

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
 Data File : BG034060.D  
 Acq On : 28 Apr 2018 5:48  
 Operator : SJ/JU  
 Sample : J2571-13  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 042418-JETB-E-D-M

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:\HPCHEM1\BNA\_G\METHODS\8270-BG041918.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         | 4.893       | 267           | 273         | 286          | rBV      | 52526          | 107115        | 1.75%           | 0.297%        |
| 2         | 5.375       | 349           | 355         | 372          | rBV2     | 60285          | 267307        | 4.37%           | 0.742%        |
| 3         | 7.296       | 675           | 682         | 719          | rBV      | 239058         | 918664        | 15.02%          | 2.549%        |
| 4         | 7.760       | 754           | 761         | 784          | rBV2     | 77640          | 305010        | 4.99%           | 0.846%        |
| 5         | 8.066       | 806           | 813         | 842          | rBV      | 651001         | 1576656       | 25.78%          | 4.374%        |
| 6         | 8.912       | 949           | 957         | 994          | rBV      | 245370         | 1057302       | 17.29%          | 2.933%        |
| 7         | 10.539      | 1227          | 1234        | 1258         | rBV      | 119743         | 459017        | 7.51%           | 1.273%        |
| 8         | 13.013      | 1648          | 1655        | 1677         | rBV      | 1659843        | 3675903       | 60.10%          | 10.198%       |
| 9         | 13.859      | 1794          | 1799        | 1812         | rBV      | 30974          | 102786        | 1.68%           | 0.285%        |
| 10        | 14.388      | 1883          | 1889        | 1911         | rBV2     | 417379         | 892400        | 14.59%          | 2.476%        |
| 11        | 15.263      | 2034          | 2038        | 2046         | rVB2     | 40911          | 69938         | 1.14%           | 0.194%        |
| 12        | 15.886      | 2138          | 2144        | 2168         | rBV2     | 802072         | 1974376       | 32.28%          | 5.477%        |
| 13        | 16.121      | 2180          | 2184        | 2186         | rBV3     | 86926          | 118808        | 1.94%           | 0.330%        |
| 14        | 16.920      | 2316          | 2320        | 2324         | rBV      | 71170          | 98791         | 1.62%           | 0.274%        |
| 15        | 17.055      | 2339          | 2343        | 2352         | rBV      | 168092         | 243141        | 3.98%           | 0.675%        |
| 16        | 17.149      | 2353          | 2359        | 2367         | rBV2     | 504335         | 1136087       | 18.58%          | 3.152%        |
| 17        | 17.320      | 2385          | 2388        | 2397         | rVB      | 74448          | 109066        | 1.78%           | 0.303%        |
| 18        | 17.519      | 2418          | 2422        | 2427         | rBV      | 124859         | 175610        | 2.87%           | 0.487%        |
| 19        | 17.654      | 2443          | 2445        | 2450         | rVB2     | 70095          | 86314         | 1.41%           | 0.239%        |
| 20        | 17.819      | 2468          | 2473        | 2479         | rBV      | 407619         | 625401        | 10.23%          | 1.735%        |
| 21        | 18.054      | 2510          | 2513        | 2518         | rBV      | 47666          | 70062         | 1.15%           | 0.194%        |
| 22        | 18.354      | 2561          | 2564        | 2569         | rVB      | 64130          | 74965         | 1.23%           | 0.208%        |
| 23        | 18.530      | 2590          | 2594        | 2598         | rBV      | 134054         | 167353        | 2.74%           | 0.464%        |
| 24        | 18.976      | 2666          | 2670        | 2673         | rBV2     | 165474         | 251349        | 4.11%           | 0.697%        |
| 25        | 19.006      | 2673          | 2675        | 2680         | rVB3     | 146974         | 194080        | 3.17%           | 0.538%        |
| 26        | 19.699      | 2789          | 2793        | 2796         | rBV5     | 56848          | 82744         | 1.35%           | 0.230%        |
| 27        | 19.781      | 2801          | 2807        | 2823         | rBV      | 3745008        | 6115855       | 100.00%         | 16.967%       |
| 28        | 19.952      | 2834          | 2836        | 2842         | rVB6     | 58335          | 70300         | 1.15%           | 0.195%        |
| 29        | 20.016      | 2845          | 2847        | 2850         | rVB3     | 80609          | 84146         | 1.38%           | 0.233%        |
| 30        | 20.140      | 2865          | 2868        | 2870         | rBV4     | 69498          | 80920         | 1.32%           | 0.224%        |
| 31        | 20.204      | 2877          | 2879        | 2882         | rVB3     | 67946          | 64564         | 1.06%           | 0.179%        |
| 32        | 20.340      | 2898          | 2902        | 2905         | rBV5     | 71843          | 113434        | 1.85%           | 0.315%        |
| 33        | 20.375      | 2905          | 2908        | 2911         | rBV5     | 57273          | 69171         | 1.13%           | 0.192%        |
| 34        | 20.416      | 2911          | 2915        | 2921         | rVB6     | 115171         | 222332        | 3.64%           | 0.617%        |

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 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 042418-JETB-E-D-M

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:\HPCHEM1\BNA\_G\METHODS\8270-BG041918.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 35 | 20.475 | 2921 | 2925 | 2927 | rBV2 | 182934 | 238760  | 3.90%  | 0.662% |
| 36 | 20.616 | 2947 | 2949 | 2952 | rBV4 | 118419 | 125672  | 2.05%  | 0.349% |
| 37 | 20.733 | 2967 | 2969 | 2973 | rBV3 | 101628 | 105380  | 1.72%  | 0.292% |
| 38 | 20.804 | 2978 | 2981 | 2984 | rBV5 | 83077  | 106438  | 1.74%  | 0.295% |
| 39 | 20.845 | 2984 | 2988 | 2992 | rBV2 | 219583 | 345634  | 5.65%  | 0.959% |
| 40 | 20.915 | 2995 | 3000 | 3002 | rBV5 | 222358 | 363190  | 5.94%  | 1.008% |
| 41 | 21.021 | 3012 | 3018 | 3021 | rBV5 | 239624 | 466466  | 7.63%  | 1.294% |
| 42 | 21.115 | 3030 | 3034 | 3042 | rVB6 | 297799 | 560588  | 9.17%  | 1.555% |
| 43 | 21.186 | 3043 | 3046 | 3048 | rBV2 | 208949 | 239433  | 3.91%  | 0.664% |
| 44 | 21.256 | 3055 | 3058 | 3060 | rBV3 | 96670  | 130192  | 2.13%  | 0.361% |
| 45 | 21.327 | 3066 | 3070 | 3072 | rBV2 | 169788 | 233787  | 3.82%  | 0.649% |
| 46 | 21.362 | 3073 | 3076 | 3077 | rVV3 | 135541 | 160793  | 2.63%  | 0.446% |
| 47 | 21.397 | 3077 | 3082 | 3089 | rVB3 | 868792 | 1845630 | 30.18% | 5.120% |
| 48 | 21.485 | 3093 | 3097 | 3098 | rBV  | 135445 | 191710  | 3.13%  | 0.532% |
| 49 | 21.620 | 3115 | 3120 | 3129 | rVB3 | 442066 | 864375  | 14.13% | 2.398% |
| 50 | 21.714 | 3130 | 3136 | 3141 | rBV4 | 450708 | 926051  | 15.14% | 2.569% |
| 51 | 21.861 | 3157 | 3161 | 3163 | rBV3 | 153250 | 220375  | 3.60%  | 0.611% |
| 52 | 21.938 | 3170 | 3174 | 3182 | rVB3 | 401079 | 784284  | 12.82% | 2.176% |
| 53 | 22.038 | 3186 | 3191 | 3202 | rVV5 | 353745 | 988804  | 16.17% | 2.743% |
| 54 | 22.120 | 3202 | 3205 | 3209 | rVB4 | 97225  | 135732  | 2.22%  | 0.377% |
| 55 | 22.184 | 3210 | 3216 | 3225 | rVB3 | 268238 | 616025  | 10.07% | 1.709% |
| 56 | 22.284 | 3227 | 3233 | 3239 | rBV3 | 420352 | 817336  | 13.36% | 2.268% |
| 57 | 22.449 | 3256 | 3261 | 3266 | rBV7 | 126283 | 297199  | 4.86%  | 0.825% |
| 58 | 22.537 | 3273 | 3276 | 3290 | rVB4 | 202123 | 472206  | 7.72%  | 1.310% |
| 59 | 22.660 | 3291 | 3297 | 3306 | rVV2 | 312523 | 698481  | 11.42% | 1.938% |
| 60 | 22.819 | 3321 | 3324 | 3331 | rVB3 | 104797 | 192827  | 3.15%  | 0.535% |
| 61 | 22.936 | 3338 | 3344 | 3352 | rBV2 | 192134 | 389075  | 6.36%  | 1.079% |
| 62 | 23.377 | 3413 | 3419 | 3434 | rVB3 | 169461 | 419103  | 6.85%  | 1.163% |
| 63 | 24.211 | 3557 | 3561 | 3570 | rVB7 | 49411  | 113702  | 1.86%  | 0.315% |
| 64 | 24.388 | 3582 | 3591 | 3614 | rBV2 | 385961 | 1335043 | 21.83% | 3.704% |

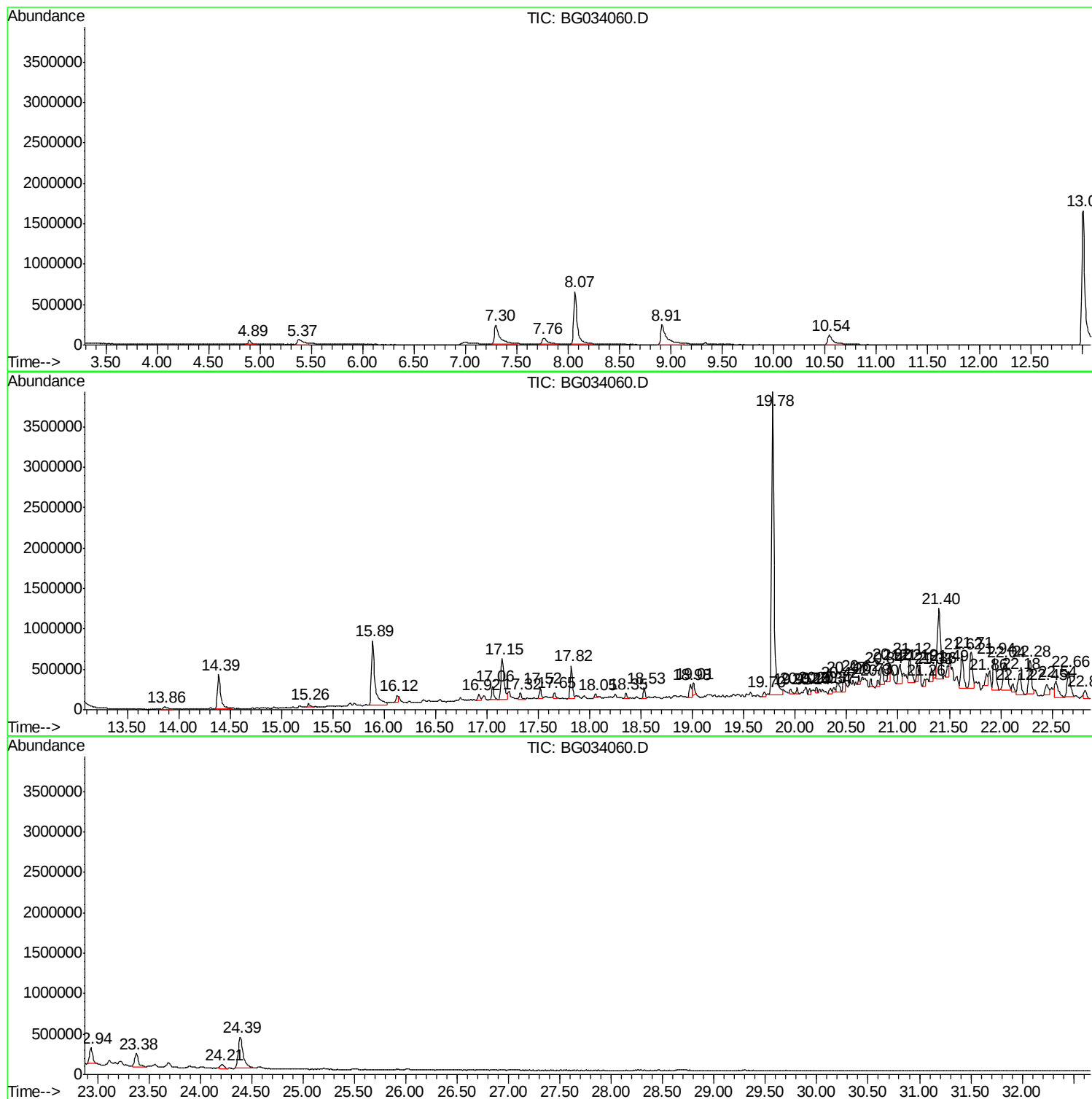
Sum of corrected areas: 36045258

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Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.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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Data File : BG034060.D  
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Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

