

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
 Data File : BG041355.D  
 Acq On : 20 Jun 2019 13:55  
 Operator : HP/JU  
 Sample : K3158-03 2X  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OK-01-061719

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\_G\METHODS\8270-BG061719.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         | 3.146       | 5             | 10          | 19           | rVB      | 91946          | 135088        | 10.41%          | 1.062%        |
| 2         | 4.609       | 255           | 259         | 265          | rBV2     | 9571           | 15073         | 1.16%           | 0.118%        |
| 3         | 5.614       | 423           | 430         | 439          | rBV      | 352769         | 571746        | 44.06%          | 4.495%        |
| 4         | 7.230       | 698           | 705         | 718          | rVB      | 402907         | 695032        | 53.56%          | 5.464%        |
| 5         | 7.600       | 761           | 768         | 781          | rVB      | 372956         | 663869        | 51.16%          | 5.219%        |
| 6         | 8.070       | 843           | 848         | 856          | rBV      | 144402         | 247124        | 19.05%          | 1.943%        |
| 7         | 8.399       | 897           | 904         | 914          | rVB      | 282202         | 492017        | 37.92%          | 3.868%        |
| 8         | 9.251       | 1041          | 1049        | 1059         | rBV      | 237022         | 440405        | 33.94%          | 3.462%        |
| 9         | 10.896      | 1321          | 1329        | 1345         | rBV      | 200248         | 363034        | 27.98%          | 2.854%        |
| 10        | 13.328      | 1737          | 1743        | 1750         | rVB      | 635439         | 968492        | 74.64%          | 7.614%        |
| 11        | 14.033      | 1858          | 1863        | 1867         | rBV6     | 11626          | 20167         | 1.55%           | 0.159%        |
| 12        | 14.151      | 1879          | 1883        | 1889         | rBV      | 56943          | 90804         | 7.00%           | 0.714%        |
| 13        | 14.269      | 1898          | 1903        | 1908         | rBV7     | 7071           | 14552         | 1.12%           | 0.114%        |
| 14        | 14.439      | 1928          | 1932        | 1936         | rBV3     | 14598          | 25688         | 1.98%           | 0.202%        |
| 15        | 14.521      | 1943          | 1946        | 1949         | rBV5     | 8386           | 14142         | 1.09%           | 0.111%        |
| 16        | 14.709      | 1972          | 1978        | 1986         | rVB2     | 303740         | 472800        | 36.44%          | 3.717%        |
| 17        | 15.173      | 2053          | 2057        | 2061         | rVB6     | 14712          | 24788         | 1.91%           | 0.195%        |
| 18        | 15.314      | 2075          | 2081        | 2086         | rBV8     | 15426          | 38595         | 2.97%           | 0.303%        |
| 19        | 15.461      | 2103          | 2106        | 2110         | rBV6     | 8187           | 13156         | 1.01%           | 0.103%        |
| 20        | 15.778      | 2152          | 2160        | 2161         | rBV7     | 36255          | 78456         | 6.05%           | 0.617%        |
| 21        | 16.202      | 2226          | 2232        | 2239         | rVB      | 473081         | 729377        | 56.21%          | 5.734%        |
| 22        | 16.860      | 2337          | 2344        | 2345         | rBV7     | 140553         | 278206        | 21.44%          | 2.187%        |
| 23        | 17.288      | 2412          | 2417        | 2419         | rBV4     | 187244         | 343352        | 26.46%          | 2.699%        |
| 24        | 17.459      | 2441          | 2446        | 2450         | rVB      | 404005         | 599641        | 46.21%          | 4.714%        |
| 25        | 18.035      | 2542          | 2544        | 2550         | rVB2     | 23953          | 29175         | 2.25%           | 0.229%        |
| 26        | 18.311      | 2586          | 2591        | 2598         | rBV2     | 116307         | 255914        | 19.72%          | 2.012%        |
| 27        | 18.375      | 2598          | 2602        | 2609         | rVB      | 78755          | 122048        | 9.41%           | 0.959%        |
| 28        | 18.516      | 2621          | 2626        | 2631         | rBV2     | 76563          | 176336        | 13.59%          | 1.386%        |
| 29        | 18.822      | 2674          | 2678        | 2681         | rBV2     | 42386          | 62100         | 4.79%           | 0.488%        |
| 30        | 19.034      | 2712          | 2714        | 2715         | rBV2     | 21427          | 18225         | 1.40%           | 0.143%        |
| 31        | 19.116      | 2725          | 2728        | 2731         | rVB      | 45941          | 51126         | 3.94%           | 0.402%        |
| 32        | 19.233      | 2742          | 2748        | 2753         | rBV2     | 124565         | 292636        | 22.55%          | 2.301%        |
| 33        | 19.504      | 2789          | 2794        | 2799         | rBV      | 439680         | 616393        | 47.50%          | 4.846%        |
| 34        | 19.833      | 2846          | 2850        | 2852         | rBV2     | 69009          | 107099        | 8.25%           | 0.842%        |

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Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

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\_G\METHODS\8270-BG061719.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |         |
|----|--------|------|------|------|------|---------|---------|---------|---------|
| 35 | 19.868 | 2852 | 2856 | 2862 | rVB  | 526093  | 747591  | 57.61%  | 5.877%  |
| 36 | 19.932 | 2863 | 2867 | 2870 | rBV  | 75708   | 119192  | 9.19%   | 0.937%  |
| 37 | 20.050 | 2883 | 2887 | 2892 | rVB  | 1029334 | 1297577 | 100.00% | 10.201% |
| 38 | 20.320 | 2931 | 2933 | 2936 | rBV3 | 52963   | 61077   | 4.71%   | 0.480%  |
| 39 | 20.649 | 2987 | 2989 | 2993 | rVB  | 75742   | 77018   | 5.94%   | 0.605%  |
| 40 | 21.730 | 3166 | 3173 | 3178 | rBV3 | 416654  | 1014754 | 78.20%  | 7.978%  |
| 41 | 21.777 | 3179 | 3181 | 3190 | rVB  | 227490  | 336159  | 25.91%  | 2.643%  |

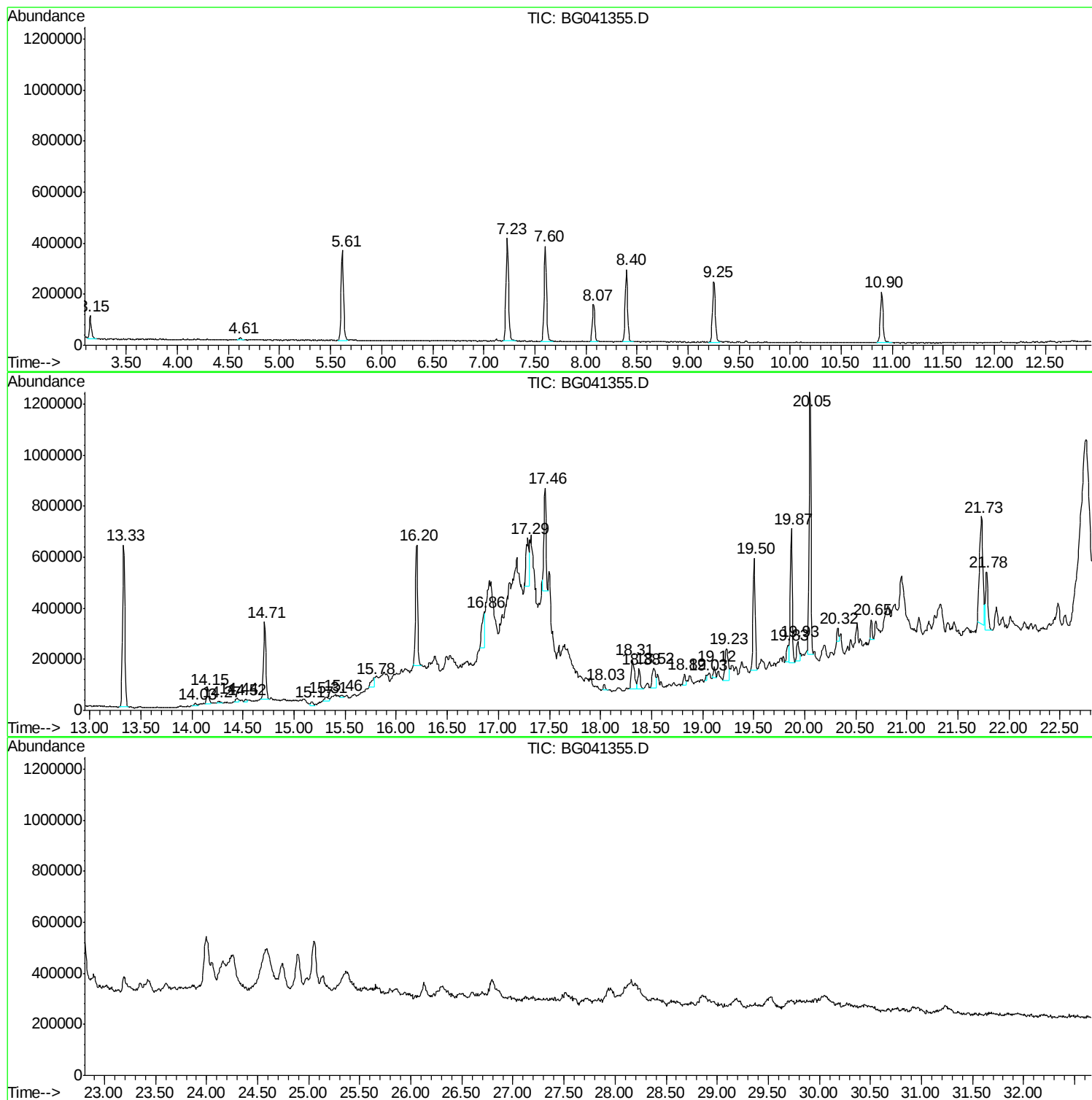
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
OK-01-061719

