

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
 Data File : BP002790.D  
 Acq On : 17 Jul 2020 14:03  
 Operator : CG/JU  
 Sample : L3324-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 10:1-COMP

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP071020.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.717       | 306           | 311         | 318          | rVB2     | 36689          | 59866         | 1.16%           | 0.110%        |
| 2         | 5.187       | 385           | 391         | 415          | rVB      | 2490867        | 3798843       | 73.76%          | 6.974%        |
| 3         | 6.763       | 652           | 659         | 676          | rBV      | 2374661        | 4058631       | 78.81%          | 7.451%        |
| 4         | 7.175       | 724           | 729         | 742          | rBV      | 74335          | 160181        | 3.11%           | 0.294%        |
| 5         | 7.558       | 786           | 794         | 804          | rBV      | 871665         | 1388375       | 26.96%          | 2.549%        |
| 6         | 8.705       | 975           | 989         | 1004         | rBV      | 1581914        | 2750712       | 53.41%          | 5.050%        |
| 7         | 10.328      | 1256          | 1265        | 1278         | rBV      | 1141117        | 1994008       | 38.72%          | 3.661%        |
| 8         | 11.028      | 1379          | 1384        | 1391         | rVB6     | 26446          | 64863         | 1.26%           | 0.119%        |
| 9         | 11.099      | 1391          | 1396        | 1404         | rBV7     | 34426          | 75625         | 1.47%           | 0.139%        |
| 10        | 11.393      | 1440          | 1446        | 1457         | rBV6     | 72128          | 193368        | 3.75%           | 0.355%        |
| 11        | 11.616      | 1479          | 1484        | 1491         | rBV5     | 36650          | 104963        | 2.04%           | 0.193%        |
| 12        | 11.804      | 1511          | 1516        | 1523         | rBV5     | 49512          | 112864        | 2.19%           | 0.207%        |
| 13        | 12.151      | 1572          | 1575        | 1579         | rVB      | 60227          | 72843         | 1.41%           | 0.134%        |
| 14        | 12.269      | 1591          | 1595        | 1598         | rBV2     | 65938          | 94132         | 1.83%           | 0.173%        |
| 15        | 12.634      | 1653          | 1657        | 1662         | rBV5     | 76379          | 127531        | 2.48%           | 0.234%        |
| 16        | 12.763      | 1675          | 1679        | 1682         | rBV      | 325346         | 474232        | 9.21%           | 0.871%        |
| 17        | 12.822      | 1683          | 1689        | 1699         | rVB      | 3222941        | 5150095       | 100.00%         | 9.455%        |
| 18        | 12.969      | 1710          | 1714        | 1716         | rBV3     | 105376         | 169444        | 3.29%           | 0.311%        |
| 19        | 13.028      | 1720          | 1724        | 1728         | rVV5     | 182709         | 316397        | 6.14%           | 0.581%        |
| 20        | 13.146      | 1742          | 1744        | 1749         | rVB3     | 150402         | 191312        | 3.71%           | 0.351%        |
| 21        | 13.369      | 1779          | 1782        | 1785         | rVB3     | 126416         | 150356        | 2.92%           | 0.276%        |
| 22        | 13.487      | 1798          | 1802        | 1807         | rBV4     | 435889         | 623964        | 12.12%          | 1.146%        |
| 23        | 13.575      | 1813          | 1817        | 1823         | rBV5     | 364690         | 654462        | 12.71%          | 1.202%        |
| 24        | 13.634      | 1824          | 1827        | 1830         | rBV3     | 268136         | 340132        | 6.60%           | 0.624%        |
| 25        | 13.675      | 1830          | 1834        | 1837         | rVV      | 510318         | 728935        | 14.15%          | 1.338%        |
| 26        | 13.728      | 1838          | 1843        | 1852         | rVB3     | 971747         | 1707763       | 33.16%          | 3.135%        |
| 27        | 13.940      | 1876          | 1879        | 1884         | rBV3     | 203641         | 352195        | 6.84%           | 0.647%        |
| 28        | 13.998      | 1886          | 1889        | 1892         | rVV2     | 240999         | 302957        | 5.88%           | 0.556%        |
| 29        | 14.045      | 1893          | 1897        | 1901         | rVV3     | 454969         | 690167        | 13.40%          | 1.267%        |
| 30        | 14.128      | 1908          | 1911        | 1917         | rBV6     | 212339         | 526014        | 10.21%          | 0.966%        |
| 31        | 14.210      | 1918          | 1925        | 1930         | rBV      | 1567758        | 2971343       | 57.69%          | 5.455%        |
| 32        | 14.775      | 2017          | 2021        | 2026         | rVB3     | 659958         | 934240        | 18.14%          | 1.715%        |
| 33        | 15.016      | 2059          | 2062        | 2066         | rVB3     | 496089         | 622843        | 12.09%          | 1.143%        |
| 34        | 15.522      | 2144          | 2148        | 2155         | rVB      | 1271683        | 2022468       | 39.27%          | 3.713%        |

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Instrument :  
 BNA\_P  
 ClientSampleId :  
 10:1-COMP

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON Filtering: 5  
 Sampling : 1 Min Area: 3 % of largest Peak  
 Start Thrs: 0.2 Max Peaks: 100  
 Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP071020.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 15.722 | 2177 | 2182 | 2188 | rBV  | 1474580 | 2241688 | 43.53% | 4.115% |
| 36 | 16.004 | 2226 | 2230 | 2237 | rVB  | 1954266 | 2950123 | 57.28% | 5.416% |
| 37 | 16.828 | 2366 | 2370 | 2376 | rVB  | 2105925 | 3438880 | 66.77% | 6.313% |
| 38 | 19.569 | 2833 | 2836 | 2847 | rVB  | 2197318 | 3006024 | 58.37% | 5.519% |
| 39 | 20.816 | 3046 | 3048 | 3057 | rVB2 | 915188  | 1332706 | 25.88% | 2.447% |
| 40 | 21.086 | 3090 | 3094 | 3097 | rVB  | 1059607 | 1411347 | 27.40% | 2.591% |
| 41 | 21.686 | 3192 | 3196 | 3207 | rVB2 | 355430  | 809576  | 15.72% | 1.486% |
| 42 | 22.351 | 3306 | 3309 | 3314 | rBV2 | 331878  | 406645  | 7.90%  | 0.747% |
| 43 | 22.627 | 3352 | 3356 | 3360 | rBV2 | 384848  | 627524  | 12.18% | 1.152% |
| 44 | 22.674 | 3360 | 3364 | 3375 | rVB2 | 687198  | 1267567 | 24.61% | 2.327% |
| 45 | 23.274 | 3460 | 3466 | 3472 | rBV2 | 933856  | 1669851 | 32.42% | 3.066% |
| 46 | 23.827 | 3555 | 3560 | 3568 | rVB  | 357052  | 697425  | 13.54% | 1.280% |
| 47 | 23.910 | 3570 | 3574 | 3584 | rVB2 | 283269  | 624692  | 12.13% | 1.147% |

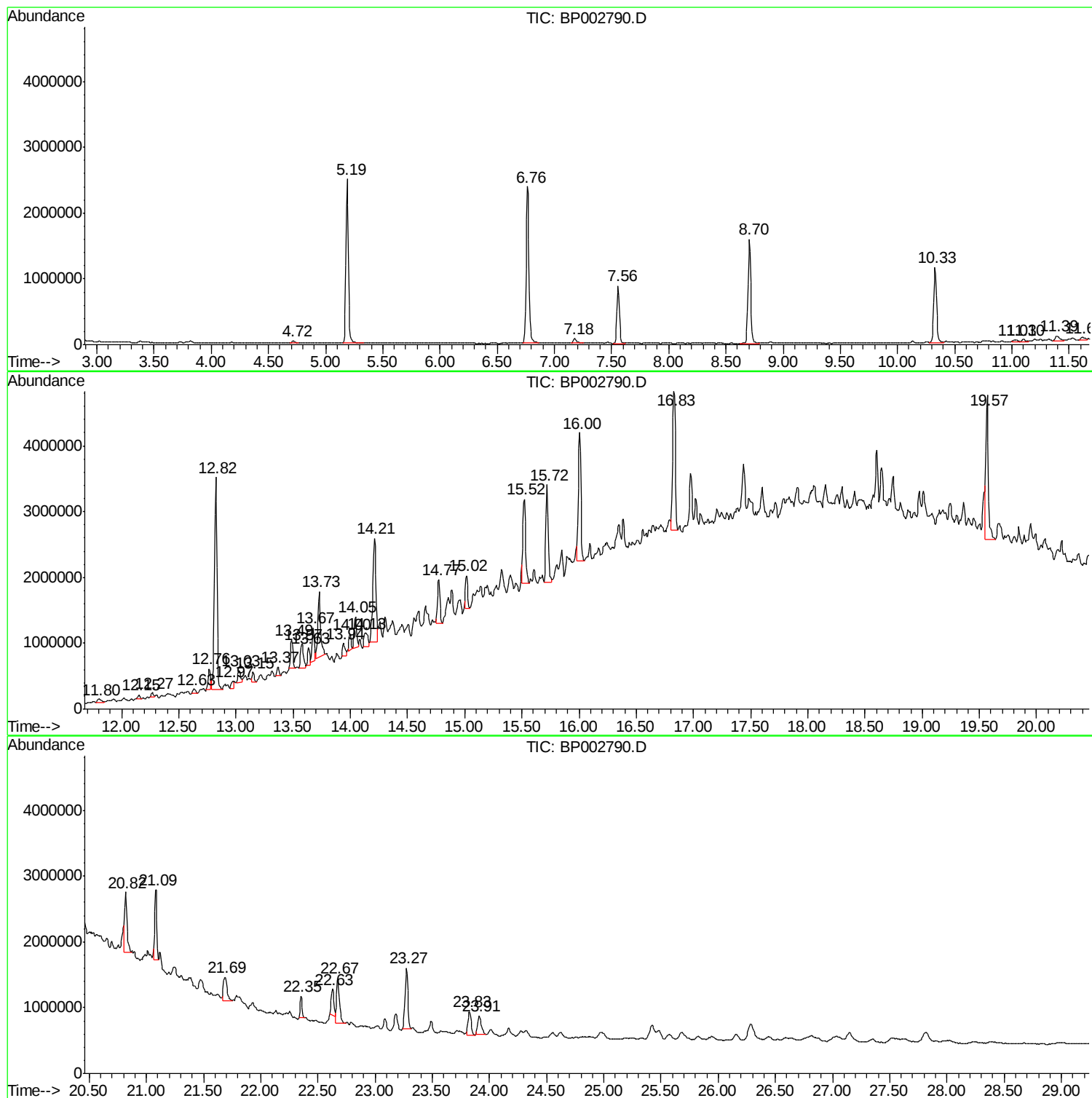
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ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
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Instrument :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

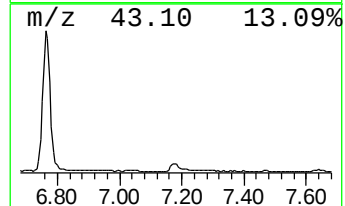
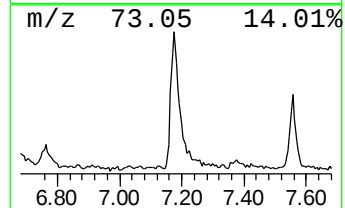
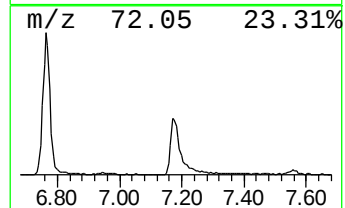
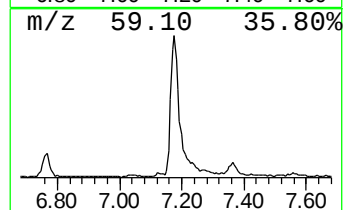
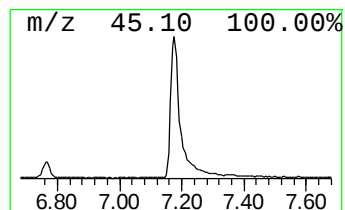
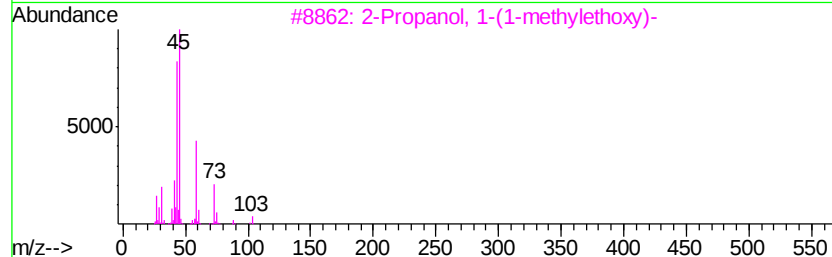
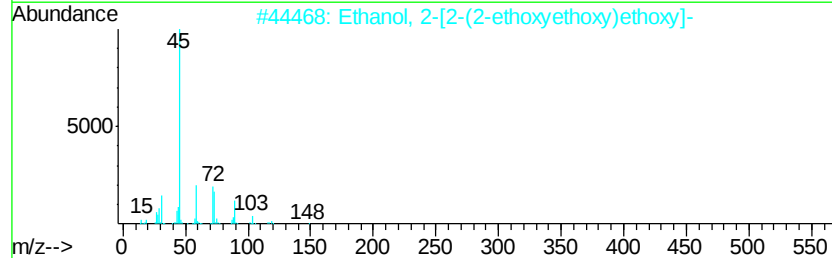
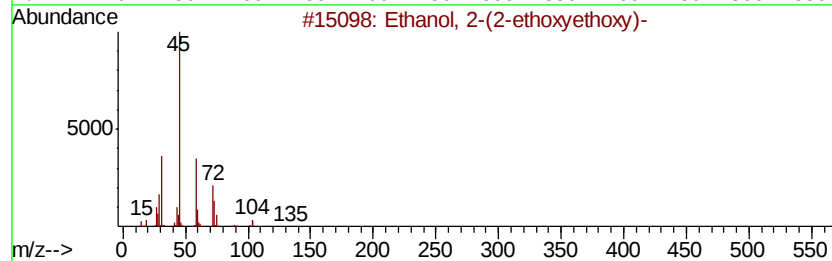
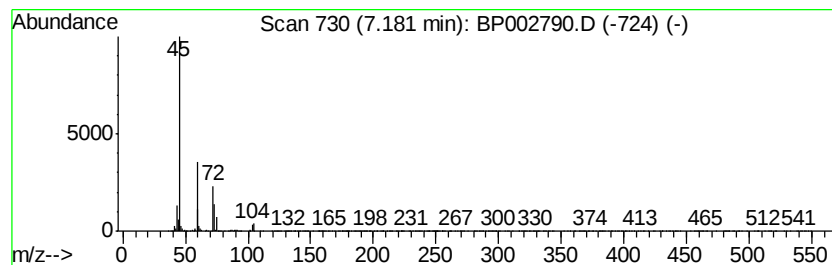
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.18 | 2.31 ng | 160181 | 1,4-Dichlorobenzene-d4 | 7.56 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Ethanol, 2-(2-ethoxyethoxy)-          | 134 | C6H14O3 | 000111-90-0 | 90   |
| 2    |    | Ethanol, 2-[2-(2-ethoxyethoxy)eth...] | 178 | C8H18O4 | 000112-50-5 | 56   |
| 3    |    | 2-Propanol, 1-(1-methoxyethoxy)-      | 118 | C6H14O2 | 003944-36-3 | 50   |
| 4    |    | Ethanol, 2-(2-methoxyethoxy)-         | 120 | C5H12O3 | 000111-77-3 | 45   |
| 5    |    | Ethanol, 2-[2-(2-propenyloxy)eth...]  | 146 | C7H14O3 | 015075-50-0 | 40   |