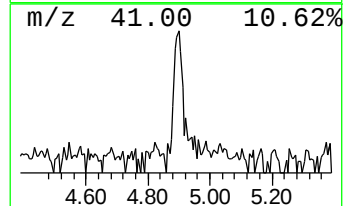
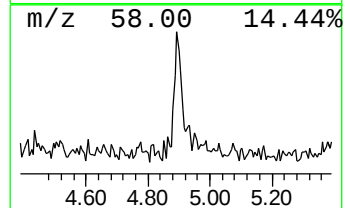
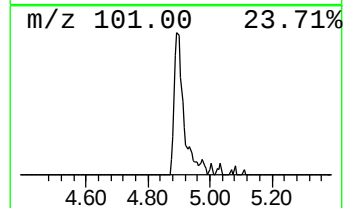
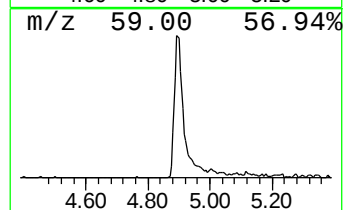
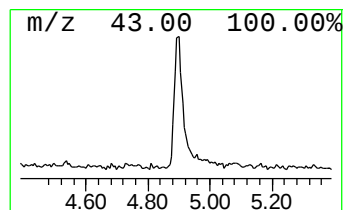
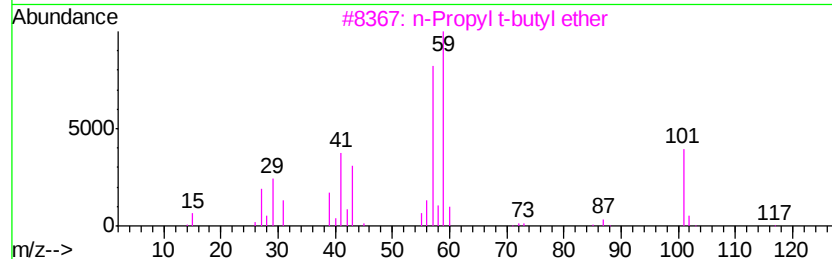
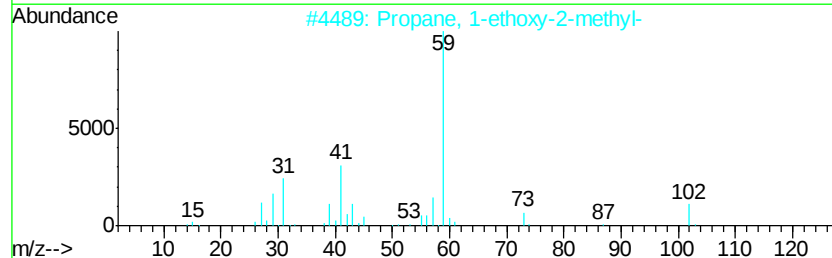
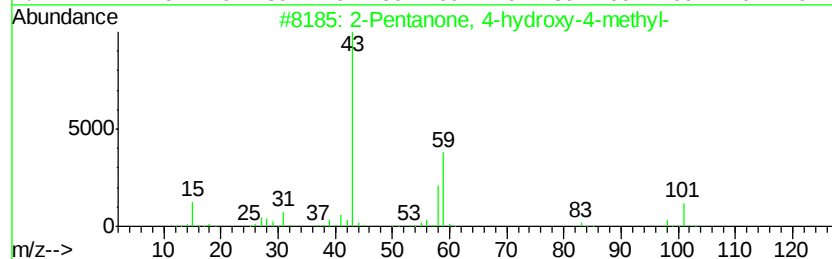
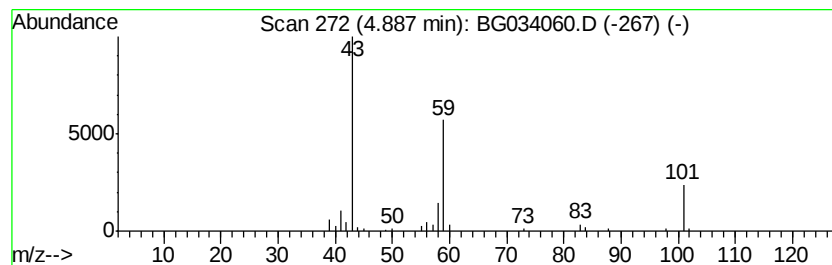
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.89 | 7.02 ng | 107115 | 1,4-Dichlorobenzene-d4 | 7.77 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#         | Qual |
|------|----|---|----------------------------------|-----|---------|--------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl- | 116 | C6H12O2 | 000123-42-2  | 50   |
| 2    |    |   | Propane, 1-ethoxy-2-methyl-      | 102 | C6H14O  | 000627-02-1  | 35   |
| 3    |    |   | n-Propyl t-butyl ether           | 116 | C7H16O  | 1000334-81-9 | 33   |
| 4    |    |   | 2-Decanol, methyl ether          | 172 | C11H24O | 1000333-83-9 | 23   |
| 5    |    |   | Morpholine, 4-methyl-            | 101 | C5H11NO | 000109-02-4  | 9    |



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Instrument :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

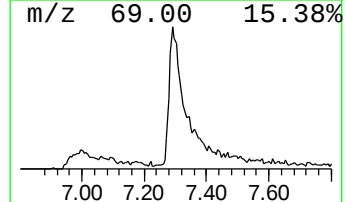
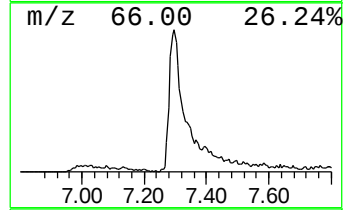
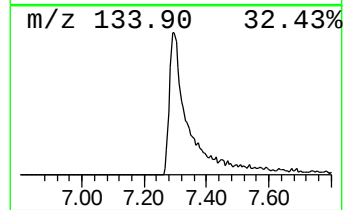
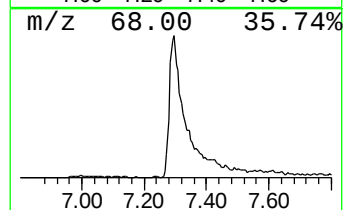
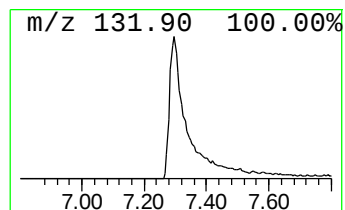
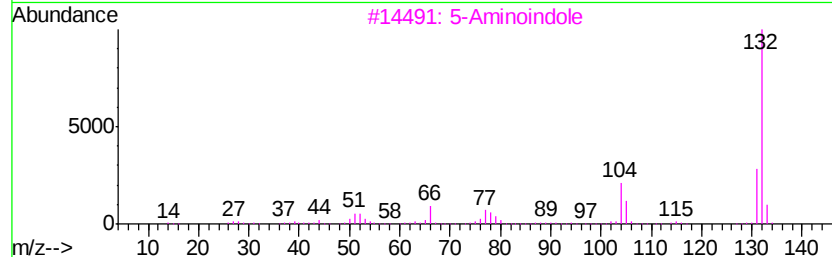
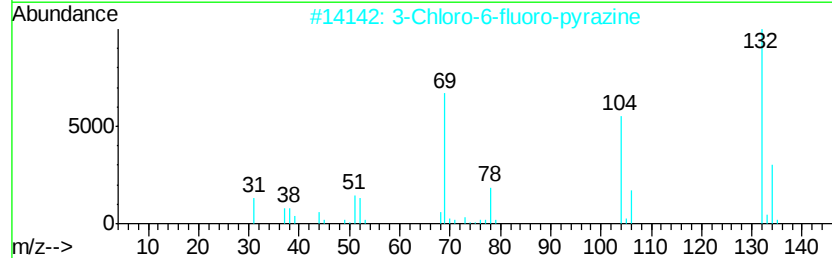
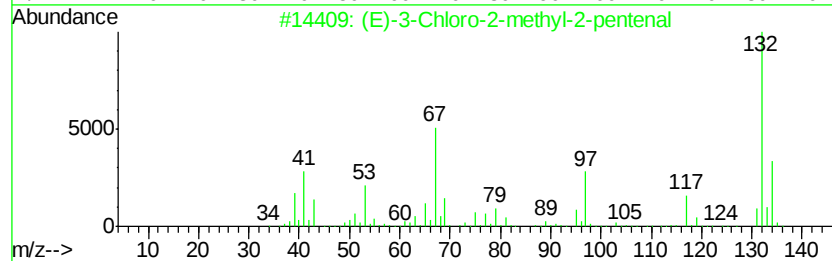
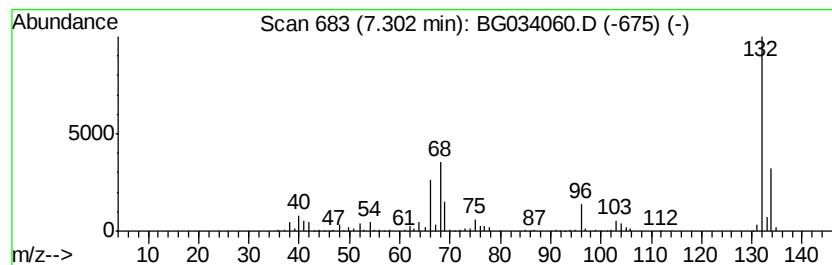
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown7.30 Concentration Rank 2

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.30 | 60.24 ng | 918664 | 1,4-Dichlorobenzene-d4 | 7.77 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 2    |    | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 16   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 10   |
| 4    |    | Amphetaminil                     | 250 | C17H18N2  | 017590-01-1  | 10   |
| 5    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |



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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

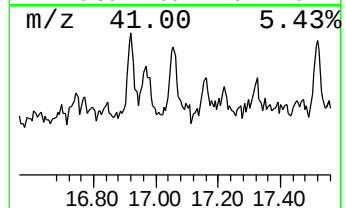
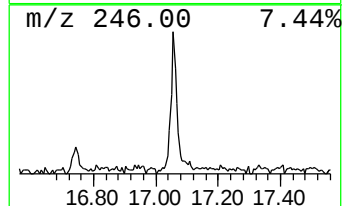
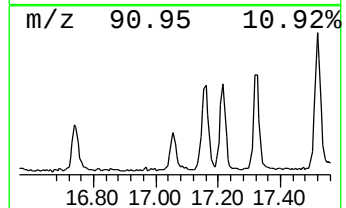
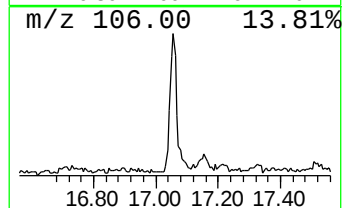
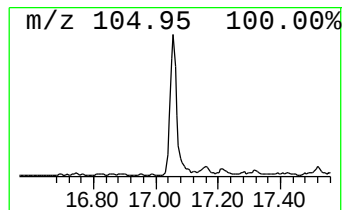
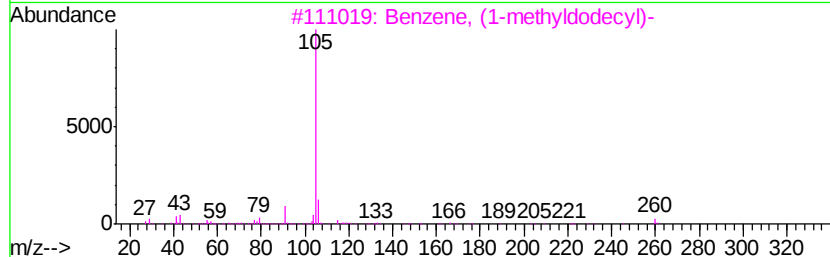
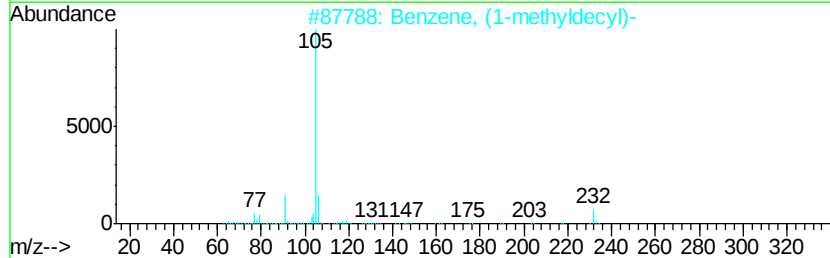
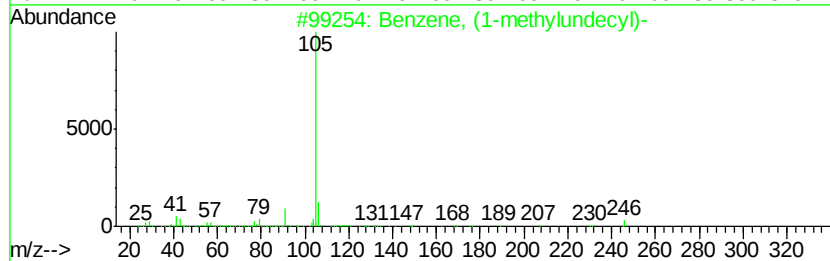
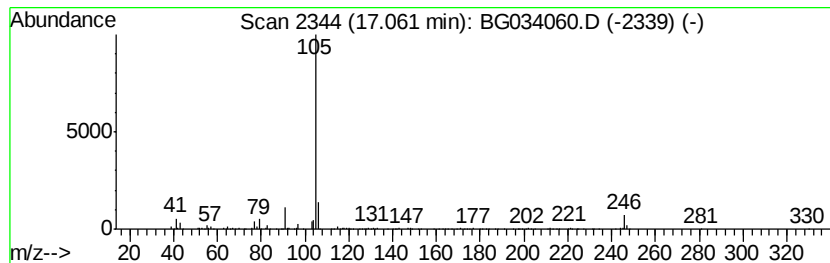
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 Benzene, (1-methylundecyl)- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.06 | 4.28 ng | 243141 | Phenanthrene-d10 | 17.14 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, (1-methylundecyl)- | 246 | C18H30  | 002719-61-1 | 93   |
| 2    |    | Benzene, (1-methyldecyl)-   | 232 | C17H28  | 004536-88-3 | 76   |
| 3    |    | Benzene, (1-methyldodecyl)- | 260 | C19H32  | 004534-53-6 | 59   |
| 4    |    | Benzene, (1-methylnonyl)-   | 218 | C16H26  | 004537-13-7 | 59   |
| 5    |    | Benzene, (1-methylheptyl)-  | 190 | C14H22  | 000777-22-0 | 59   |