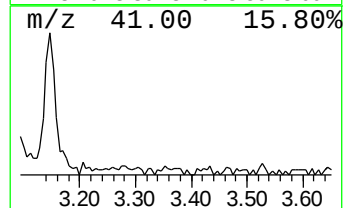
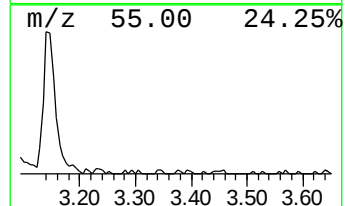
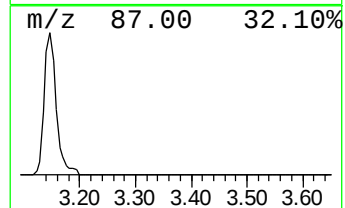
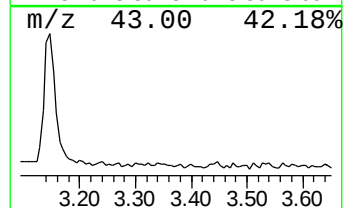
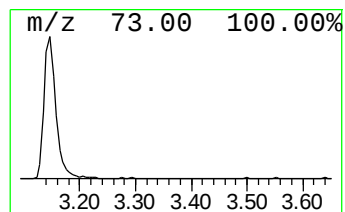
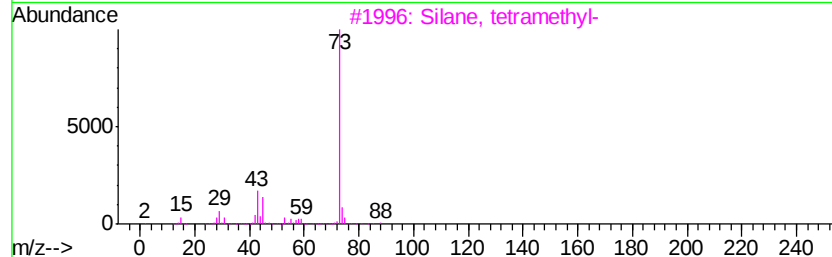
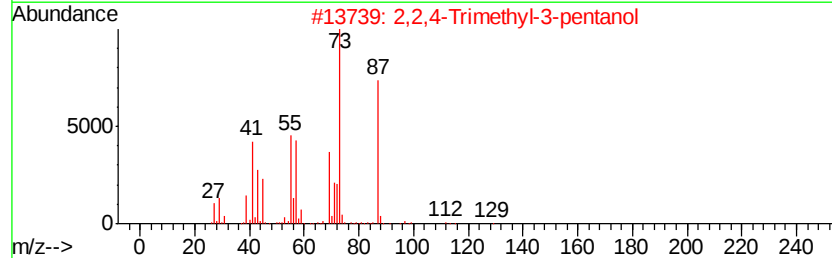
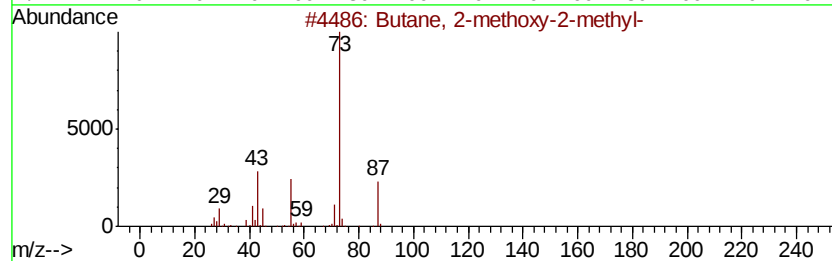
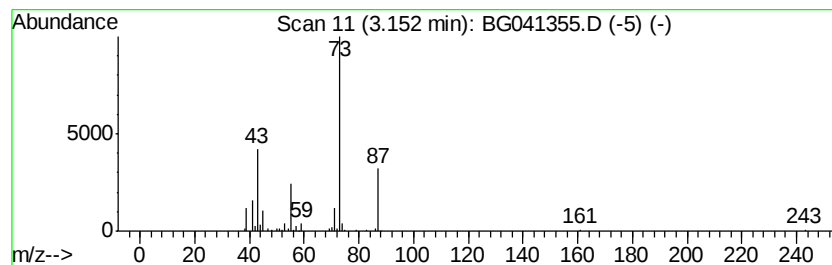
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.15 | 10.93 ng | 135088 | 1,4-Dichlorobenzene-d4 | 8.08 |

| Hit# | of | Tentative ID                            | MW  | MolForm | CAS#         | Qual |
|------|----|-----------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl-             | 102 | C6H14O  | 000994-05-8  | 78   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol              | 130 | C8H18O  | 005162-48-1  | 40   |
| 3    |    | Silane, tetramethyl-                    | 88  | C4H12Si | 000075-76-3  | 38   |
| 4    |    | Butanoic acid, 2-hydroxy-2-methyl-      | 132 | C6H12O3 | 032793-34-3  | 37   |
| 5    |    | 1-(2-Methoxyethoxy)-2-methyl-2-propanol | 162 | C8H18O3 | 1000364-24-4 | 25   |