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10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

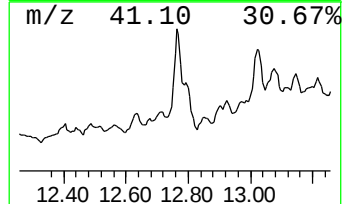
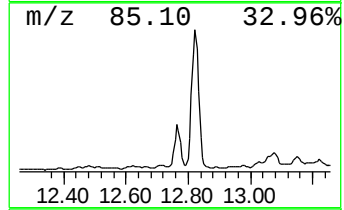
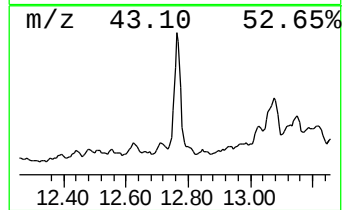
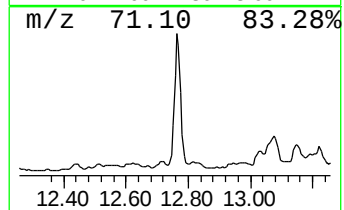
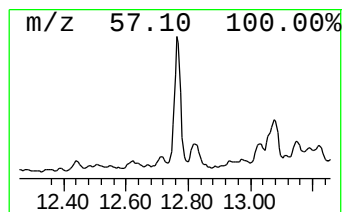
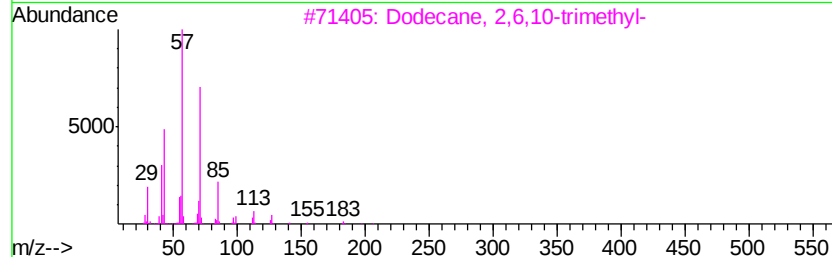
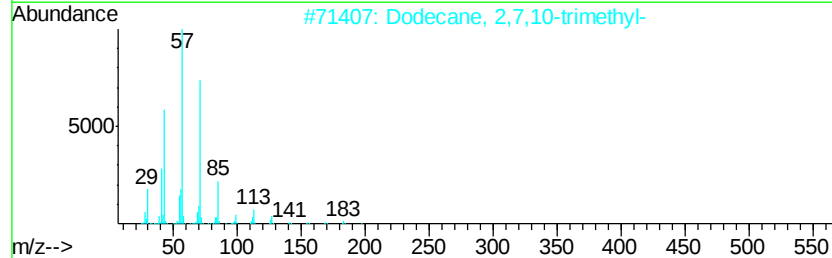
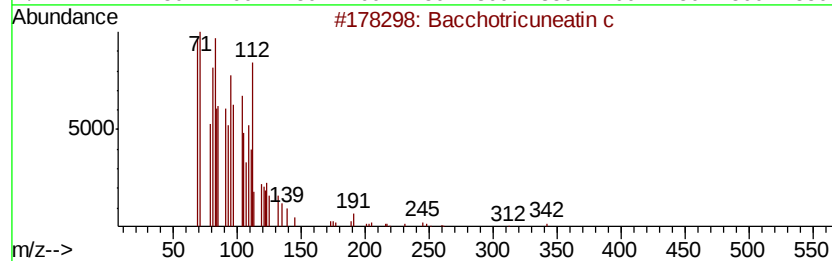
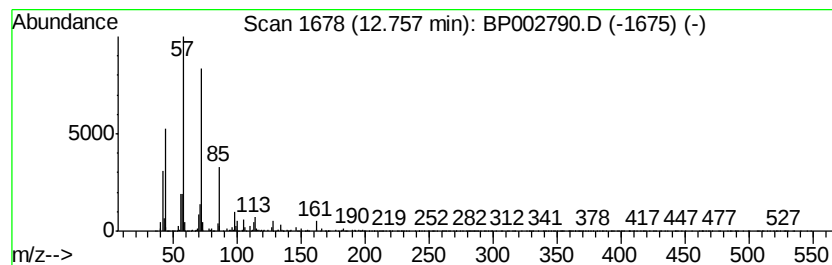
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Bacchotricuneatin c Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.76 | 3.19 ng | 474232 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------------|-----|----------|-------------|------|
| 1    | 5  | Bacchotricuneatin c         | 342 | C20H22O5 | 066563-30-2 | 96   |
| 2    |    | Dodecane, 2,7,10-trimethyl- | 212 | C15H32   | 074645-98-0 | 91   |
| 3    |    | Dodecane, 2,6,10-trimethyl- | 212 | C15H32   | 003891-98-3 | 90   |
| 4    |    | Dodecane, 2,6,11-trimethyl- | 212 | C15H32   | 031295-56-4 | 78   |
| 5    |    | Heptadecane                 | 240 | C17H36   | 000629-78-7 | 72   |



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Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

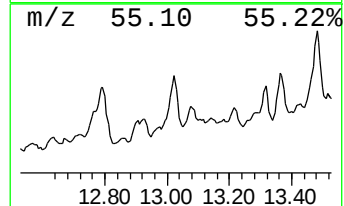
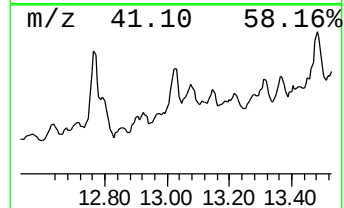
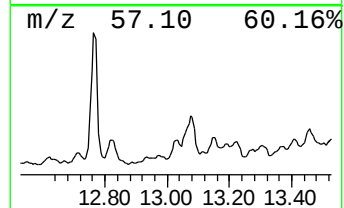
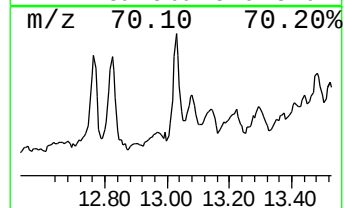
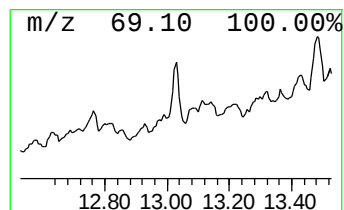
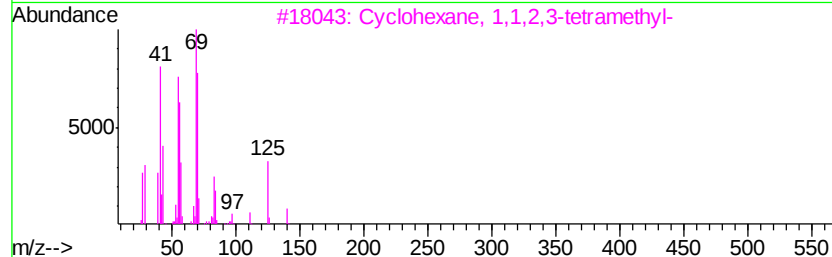
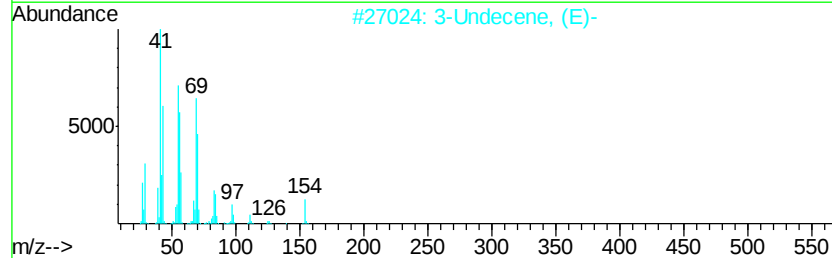
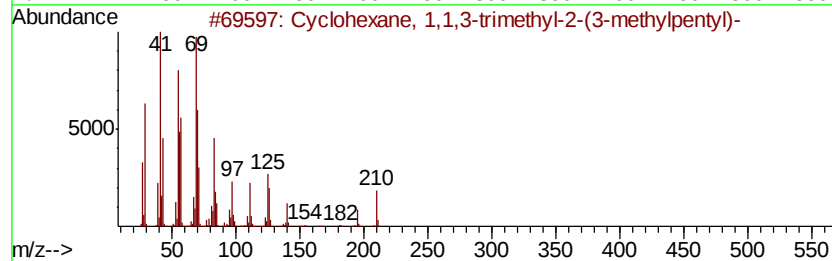
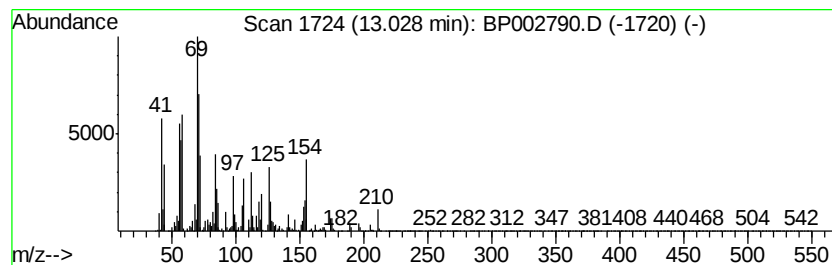
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Cyclohexane, 1,1,3-trimethy... Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.03 | 2.13 ng | 316397 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, 1,1,3-trimethyl-2-(... | 210 | C15H30  | 054965-05-8 | 74   |
| 2    |    | 3-Undecene, (E)-                    | 154 | C11H22  | 001002-68-2 | 64   |
| 3    |    | Cyclohexane, 1,1,2,3-tetramethyl-   | 140 | C10H20  | 006783-92-2 | 58   |
| 4    |    | 5-Undecene                          | 154 | C11H22  | 004941-53-1 | 58   |
| 5    |    | 5-Undecene, (E)-                    | 154 | C11H22  | 000764-97-6 | 43   |



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ALS Vial : 7 Sample Multiplier: 1

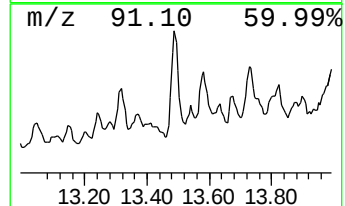
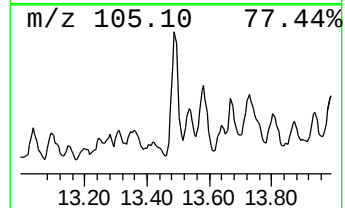
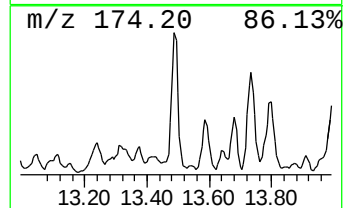
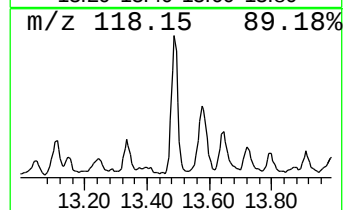
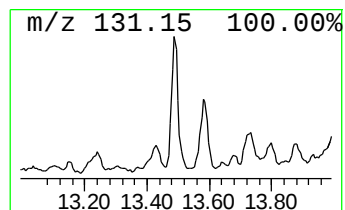
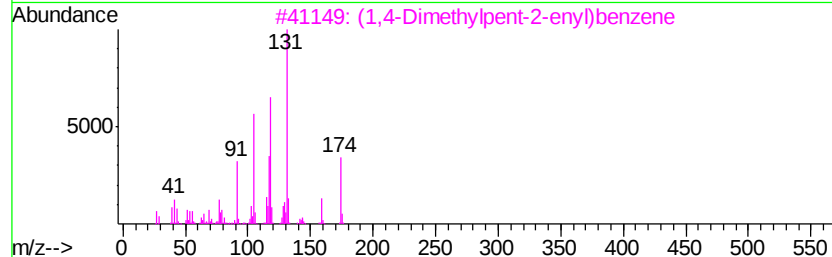
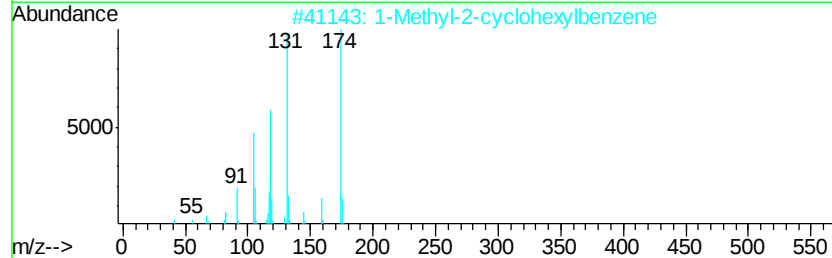
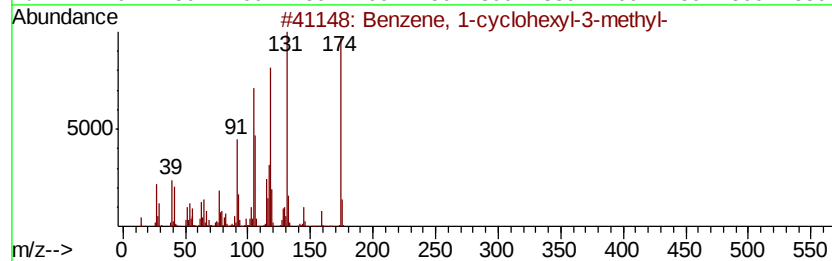
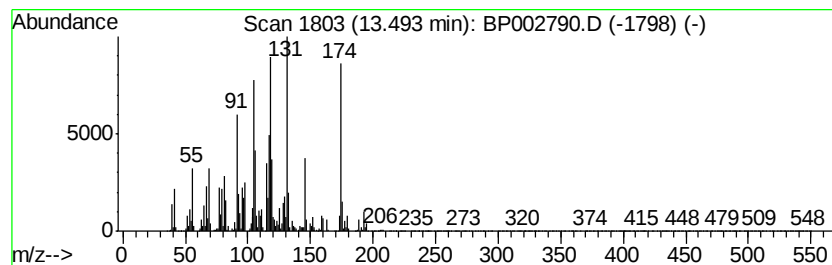
Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Benzene, 1-cyclohexyl-3-met... Concentration Rank 14

