629-50-5 | 91   |
| 4    |    | Tridecane                          | 184 | C13H28  | 000629-50-5 | 87   |
| 5    |    | Tetracosane                        | 338 | C24H50  | 000646-31-1 | 86   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

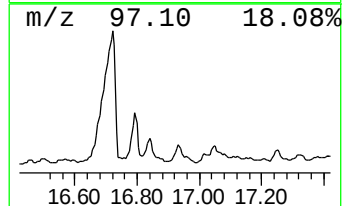
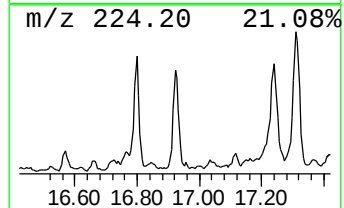
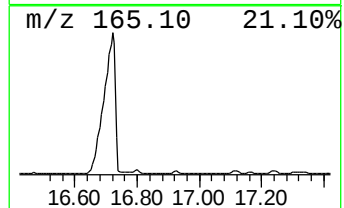
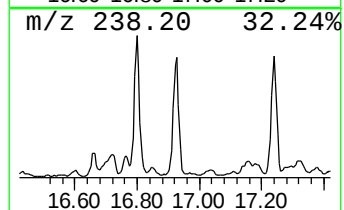
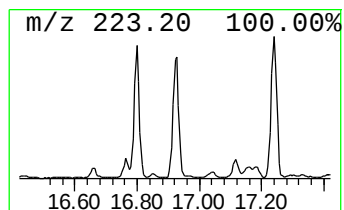
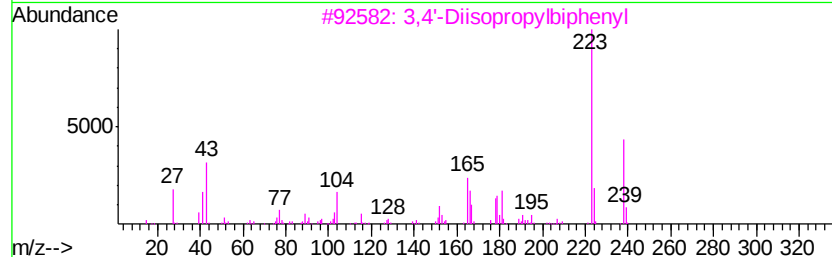
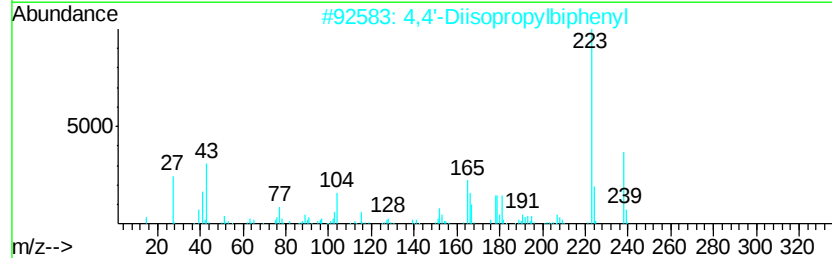
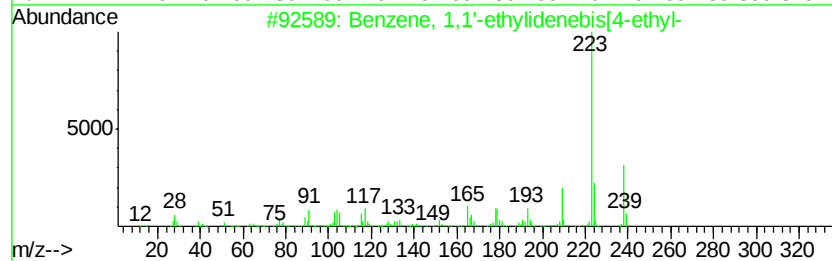
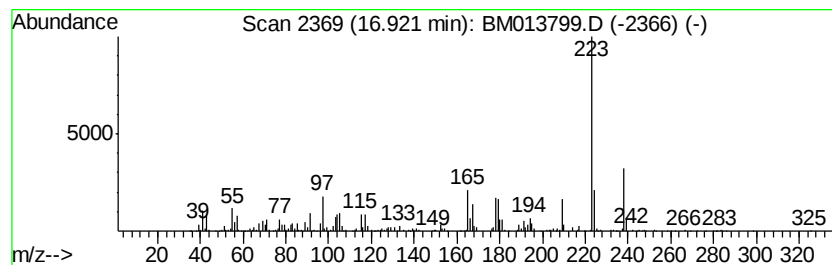
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TIC Integration Parameters: LSCINT.P

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Peak Number 21 Benzene, 1,1'-ethylidenebis... Concentration Rank 31

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.92 | 9.53 ng/ul | 3360940 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID                            | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Benzene, 1,1'-ethylidenebis[4-ethyl-... | 238 | C18H22   | 010224-91-6 | 89   |
| 2    |    | 4,4'-Diisopropylbiphenyl                | 238 | C18H22   | 018970-30-4 | 87   |
| 3    |    | 3,4'-Diisopropylbiphenyl                | 238 | C18H22   | 061434-46-6 | 80   |
| 4    |    | Benzene, 1,1'-ethylidenebis[3,4-...     | 238 | C18H22   | 001742-14-9 | 58   |
| 5    |    | 9-Acridinecarboxylic acid               | 223 | C14H9NO2 | 005336-90-3 | 53   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

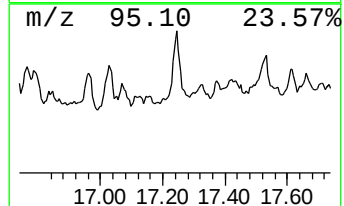
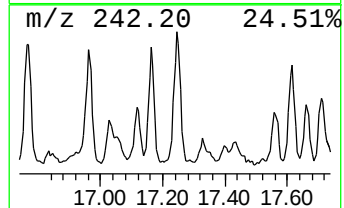
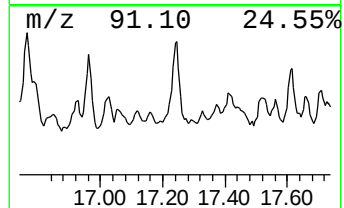
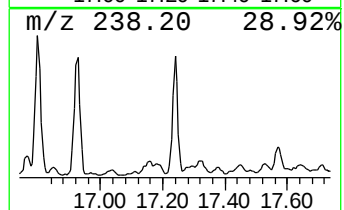
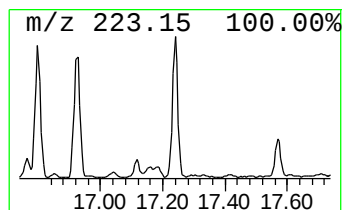
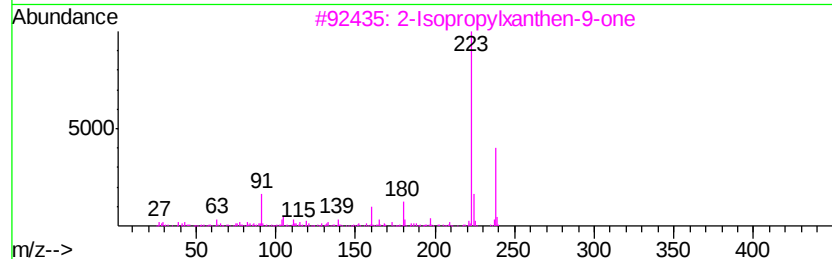
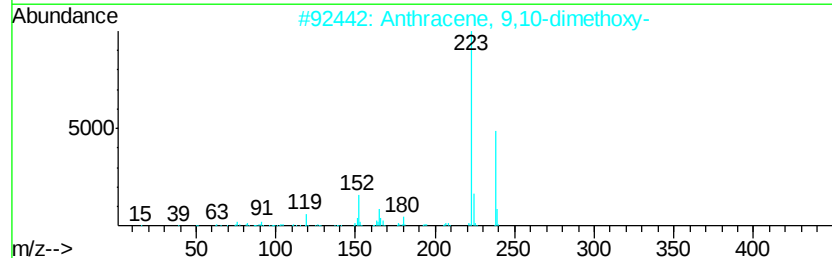
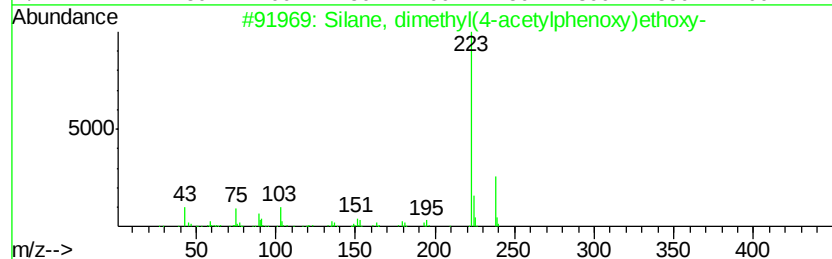
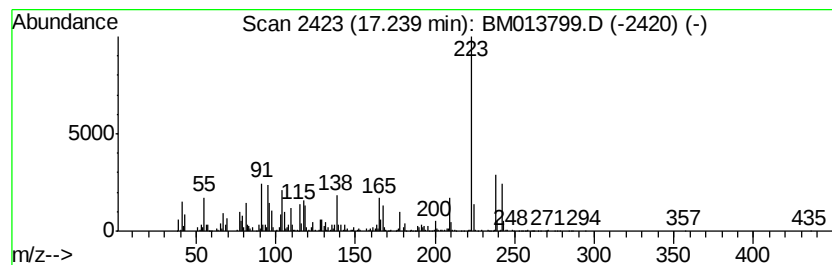
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 unknown-01 Concentration Rank 20

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.24 | 22.38 ng/ul | 7894580 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID                                                      | MW  | MolForm      | CAS#         | Qual |
|------|----|---|-------------------------------------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | Silane, dimethyl(4-acetylphenoxy)ethoxy-                          | 238 | C12H18O3Si   | 1000347-30-7 | 42   |
| 2    |    |   | Anthracene, 9,10-dimethoxy-                                       | 238 | C16H14O2     | 002395-97-3  | 30   |
| 3    |    |   | 2-Isopropylxanthen-9-one                                          | 238 | C16H14O2     | 069737-66-2  | 27   |
| 4    |    |   | 5-Acetonyl-5-(2-bromo-2-propenyl)-2-phenyl-1H-isobenzofuran-3-one | 302 | C10H11BrN2O4 | 021248-31-7  | 22   |
| 5    |    |   | 1H-Isoindole-1,3(2H)-dione, 2-phenyl-                             | 223 | C14H9N2O2    | 000520-03-6  | 22   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