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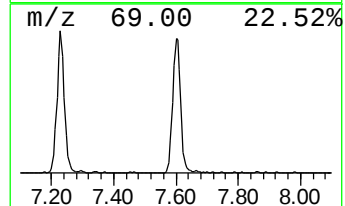
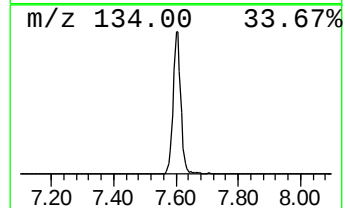
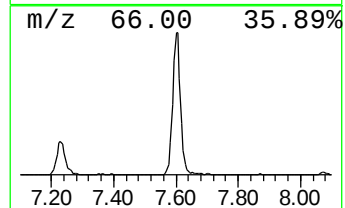
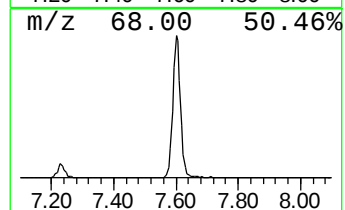
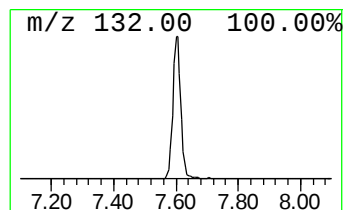
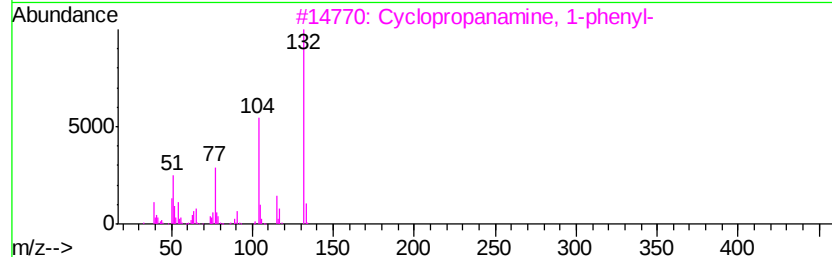
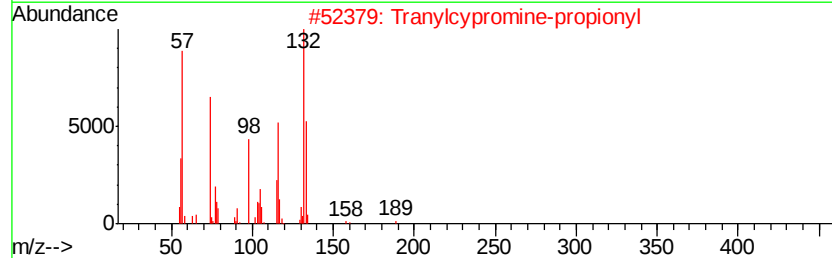
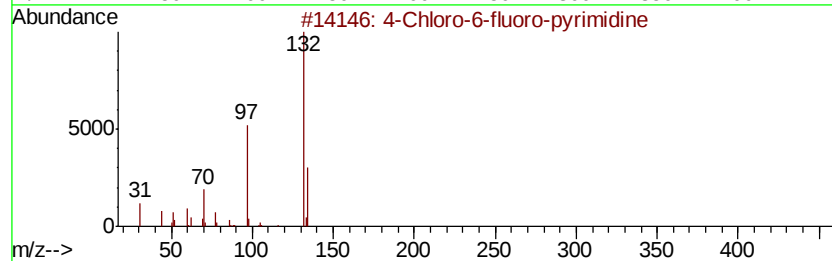
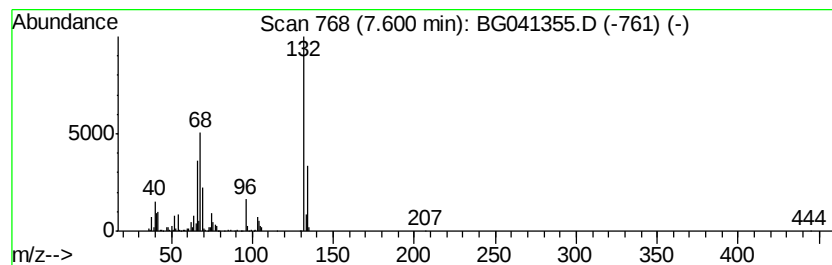
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown7.60 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.60 | 53.73 ng | 663869 | 1,4-Dichlorobenzene-d4 | 8.08 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | Tranlycypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 22   |
| 3    |    | Cyclopropanamine, 1-phenyl-         | 133 | C9H11N    | 1000351-26-2 | 22   |
| 4    |    | 1H-Benzimidazole, 2-methyl-         | 132 | C8H8N2    | 000615-15-6  | 22   |
| 5    |    | 2H-Cyclopenta[d]pyridazine, 2-me... | 132 | C8H8N2    | 022291-85-6  | 16   |



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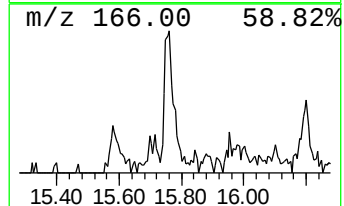
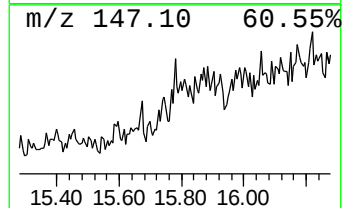
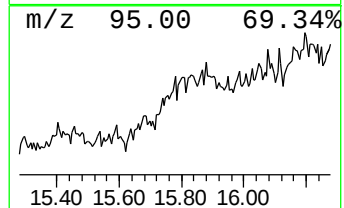
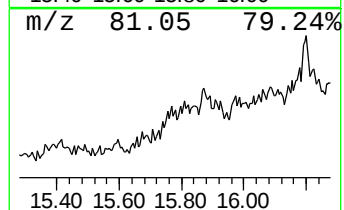
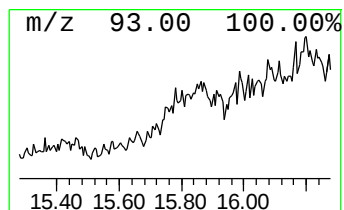
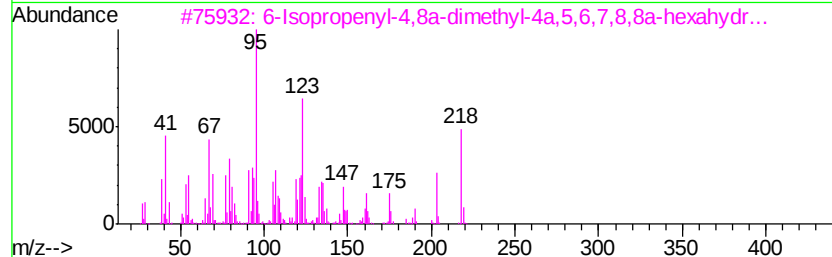
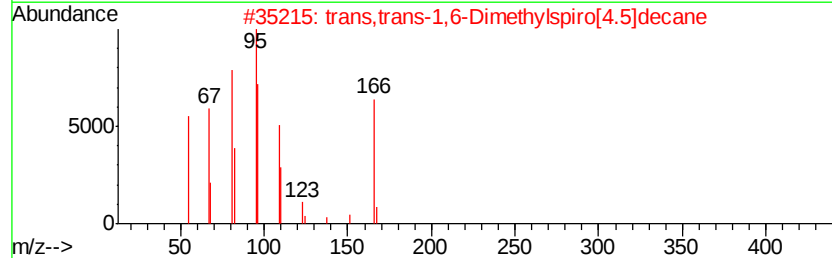
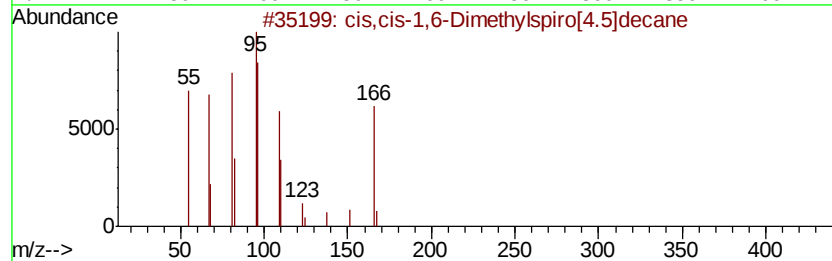
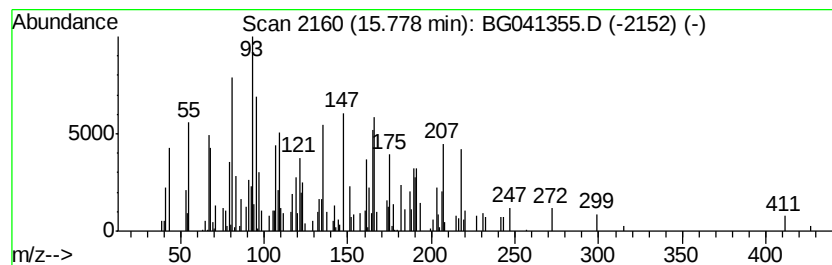
Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 unknown15.78 Concentration Rank 10

| R.T.    | EstConc | Area                                    | Relative to ISTD | R.T.            |
|---------|---------|-----------------------------------------|------------------|-----------------|
| 15.78   | 3.32 ng | 78456                                   | Acenaphthene-d10 | 14.71           |
| Hit# of | 5       | Tentative ID                            | MW MolForm       | CAS# Qual       |
| 1       |         | cis,cis-1,6-Dimethylspiro[4.5]de...     | 166 C12H22       | 1000111-72-4 25 |
| 2       |         | trans,trans-1,6-Dimethylspiro[4.5]de... | 166 C12H22       | 1000111-72-1 25 |
| 3       |         | 6-Isopropenyl-4,8a-dimethyl-4a,5...     | 218 C15H22O      | 086917-79-5 15  |
| 4       |         | 1,4-Methanoazulen-7(1H)-one, oct...     | 220 C15H24O      | 018319-28-3 11  |
| 5       |         | Ethionamide                             | 166 C8H10N2S     | 000536-33-4 11  |