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 13.49   | 4.20 ng | 623964                              | Acenaphthene-d10 | 14.21           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | Benzene, 1-cvclohexvl-3-methyl-     | 174 C13H18       | 004575-46-6 96  |
| 2       |         | 1-Methyl-2-cyclohexylbenzene        | 174 C13H18       | 004501-35-3 86  |
| 3       |         | (1,4-Dimethylpent-2-enyl)benzene    | 174 C13H18       | 1000184-97-9 64 |
| 4       |         | 1H-1,3,2-Benzodiazaborole, 2-but... | 174 C10H15BN2    | 031748-14-8 46  |
| 5       |         | 2-Pentanone, 3-(phenylmethylene)-   | 174 C12H14O      | 003437-89-6 22  |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

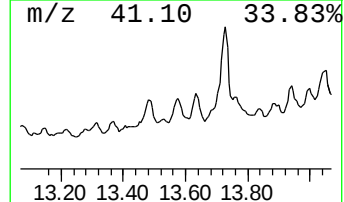
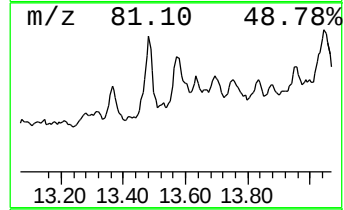
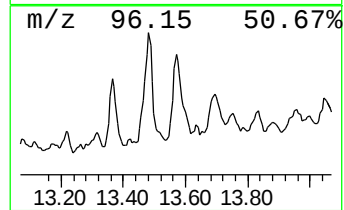
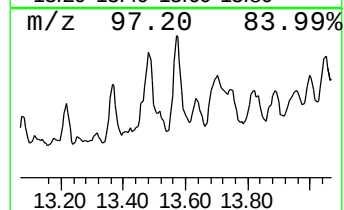
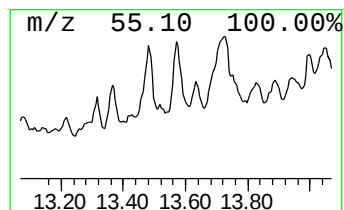
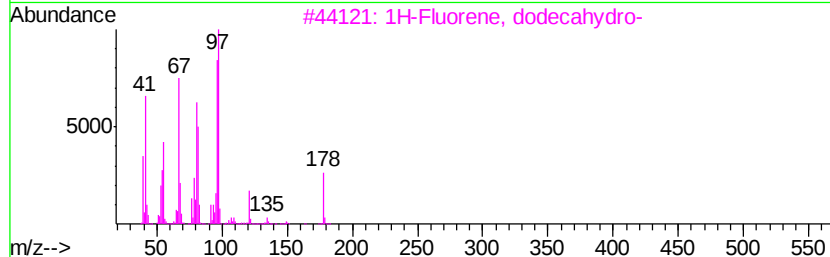
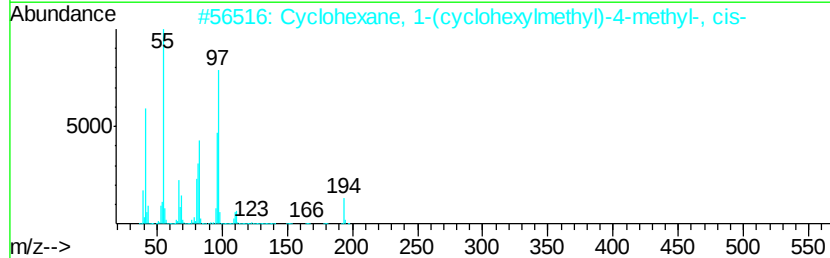
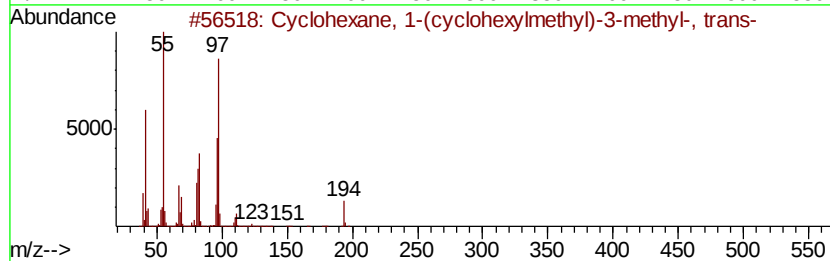
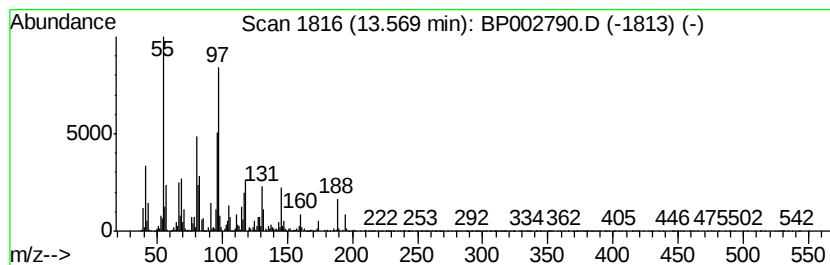
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown13.57 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.57 | 4.41 ng | 654462 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|--------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Cyclohexane, 1-(cyclohexylmethyl)... | 194 | C14H26   | 054823-95-9  | 46   |
| 2    |    | Cyclohexane, 1-(cyclohexylmethyl)... | 194 | C14H26   | 054823-97-1  | 46   |
| 3    |    | 1H-Fluorene, dodecahydro-            | 178 | C13H22   | 005744-03-6  | 38   |
| 4    |    | Oxalic acid, di(cyclohexylmethyl)... | 282 | C16H26O4 | 1000309-68-5 | 35   |
| 5    |    | Cyclohexane, 1-methyl-4-(1-methy...  | 140 | C10H20   | 006069-98-3  | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

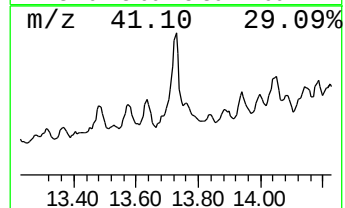
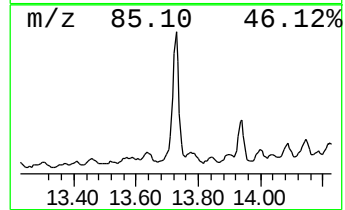
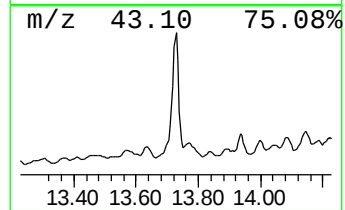
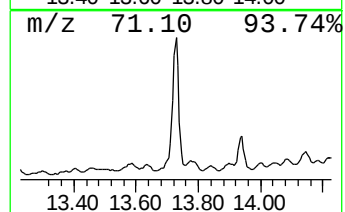
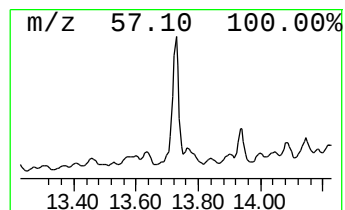
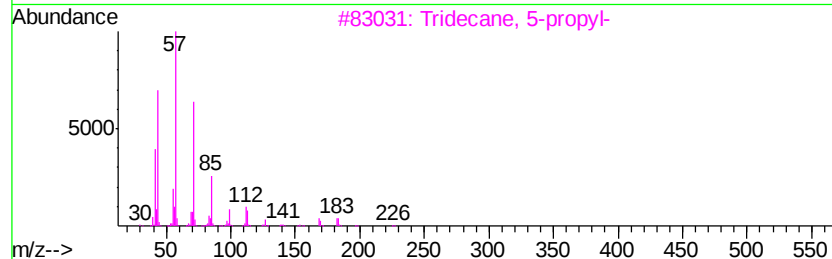
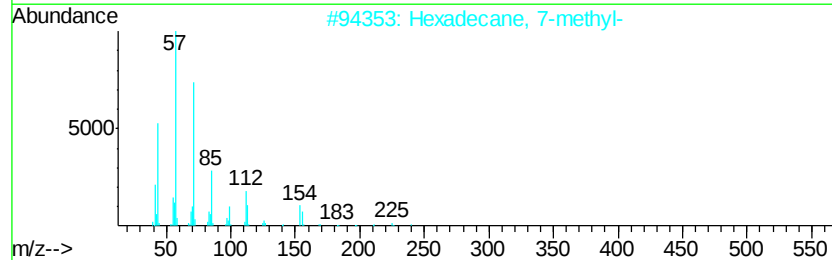
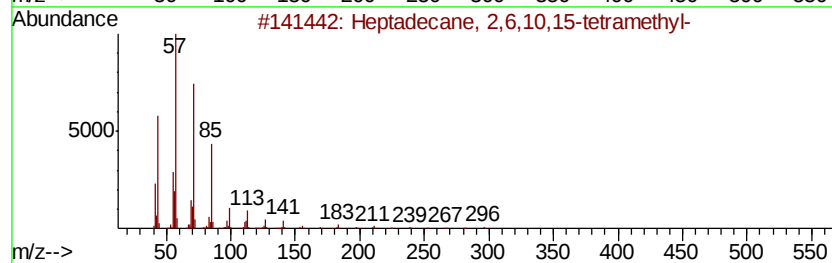
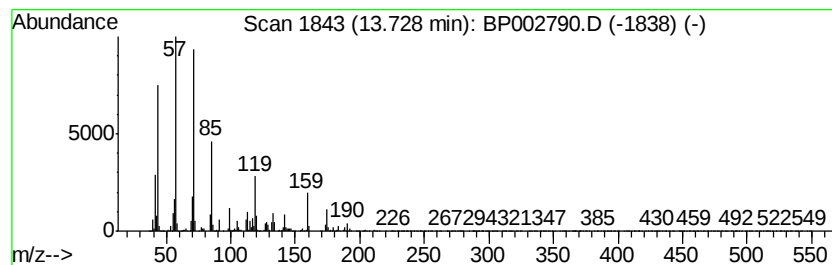
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Hexadecane, 7-methyl- Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 13.73 | 11.49 ng | 1707760 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44      | 054833-48-6 | 58   |
| 2    |    | Hexadecane, 7-methyl-               | 240 | C17H36      | 026730-20-1 | 58   |
| 3    |    | Tridecane, 5-propyl-                | 226 | C16H34      | 055045-11-9 | 53   |
| 4    |    | Silane, trichlorooctadecyl-         | 386 | C18H37Cl3Si | 000112-04-9 | 50   |
| 5    |    | Octane, 5-ethyl-2-methyl-           | 156 | C11H24      | 062016-18-6 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