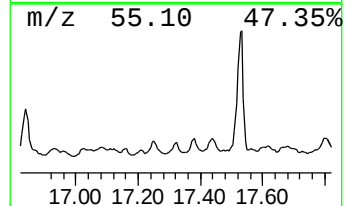
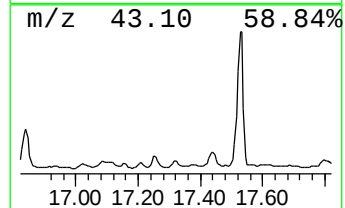
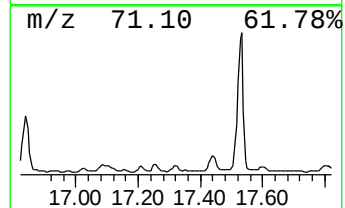
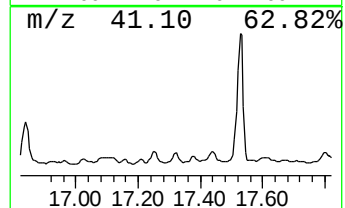
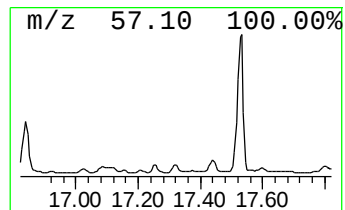
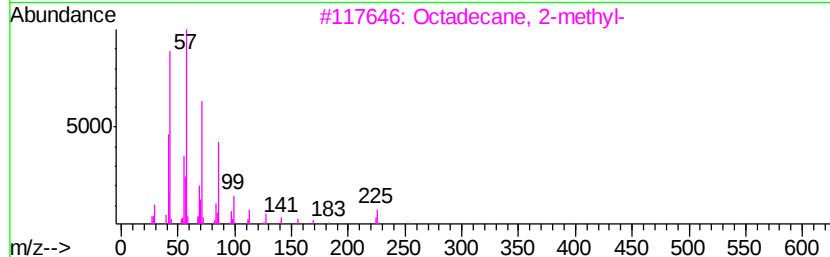
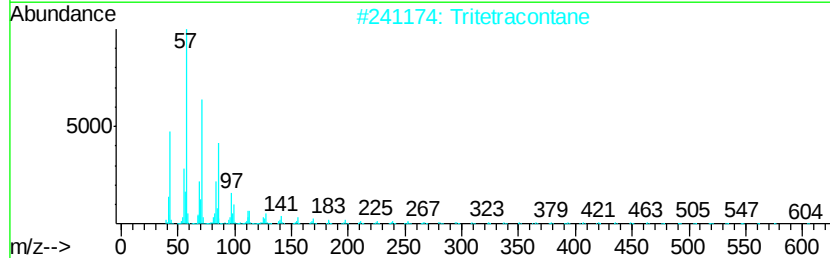
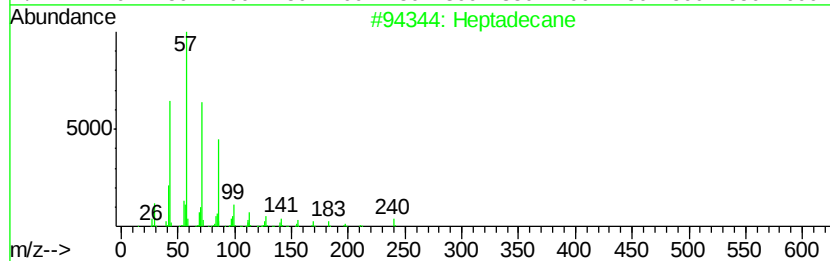
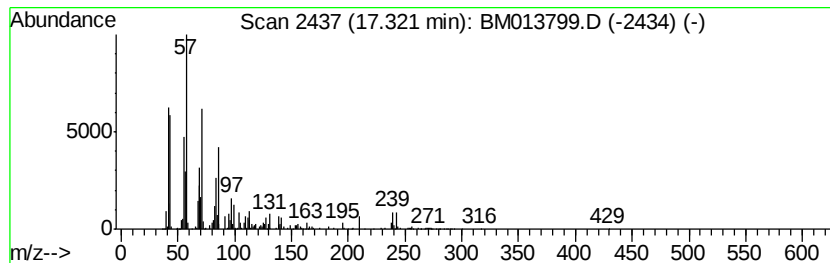
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.32 | 10.80 ng/ul | 3810670 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------|-----|---------|--------------|------|
| 1    |    |   | Heptadecane           | 240 | C17H36  | 000629-78-7  | 83   |
| 2    |    |   | Tritetracontane       | 605 | C43H88  | 007098-21-7  | 64   |
| 3    |    |   | Octadecane, 2-methyl- | 268 | C19H40  | 001560-88-9  | 64   |
| 4    |    |   | Tetratetracontane     | 619 | C44H90  | 007098-22-8  | 64   |
| 5    |    |   | 2-methyltetracosane   | 352 | C25H52  | 1000376-72-6 | 62   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

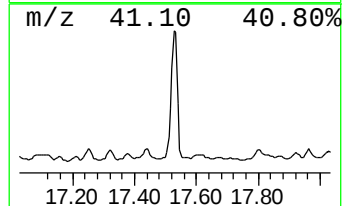
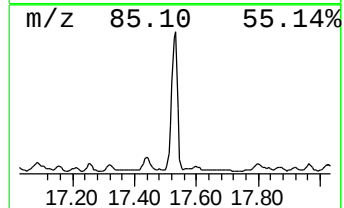
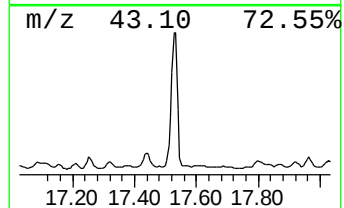
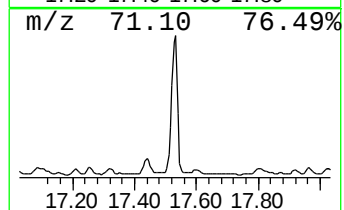
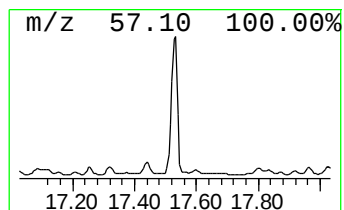
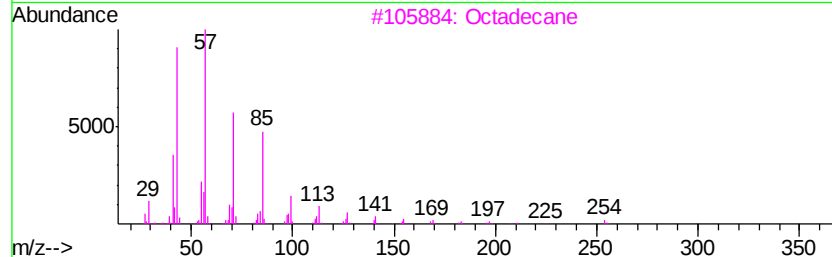
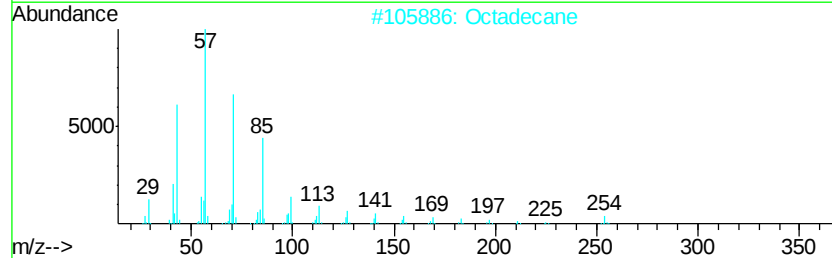
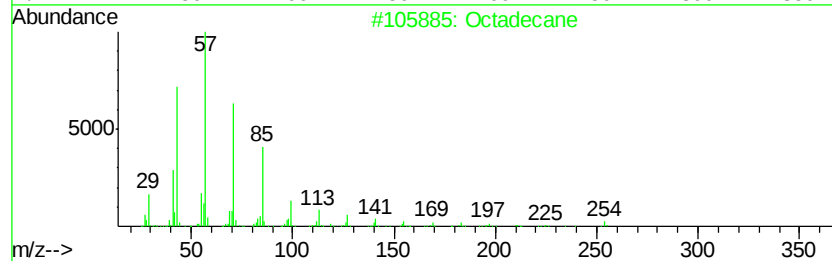
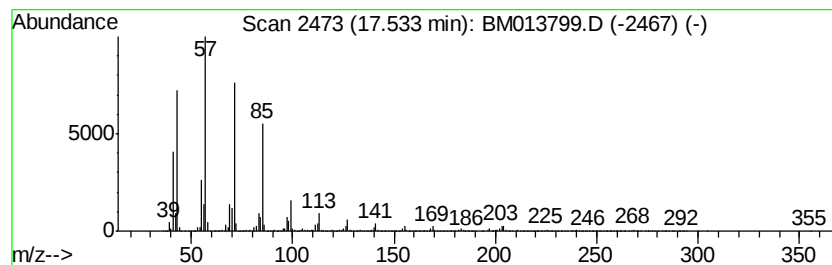
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 17.53 | 130.86 ng/ul | 46166800 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Octadecane   | 254 | C18H38  | 000593-45-3 | 97   |
| 2    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 97   |
| 3    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 94   |
| 4    |    | Tetradecane  | 198 | C14H30  | 000629-59-4 | 91   |
| 5    |    | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