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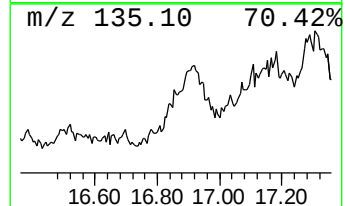
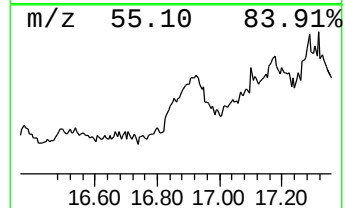
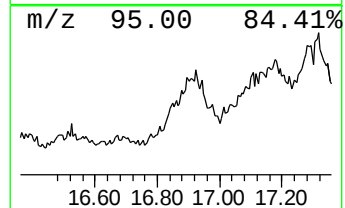
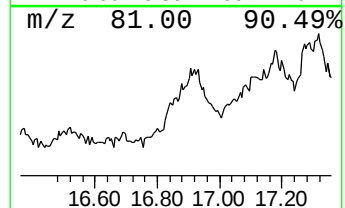
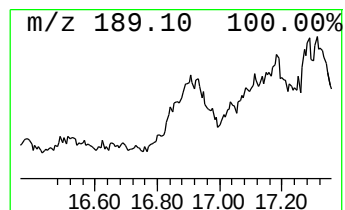
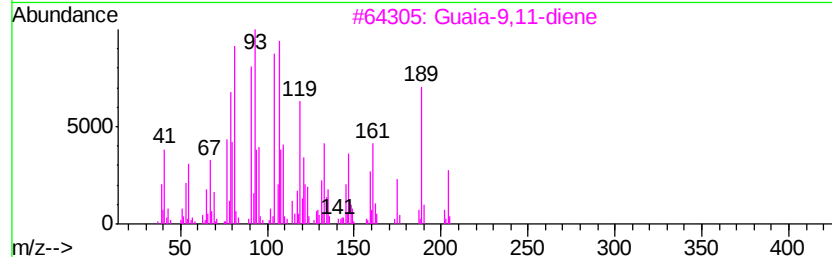
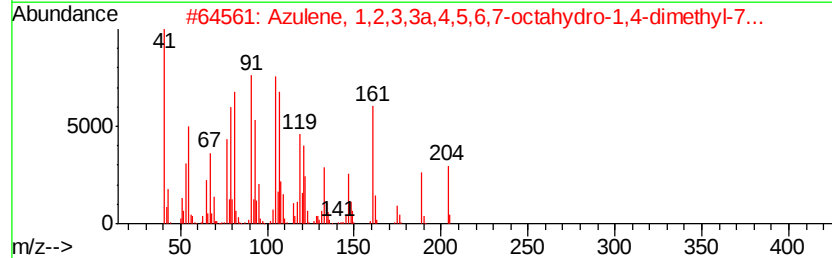
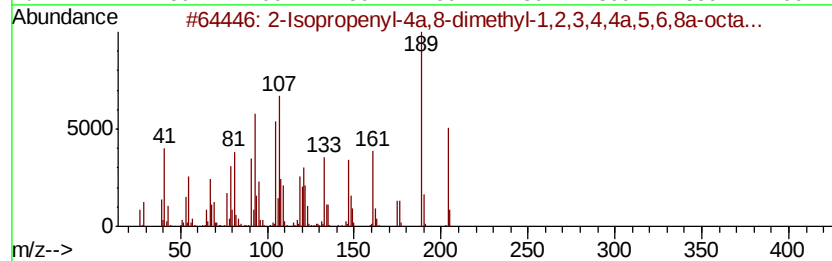
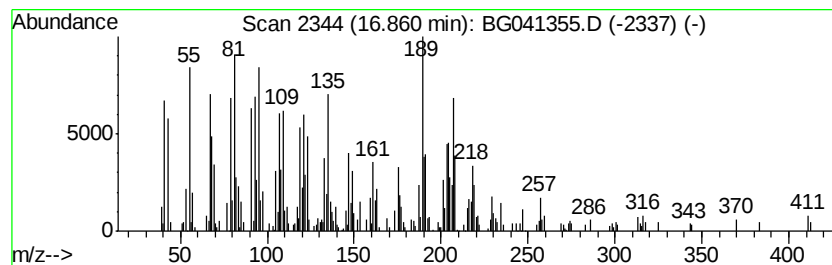
Instrument :  
BNA\_G  
ClientSampleID :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 2-Isopropenyl-4a,8-dimethyl... Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 16.86     | 9.28 ng                             | 278206 | Phenanthrene-d10 | 17.46        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2-Isopropenyl-4a,8-dimethyl-1,2,... | 204    | C15H24           | 1000193-57-0 | 86   |
| 2         | Azulene, 1,2,3,3a,4,5,6,7-octahy... | 204    | C15H24           | 022567-17-5  | 53   |
| 3         | Guaia-9,11-diene                    | 204    | C15H24           | 1000374-19-8 | 50   |
| 4         | Naphthalene, 2,3,4,4a,5,6-hexahy... | 204    | C15H24           | 000473-14-3  | 44   |
| 5         | Naphthalene, decahydro-4a-methyl... | 204    | C15H24           | 017066-67-0  | 44   |



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Data File : BG041355.D  
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Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

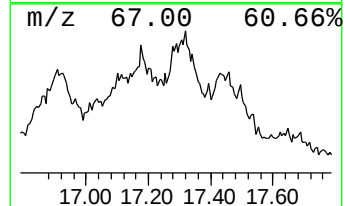
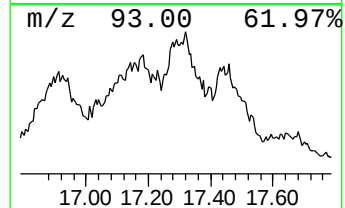
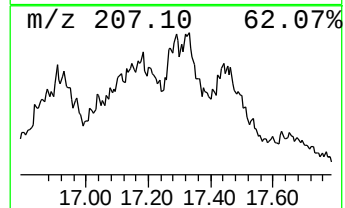
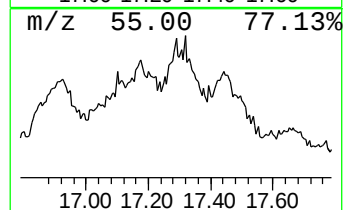
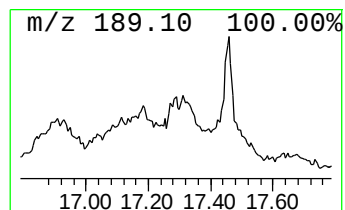
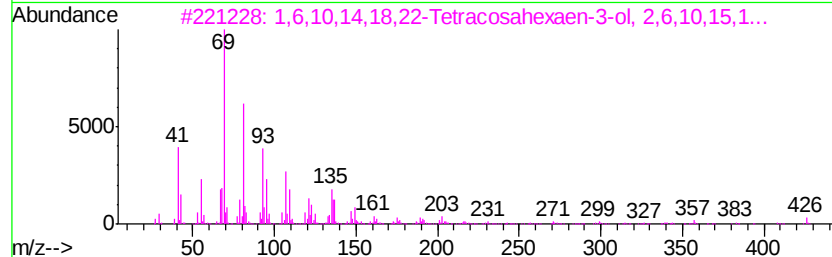
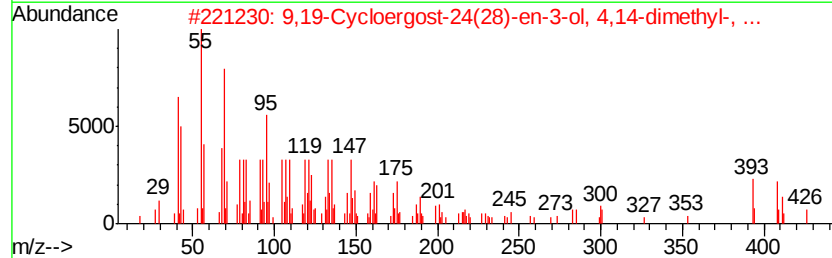
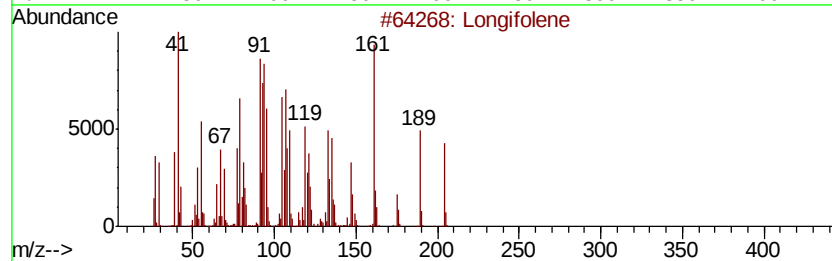
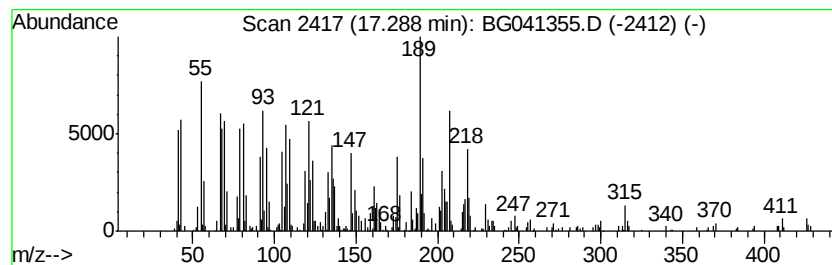
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ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 unknown17.29 Concentration Rank 3