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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

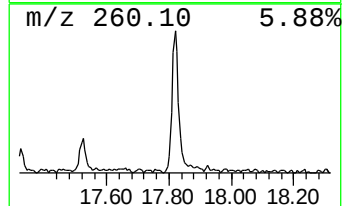
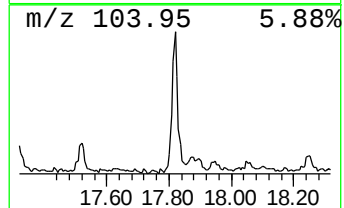
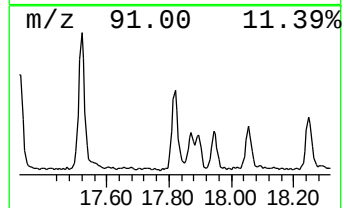
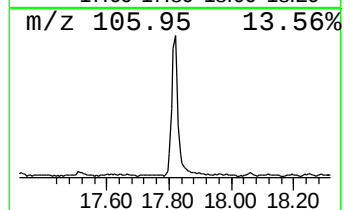
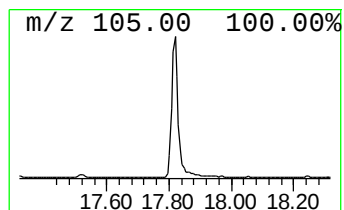
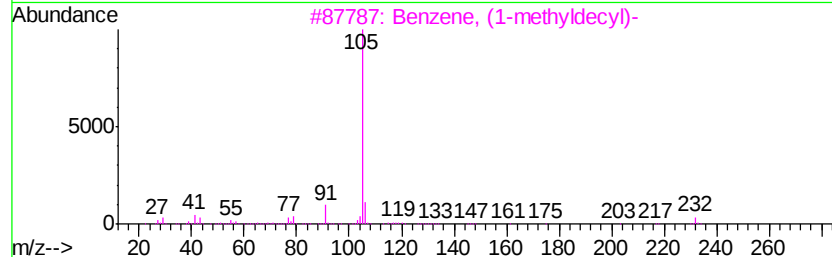
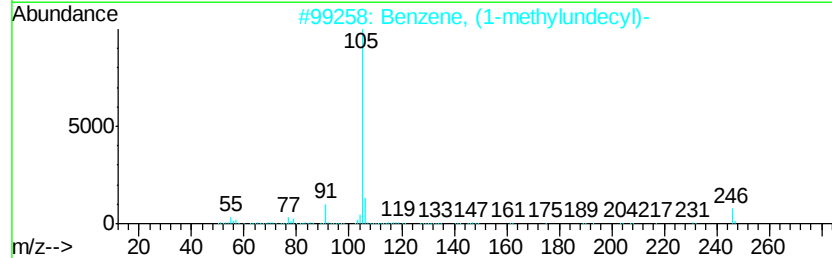
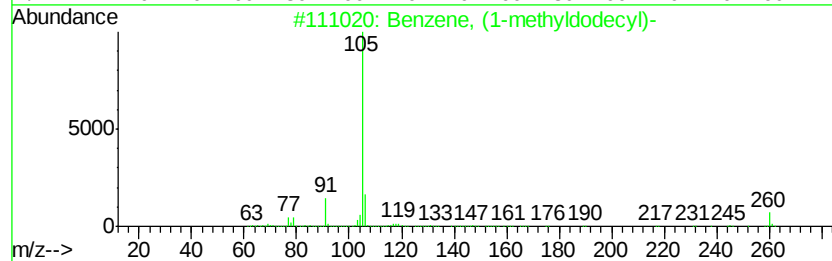
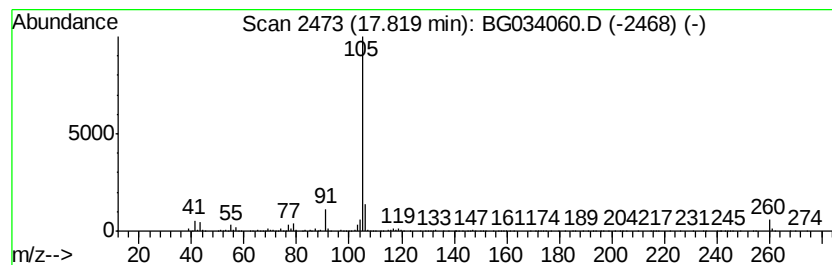
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Benzene, (1-methyldodecyl)- Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 17.82 | 11.01 ng | 625401 | Phenanthrene-d10 | 17.14 |

| Hit# of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------------|-----|---------|-------------|------|
| 1       |   | Benzene, (1-methyldodecyl)-   | 260 | C19H32  | 004534-53-6 | 94   |
| 2       |   | Benzene, (1-methylundecyl)-   | 246 | C18H30  | 002719-61-1 | 74   |
| 3       |   | Benzene, (1-methyldecyl)-     | 232 | C17H28  | 004536-88-3 | 72   |
| 4       |   | Benzene, (1-methylnonyl)-     | 218 | C16H26  | 004537-13-7 | 64   |
| 5       |   | Benzene, (1-methylnonadecyl)- | 358 | C26H46  | 002398-66-5 | 53   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

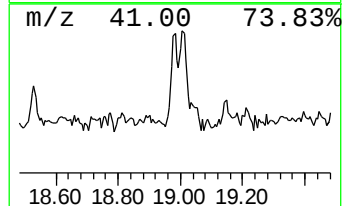
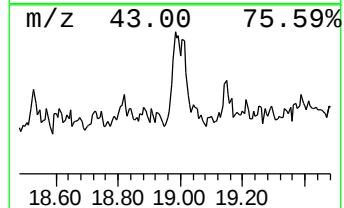
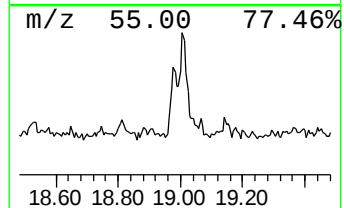
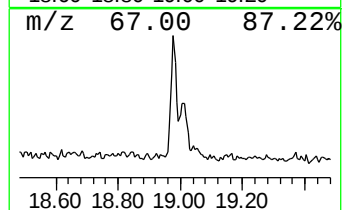
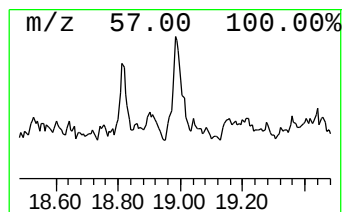
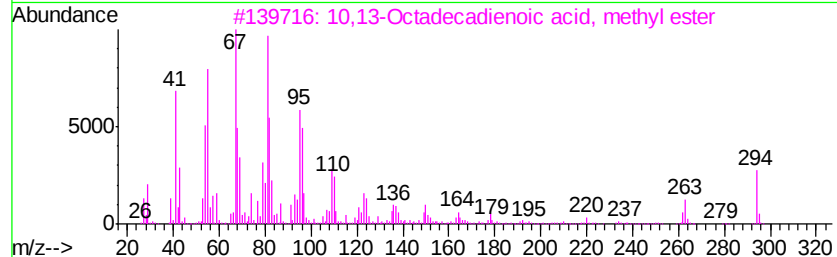
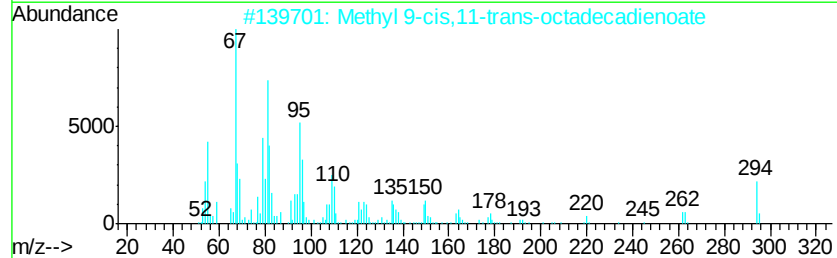
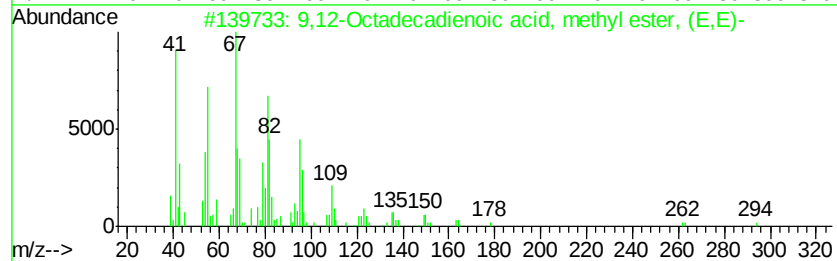
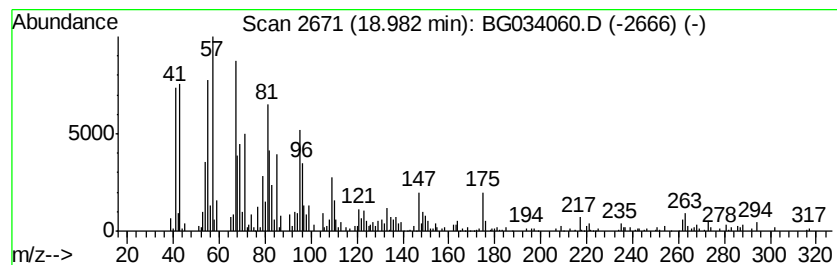
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 9,12-Octadecadienoic acid, ... Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.98 | 4.42 ng | 251349 | Phenanthrene-d10 | 17.14 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4  | 99   |
| 2    |    |   | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 1000336-44-0 | 97   |
| 3    |    |   | 10,13-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-62-2  | 97   |
| 4    |    |   | Methyl 10-trans,12-cis-octadecad... | 294 | C19H34O2 | 1000336-44-2 | 95   |
| 5    |    |   | 9,12-Octadecadienoic acid (Z,Z)-    | 280 | C18H32O2 | 000060-33-3  | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