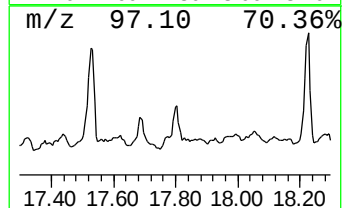
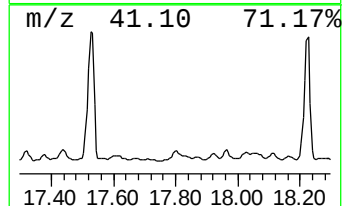
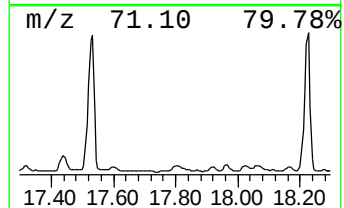
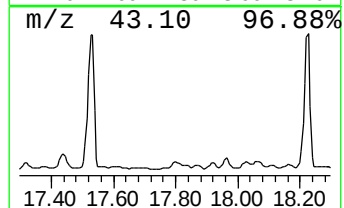
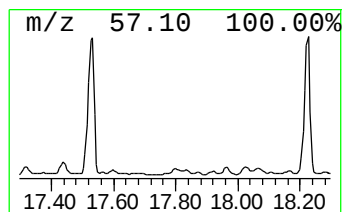
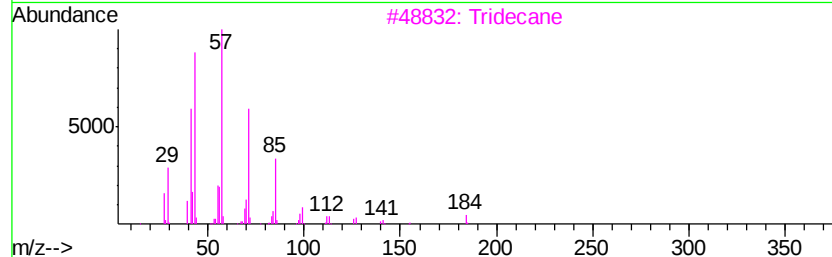
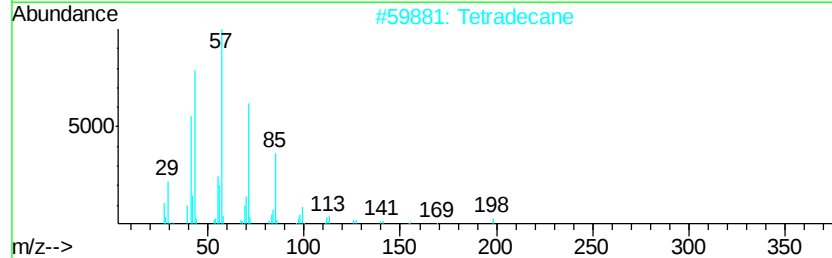
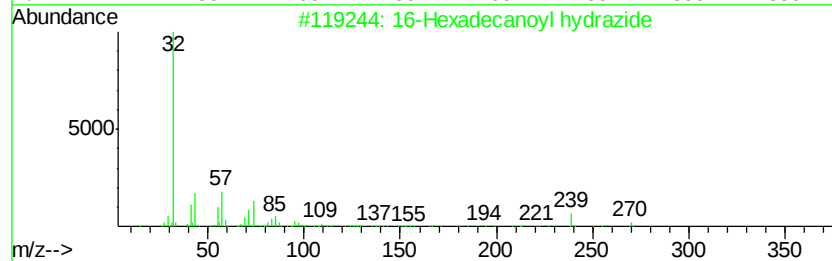
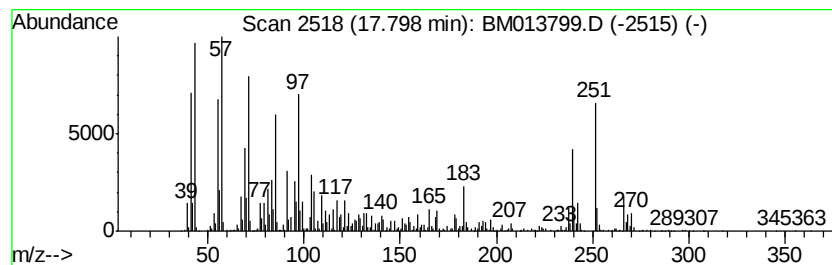
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 25 16-Hexadecanoyl hydrazide Concentration Rank 21

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 17.80 | 20.96 ng/ul | 7395560 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID              | MW  | MolForm   | CAS#        | Qual |
|------|----|---------------------------|-----|-----------|-------------|------|
| 1    | 5  | 16-Hexadecanovl hydrazide | 270 | C16H34N2O | 002619-88-7 | 64   |
| 2    |    | Tetradecane               | 198 | C14H30    | 000629-59-4 | 56   |
| 3    |    | Tridecane                 | 184 | C13H28    | 000629-50-5 | 53   |
| 4    |    | Nonadecane                | 268 | C19H40    | 000629-92-5 | 47   |
| 5    |    | Tridecane                 | 184 | C13H28    | 000629-50-5 | 47   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

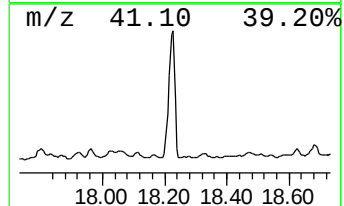
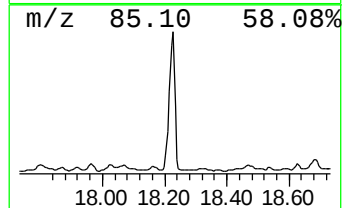
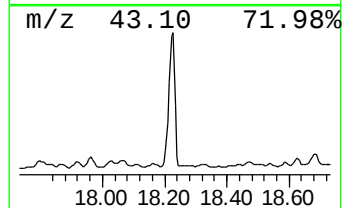
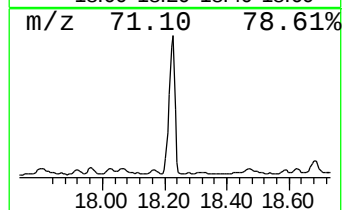
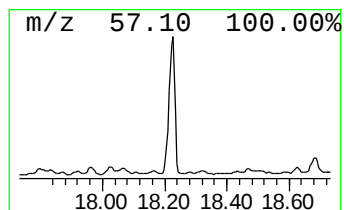
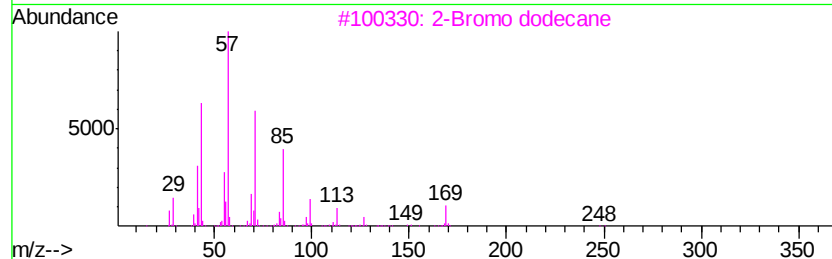
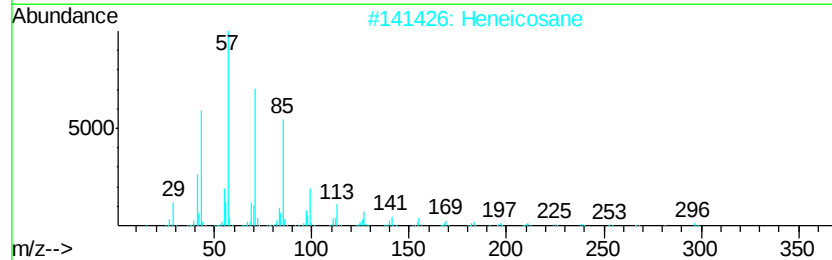
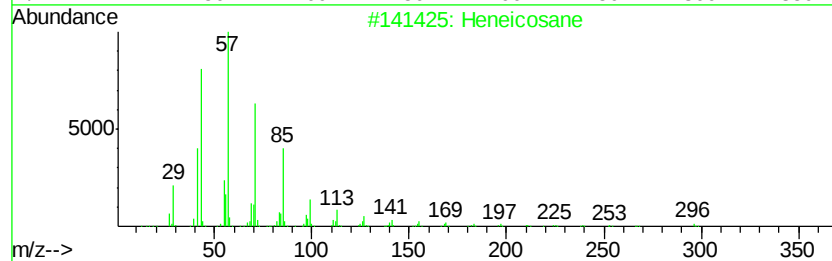
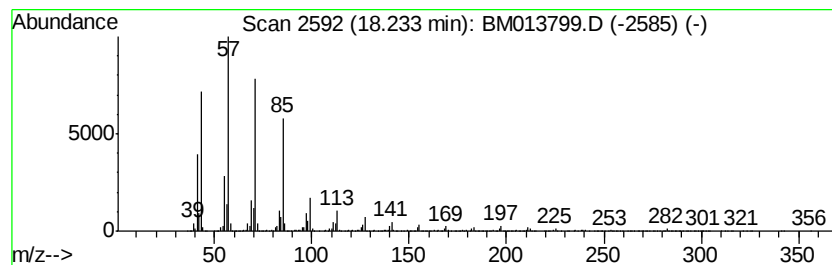
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 18.23 | 126.64 ng/ul | 44678400 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | Heneicosane           | 296 | C21H44   | 000629-94-7 | 97   |
| 2    |    | Heneicosane           | 296 | C21H44   | 000629-94-7 | 96   |
| 3    |    | 2-Bromo dodecane      | 248 | C12H25Br | 013187-99-0 | 94   |
| 4    |    | Hexacosane            | 366 | C26H54   | 000630-01-3 | 94   |
| 5    |    | Pentadecane, 8-hexyl- | 296 | C21H44   | 013475-75-7 | 94   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

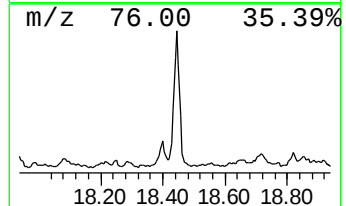
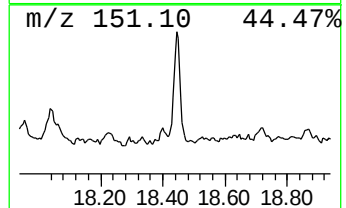
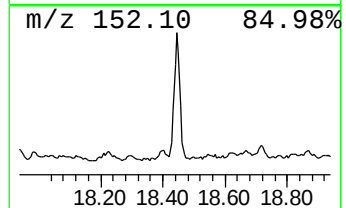
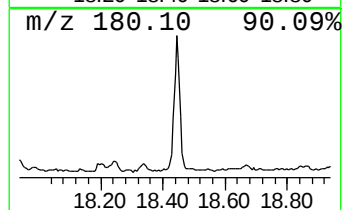
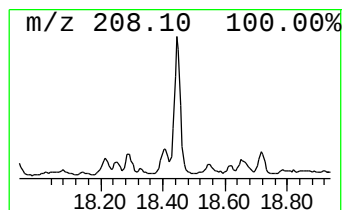
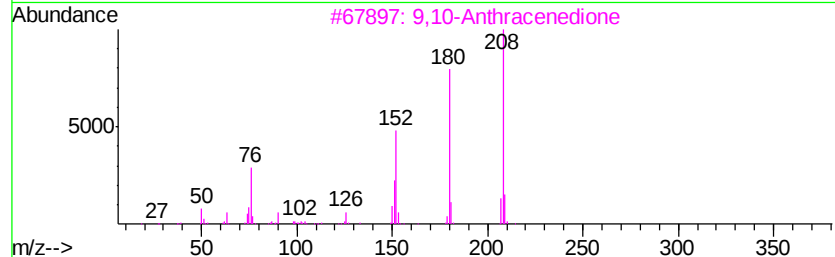
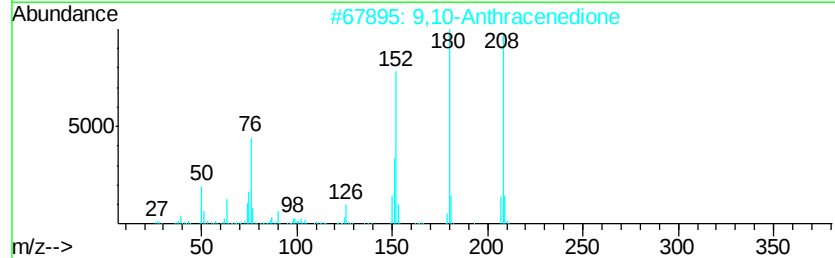
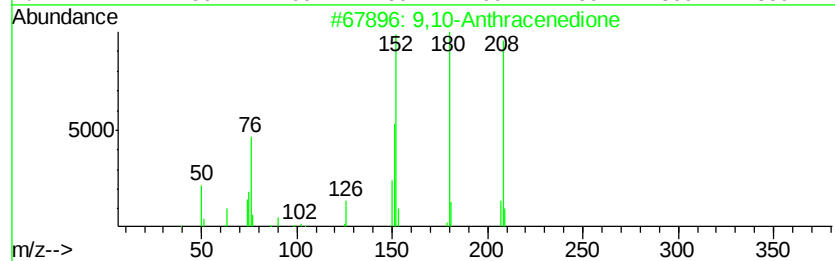
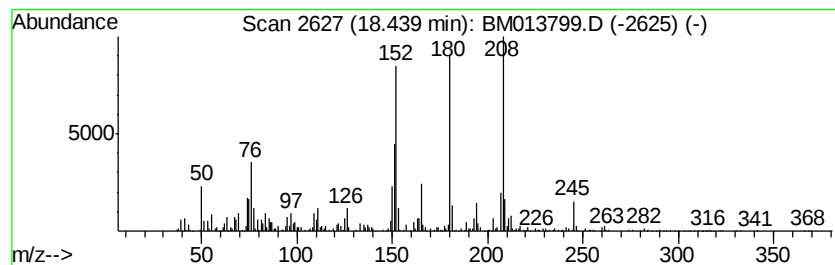
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 27 9,10-Anthracenedione Concentration Rank 30