| R.T.    | EstConc  | Area                                 | Relative to ISTD | R.T.            |
|---------|----------|--------------------------------------|------------------|-----------------|
| 17.29   | 11.45 ng | 343352                               | Phenanthrene-d10 | 17.46           |
| Hit# of | 5        | Tentative ID                         | MW MolForm       | CAS# Qual       |
| 1       |          | Longifolene                          | 204 C15H24       | 000475-20-7 47  |
| 2       |          | 9,19-Cycloergost-24(28)-en-3-ol, ... | 426 C30H500      | 000469-39-6 44  |
| 3       |          | 1,6,10,14,18,22-Tetracosahexaen-...  | 426 C30H500      | 054159-46-5 42  |
| 4       |          | Lupeol                               | 426 C30H500      | 000545-47-1 38  |
| 5       |          | Diepicedrene-1-oxide                 | 220 C15H240      | 1000156-11-0 35 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

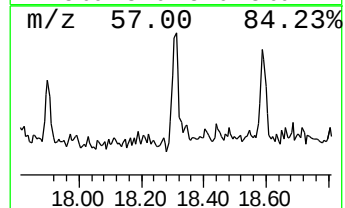
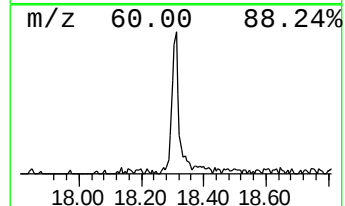
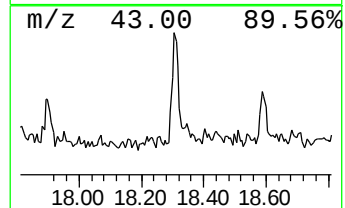
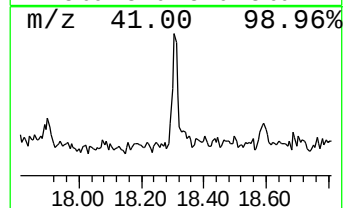
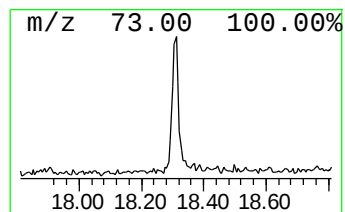
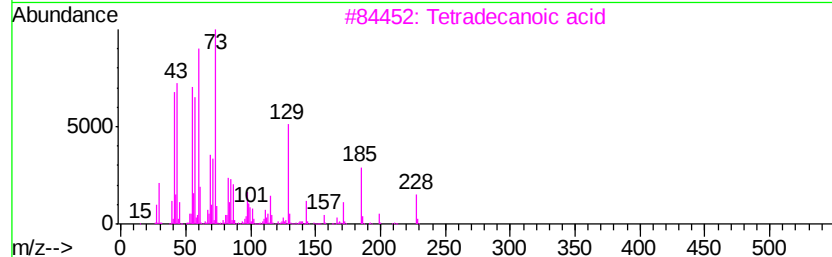
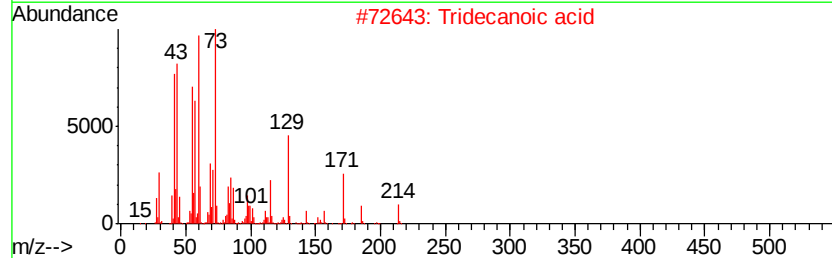
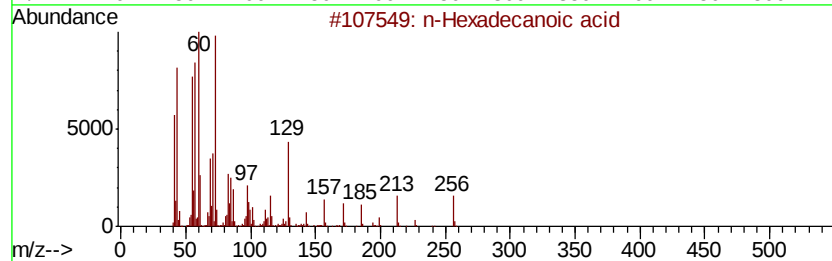
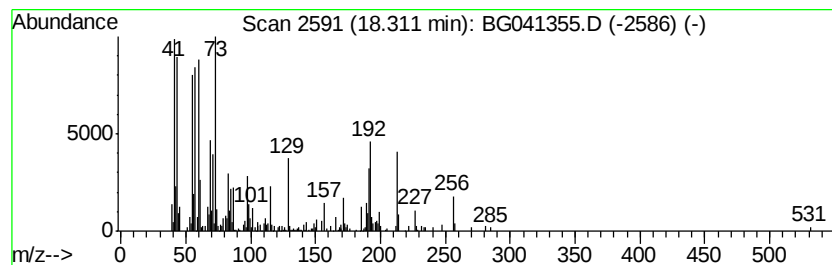
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 n-Hexadecanoic acid Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.31 | 8.54 ng | 255914 | Phenanthrene-d10 | 17.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid                 | 256 | C16H32O2   | 000057-10-3 | 97   |
| 2    |    | Tridecanoic acid                    | 214 | C13H26O2   | 000638-53-9 | 78   |
| 3    |    | Tetradecanoic acid                  | 228 | C14H28O2   | 000544-63-8 | 78   |
| 4    |    | Pentadecanoic acid                  | 242 | C15H30O2   | 001002-84-2 | 60   |
| 5    |    | 1-Phenazinecarboxylic acid, 6-[1... | 506 | C31H42N2O4 | 120464-92-8 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