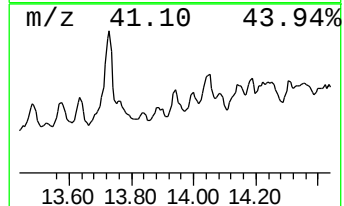
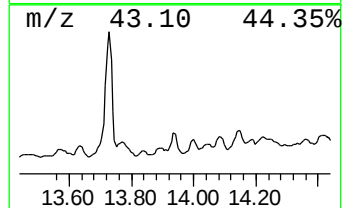
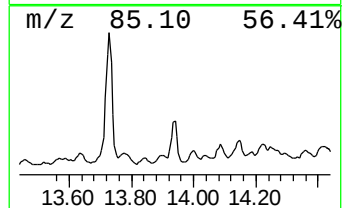
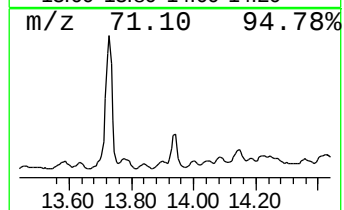
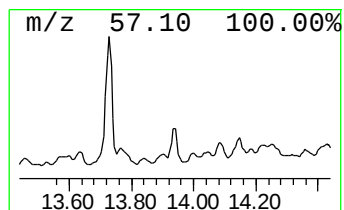
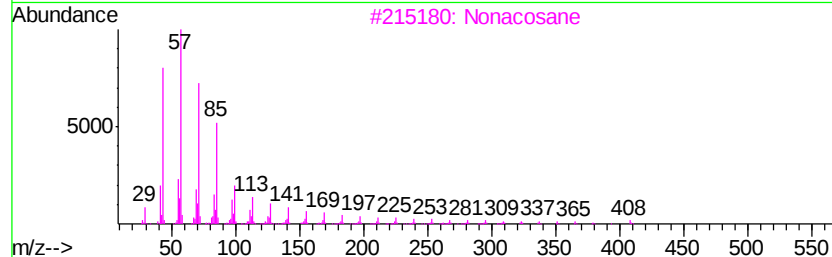
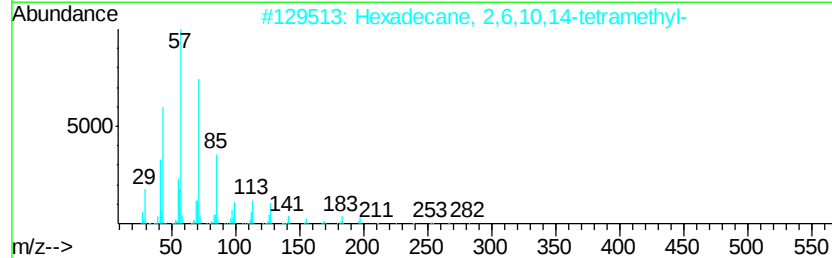
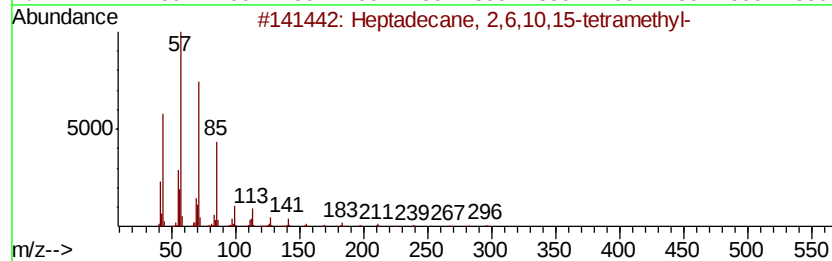
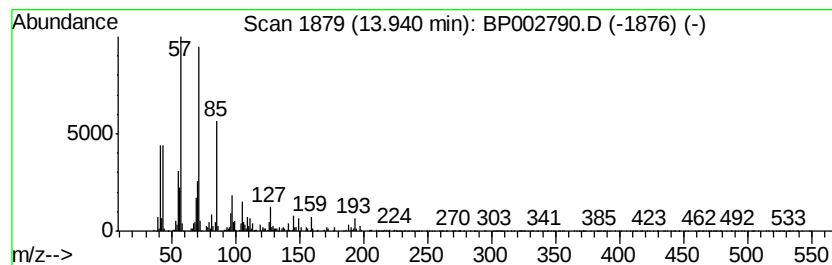
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Heptadecane, 2,6,10,15-tetr... Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.94 | 2.37 ng | 352195 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44    | 054833-48-6  | 64   |
| 2    |    | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42    | 000638-36-8  | 59   |
| 3    |    | Nonacosane                          | 408 | C29H60    | 000630-03-5  | 53   |
| 4    |    | Tetracontane                        | 563 | C40H82    | 004181-95-7  | 53   |
| 5    |    | Sulfurous acid, 2-ethylhexyl hep... | 432 | C25H52O3S | 1000309-20-0 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

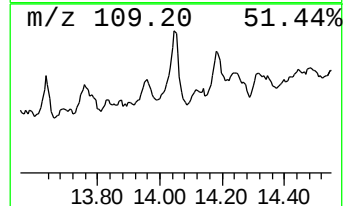
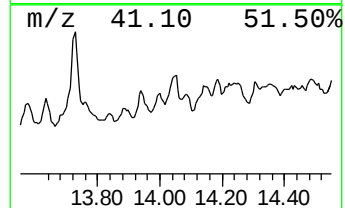
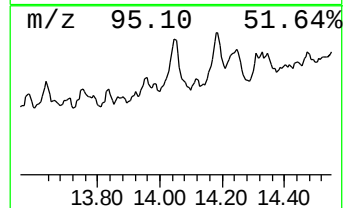
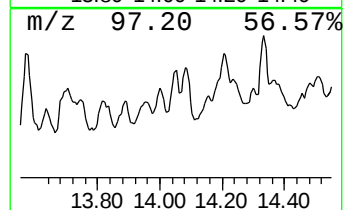
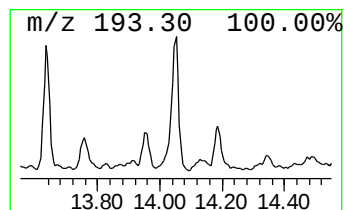
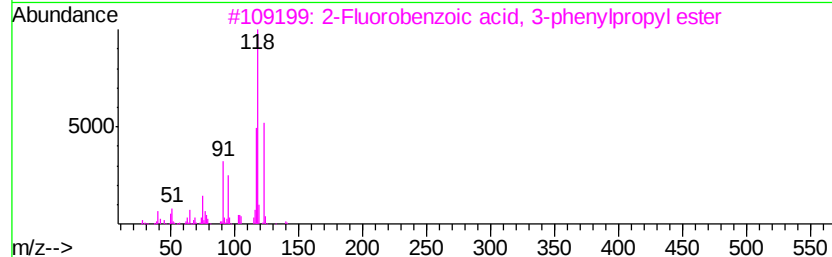
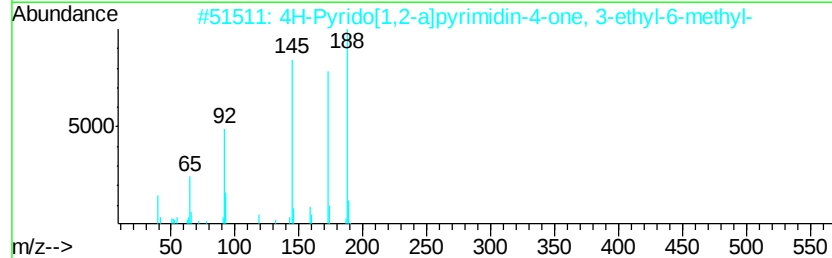
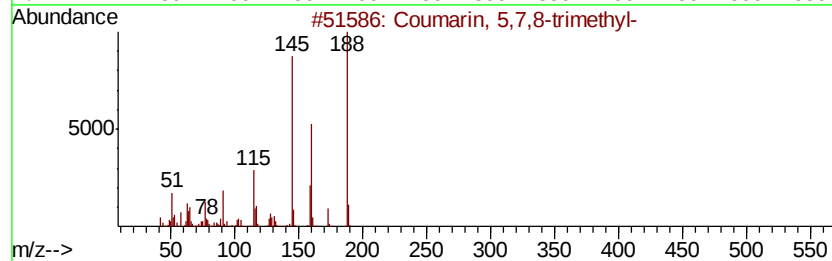
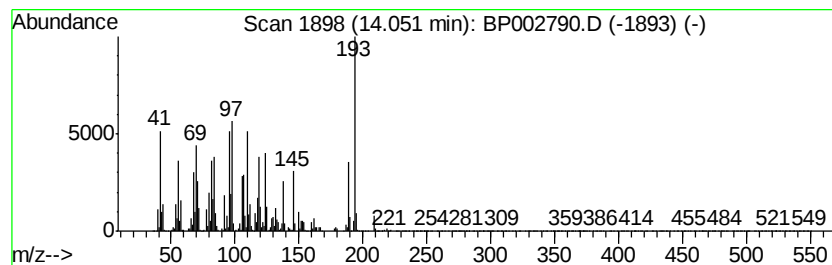
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 unknown14.05 Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.05 | 4.65 ng | 690167 | Acenaphthene-d10 | 14.21 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Coumarin, 5,7,8-trimethyl-          | 188 | C12H12O2  | 040662-19-9  | 25   |
| 2    |    |   | 4H-Pyrido[1,2-a]pyrimidin-4-one,... | 188 | C11H12N2O | 057773-19-0  | 15   |
| 3    |    |   | 2-Fluorobenzoic acid, 3-phenylpr... | 258 | C16H15FO2 | 1000278-98-4 | 14   |
| 4    |    |   | Carbonic acid, monoamide, N-meth... | 193 | C11H15NO2 | 1000340-19-0 | 11   |
| 5    |    |   | Decahydro-4,4,8,9,10-pentamethyl... | 208 | C15H28    | 080655-44-3  | 11   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

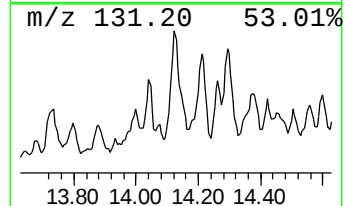
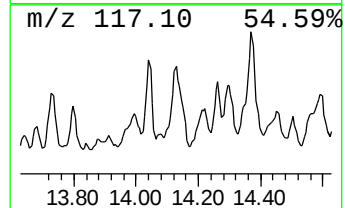
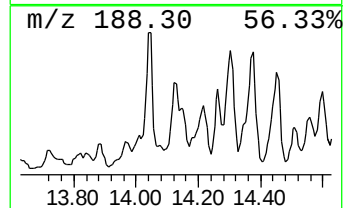
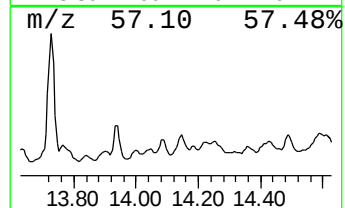
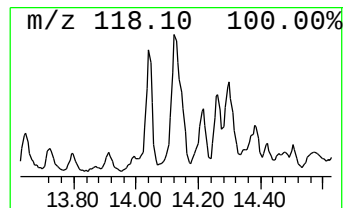
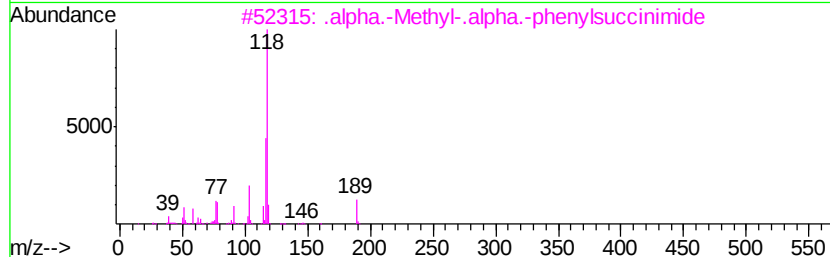
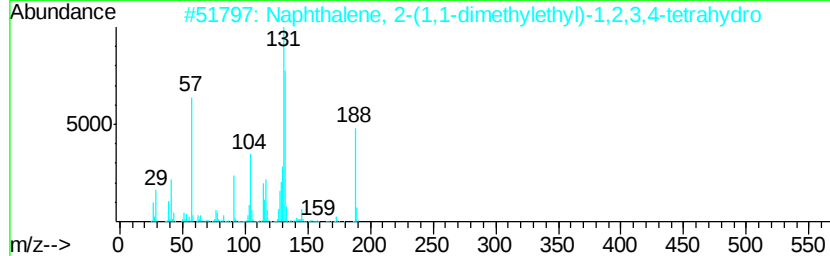
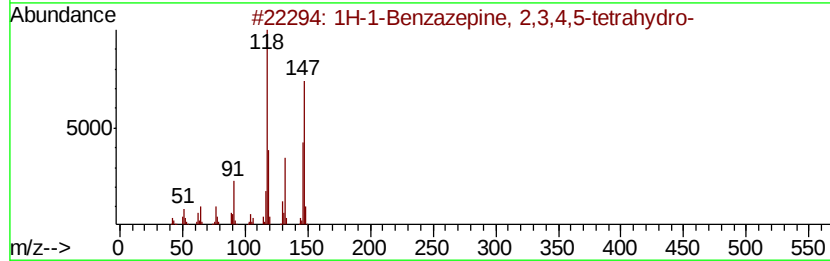
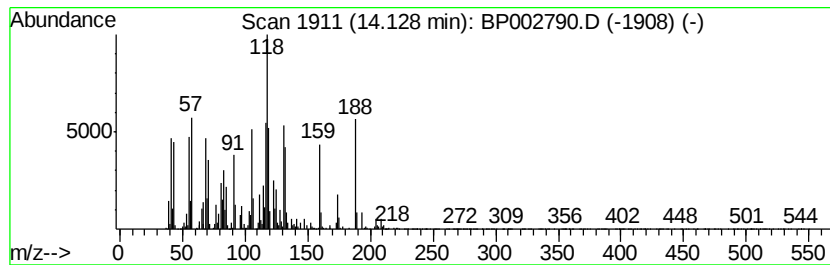
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown14.13 Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.13 | 3.54 ng | 526014 | Acenaphthene-d10 | 14.21 |

| Hit# | of | 5 | Tentative ID                             | MW  | MolForm    | CAS#         | Qual |
|------|----|---|------------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1H-1-Benzazepine, 2,3,4,5-tetrahydro-    | 147 | C10H13N    | 001701-57-1  | 35   |
| 2    |    |   | Naphthalene, 2-(1,1-dimethylethyl)-      | 188 | C14H20     | 042044-22-4  | 30   |
| 3    |    |   | .alpha.-Methyl-.alpha.-phenylsuccinimide | 189 | C11H11NO2  | 001497-17-2  | 25   |
| 4    |    |   | 4-Trifluoromethylbenzoic acid, 3-methyl- | 308 | C17H15F3O2 | 150272-46-1  | 22   |
| 5    |    |   | Dimethylmalonic acid, monochloride       | 268 | C14H17ClO3 | 1000361-84-8 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
 Data File : BP002790.D  
 Acq On : 17 Jul 2020 14:03  
 Operator : CG/JU  
 Sample : L3324-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