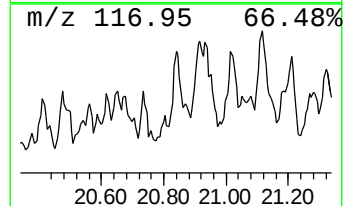
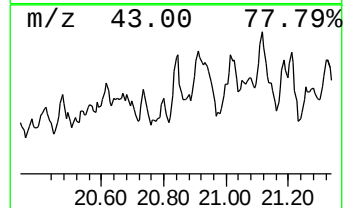
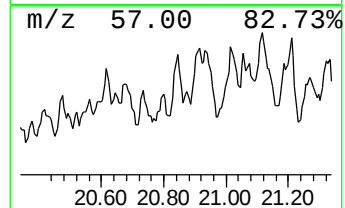
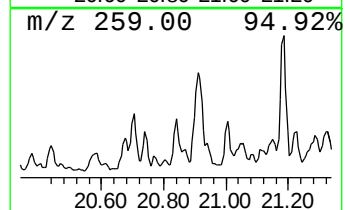
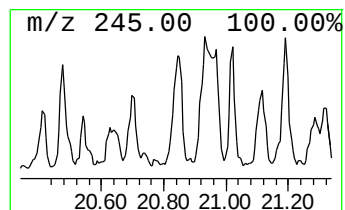
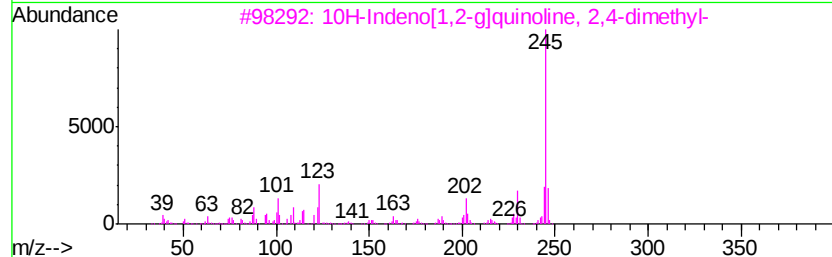
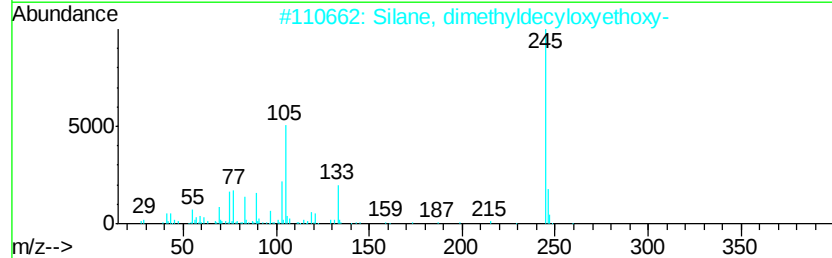
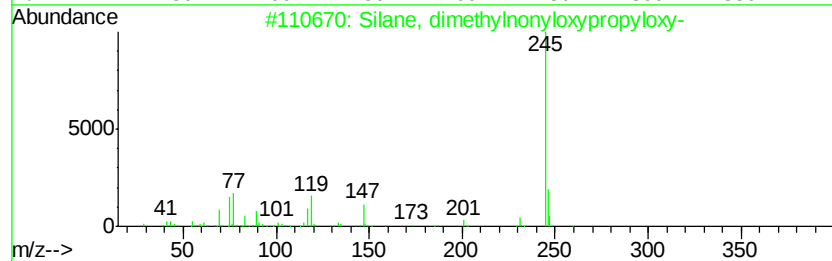
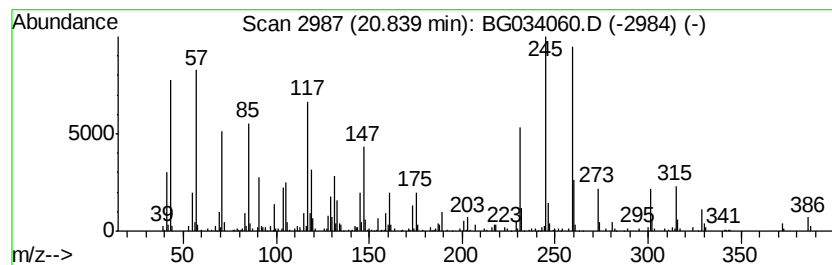
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 unknown20.84 Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.84 | 3.75 ng | 345634 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                              | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Silane, dimethylnonyloxypropyloxy-        | 260 | C14H32O2Si | 1000356-04-5 | 32   |
| 2    |    | Silane, dimethyldecyloxyethoxy-           | 260 | C14H32O2Si | 1000346-91-9 | 10   |
| 3    |    | 10H-Indeno[1,2-a]quinoline, 2,4-dimethyl- | 245 | C18H15N    | 050519-37-4  | 10   |
| 4    |    | Silane, dimethyl(2-heptyloxy)non-         | 316 | C18H40O2Si | 1000347-65-9 | 10   |
| 5    |    | Furo[3,2-c]quinoline, 2,3-dihydro-        | 245 | C14H15NO3  | 1000264-83-9 | 10   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

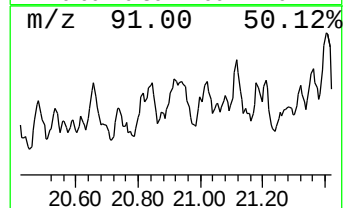
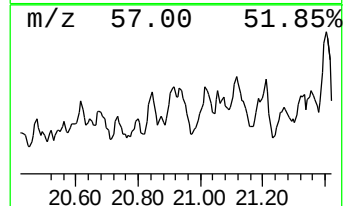
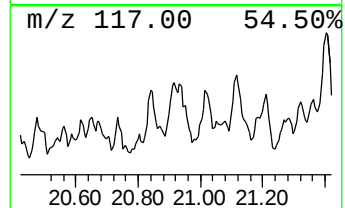
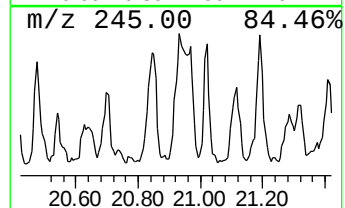
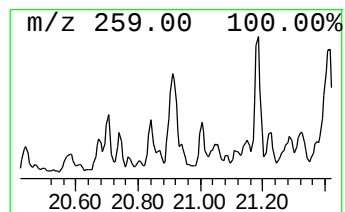
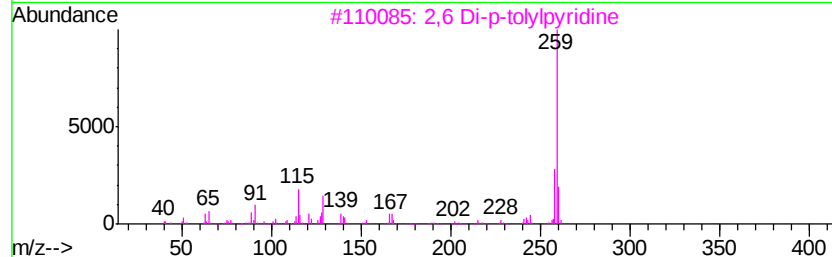
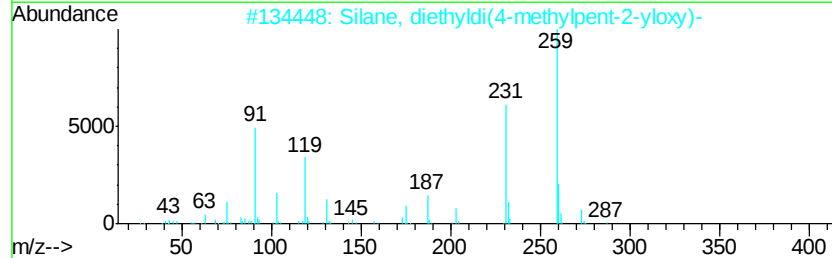
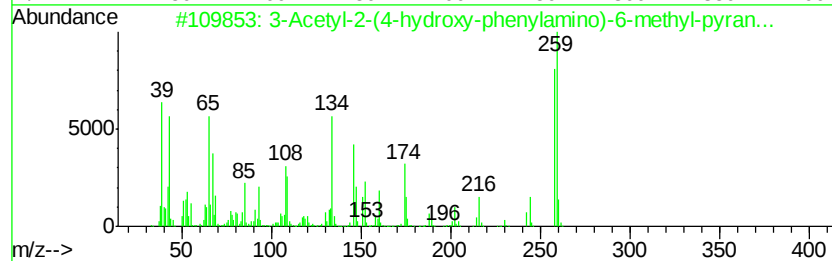
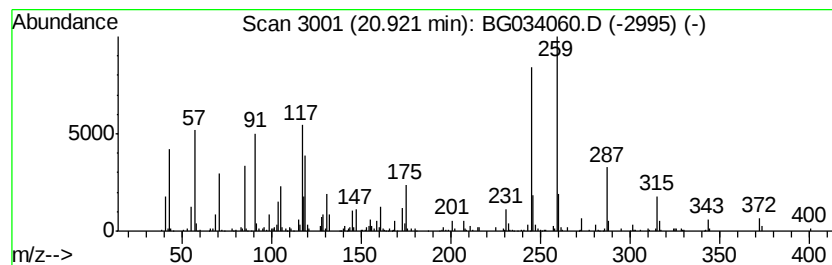
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 3-Acetyl-2-(4-hydroxy-pheny... Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.92 | 3.94 ng | 363190 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 3-Acetyl-2-(4-hydroxy-phenylamin... | 259 | C14H13NO4  | 1000300-99-5 | 80   |
| 2    |    | Silane, diethyldi(4-methylpent-2... | 288 | C16H36O2Si | 1000363-44-3 | 25   |
| 3    |    | 2,6 Di-p-tolylpyridine              | 259 | C19H17N    | 014435-88-2  | 22   |
| 4    |    | Cyclohex-3-ene-1-carboxylic acid... | 259 | C15H17NO3  | 311316-40-2  | 22   |
| 5    |    | Silane, dimethyl(4-heptyloxy)oct... | 302 | C17H38O2Si | 1000347-60-3 | 17   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

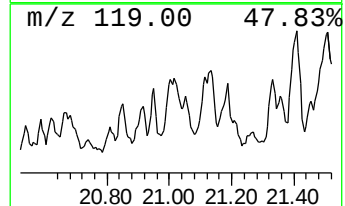
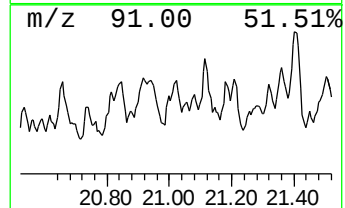
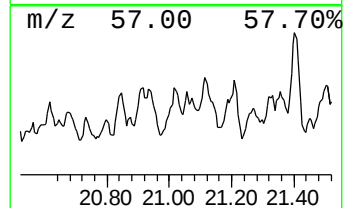
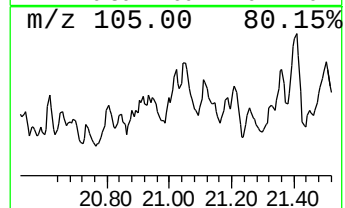
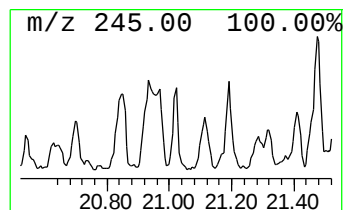
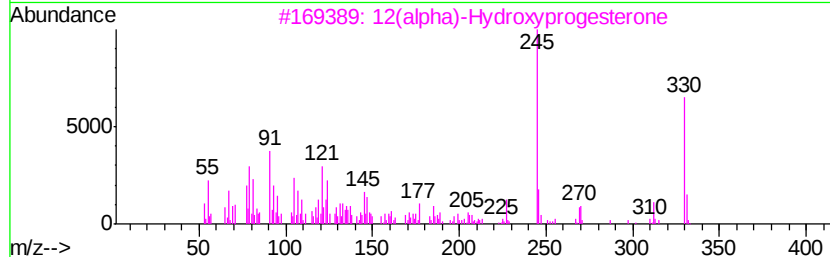
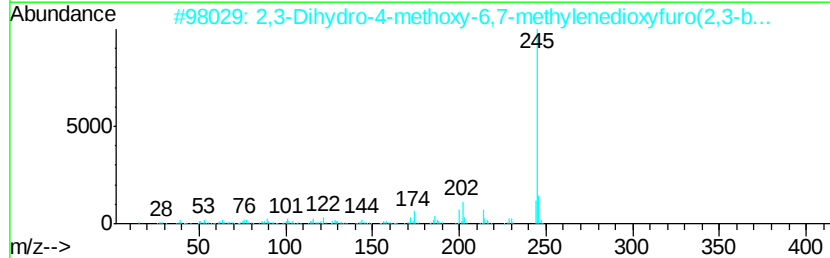
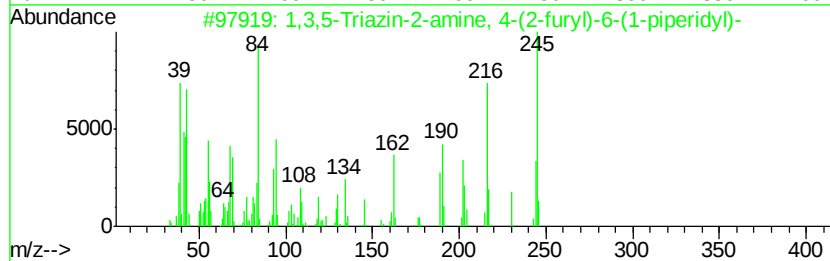
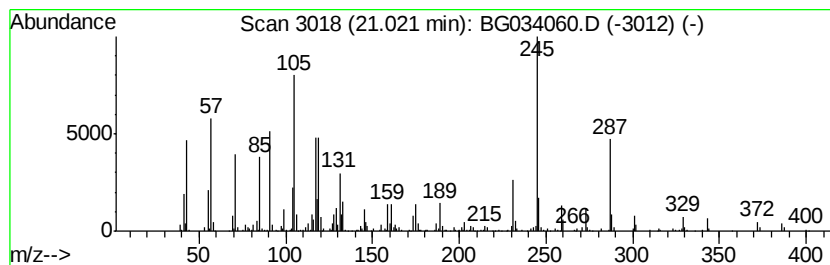
Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 1,3,5-Triazin-2-amine, 4-(2... Concentration Rank 16

| R.T.      | EstConc                              | Area   | Relative to ISTD | R.T.         |      |
|-----------|--------------------------------------|--------|------------------|--------------|------|
| 21.02     | 5.05 ng                              | 466466 | Chrysene-d12     | 21.39        |      |
| Hit# of 5 | Tentative ID                         | MW     | MolForm          | CAS#         | Qual |
| 1         | 1,3,5-Triazin-2-amine, 4-(2-furyl... | 245    | C12H15N5O        | 148403-53-6  | 53   |
| 2         | 2,3-Dihydro-4-methoxy-6,7-methyl...  | 245    | C13H11N04        | 094521-55-8  | 15   |
| 3         | 12(alpha)-Hvdroxvprodesterone        | 330    | C21H30O3         | 019897-02-0  | 14   |
| 4         | 2,3-Dimethyl-5-oxo-1-phenyl-3-pv...  | 245    | C12H11N3OS       | 091397-05-6  | 14   |
| 5         | .beta.-Methylfentanyl                | 350    | C23H30N2O        | 1000248-46-0 | 14   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