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 18.44 | 9.81 ng/ul | 3460280 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 91   |
| 2    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 89   |
| 3    |    |   | 9,10-Anthracenedione | 208 | C14H8O2 | 000084-65-1 | 86   |
| 4    |    |   | Benzo[c]cinnoline    | 180 | C12H8N2 | 000230-17-1 | 78   |
| 5    |    |   | Benzo[c]cinnoline    | 180 | C12H8N2 | 000230-17-1 | 60   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

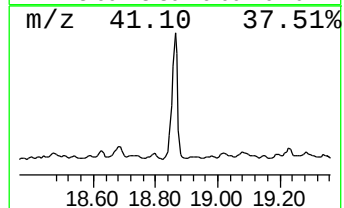
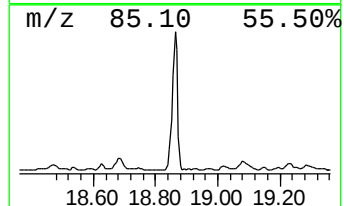
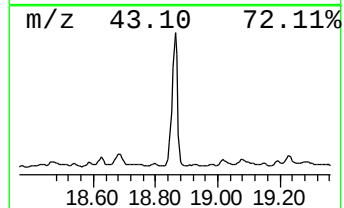
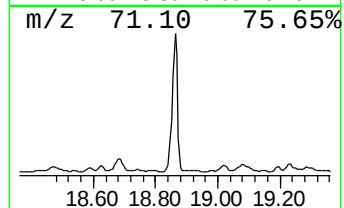
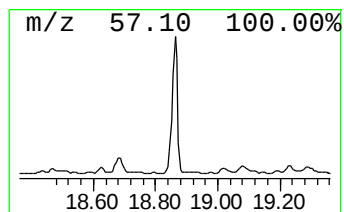
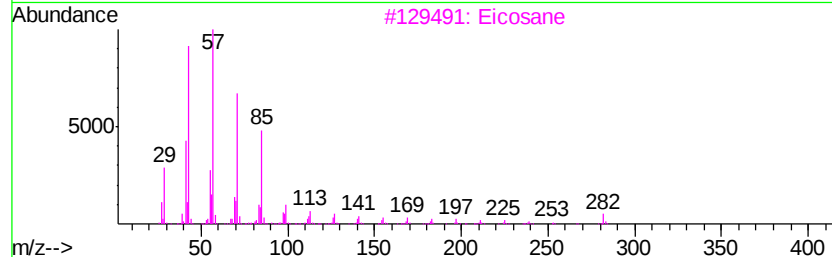
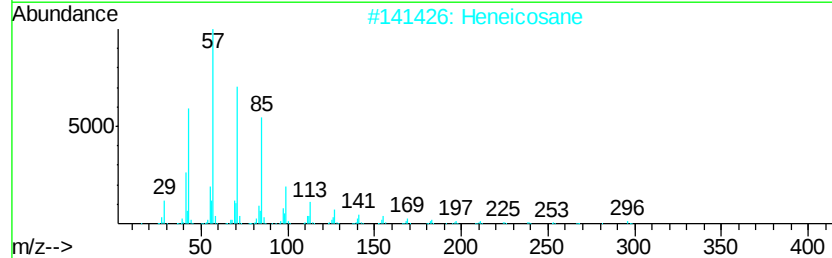
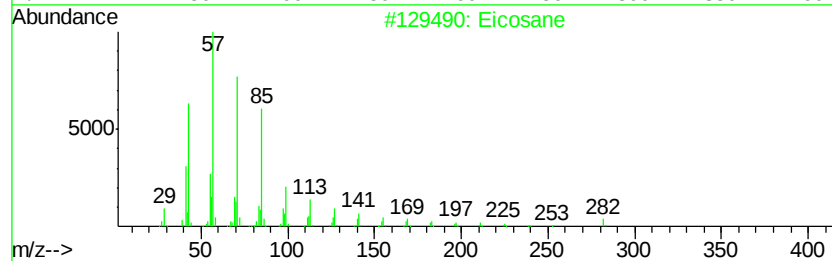
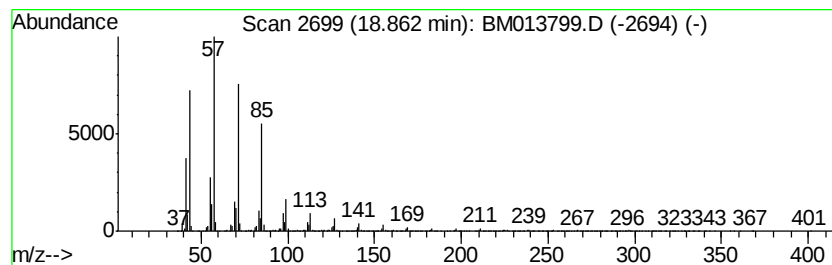
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 4

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 18.86 | 127.49 ng/ul | 44977500 | Phenanthrene-d10 | 17.03 |

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



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

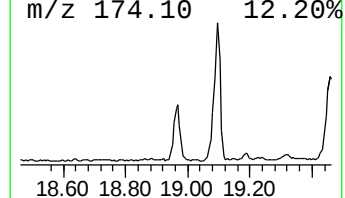
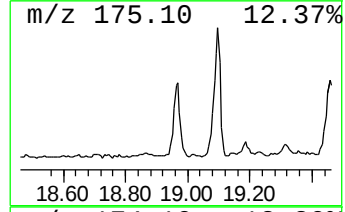
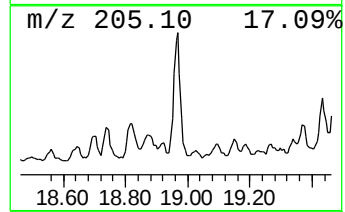
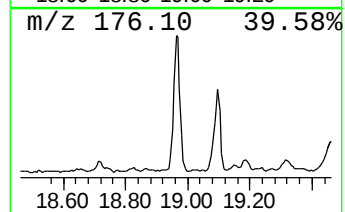
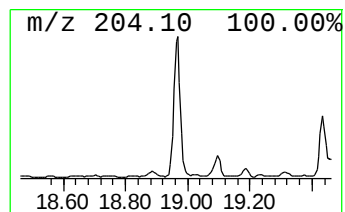
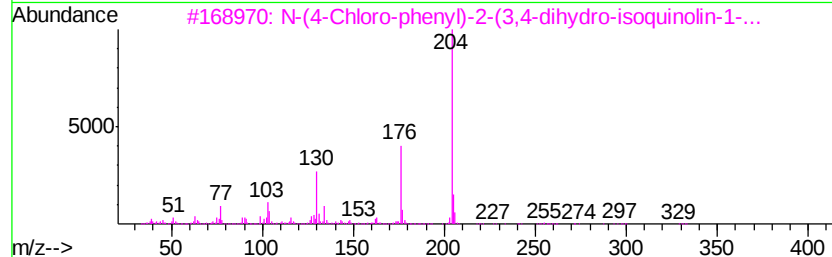
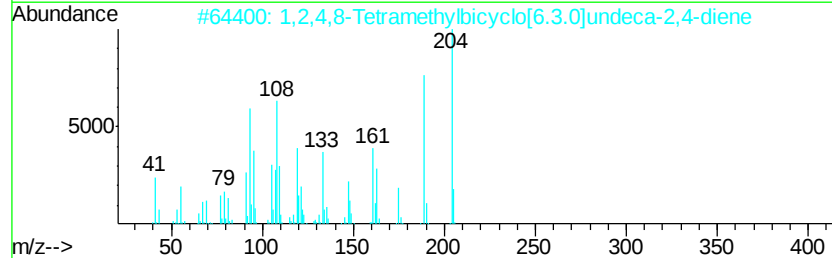
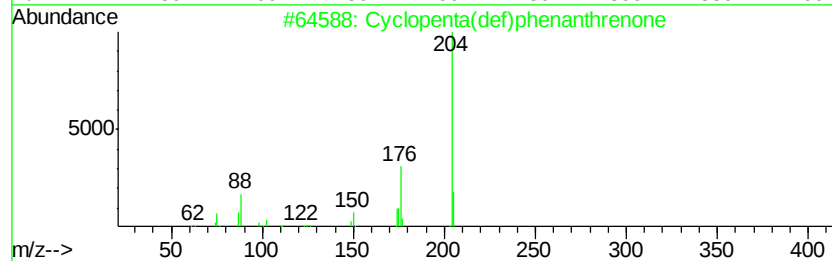
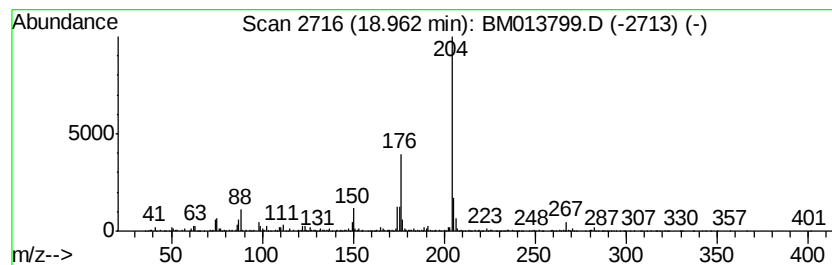
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 29 Cyclopenta(def)phenanthrenone Concentration Rank 24