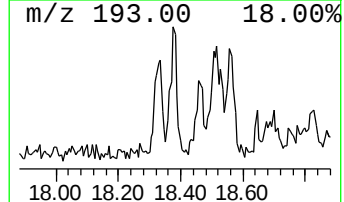
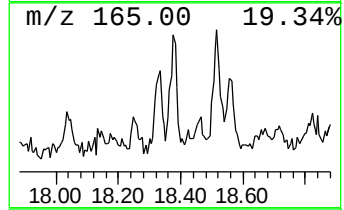
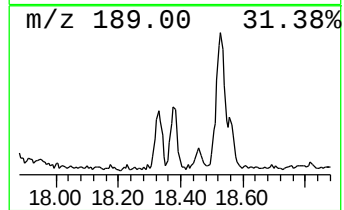
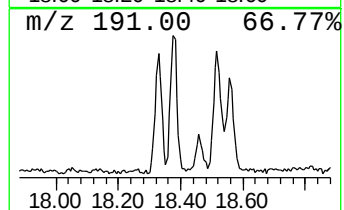
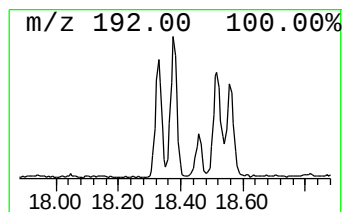
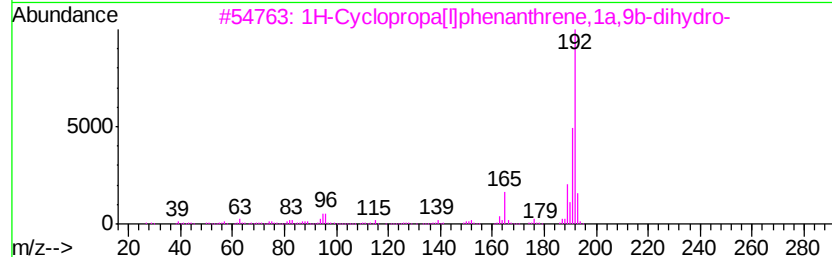
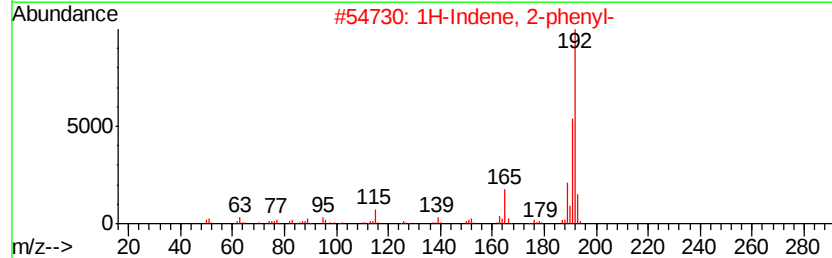
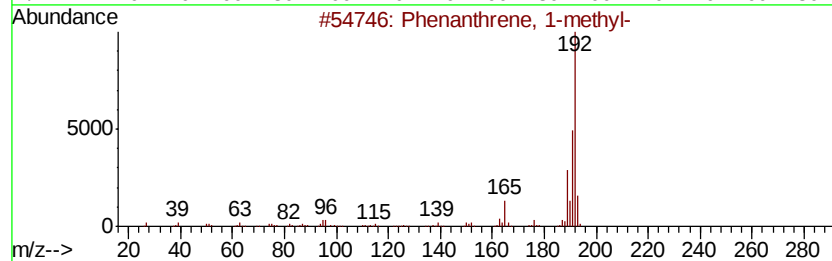
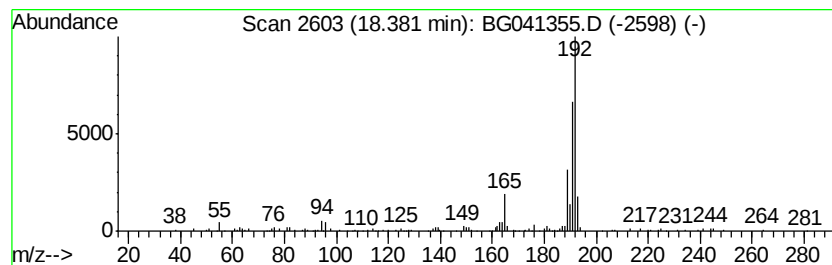
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.38 | 4.07 ng | 122048 | Phenanthrene-d10 | 17.46 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 1-methyl-              | 192 | C15H12  | 000832-69-9 | 94   |
| 2    |    | 1H-Indene, 2-phenyl-                 | 192 | C15H12  | 004505-48-0 | 94   |
| 3    |    | 1H-Cyclopropa[1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 91   |
| 4    |    | Anthracene, 2-methyl-                | 192 | C15H12  | 000613-12-7 | 91   |
| 5    |    | 1H-Indene, 1-phenyl-                 | 192 | C15H12  | 001961-96-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

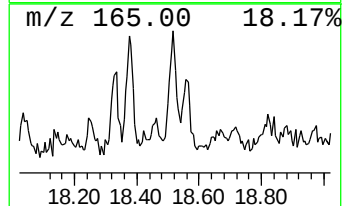
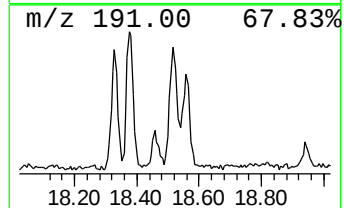
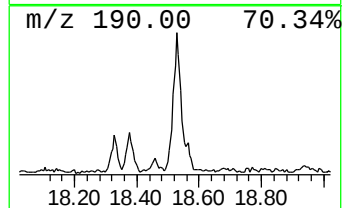
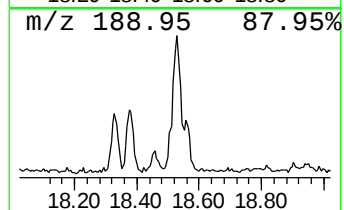
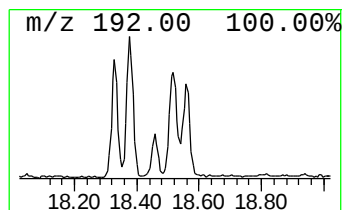
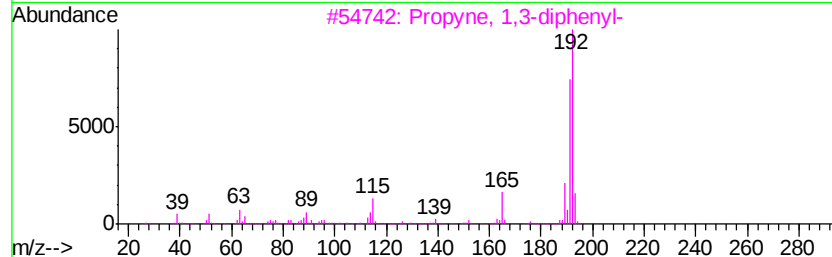
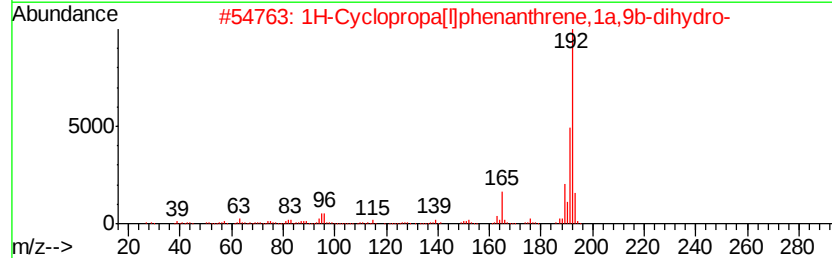
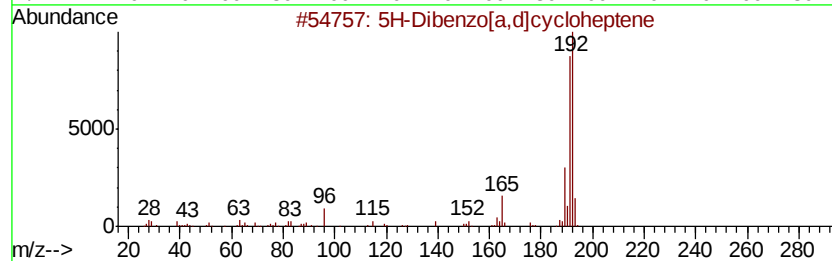
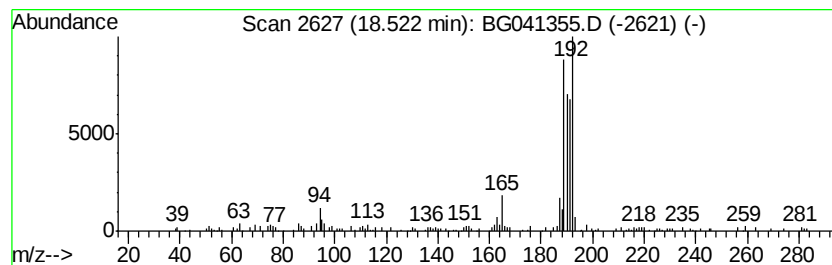
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 5H-Dibenzo[a,d]cycloheptene Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.52 | 5.88 ng | 176336 | Phenanthrene-d10 | 17.46 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 5H-Dibenzo[a,d]cycloheptene         | 192 | C15H12  | 000256-81-5 | 60   |
| 2    |    | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 53   |
| 3    |    | Propyne, 1,3-diphenyl-              | 192 | C15H12  | 004980-70-5 | 53   |
| 4    |    | 1a,9b-Dihydro-1H-cyclopropa[a]an... | 192 | C15H12  | 006109-24-6 | 53   |
| 5    |    | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 52   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
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Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