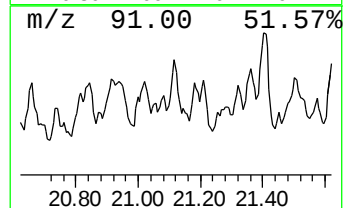
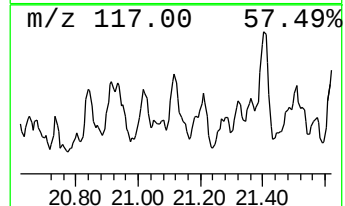
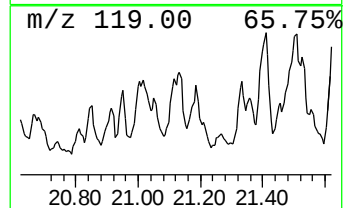
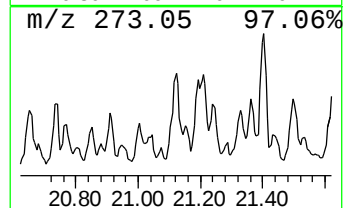
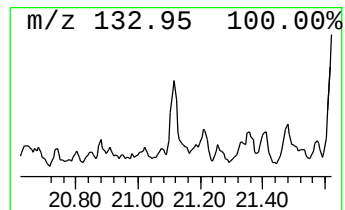
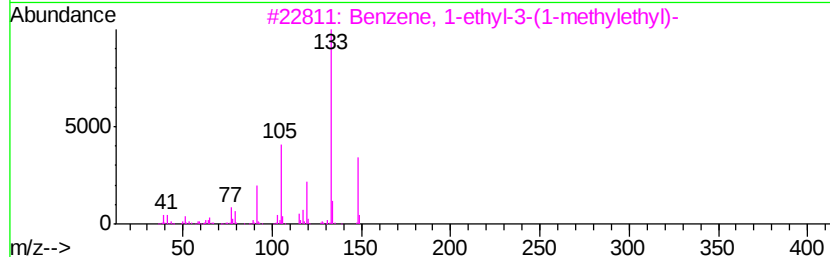
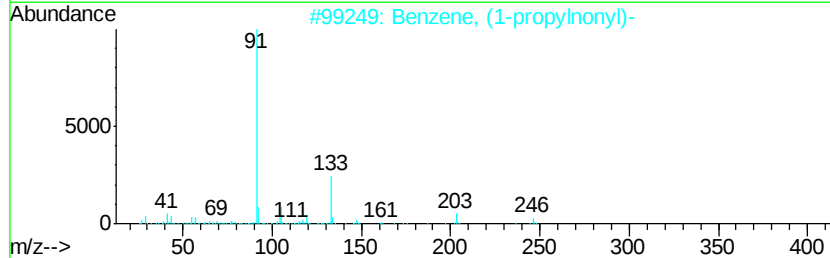
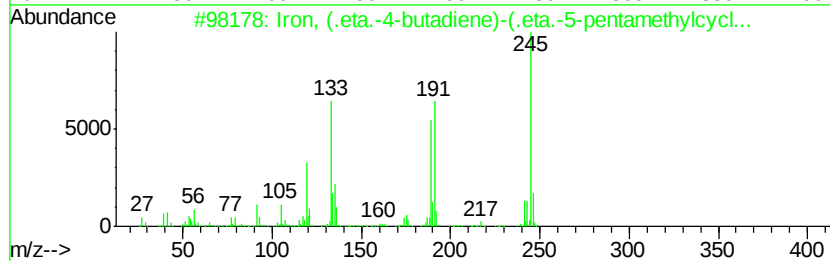
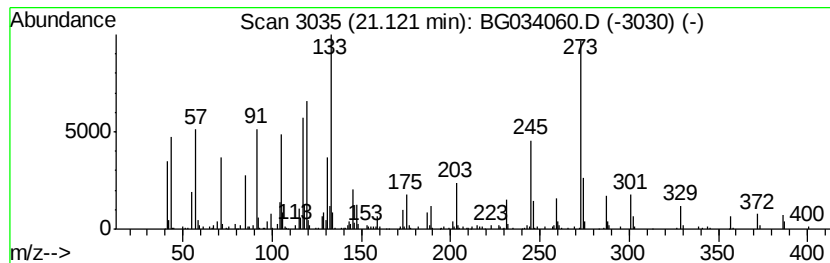
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown21.12 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.12 | 6.07 ng | 560588 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Iron, (.eta.-4-butadiene)-(.eta.... | 245 | C14H21Fe   | 1000161-69-5 | 38   |
| 2    |    | Benzene, (1-propylnonyl)-           | 246 | C18H30     | 002719-64-4  | 18   |
| 3    |    | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16     | 004920-99-4  | 14   |
| 4    |    | 1-Bromo-1,4,4a,5,8,8a-hexahydron... | 212 | C10H13Br   | 1000192-97-4 | 14   |
| 5    |    | 2-Methyl-4-(4-chlorophenyl)-2,3-... | 287 | C16H14ClNS | 078031-24-0  | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
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Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

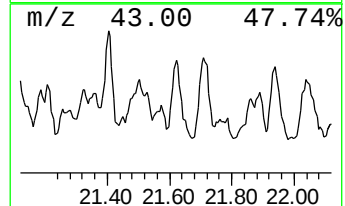
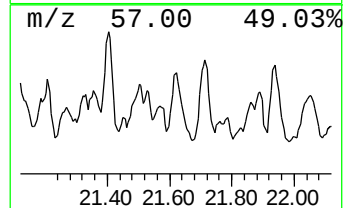
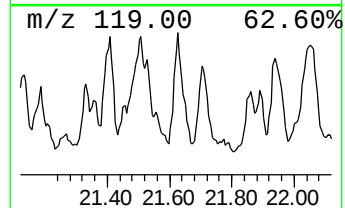
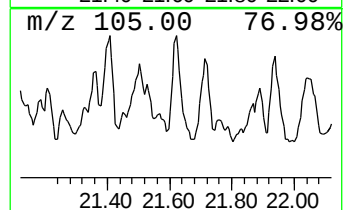
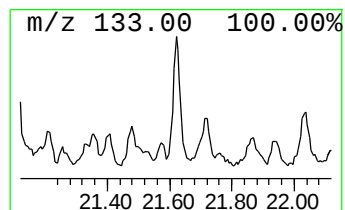
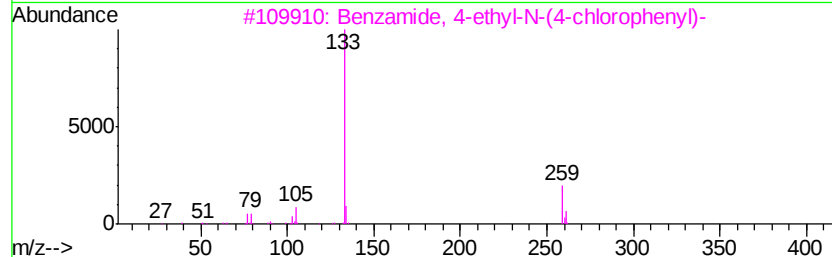
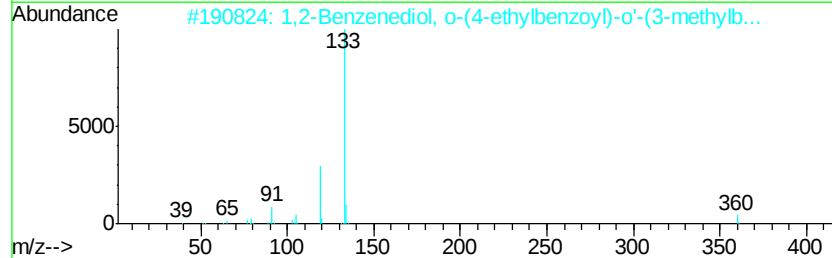
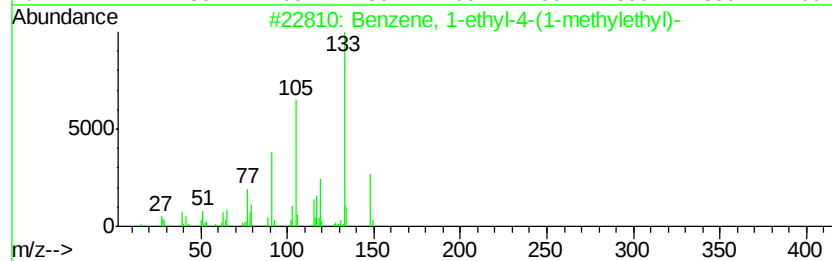
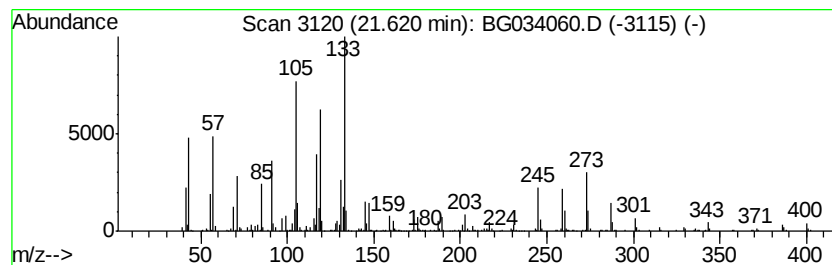
Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 unknown21.62 Concentration Rank 6

| R.T.    | EstConc | Area                                | Relative to ISTD | R.T.            |
|---------|---------|-------------------------------------|------------------|-----------------|
| 21.62   | 9.37 ng | 864375                              | Chrysene-d12     | 21.39           |
| Hit# of | 5       | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |         | Benzene, 1-ethyl-4-(1-methylethyl)- | 148 C11H16       | 004218-48-8 35  |
| 2       |         | 1,2-Benzenediol, o-(4-ethylbenzo... | 360 C23H20O4     | 1000325-97-4 35 |
| 3       |         | Benzamide, 4-ethyl-N-(4-chloroph... | 259 C15H14ClNO   | 1000340-35-1 22 |
| 4       |         | 1H-Inden-2-amine, 2,3-dihydro-      | 133 C9H11N       | 002975-41-9 22  |
| 5       |         | 1-Bromo-1,4,4a,5,8,8a-hexahydron... | 212 C10H13Br     | 1000192-97-4 22 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