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 18.96 | 15.71 ng/ul | 5542640 | Phenanthrene-d10 | 17.03 |

| Hit# | of | Tentative ID                                                        | MW  | MolForm      | CAS#         | Qual |
|------|----|---------------------------------------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Cyclopenta(def)phenanthrenone                                       | 204 | C15H8O       | 005737-13-3  | 91   |
| 2    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0]undeca-2,4-diene                   | 204 | C15H24       | 137235-51-9  | 89   |
| 3    |    | N-(4-Chloro-phenyl)-2-(3,4-dihydroisoquinolin-1-ylidene)ethan-1-one | 330 | C17H15ClN2OS | 1000296-34-3 | 59   |
| 4    |    | [1,1'-Biphenyl]-4,4'-dicarbonitrile                                 | 204 | C14H8N2      | 001591-30-6  | 53   |
| 5    |    | 6-Phenyl-2-thiouracil                                               | 204 | C10H8N2OS    | 005590-94-3  | 50   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

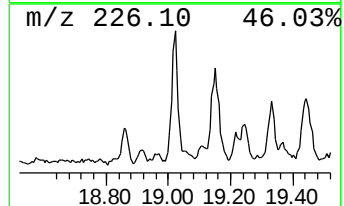
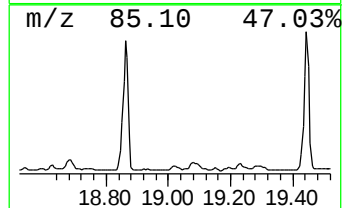
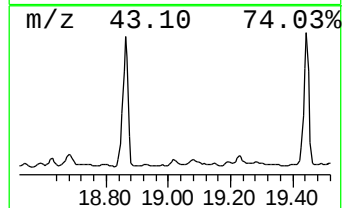
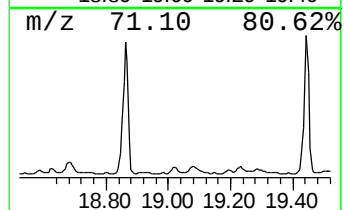
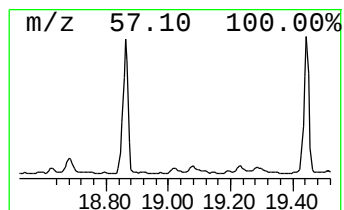
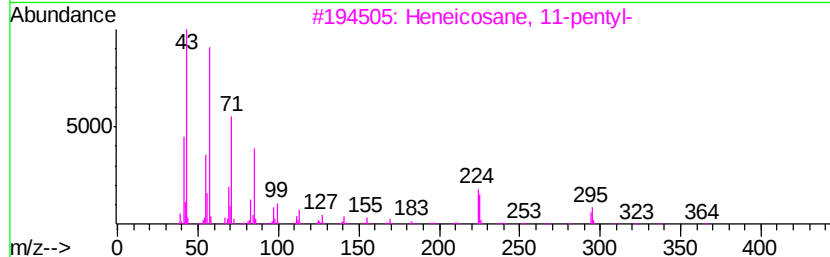
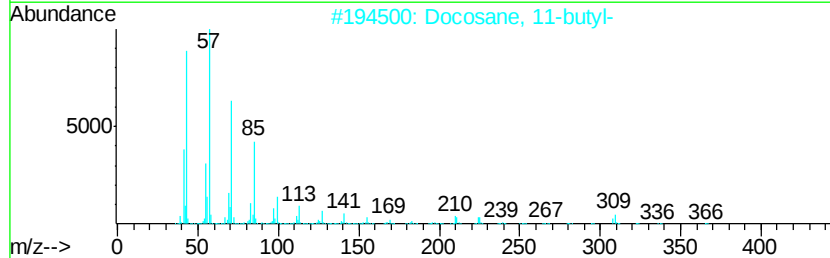
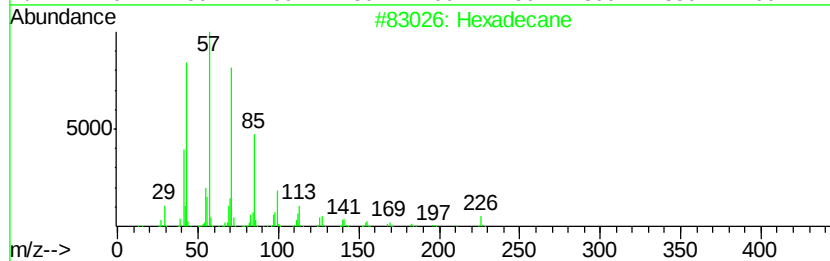
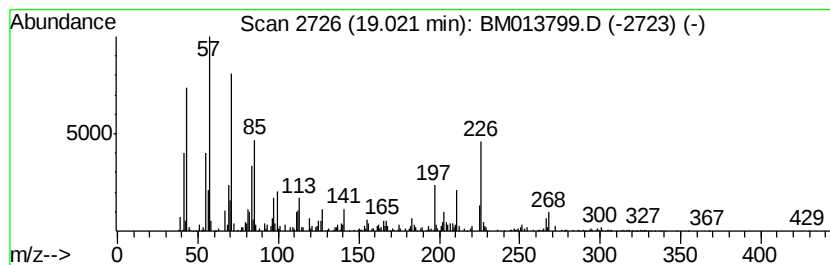
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.02 | 10.63 ng/ul | 3749830 | Phenanthrene-d10 | 17.03 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane              | 226 | C16H34  | 000544-76-3 | 78   |
| 2    |    |   | Docosane, 11-butyl-     | 366 | C26H54  | 013475-76-8 | 49   |
| 3    |    |   | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1 | 46   |
| 4    |    |   | Hexadecane, 2-methyl-   | 240 | C17H36  | 001560-92-5 | 45   |
| 5    |    |   | Octadecane              | 254 | C18H38  | 000593-45-3 | 45   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

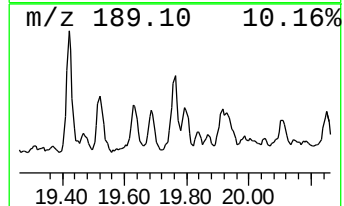
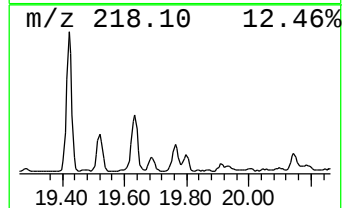
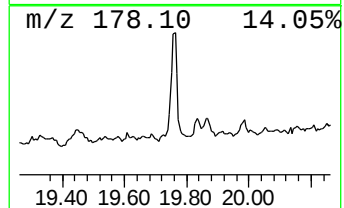
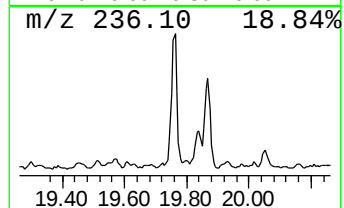
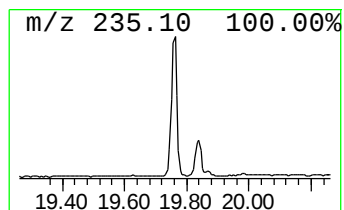
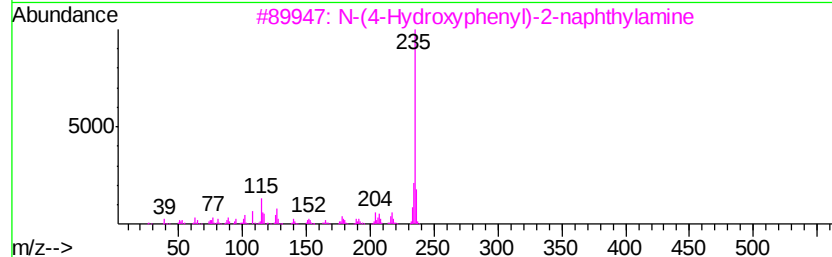
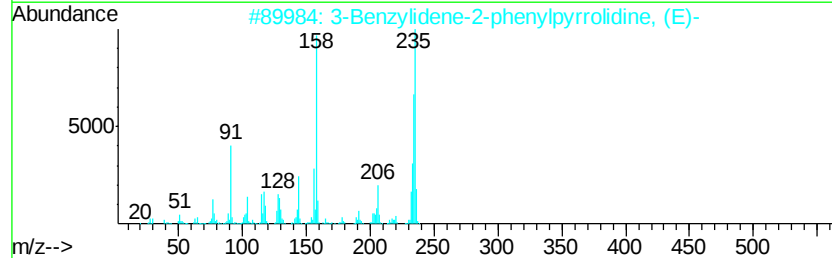
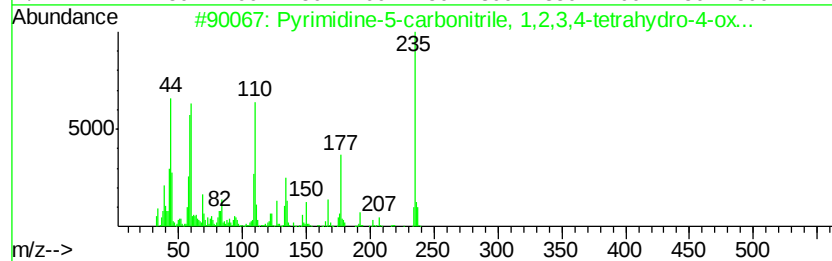
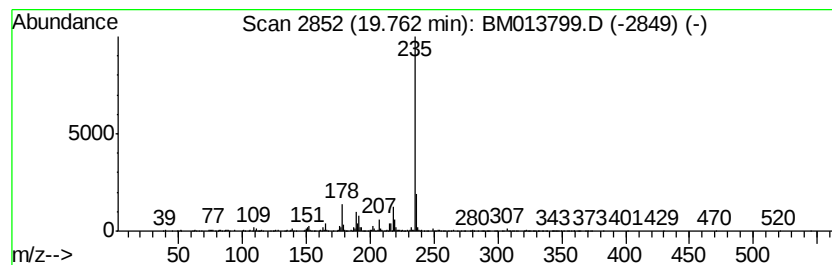
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 Pyrimidine-5-carbonitrile, ... Concentration Rank 17

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 19.76 | 30.98 ng/ul | 10666900 | Chrysene-d12     | 21.22 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Pyrimidine-5-carbonitrile, 1,2,3... | 235 | C9H5N3OS2  | 109532-65-2 | 53   |
| 2    |    | 3-Benzylidene-2-phenylpyrrolidin... | 235 | C17H17N    | 122949-15-9 | 53   |
| 3    |    | N-(4-Hydroxyphenyl)-2-naphthylamine | 235 | C16H13NO   | 000093-45-8 | 45   |
| 4    |    | Acetamide, N-[3-[(2-cyanoethyl)e... | 275 | C15H21N3O2 | 067905-64-0 | 45   |
| 5    |    | 4-Hydroxy-3-methyl-beta-phenylci... | 235 | C16H13NO   | 016299-28-8 | 45   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