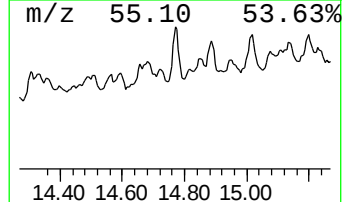
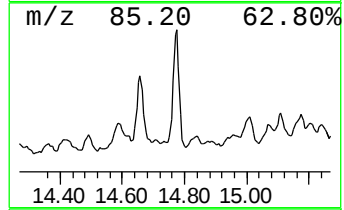
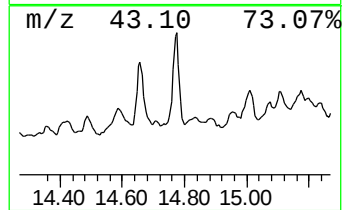
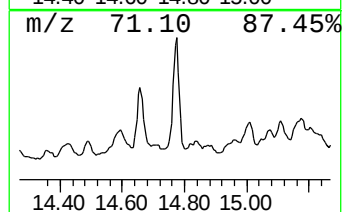
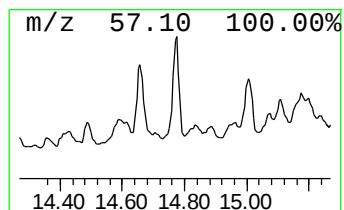
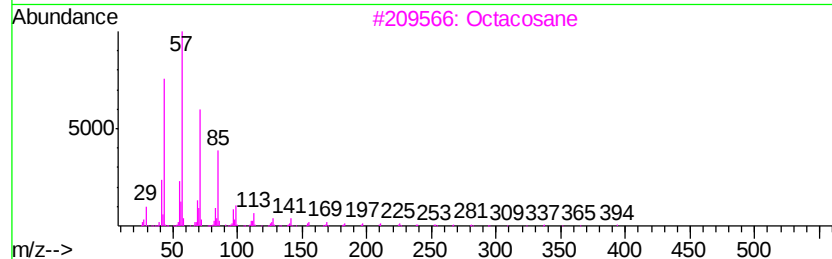
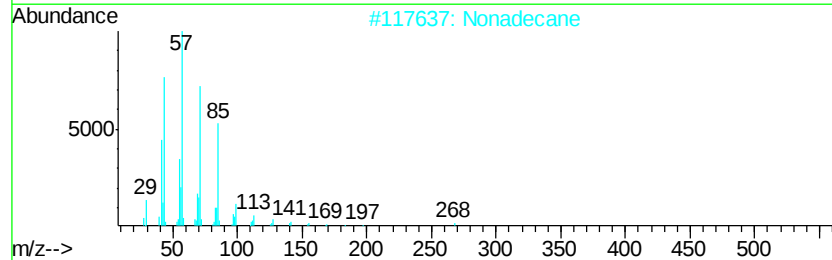
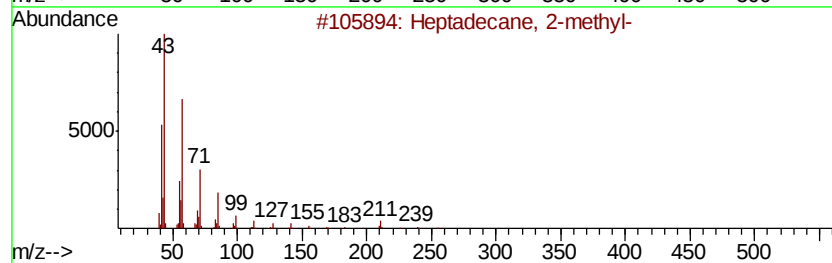
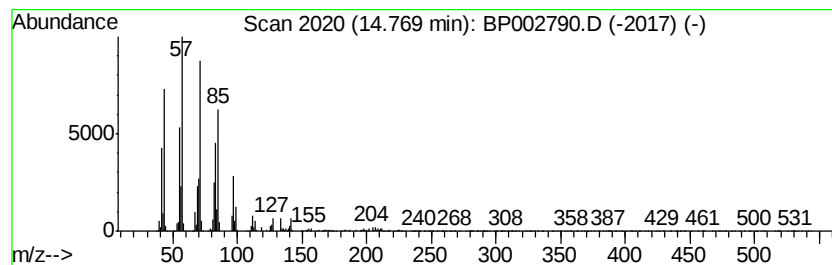
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 10 Heptadecane, 2-methyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.77 | 6.29 ng | 934240 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 2-methyl-   | 254 | C18H38  | 001560-89-0 | 60   |
| 2    |    | Nonadecane               | 268 | C19H40  | 000629-92-5 | 58   |
| 3    |    | Octacosane               | 394 | C28H58  | 000630-02-4 | 58   |
| 4    |    | Tetracosane              | 338 | C24H50  | 000646-31-1 | 53   |
| 5    |    | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 52   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

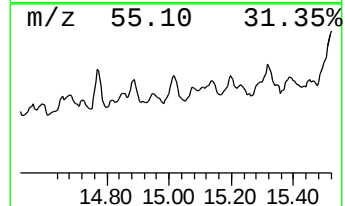
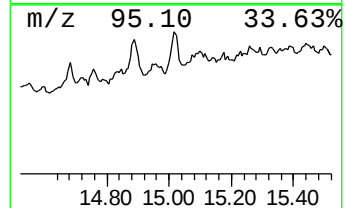
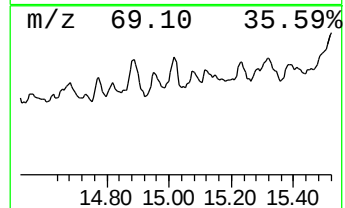
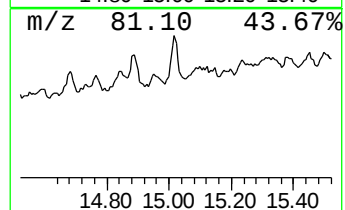
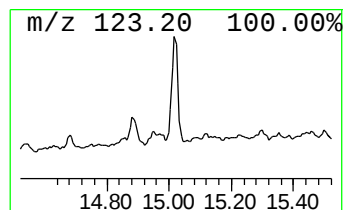
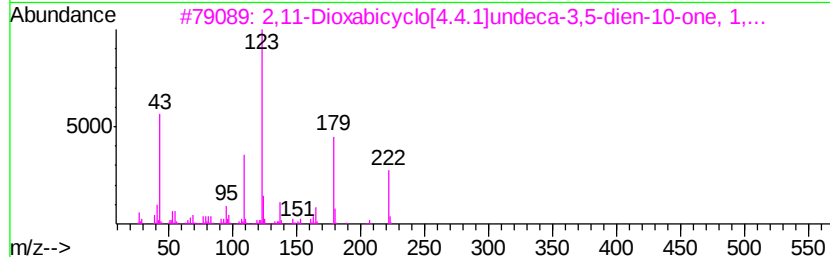
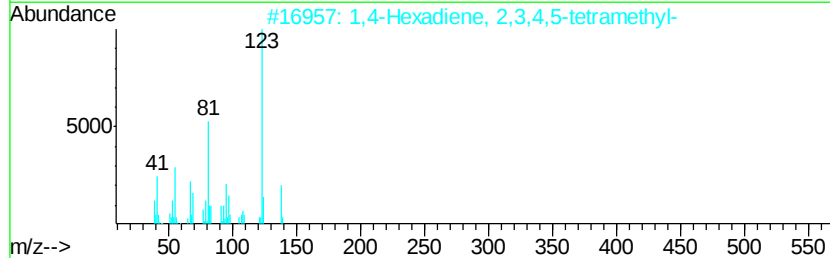
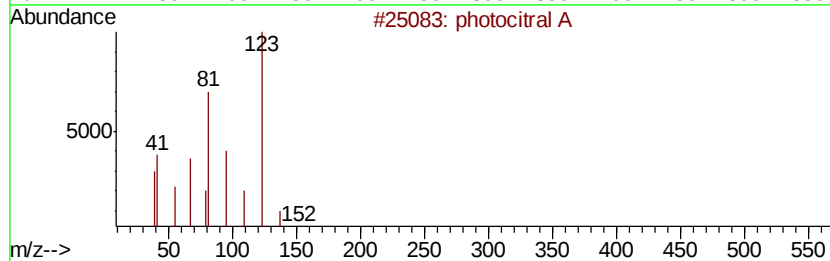
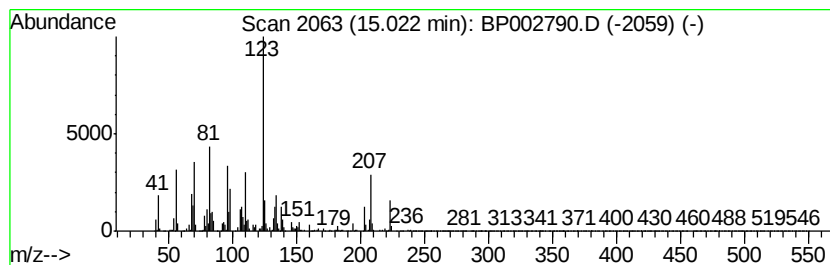
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown15.02 Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.02 | 4.19 ng | 622843 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | photocitral A                       | 152 | C10H16O   | 055253-28-6  | 49   |
| 2    |    | 1,4-Hexadiene, 2,3,4,5-tetramethyl- | 138 | C10H18    | 051504-54-2  | 47   |
| 3    |    | 2,11-Dioxabicyclo[4.4.1]undeca-3... | 222 | C13H18O3  | 070412-52-1  | 47   |
| 4    |    | 4-Fluorobenzoic acid, 3-pentadec... | 350 | C22H35FO2 | 1000280-68-4 | 38   |
| 5    |    | 1-Trimethylsilylpent-1-en-4-yne     | 138 | C8H14Si   | 079516-25-9  | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

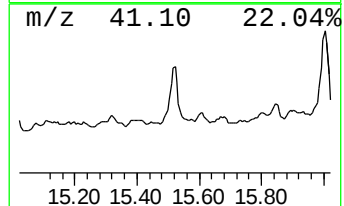
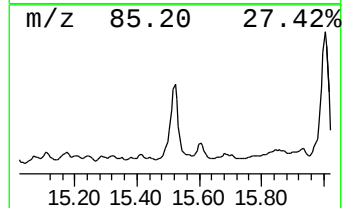
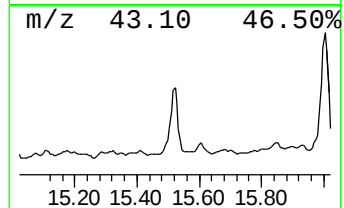
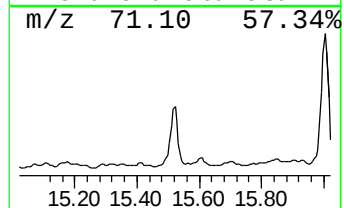
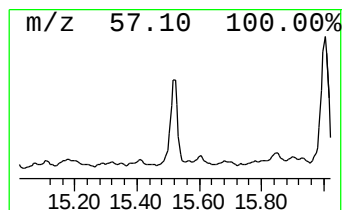
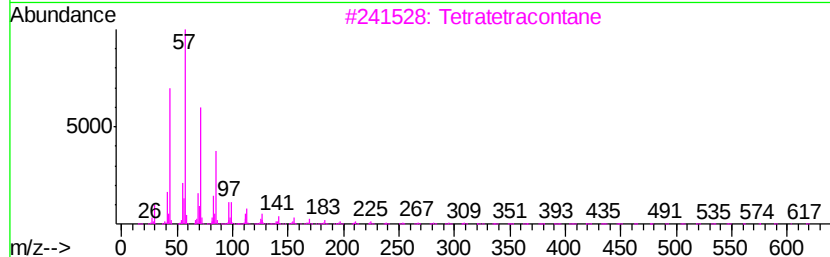
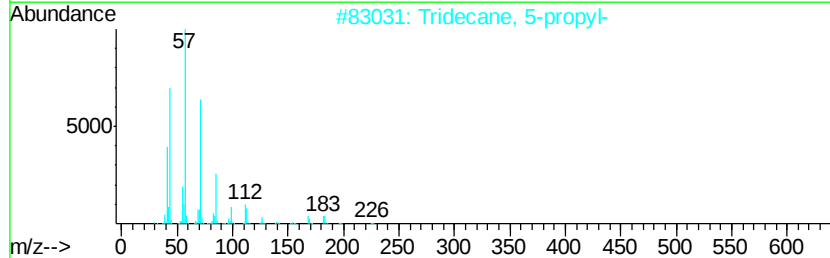
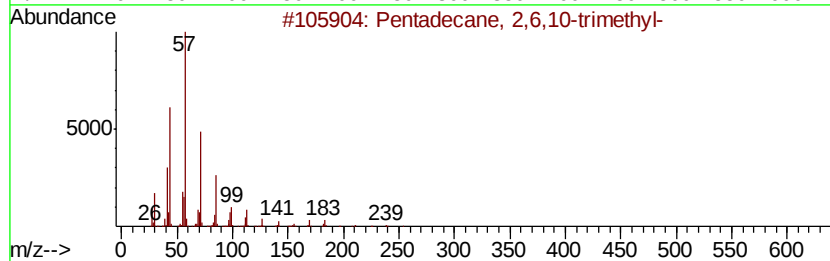
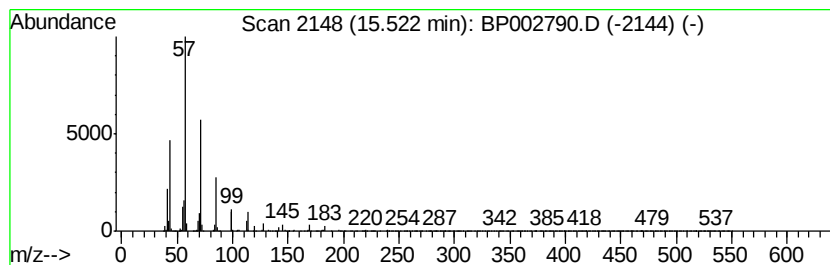
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Pentadecane, 2,6,10-trimethyl- Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 15.52 | 13.61 ng | 2022470 | Acenaphthene-d10 | 14.21 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38  | 003892-00-0 | 93   |
| 2    |    | Tridecane, 5-propyl-           | 226 | C16H34  | 055045-11-9 | 87   |
| 3    |    | Tetratetracontane              | 619 | C44H90  | 007098-22-8 | 64   |
| 4    |    | Nonane, 2-methyl-5-propyl-     | 184 | C13H28  | 031081-17-1 | 59   |
| 5    |    | Dodecane, 2,7,10-trimethyl-    | 212 | C15H32  | 074645-98-0 | 59   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