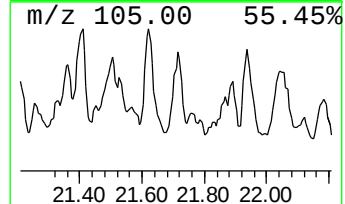
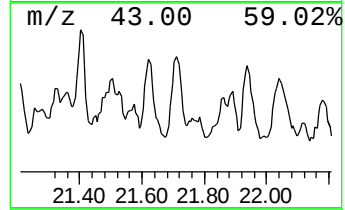
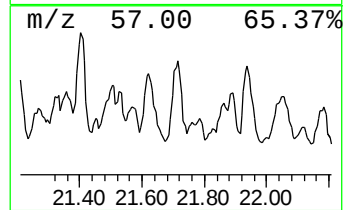
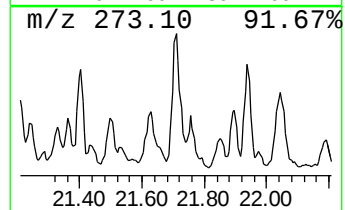
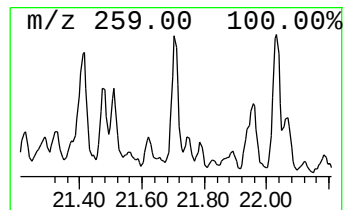
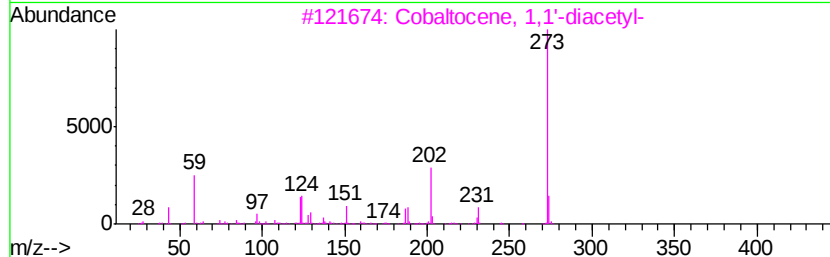
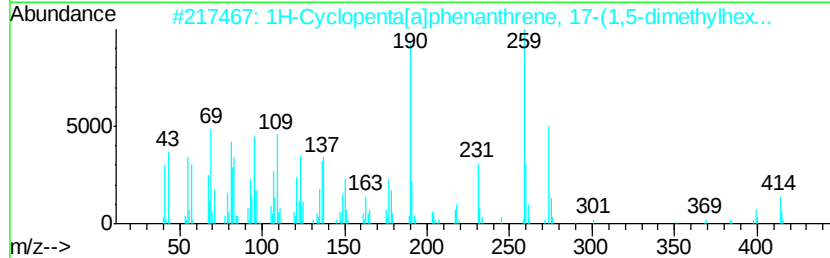
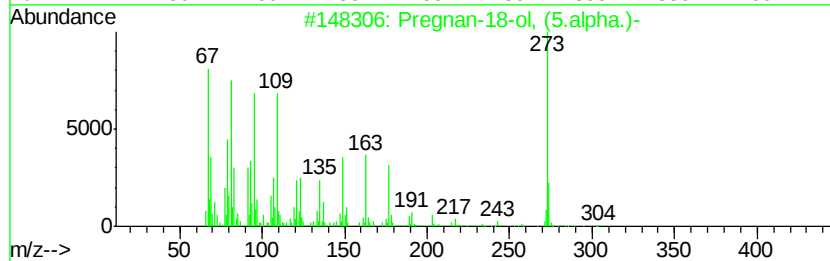
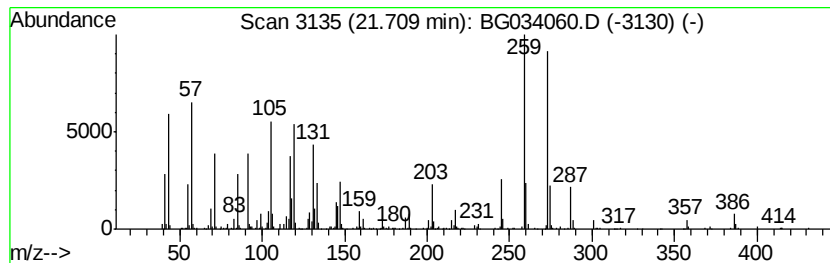
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Pregnan-18-ol, (5.alpha.)- Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 21.71 | 10.04 ng | 926051 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                                 | MW  | MolForm    | CAS#         | Qual |
|------|----|----------------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Pregnan-18-ol, (5.alpha.)-                   | 304 | C21H36O    | 056053-09-9  | 56   |
| 2    |    | 1H-Cyclopenta[ <i>a</i> ]phenanthrene, 17... | 414 | C30H54     | 000474-20-4  | 15   |
| 3    |    | Cobaltocene, 1,1'-diacetyl-                  | 273 | C14H14CoO2 | 073249-54-4  | 15   |
| 4    |    | 1H-1,3-Benzimidazole-1-acetamide...          | 287 | C17H25N3O  | 1000351-15-0 | 15   |
| 5    |    | Tolcapone                                    | 273 | C14H11NO5  | 134308-13-7  | 14   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

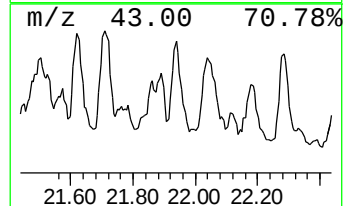
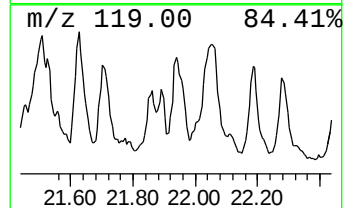
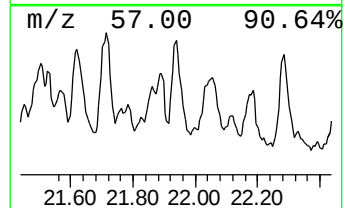
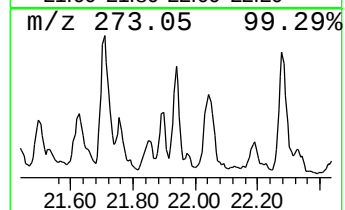
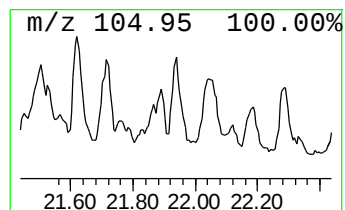
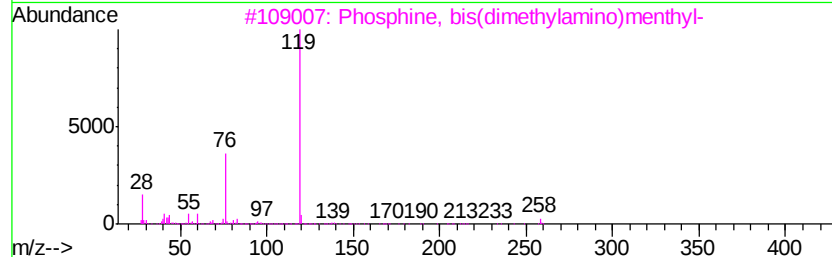
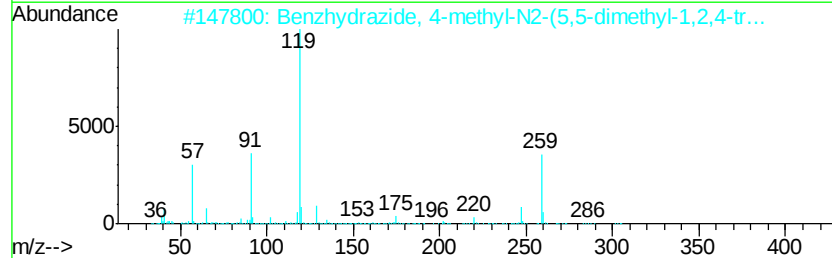
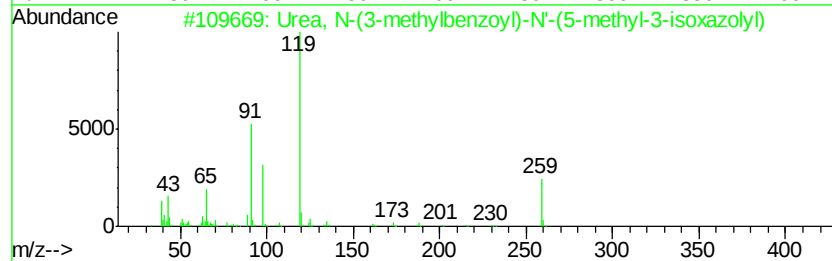
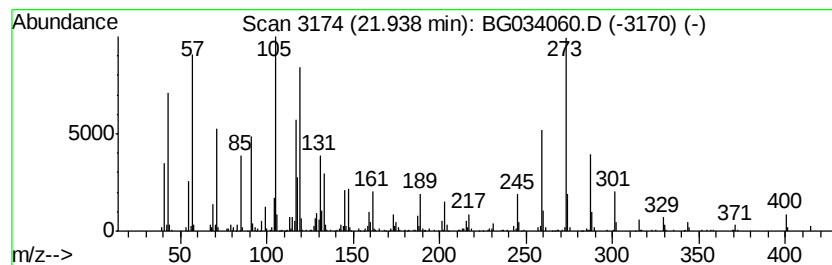
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 unknown21.94 Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.94 | 8.50 ng | 784284 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Urea, N-(3-methylbenzoyl)-N'-(5-... | 259 | C13H13N3O3 | 1000320-00-2 | 14   |
| 2    |    | Benzhydrazide, 4-methyl-N2-(5,5-... | 304 | C16H20N2O4 | 1000264-15-6 | 14   |
| 3    |    | Phosphine, bis(dimethylamino)men... | 258 | C14H31N2P  | 1000150-77-6 | 11   |
| 4    |    | Cobaltocene, 1,1'-diacetyl-         | 273 | C14H14CoO2 | 073249-54-4  | 11   |
| 5    |    | N-Methylpiperidino[2,4-b,c]1,2,3... | 273 | C18H27NO   | 1000128-54-7 | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

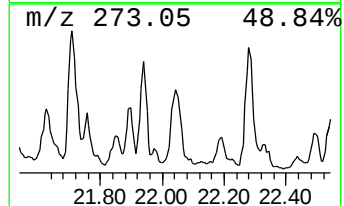
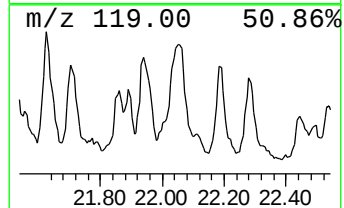
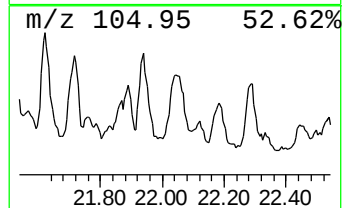
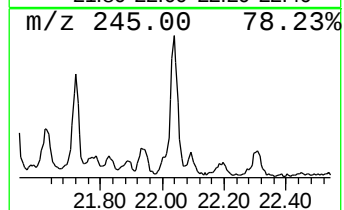
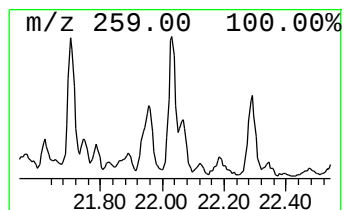
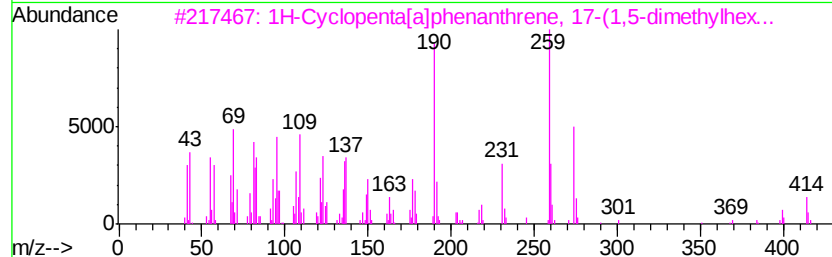
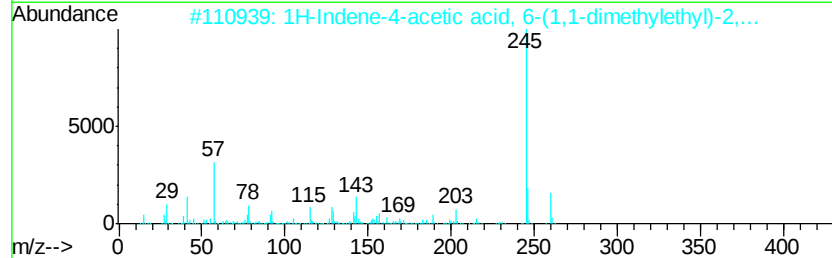
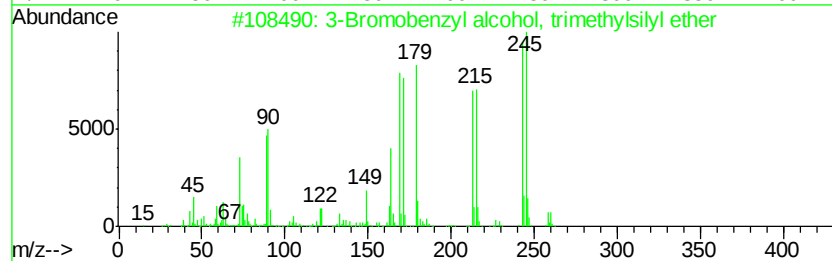
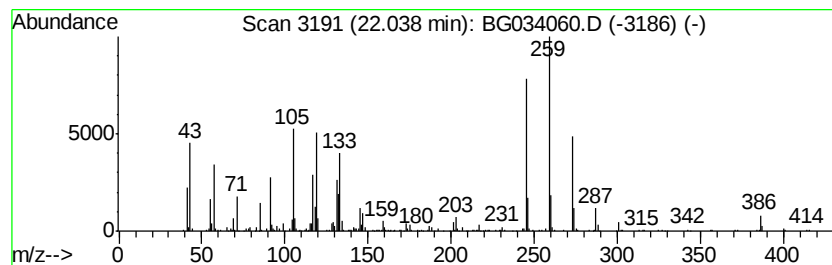
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 unknown22.04 Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 22.04 | 10.72 ng | 988804 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 3-Bromobenzyl alcohol, trimethyl... | 258 | C10H15BrOSi | 1000364-25-3 | 25   |
| 2    |    | 1H-Indene-4-acetic acid, 6-(1,1-... | 260 | C17H24O2    | 055591-05-4  | 18   |
| 3    |    | 1H-Cyclopentafalphenanthrene, 17... | 414 | C30H54      | 000474-20-4  | 15   |
| 4    |    | Propanamide, N-(1-ethyl-1,2,3,4-... | 274 | C17H26N2O   | 063134-09-8  | 15   |
| 5    |    | 2-Phenoxy-2,2'-(3H,3'H)-spiro[1,... | 338 | C18H15N2O3P | 052961-94-1  | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