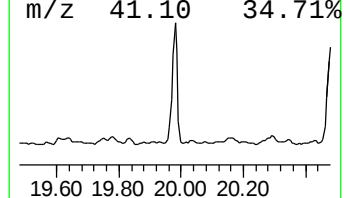
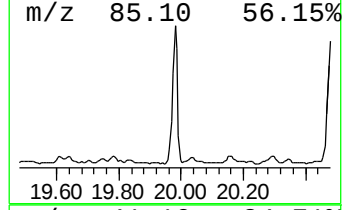
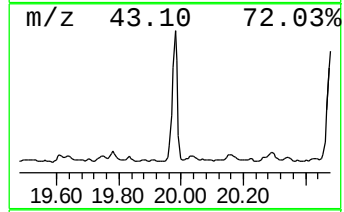
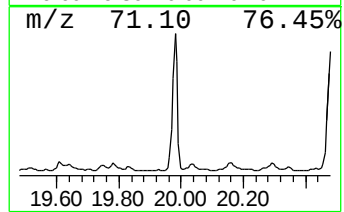
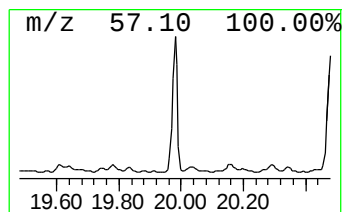
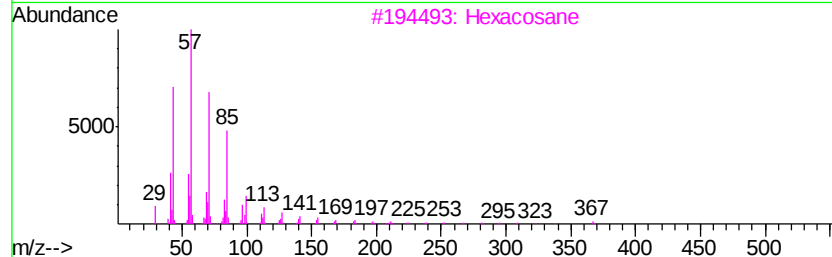
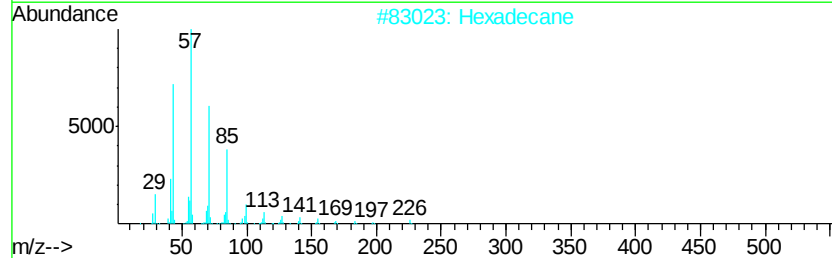
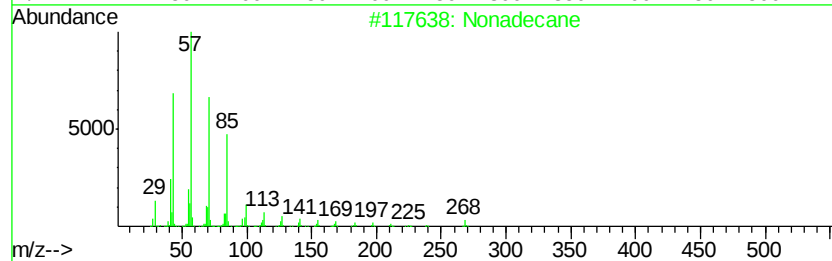
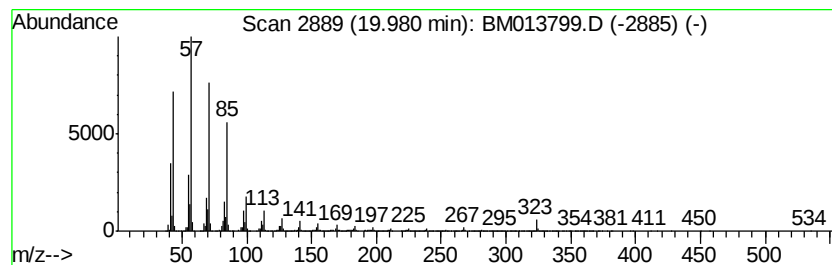
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc      | Area     | Relative to ISTD | R.T.  |
|-------|--------------|----------|------------------|-------|
| 19.98 | 109.31 ng/ul | 37641200 | Chrysene-d12     | 21.22 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane   | 268 | C19H40  | 000629-92-5 | 94   |
| 2    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 3    |    | Hexacosane   | 366 | C26H54  | 000630-01-3 | 93   |
| 4    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 93   |
| 5    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

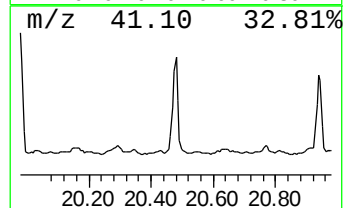
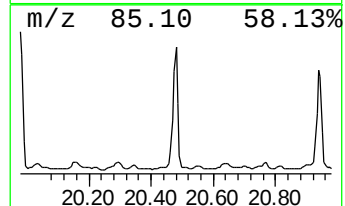
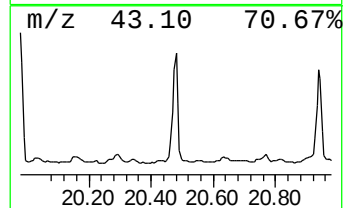
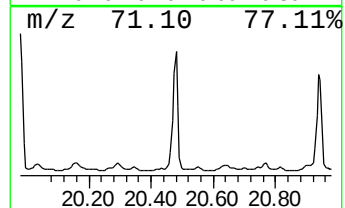
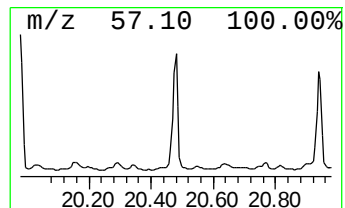
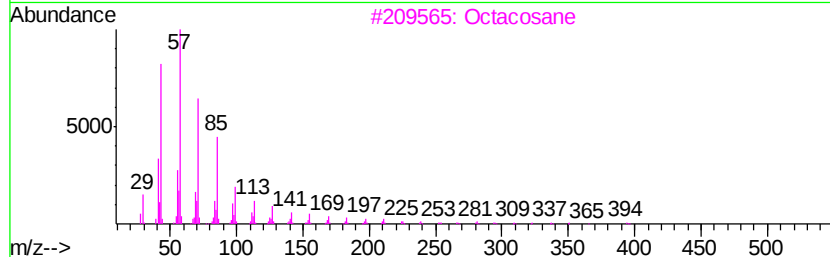
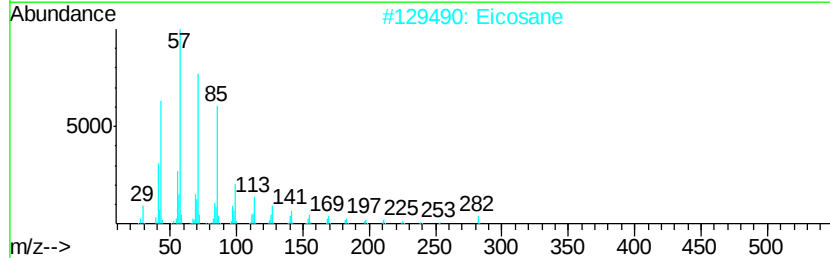
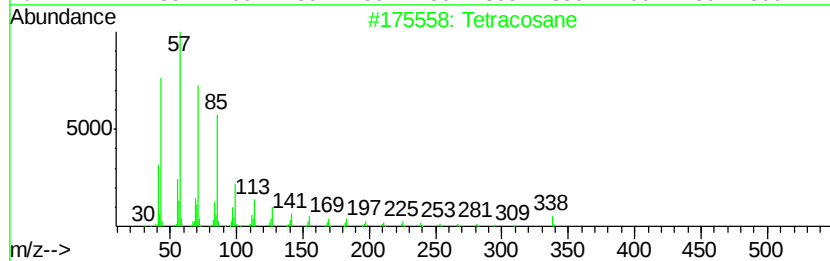
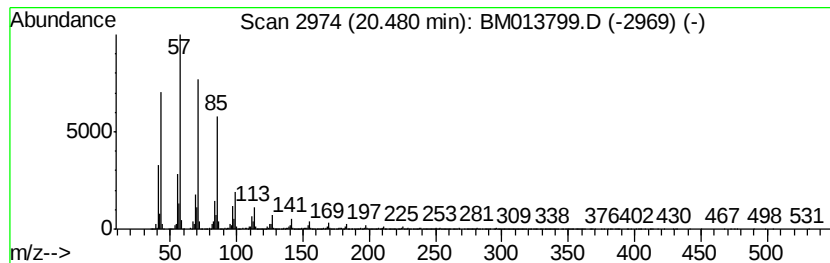
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 20.48 | 90.97 ng/ul | 31326300 | Chrysene-d12     | 21.22 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane             | 338 | C24H50  | 000646-31-1 | 98   |
| 2    |    | Eicosane                | 282 | C20H42  | 000112-95-8 | 98   |
| 3    |    | Octacosane              | 394 | C28H58  | 000630-02-4 | 97   |
| 4    |    | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1 | 93   |
| 5    |    | Hexadecane              | 226 | C16H34  | 000544-76-3 | 93   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