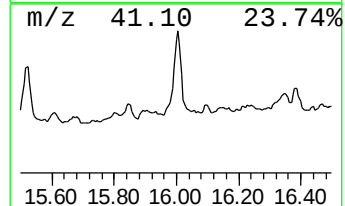
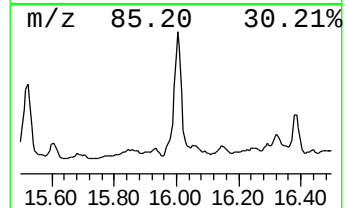
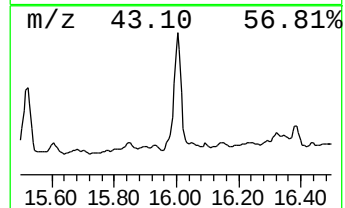
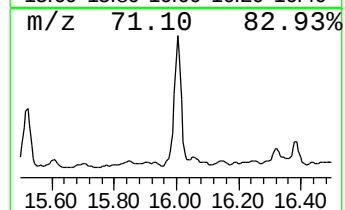
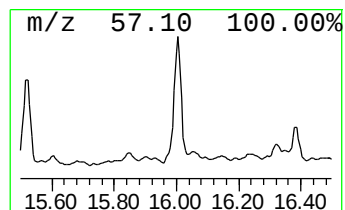
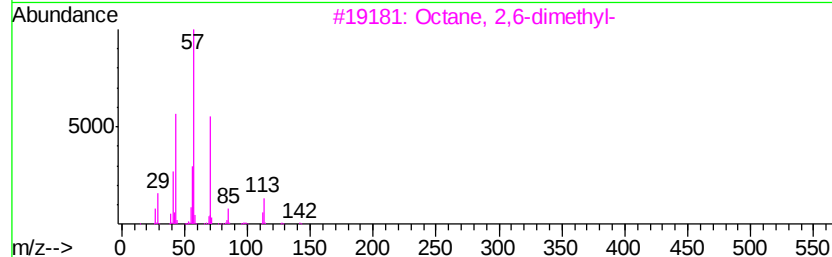
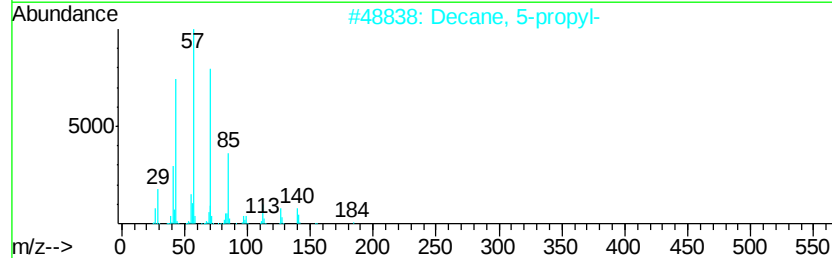
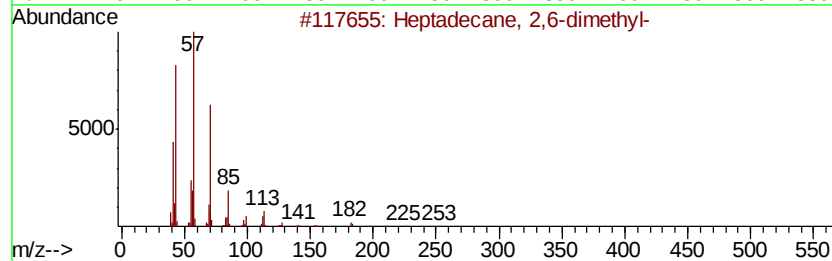
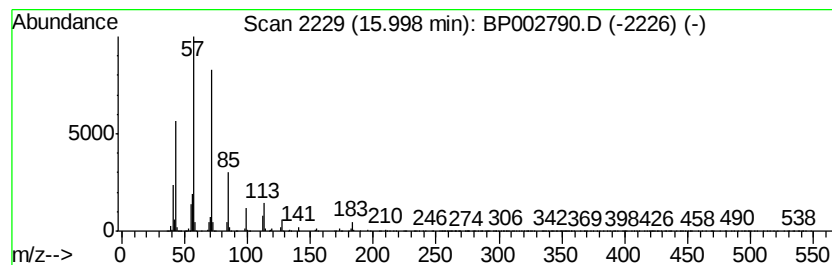
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Heptadecane, 2,6-dimethyl- Concentration Rank 3

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 16.00 | 17.16 ng | 2950120 | Phenanthrene-d10 | 16.97 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptadecane, 2,6-dimethyl- | 268 | C19H40  | 054105-67-8 | 86   |
| 2    |    | Decane, 5-propyl-          | 184 | C13H28  | 017312-62-8 | 80   |
| 3    |    | Octane, 2,6-dimethyl-      | 142 | C10H22  | 002051-30-1 | 72   |
| 4    |    | Undecane, 6-ethyl-         | 184 | C13H28  | 017312-60-6 | 72   |
| 5    |    | Tridecane, 5-propyl-       | 226 | C16H34  | 055045-11-9 | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

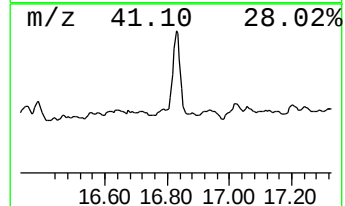
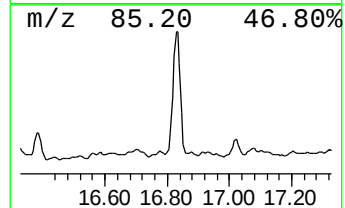
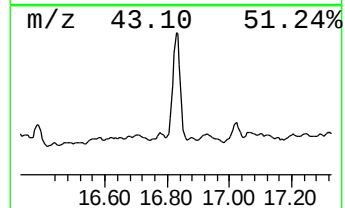
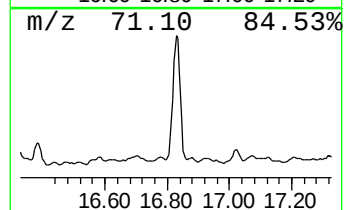
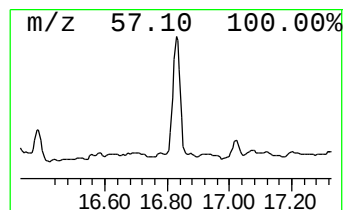
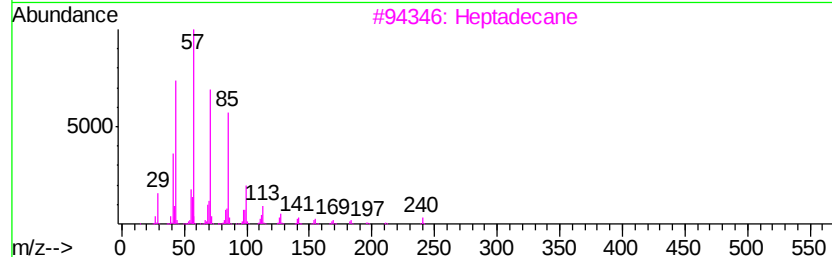
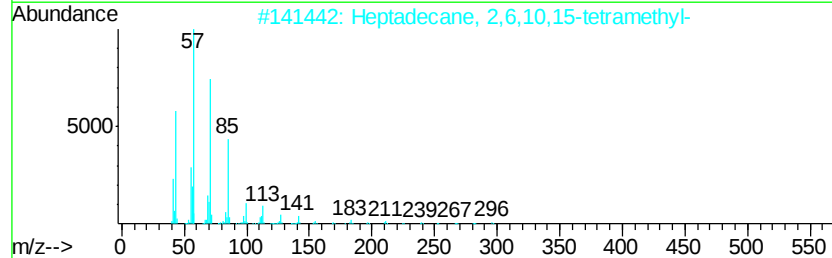
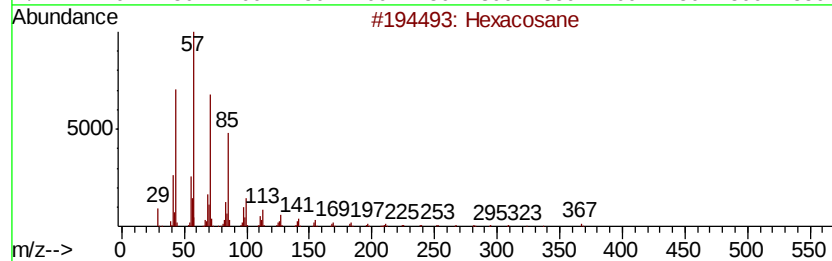
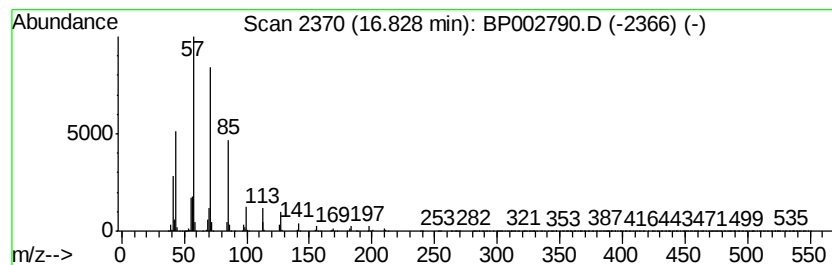
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Hexacosane Concentration Rank 1

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 16.83 | 20.00 ng | 3438880 | Phenanthrene-d10 | 16.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane                          | 366 | C26H54  | 000630-01-3 | 90   |
| 2    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 90   |
| 3    |    | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 87   |
| 4    |    | Octacosane                          | 394 | C28H58  | 000630-02-4 | 86   |
| 5    |    | Tritetracontane                     | 605 | C43H88  | 007098-21-7 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

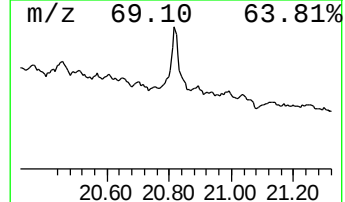
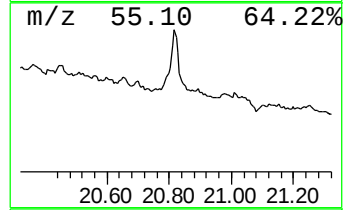
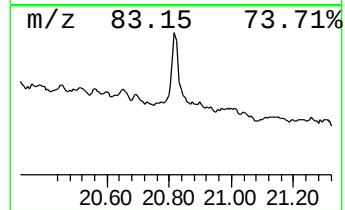
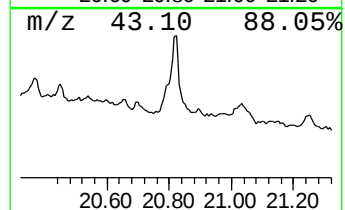
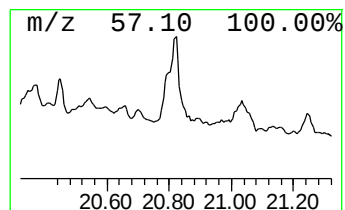
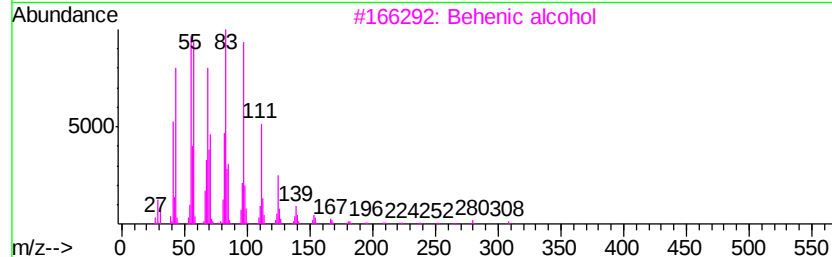
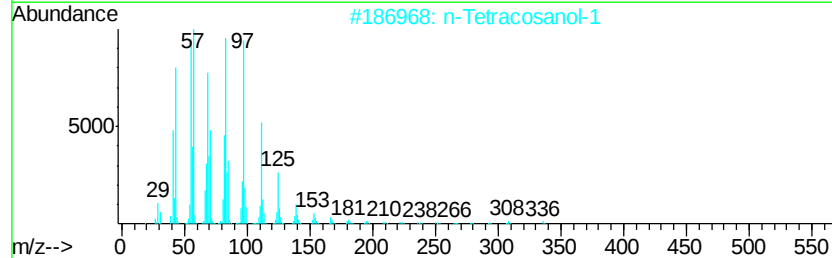
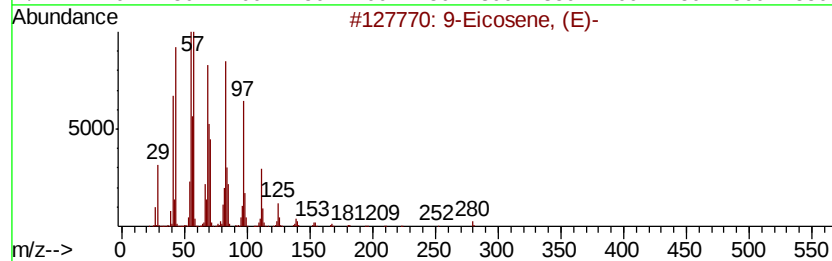
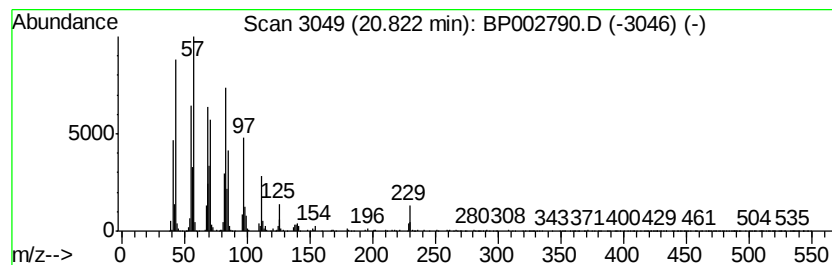
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 9-Eicosene, (E)- Concentration Rank 2