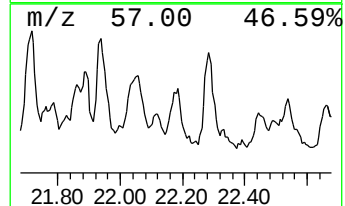
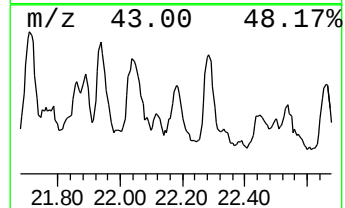
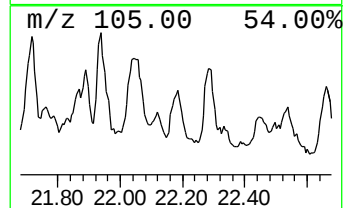
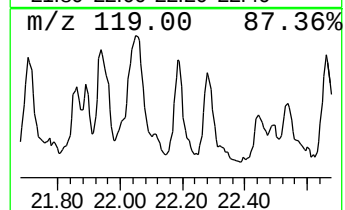
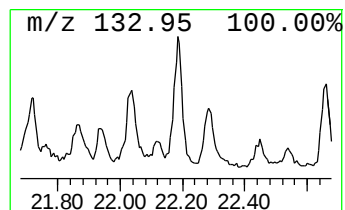
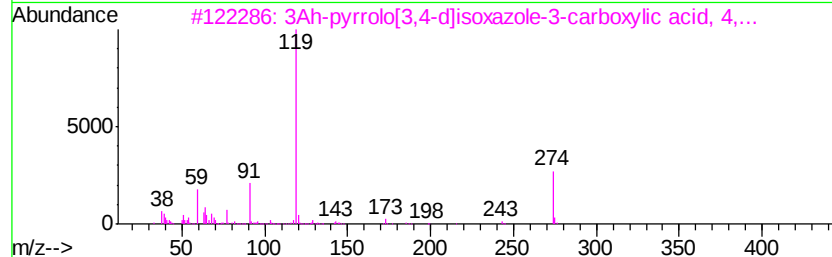
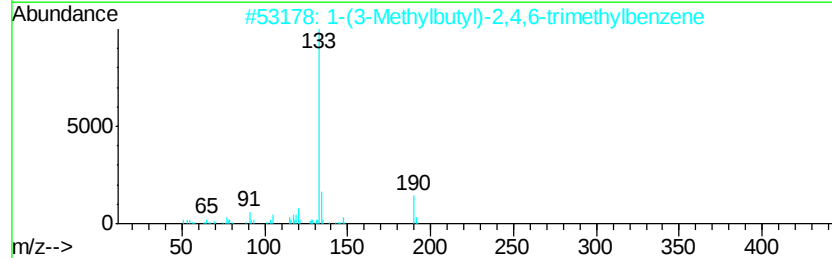
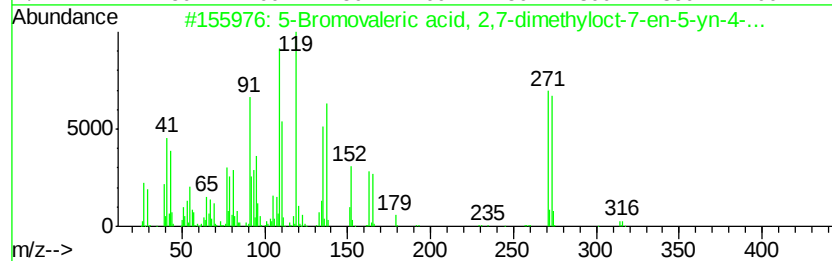
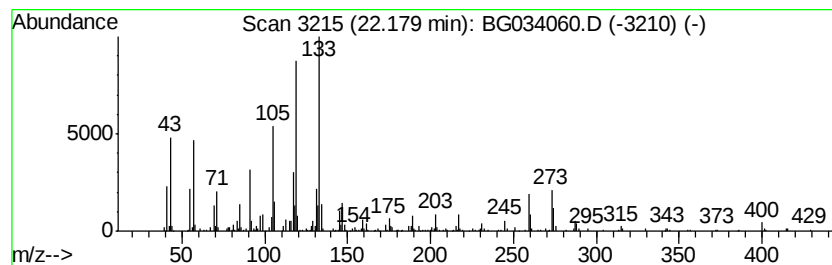
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 unknown22.18 Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.18 | 6.68 ng | 616025 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 5-Bromovaleric acid, 2,7-dimethy... | 314 | C15H23BrO2 | 1000292-56-8 | 30   |
| 2    |    | 1-(3-Methylbutyl)-2,4,6-trimethy... | 190 | C14H22     | 016204-65-2  | 27   |
| 3    |    | 3Ah-pyrrolo[3,4-d]isoxazole-3-ca... | 274 | C13H10N2O5 | 1000317-17-8 | 22   |
| 4    |    | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14     | 000874-41-9  | 18   |
| 5    |    | Benzene, (1,1,4,6,6-pentamethylh... | 246 | C18H30     | 055134-07-1  | 18   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

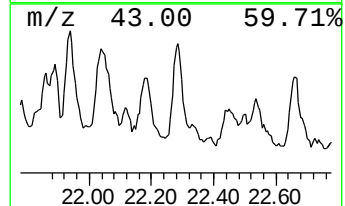
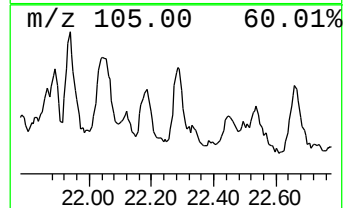
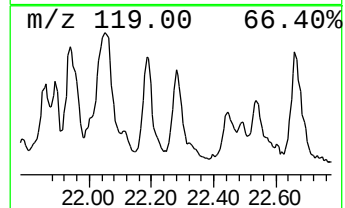
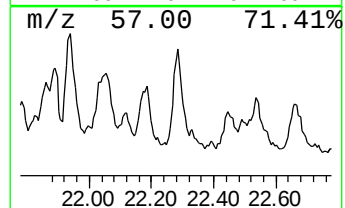
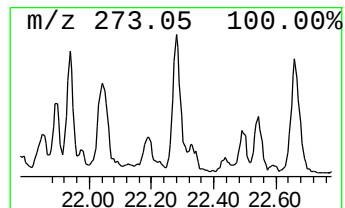
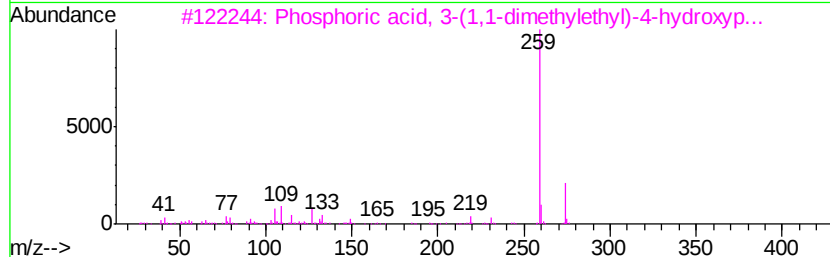
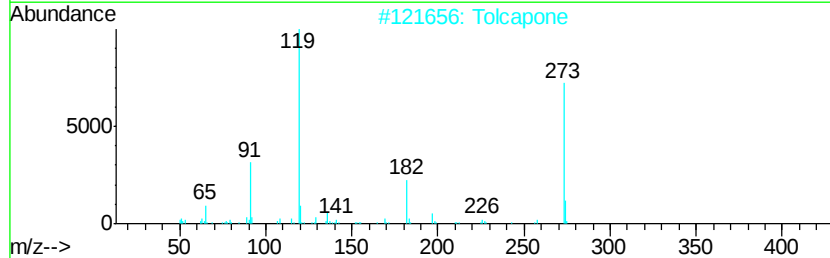
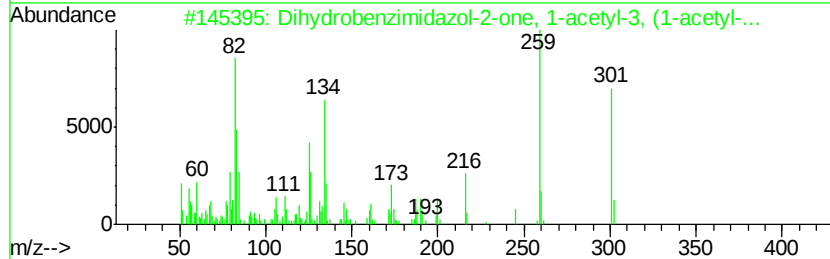
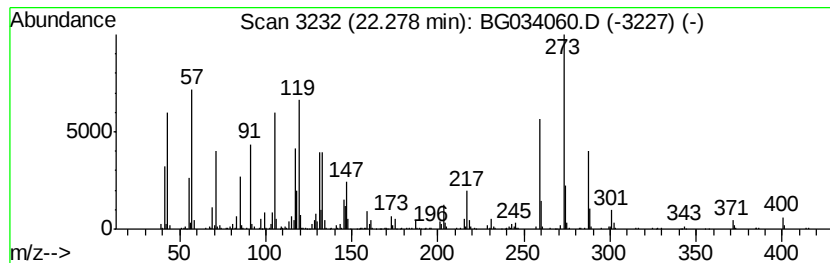
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 unknown22.28 Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.28 | 8.86 ng | 817336 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Dihydrobenzimidazol-2-one, 1-ace... | 301 | C16H19N3O3 | 1000280-84-2 | 25   |
| 2    |    | Tolcapone                           | 273 | C14H11NO5  | 134308-13-7  | 22   |
| 3    |    | Phosphoric acid, 3-(1,1-dimethyl... | 274 | C12H19O5P  | 094420-61-8  | 11   |
| 4    |    | 2,6 Di-p-tolylpyridine              | 259 | C19H17N    | 014435-88-2  | 11   |
| 5    |    | 1H-Cyclopenta[a]phenanthrene, 17... | 414 | C30H54     | 000474-20-4  | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