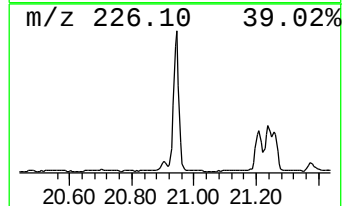
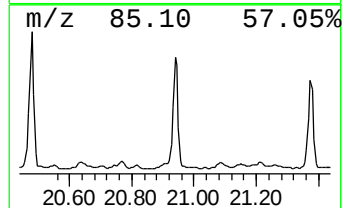
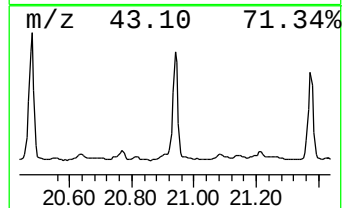
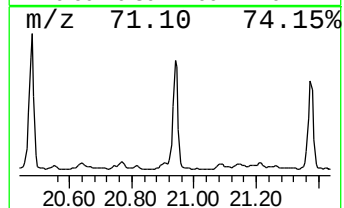
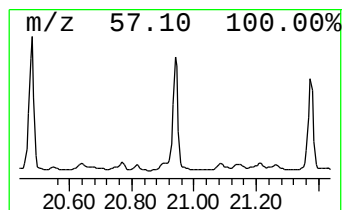
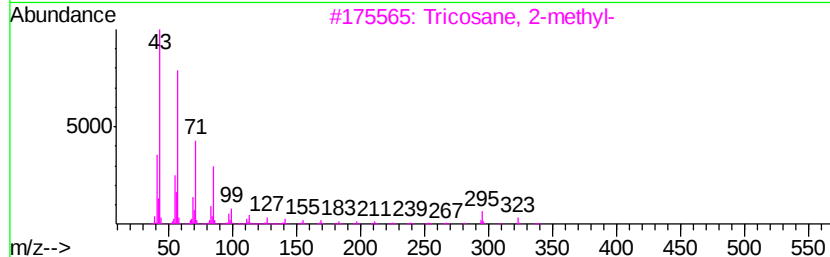
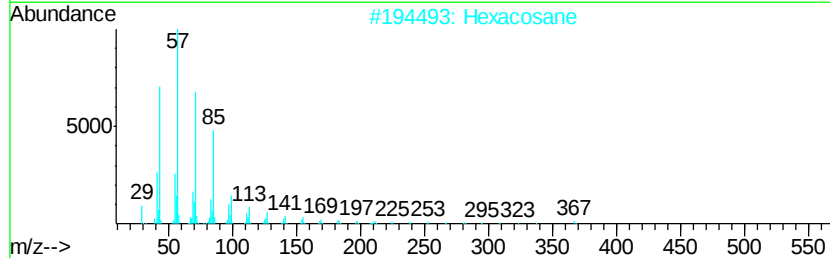
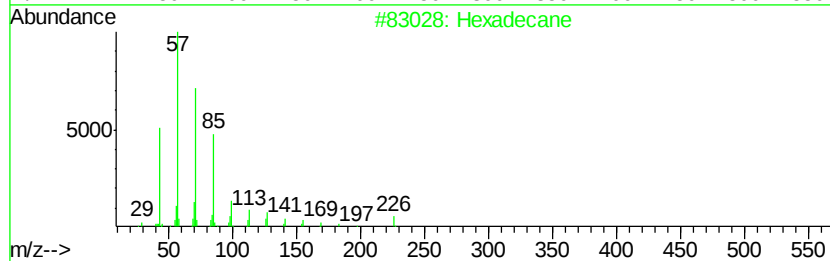
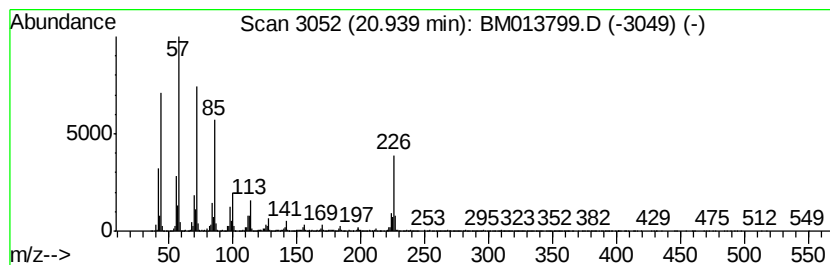
Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.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 9

| R.T.    | EstConc                 | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------|--------------|------------------|----------------|
| 20.94   | 87.41 ng/ul             | 30100000     | Chrysene-d12     | 21.22          |
| Hit# of | 5                       | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Hexadecane              |              | 226 C16H34       | 000544-76-3 83 |
| 2       | Hexacosane              |              | 366 C26H54       | 000630-01-3 76 |
| 3       | Tricosane, 2-methyl-    |              | 338 C24H50       | 001928-30-9 76 |
| 4       | 1-Iodo-2-methylundecane |              | 296 C12H25I      | 073105-67-6 76 |
| 5       | Heptadecane             |              | 240 C17H36       | 000629-78-7 70 |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

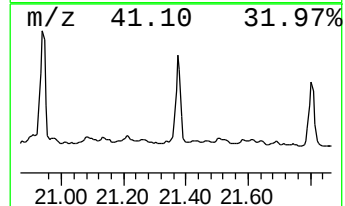
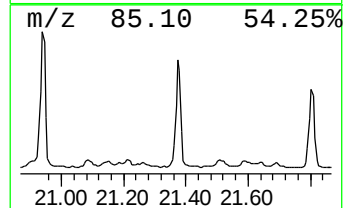
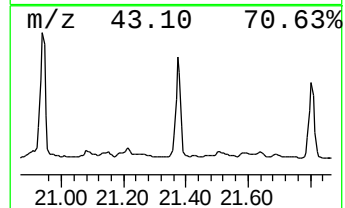
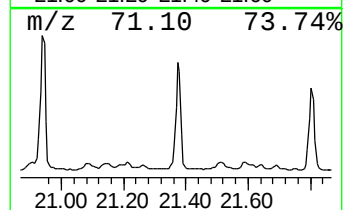
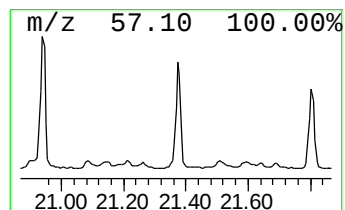
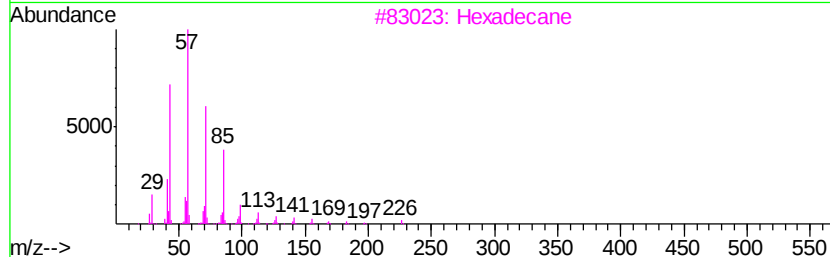
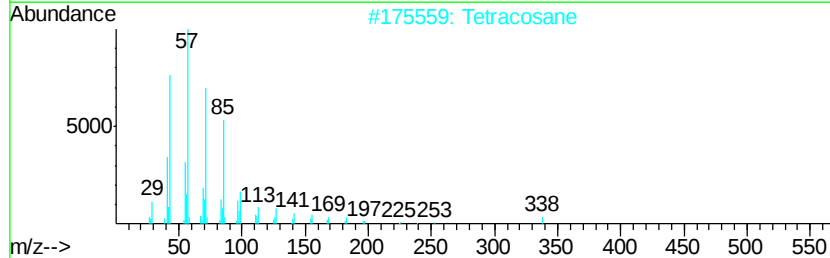
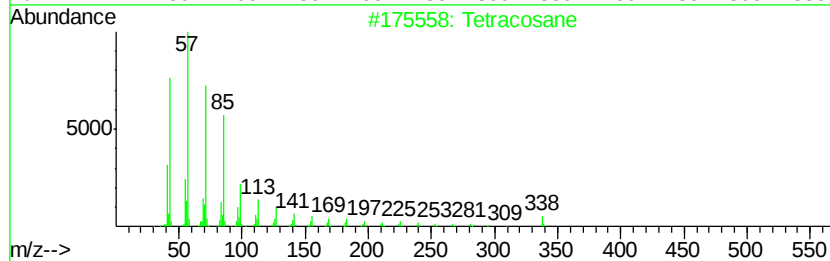
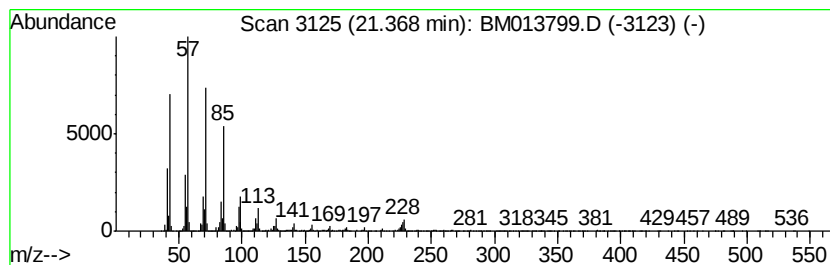
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 21.37 | 59.52 ng/ul | 20496200 | Chrysene-d12     | 21.22 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane  | 338 | C24H50  | 000646-31-1 | 98   |
| 2    |    | Tetracosane  | 338 | C24H50  | 000646-31-1 | 96   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 96   |
| 4    |    | Octacosane   | 394 | C28H58  | 000630-02-4 | 96   |
| 5    |    | Heptadecane  | 240 | C17H36  | 000629-78-7 | 93   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

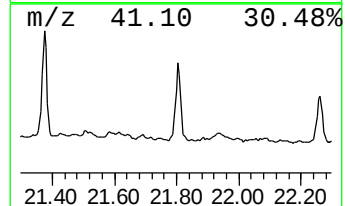
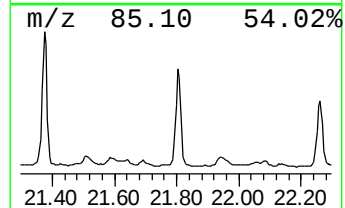
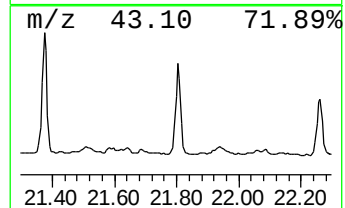
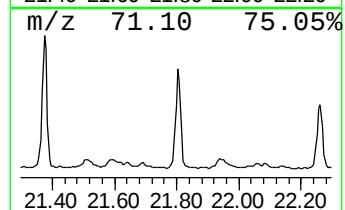
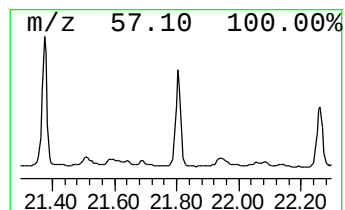
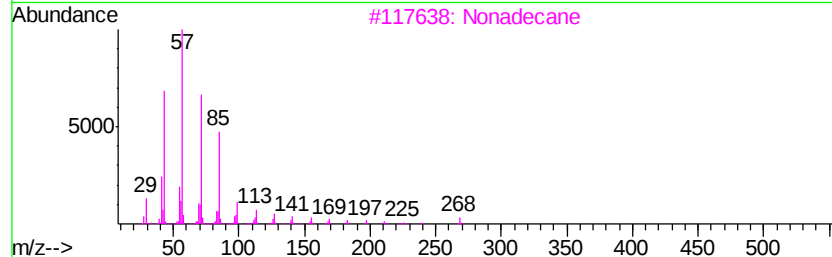
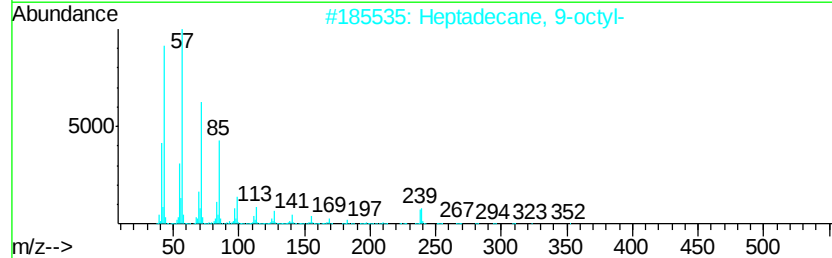
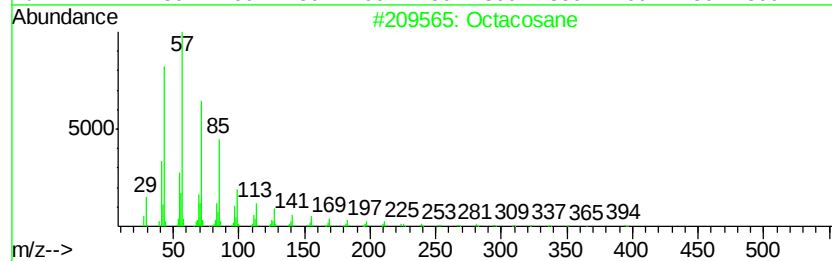
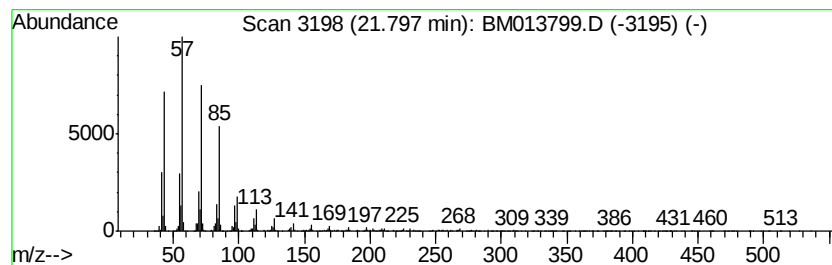
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 12