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 20.82 | 18.89 ng | 1332710 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | 9-Eicosene, (E)-                    | 280 | C20H40      | 074685-29-3 | 93   |
| 2    |    | n-Tetracosanol-1                    | 354 | C24H50O     | 000506-51-4 | 91   |
| 3    |    | Behenic alcohol                     | 326 | C22H46O     | 000661-19-8 | 91   |
| 4    |    | Trichloroacetic acid, tetradecyl... | 358 | C16H29Cl3O2 | 074339-52-9 | 91   |
| 5    |    | 17-Pentatriacontene                 | 491 | C35H70      | 006971-40-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

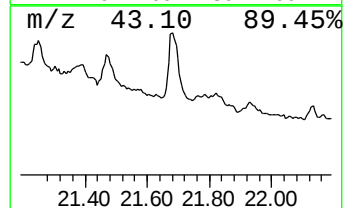
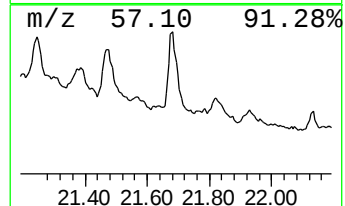
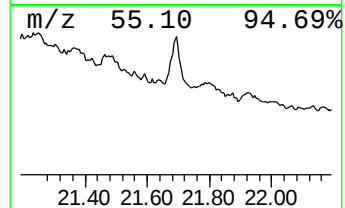
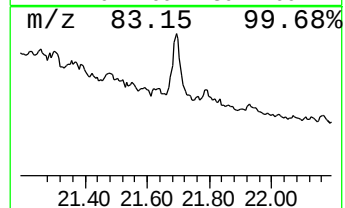
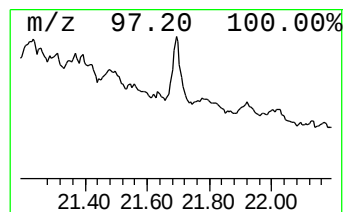
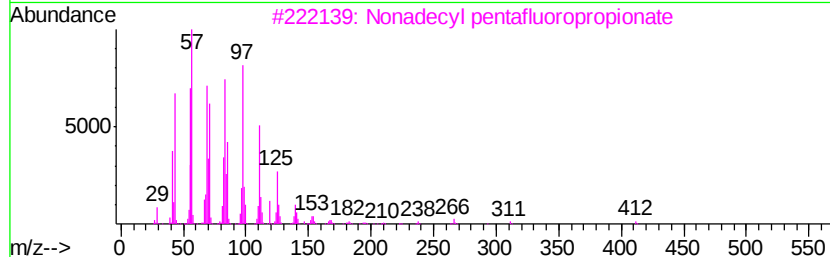
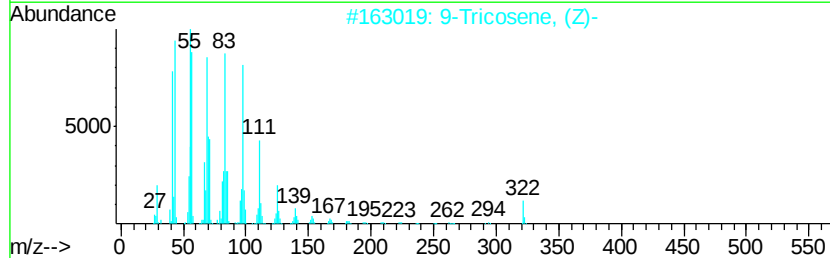
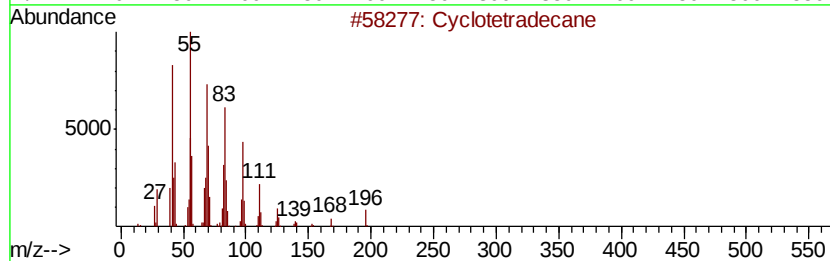
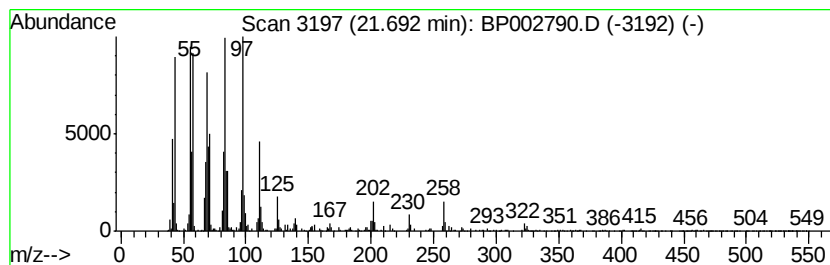
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Cyclotetradecane Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 21.69 | 11.47 ng | 809576 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Cyclotetradecane                    | 196 | C14H28      | 000295-17-0  | 95   |
| 2    |    | 9-Tricosene, (Z)-                   | 322 | C23H46      | 027519-02-4  | 78   |
| 3    |    | Nonadecyl pentafluoropropionate     | 430 | C22H39F5O2  | 1000351-88-8 | 76   |
| 4    |    | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 70   |
| 5    |    | 11-Tricosene                        | 322 | C23H46      | 052078-56-5  | 70   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

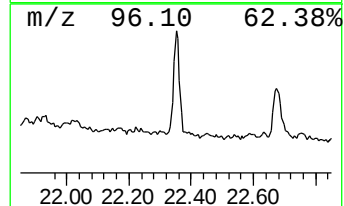
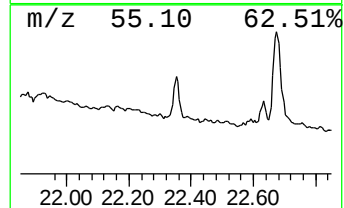
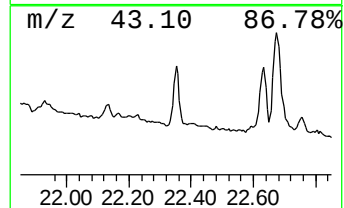
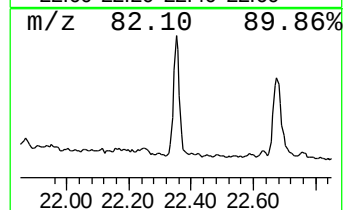
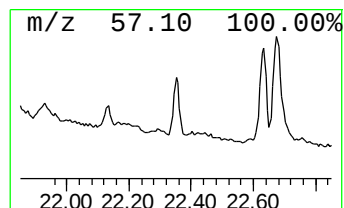
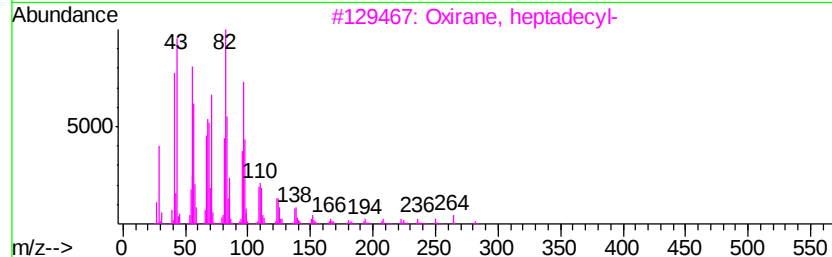
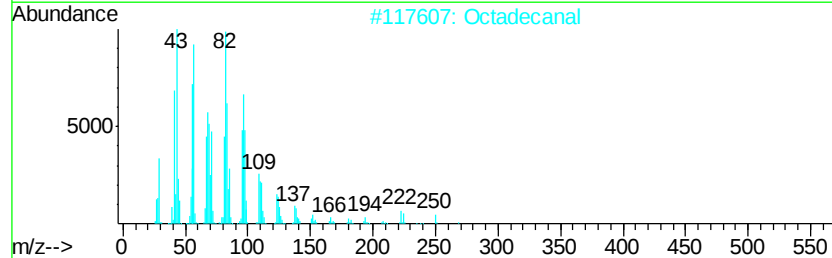
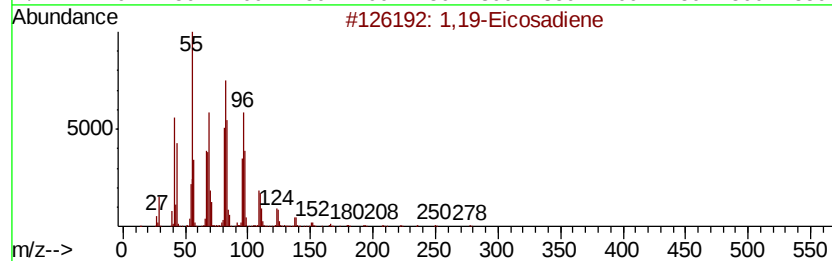
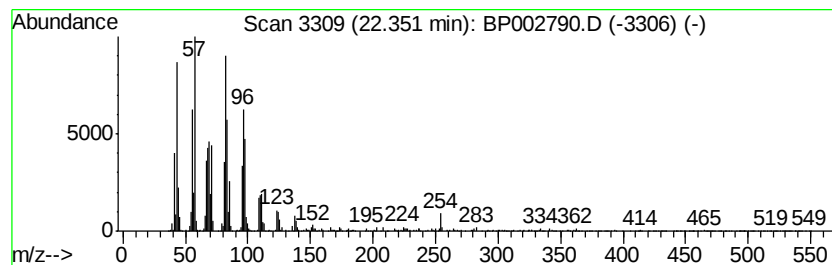
Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 1,19-Eicosadiene Concentration Rank 11

| R.T.      | EstConc                   | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------|--------|------------------|-------------|------|
| 22.35     | 4.87 ng                   | 406645 | Perylene-d12     | 23.27       |      |
| Hit# of 5 | Tentative ID              | MW     | MolForm          | CAS#        | Qual |
| 1         | 1,19-Eicosadiene          | 278    | C20H38           | 014811-95-1 | 95   |
| 2         | Octadecanal               | 268    | C18H36O          | 000638-66-4 | 91   |
| 3         | Oxirane, heptadecyl-      | 282    | C19H38O          | 067860-04-2 | 90   |
| 4         | (Z)-14-Tricosenyl formate | 366    | C24H46O2         | 077899-10-6 | 90   |
| 5         | 9-Octadecen-1-ol, (Z)-    | 268    | C18H36O          | 000143-28-2 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

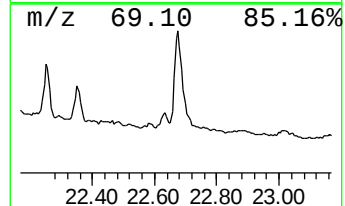
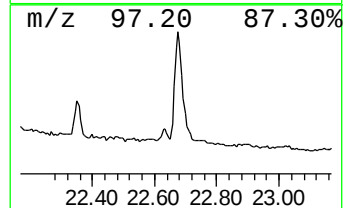
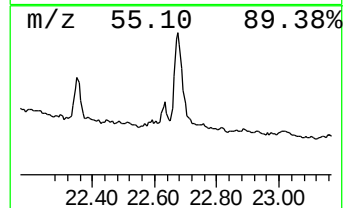
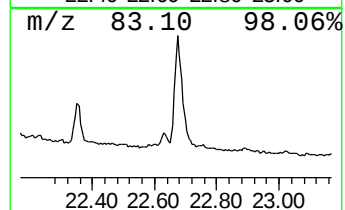
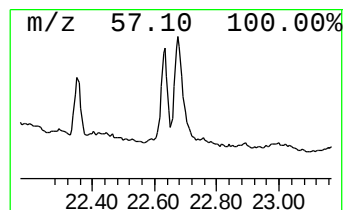
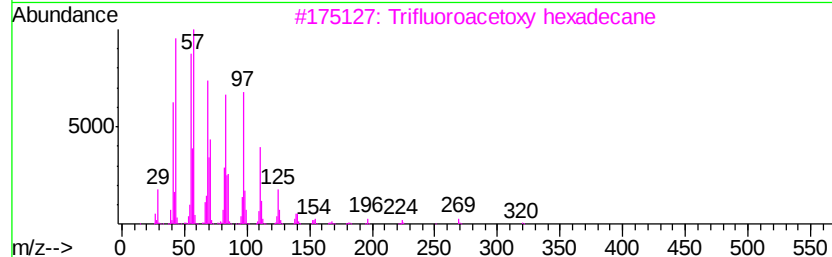
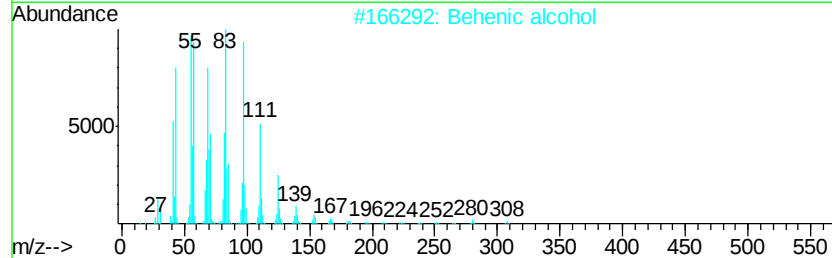
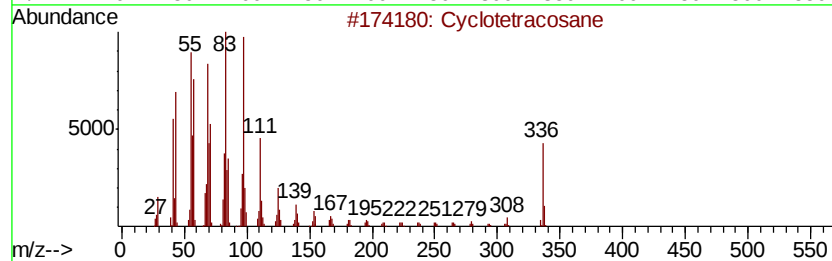
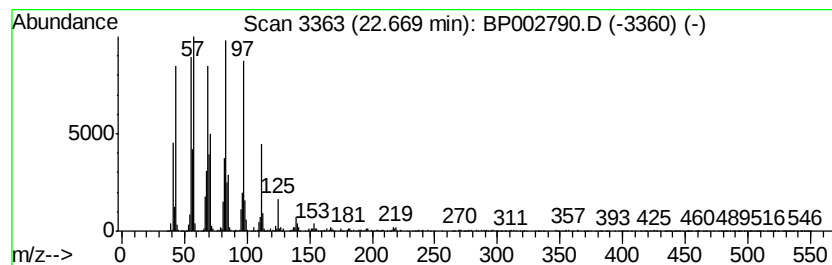
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Cyclotetracosane Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 22.67 | 15.18 ng | 1267570 | Perylene-d12     | 23.27 |

| Hit# | of | Tentative ID                | MW  | MolForm    | CAS#         | Qual |
|------|----|-----------------------------|-----|------------|--------------|------|
| 1    | 5  | Cyclotetracosane            | 336 | C24H48     | 000297-03-0  | 98   |
| 2    |    | Behenic alcohol             | 326 | C22H46O    | 000661-19-8  | 91   |
| 3    |    | Trifluoroacetoxy hexadecane | 338 | C18H33F3O2 | 006222-03-3  | 83   |
| 4    |    | Octadecane, 1-(ethenyl)-    | 296 | C20H40     | 000930-02-9  | 83   |
| 5    |    | 1-Docosanol, methyl ether   | 340 | C23H48O    | 1000333-91-9 | 76   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