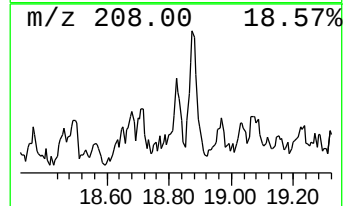
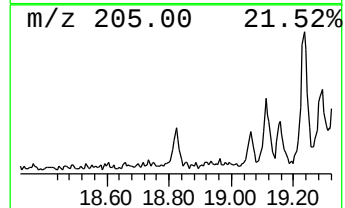
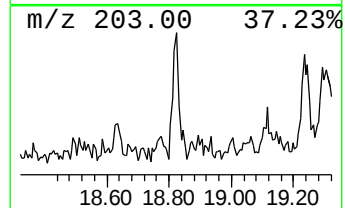
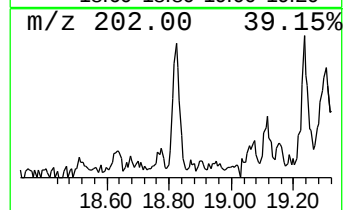
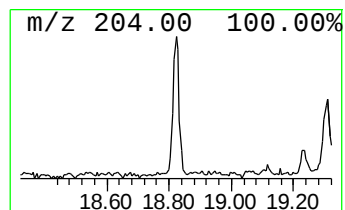
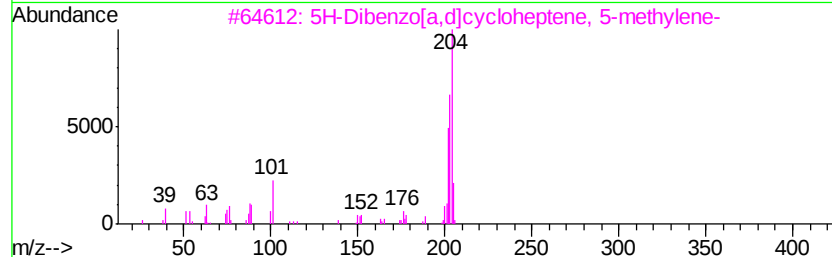
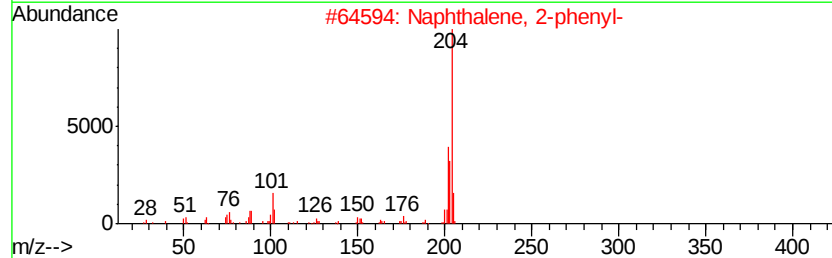
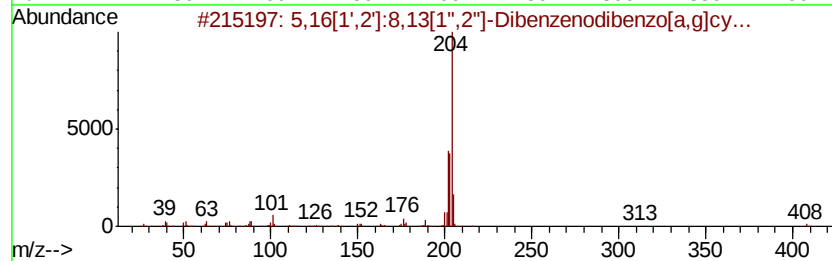
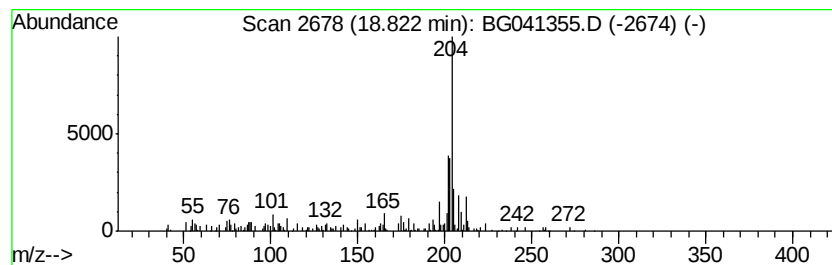
Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.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 10 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 12

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.    |             |      |
|---------|-------------------------------------|--------------|------------------|---------|-------------|------|
| 18.82   | 2.07 ng                             | 62100        | Phenanthrene-d10 | 17.46   |             |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm | CAS#        | Qual |
| 1       | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408          | C32H24           |         | 005672-97-9 | 72   |
| 2       | Naphthalene, 2-phenyl-              | 204          | C16H12           |         | 000612-94-2 | 53   |
| 3       | 5H-Dibenzo[a,d]cycloheptene, 5-m... | 204          | C16H12           |         | 002975-79-3 | 50   |
| 4       | 9,10-di(Chloromethyl)anthracene     | 274          | C16H12Cl2        |         | 010387-13-0 | 50   |
| 5       | 6-Phenylbenzocyclohepten-7-one      | 232          | C17H12O          |         | 093327-56-1 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

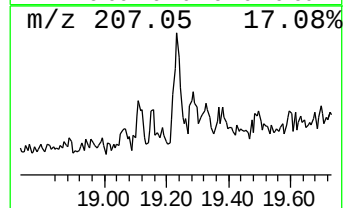
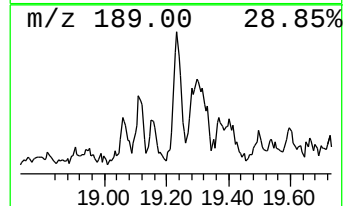
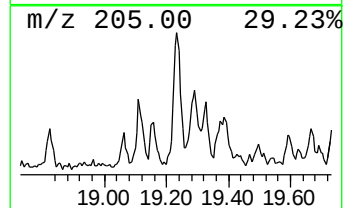
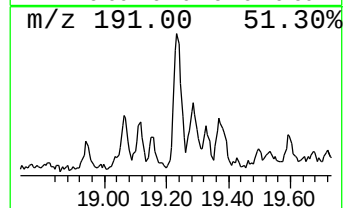
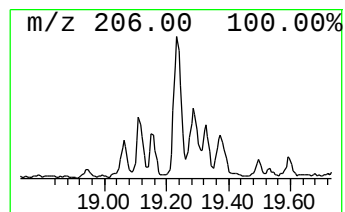
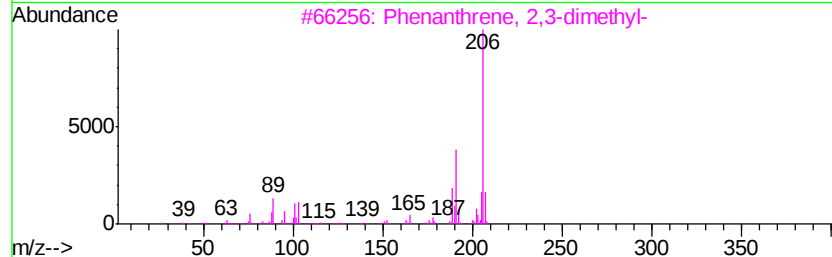
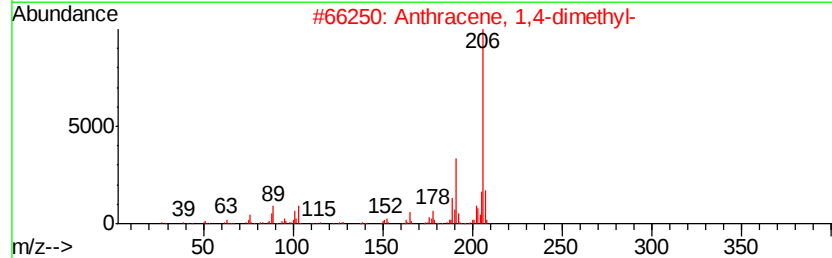
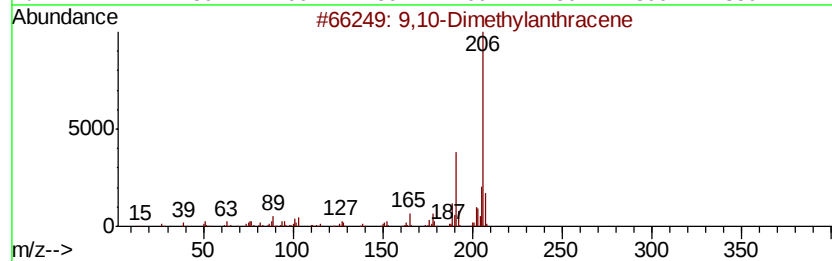
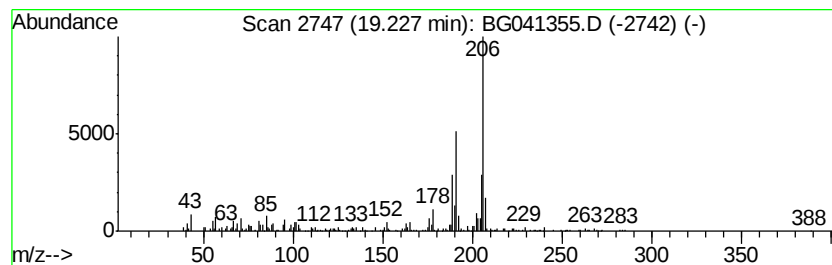
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 9,10-Dimethylantracene Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.23 | 9.76 ng | 292636 | Phenanthrene-d10 | 17.46 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 9,10-Dimethylantracene             | 206 | C16H14  | 000781-43-1 | 93   |
| 2    |    | Anthracene, 1,4-dimethyl-          | 206 | C16H14  | 000781-92-0 | 91   |
| 3    |    | Phenanthrene, 2,3-dimethyl-        | 206 | C16H14  | 003674-65-5 | 87   |
| 4    |    | Naphthalene, 1,2-dihydro-4-phenyl- | 206 | C16H14  | 007469-40-1 | 86   |
| 5    |    | Phenanthrene, 3,6-dimethyl-        | 206 | C16H14  | 001576-67-6 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