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 21.80 | 55.97 ng/ul | 19273200 | Chrysene-d12     | 21.22 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 98   |
| 2    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 95   |
| 3    |    |   | Nonadecane            | 268 | C19H40  | 000629-92-5 | 93   |
| 4    |    |   | Heptacosane           | 380 | C27H56  | 000593-49-7 | 91   |
| 5    |    |   | Pentadecane, 8-hexyl- | 296 | C21H44  | 013475-75-7 | 91   |



Data Path : Z:\HPCHEM1\BNA M\Data\BM012418\  
Data File : BM013799.D  
Acq On : 25 Jan 2018 09:43  
Operator : SJ/JU  
Sample : J1222-13  
Misc :  
ALS Vial : 33 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
F6A95

Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM011018.M  
Quant Title : SVOA CALIBRATION

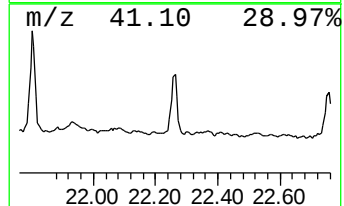
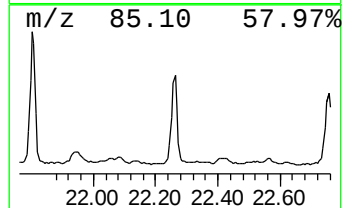
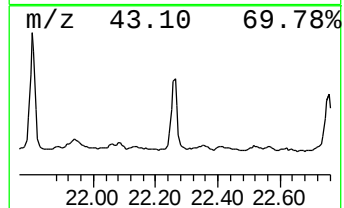
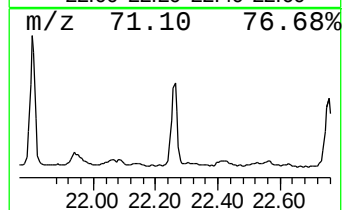
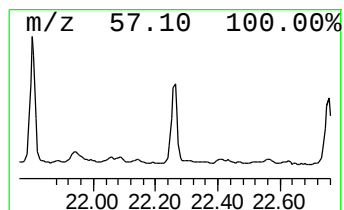
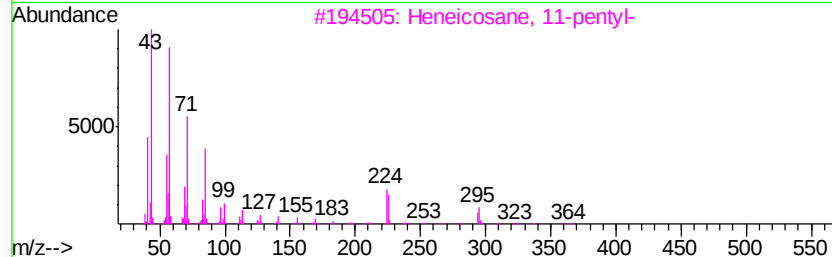
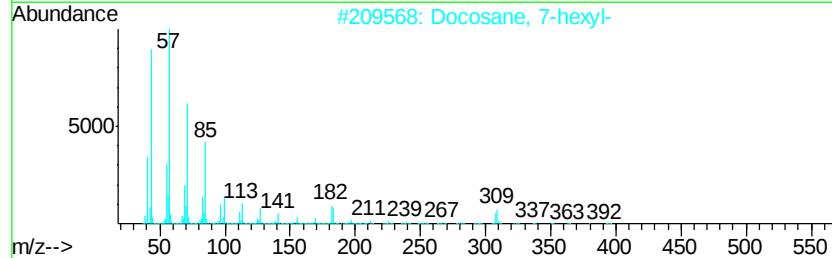
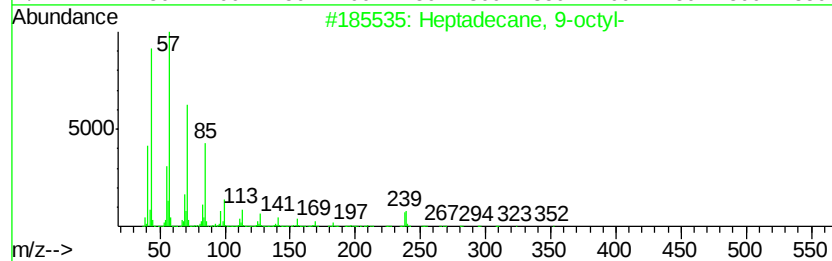
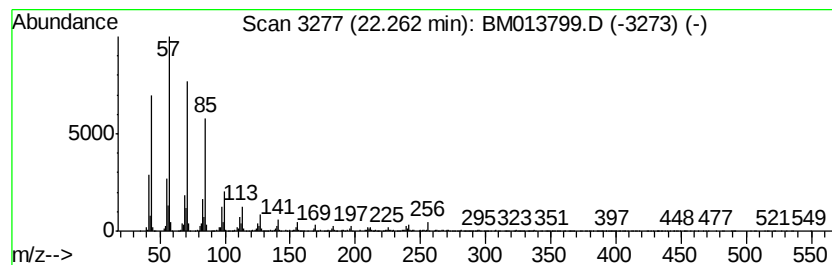
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc     | Area     | Relative to ISTD | R.T.  |
|-------|-------------|----------|------------------|-------|
| 22.26 | 33.95 ng/ul | 11690200 | Chrysene-d12     | 21.22 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------|-----|---------|--------------|------|
| 1    | 5  | Heptadecane, 9-octyl-   | 352 | C25H52  | 007225-64-1  | 93   |
| 2    |    | Docosane, 7-hexyl-      | 394 | C28H58  | 055373-86-9  | 93   |
| 3    |    | Heneicosane, 11-pentyl- | 366 | C26H54  | 014739-72-1  | 91   |
| 4    |    | 2-methyloctacosane      | 408 | C29H60  | 1000376-72-8 | 90   |
| 5    |    | Heneicosane             | 296 | C21H44  | 000629-94-7  | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\Data\BM012418\  
 Data File : BM013799.D  
 Acq On : 25 Jan 2018 09:43  
 Operator : SJ/JU  
 Sample : J1222-13  
 Misc :  
 ALS Vial : 33 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 F6A95

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011018.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 |
| (DEL) Alkane: Str...  | 8.81  | 20.0    | ng/ul | 3249440  | 1                     | 7.59  | 3249440  | 20.0 |
| Sulfurous acid, d...  | 11.40 | 11.7    | ng/ul | 4868150  | 2                     | 10.37 | 8296570  | 20.0 |
| (DEL) Alkane: Str...  | 11.80 | 25.9    | ng/ul | 10758400 | 2                     | 10.37 | 8296570  | 20.0 |
| (DEL) Alkane: Str...  | 12.77 | 7.7     | ng/ul | 4650810  | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 13.78 | 6.6     | ng/ul | 3993750  | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 14.17 | 38.4    | ng/ul | 23173200 | 3                     | 14.25 | 12072500 | 20.0 |
| 4'-(1-Pyrrolyl)ace... | 14.40 | 6.8     | ng/ul | 4113550  | 3                     | 14.25 | 12072500 | 20.0 |
| Heptacosane, 1-ch...  | 14.79 | 13.3    | ng/ul | 8021700  | 3                     | 14.25 | 12072500 | 20.0 |
| Naphthalene, 2,3,...  | 15.02 | 5.9     | ng/ul | 3577440  | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 15.13 | 49.3    | ng/ul | 29733400 | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 15.53 | 26.6    | ng/ul | 16035900 | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 15.62 | 7.0     | ng/ul | 4212920  | 3                     | 14.25 | 12072500 | 20.0 |
| (DEL) Alkane: Str...  | 15.67 | 11.4    | ng/ul | 4022740  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 16.00 | 134.4   | ng/ul | 47428300 | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 16.02 | 94.8    | ng/ul | 33435700 | 4                     | 17.03 | 7055730  | 20.0 |
| (4-Acetylphenyl)p...  | 16.05 | 17.3    | ng/ul | 6110130  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 16.50 | 9.3     | ng/ul | 3286310  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 16.79 | 149.2   | ng/ul | 52639200 | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 16.84 | 58.2    | ng/ul | 20543900 | 4                     | 17.03 | 7055730  | 20.0 |
| Benzene, 1,1'-eth...  | 16.92 | 9.5     | ng/ul | 3360940  | 4                     | 17.03 | 7055730  | 20.0 |
| unknown-01            | 17.24 | 22.4    | ng/ul | 7894580  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 17.32 | 10.8    | ng/ul | 3810670  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 17.53 | 130.9   | ng/ul | 46166800 | 4                     | 17.03 | 7055730  | 20.0 |
| 16-Hexadecanoyl h...  | 17.80 | 21.0    | ng/ul | 7395560  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 18.23 | 126.6   | ng/ul | 44678400 | 4                     | 17.03 | 7055730  | 20.0 |
| 9,10-Anthracenedione  | 18.44 | 9.8     | ng/ul | 3460280  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 18.86 | 127.5   | ng/ul | 44977500 | 4                     | 17.03 | 7055730  | 20.0 |
| Cyclopenta(def)ph...  | 18.96 | 15.7    | ng/ul | 5542640  | 4                     | 17.03 | 7055730  | 20.0 |
| (DEL) Alkane: Str...  | 19.02 | 10.6    | ng/ul | 3749830  | 4                     | 17.03 | 7055730  | 20.0 |
| Pyrimidine-5-carb...  | 19.76 | 31.0    | ng/ul | 10666900 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 19.98 | 109.3   | ng/ul | 37641200 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 20.48 | 91.0    | ng/ul | 31326300 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 20.94 | 87.4    | ng/ul | 30100000 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 21.37 | 59.5    | ng/ul | 20496200 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 21.80 | 56.0    | ng/ul | 19273200 | 5                     | 21.22 | 6887130  | 20.0 |
| (DEL) Alkane: Str...  | 22.26 | 34.0    | ng/ul | 11690200 | 5                     | 21.22 | 6887130  | 20.0 |