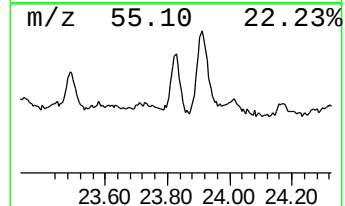
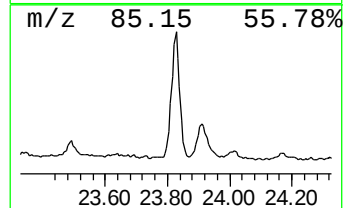
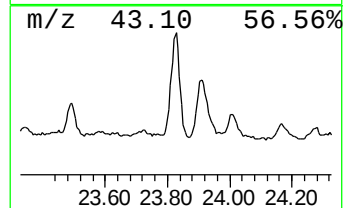
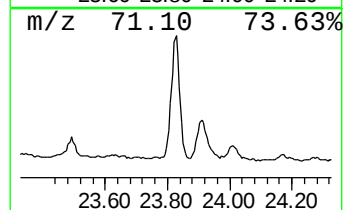
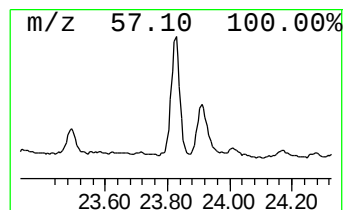
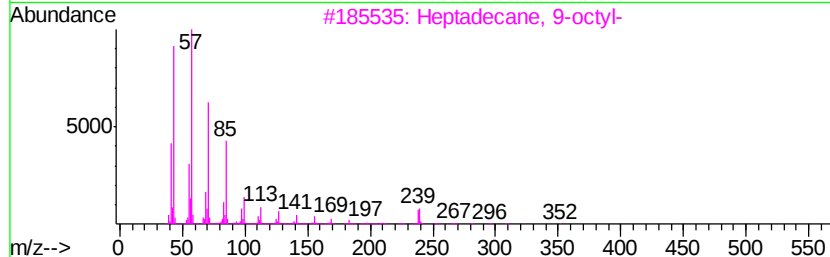
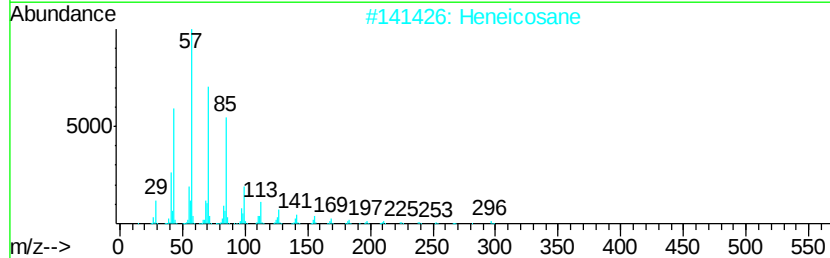
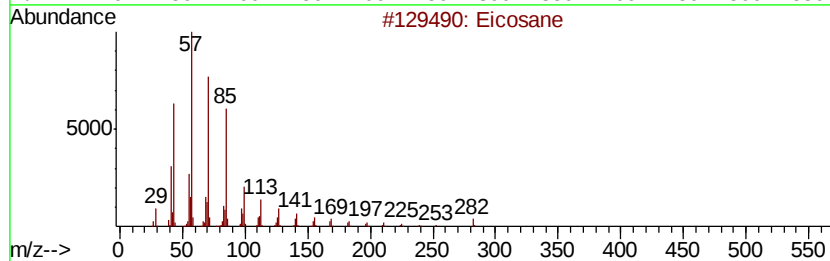
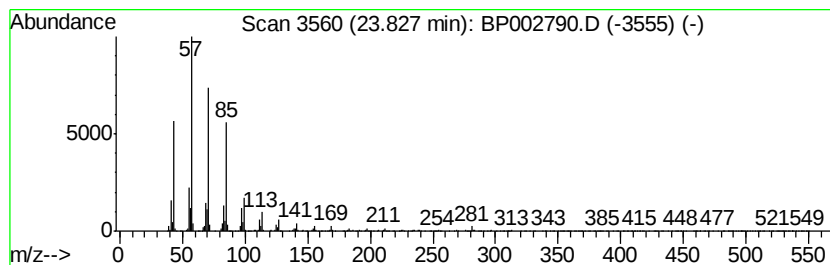
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Eicosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 23.83 | 8.35 ng | 697425 | Perylene-d12     | 23.27 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane              | 282 | C20H42  | 000112-95-8 | 94   |
| 2    |    | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |
| 4    |    | Tridecane, 6-propyl-  | 226 | C16H34  | 055045-10-8 | 90   |
| 5    |    | Nonadecane            | 268 | C19H40  | 000629-92-5 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP071720\  
Data File : BP002790.D  
Acq On : 17 Jul 2020 14:03  
Operator : CG/JU  
Sample : L3324-01  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP071020.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

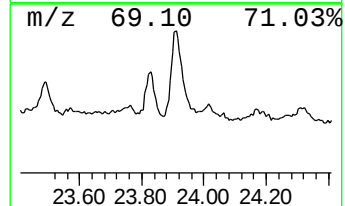
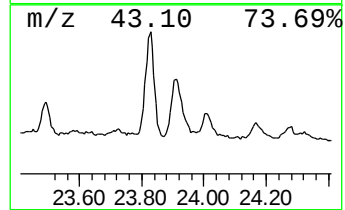
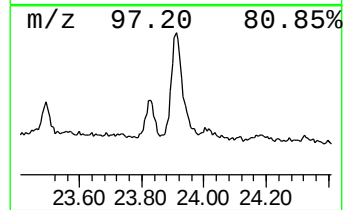
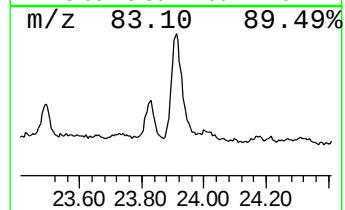
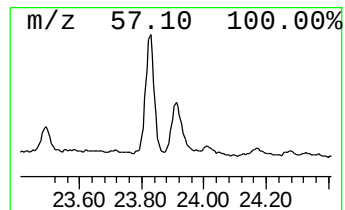
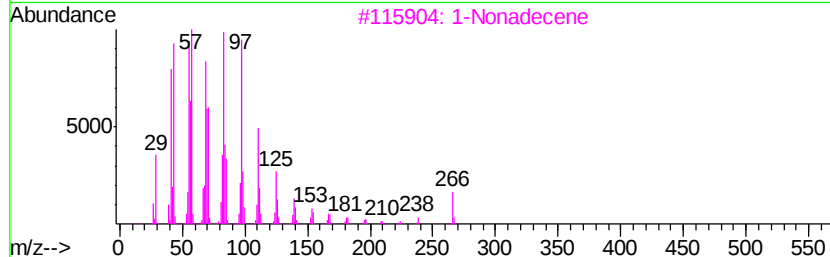
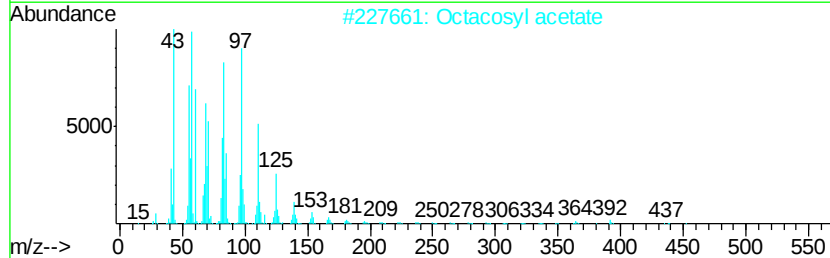
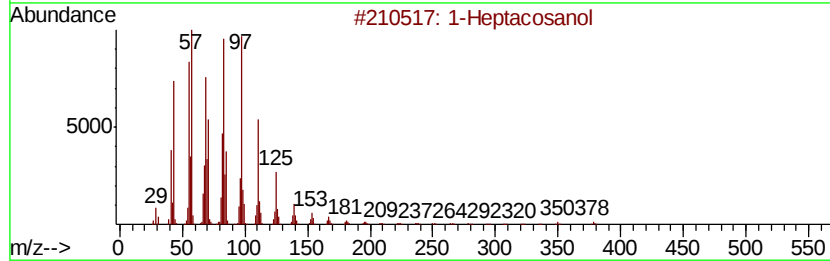
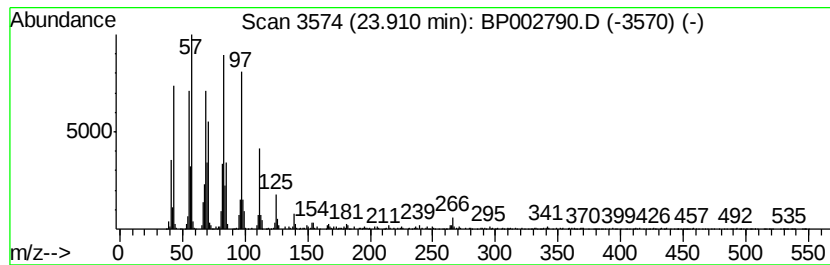
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 1-Heptacosanol Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 23.91 | 7.48 ng | 624692 | Perylene-d12     | 23.27 |

| Hit# | of | Tentative ID      | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------|-----|----------|-------------|------|
| 1    | 5  | 1-Heptacosanol    | 396 | C27H56O  | 002004-39-9 | 94   |
| 2    |    | Octacosyl acetate | 452 | C30H60O2 | 018206-97-8 | 93   |
| 3    |    | 1-Nonadecene      | 266 | C19H38   | 018435-45-5 | 87   |
| 4    |    | Behenic alcohol   | 326 | C22H46O  | 000661-19-8 | 87   |
| 5    |    | Cyclohexadecane   | 224 | C16H32   | 000295-65-8 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP071720\  
 Data File : BP002790.D  
 Acq On : 17 Jul 2020 14:03  
 Operator : CG/JU  
 Sample : L3324-01  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 10:1-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP071020.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT 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 |
| Ethanol, 2-(2-eth... | 7.18  | 2.3     | ng    | 160181   | 1                     | 7.56  | 1388380 | 20.0 |
| Bacchotricuneatin c  | 12.76 | 3.2     | ng    | 474232   | 3                     | 14.21 | 2971340 | 20.0 |
| Cyclohexane, 1,1,... | 13.03 | 2.1     | ng    | 316397   | 3                     | 14.21 | 2971340 | 20.0 |
| Benzene, 1-cycloh... | 13.49 | 4.2     | ng    | 623964   | 3                     | 14.21 | 2971340 | 20.0 |
| unknown13.57         | 13.57 | 4.4     | ng    | 654462   | 3                     | 14.21 | 2971340 | 20.0 |
| Hexadecane, 7-met... | 13.73 | 11.5    | ng    | 1707760  | 3                     | 14.21 | 2971340 | 20.0 |
| Heptadecane, 2,6,... | 13.94 | 2.4     | ng    | 352195   | 3                     | 14.21 | 2971340 | 20.0 |
| unknown14.05         | 14.05 | 4.7     | ng    | 690167   | 3                     | 14.21 | 2971340 | 20.0 |
| unknown14.13         | 14.13 | 3.5     | ng    | 526014   | 3                     | 14.21 | 2971340 | 20.0 |
| Heptadecane, 2-me... | 14.77 | 6.3     | ng    | 934240   | 3                     | 14.21 | 2971340 | 20.0 |
| unknown15.02         | 15.02 | 4.2     | ng    | 622843   | 3                     | 14.21 | 2971340 | 20.0 |
| Pentadecane, 2,6,... | 15.52 | 13.6    | ng    | 2022470  | 3                     | 14.21 | 2971340 | 20.0 |
| Heptadecane, 2,6-... | 16.00 | 17.2    | ng    | 2950120  | 4                     | 16.97 | 3438880 | 20.0 |
| Hexacosane           | 16.83 | 20.0    | ng    | 3438880  | 4                     | 16.97 | 3438880 | 20.0 |
| 9-Eicosene, (E)-     | 20.82 | 18.9    | ng    | 1332710  | 5                     | 21.09 | 1411350 | 20.0 |
| Cyclotetradecane     | 21.69 | 11.5    | ng    | 809576   | 5                     | 21.09 | 1411350 | 20.0 |
| 1,19-Eicosadiene     | 22.35 | 4.9     | ng    | 406645   | 6                     | 23.27 | 1669850 | 20.0 |
| Cyclotetracosane     | 22.67 | 15.2    | ng    | 1267570  | 6                     | 23.27 | 1669850 | 20.0 |
| Eicosane             | 23.83 | 8.3     | ng    | 697425   | 6                     | 23.27 | 1669850 | 20.0 |
| 1-Heptacosanol       | 23.91 | 7.5     | ng    | 624692   | 6                     | 23.27 | 1669850 | 20.0 |