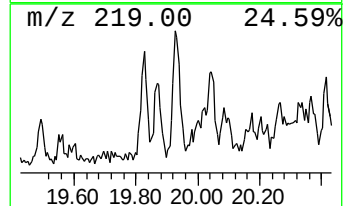
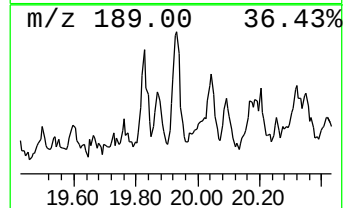
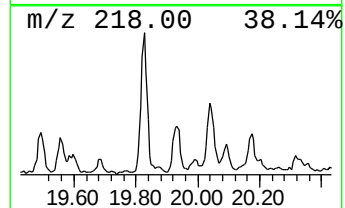
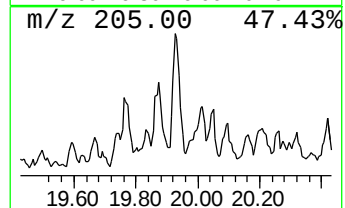
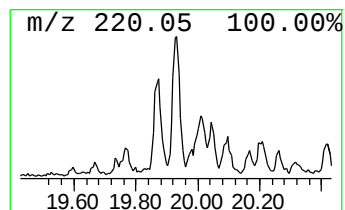
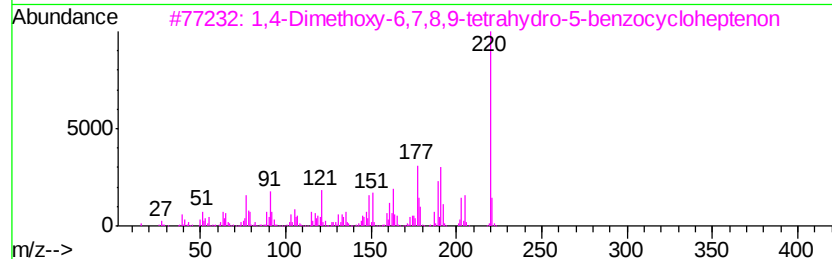
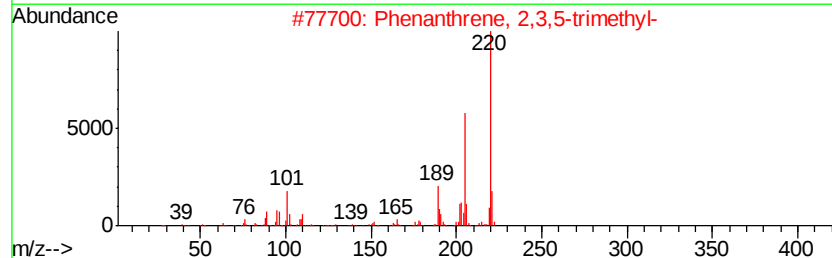
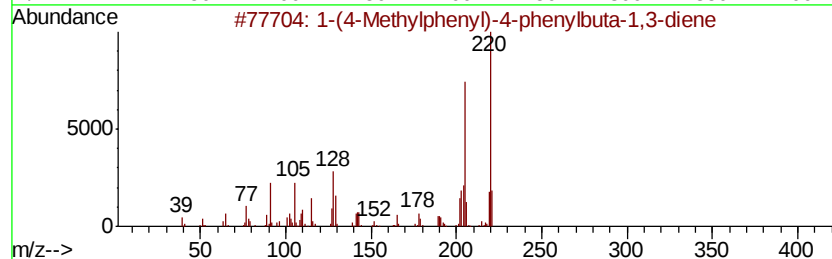
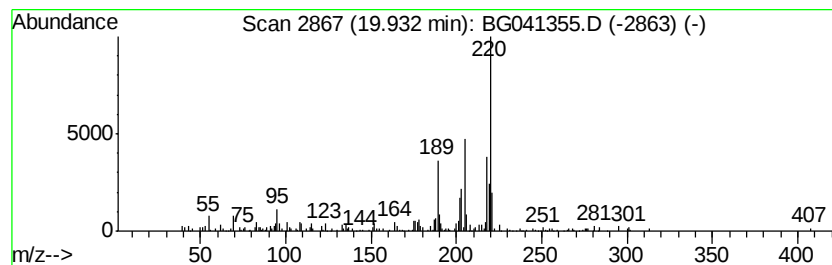
Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 1-(4-Methylphenyl)-4-phenyl... Concentration Rank 11

| R.T.    | EstConc | Area                                 | Relative to ISTD | R.T.       |              |      |
|---------|---------|--------------------------------------|------------------|------------|--------------|------|
| 19.93   | 2.35 ng | 119192                               | Chrysene-d12     | 21.74      |              |      |
| Hit# of | 5       | Tentative ID                         | MW               | MolForm    | CAS#         | Qual |
| 1       |         | 1-(4-Methylphenyl)-4-phenylbuta-...  | 220              | C17H16     | 037985-11-8  | 50   |
| 2       |         | Phenanthrene, 2,3,5-trimethyl-       | 220              | C17H16     | 003674-73-5  | 50   |
| 3       |         | 1,4-Dimethoxy-6,7,8,9-tetrahydro...  | 220              | C13H16O3   | 079381-09-2  | 47   |
| 4       |         | Phenmethylnitrile, .alpha.,.alpha... | 220              | C11H12N2O3 | 1000128-94-6 | 46   |
| 5       |         | Pyrrolo[1,2-a]thieno[2,3-d]pyrim...  | 220              | C11H12N2OS | 056929-61-4  | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_G\DATA\BG062019\  
Data File : BG041355.D  
Acq On : 20 Jun 2019 13:55  
Operator : HP/JU  
Sample : K3158-03 2X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-01-061719

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\8270-BG061719.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 |
| Butane, 2-methoxy...  | 3.15  | 10.9    | ng    | 135088   | 1                     | 8.08  | 247124  | 20.0 |
| unknown7.60           | 7.60  | 53.7    | ng    | 663869   | 1                     | 8.08  | 247124  | 20.0 |
| unknown15.78          | 15.78 | 3.3     | ng    | 78456    | 3                     | 14.71 | 472800  | 20.0 |
| 2-Isopropenyl-4a,...  | 16.86 | 9.3     | ng    | 278206   | 4                     | 17.46 | 599641  | 20.0 |
| unknown17.29          | 17.29 | 11.4    | ng    | 343352   | 4                     | 17.46 | 599641  | 20.0 |
| n-Hexadecanoic acid   | 18.31 | 8.5     | ng    | 255914   | 4                     | 17.46 | 599641  | 20.0 |
| Phenanthrene, 1-m...  | 18.38 | 4.1     | ng    | 122048   | 4                     | 17.46 | 599641  | 20.0 |
| 5H-Dibenzo[a,d]cy...  | 18.52 | 5.9     | ng    | 176336   | 4                     | 17.46 | 599641  | 20.0 |
| 5,16[1',2']:8,13[...  | 18.82 | 2.1     | ng    | 62100    | 4                     | 17.46 | 599641  | 20.0 |
| 9,10-Dimethylanthe... | 19.23 | 9.8     | ng    | 292636   | 4                     | 17.46 | 599641  | 20.0 |
| 1-(4-Methylphenyl...  | 19.93 | 2.4     | ng    | 119192   | 5                     | 21.74 | 1014750 | 20.0 |