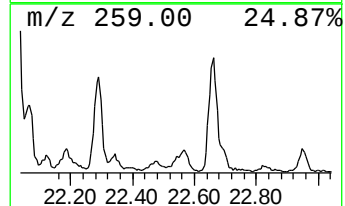
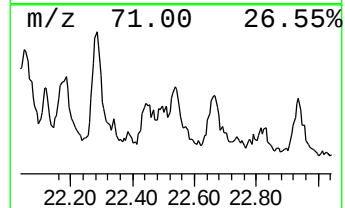
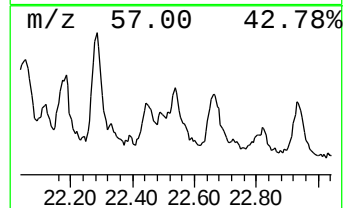
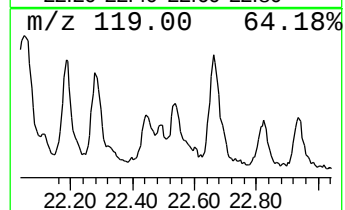
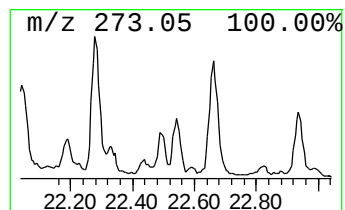
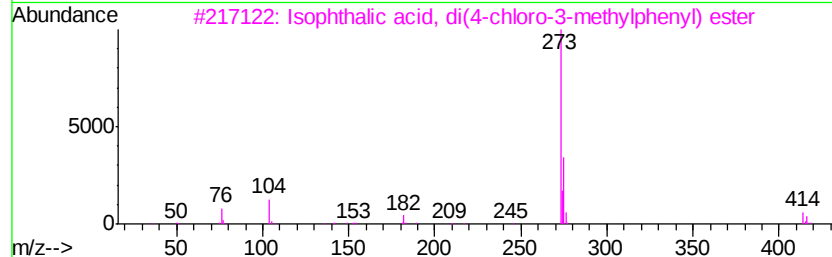
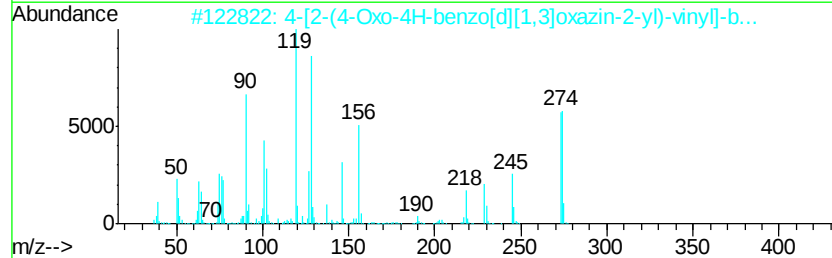
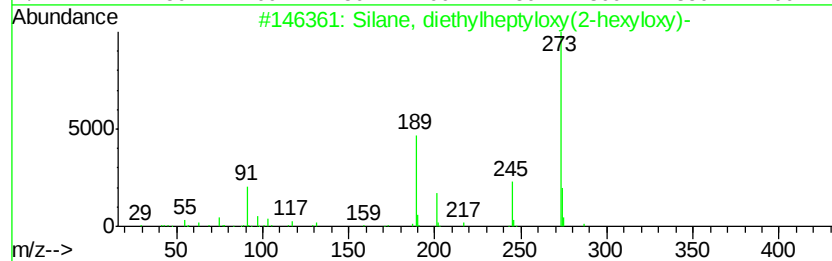
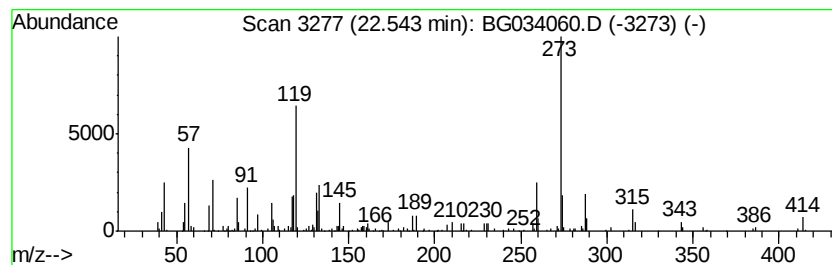
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 unknown22.54 Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.54 | 5.12 ng | 472206 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|--------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Silane, diethylheptyloxy(2-hexyl...  | 302 | C17H38O2Si  | 1000363-50-2 | 27   |
| 2    |    | 4-[2-(4-Oxo-4H-benzo[d][1,3]oxaz...  | 274 | C17H10N2O2  | 1000297-20-9 | 27   |
| 3    |    | Isophthalic acid, di(4-chloro-3-...  | 414 | C22H16Cl2O4 | 1000344-59-8 | 22   |
| 4    |    | 1H-Pyrazolo[3,4-b]quinoline, 3,6...  | 273 | C18H15N3    | 1000302-05-0 | 18   |
| 5    |    | Octahydroxanthene-1,9-dione, 3,3,... | 358 | C23H34O3    | 071827-88-8  | 18   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

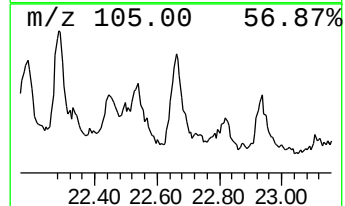
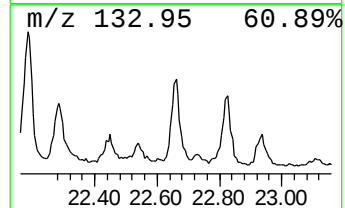
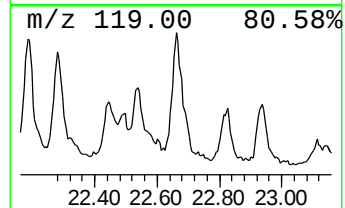
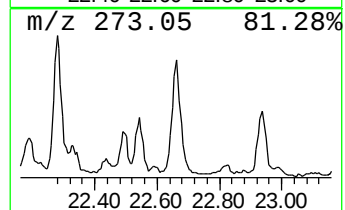
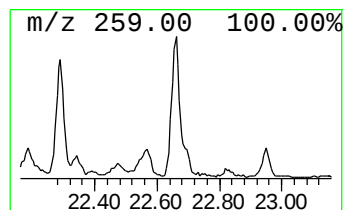
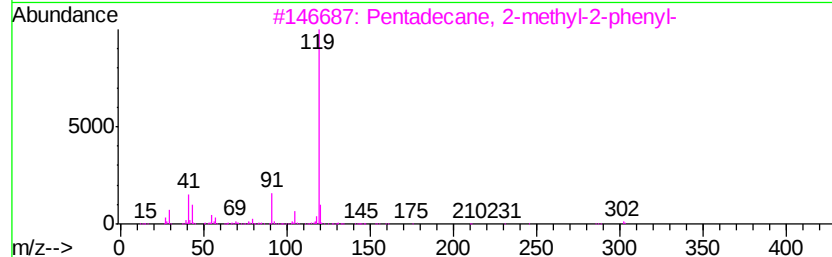
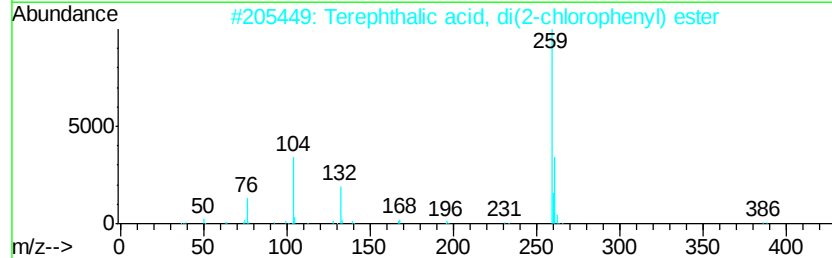
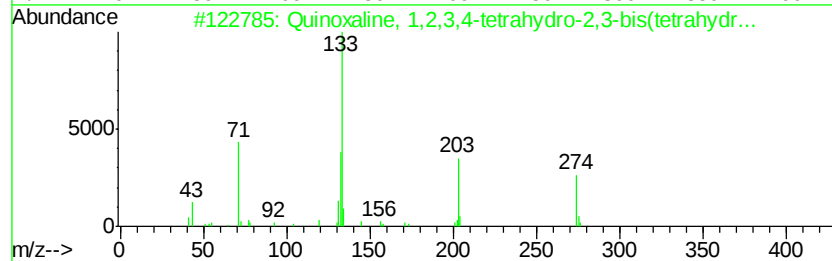
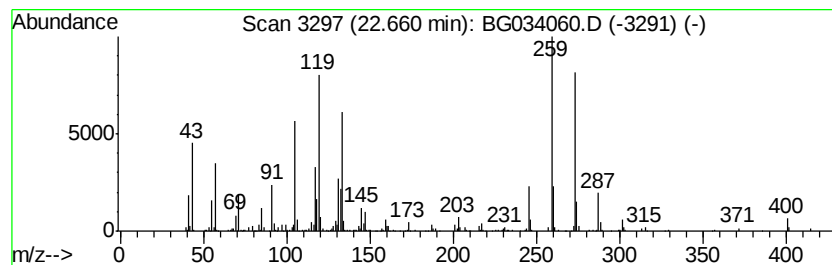
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown22.66 Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.66 | 7.57 ng | 698481 | Chrysene-d12     | 21.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Quinoxaline, 1,2,3,4-tetrahydro-... | 274 | C16H22N2O2  | 018503-40-7  | 15   |
| 2    |    | Terephthalic acid, di(2-chloroph... | 386 | C20H12Cl2O4 | 1000324-04-3 | 14   |
| 3    |    | Pentadecane, 2-methyl-2-phenyl-     | 302 | C22H38      | 029138-94-1  | 11   |
| 4    |    | Benz[clacridine-7-carboxylic acid   | 273 | C18H11NO2   | 034623-43-3  | 11   |
| 5    |    | Benzamide, N-(3-chlorophenyl)-3-... | 245 | C14H12ClNO  | 1000307-11-4 | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

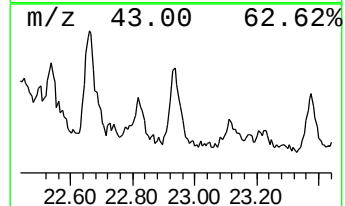
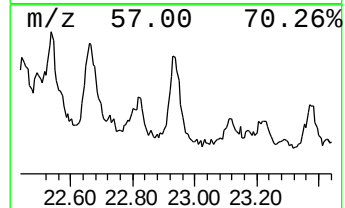
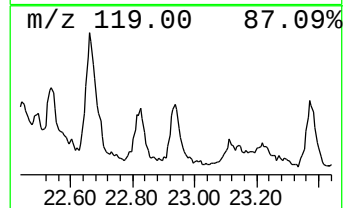
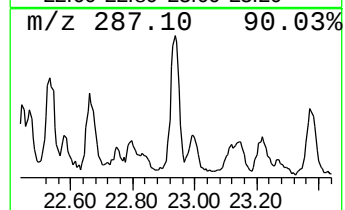
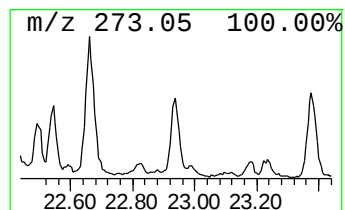
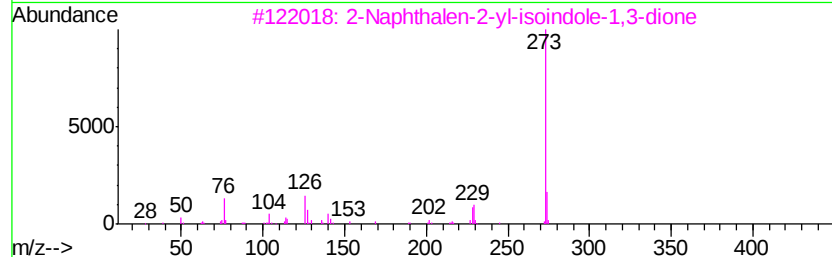
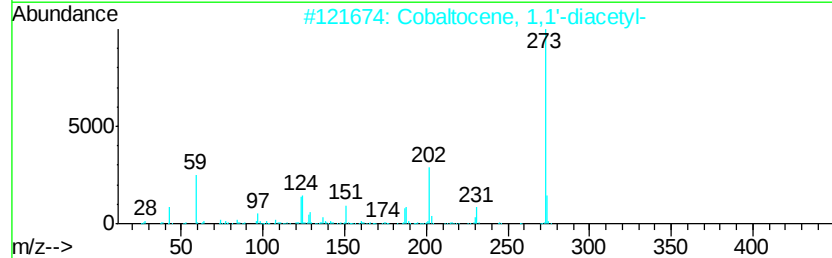
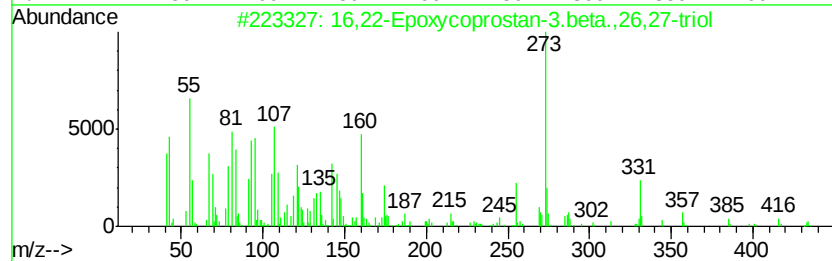
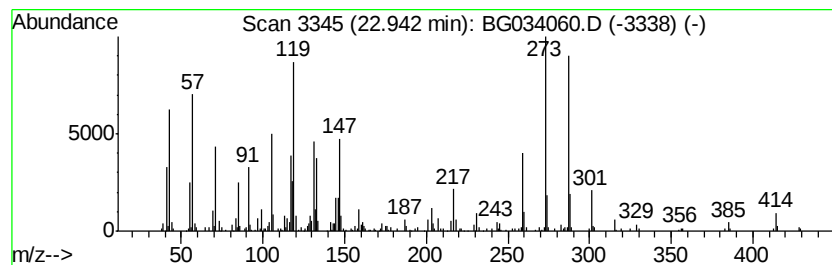
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown22.94 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 22.94 | 5.83 ng | 389075 | Perylene-d12     | 24.39 |

| Hit# | of                 | Tentative ID             | MW  | MolForm    | CAS#         | Qual |
|------|--------------------|--------------------------|-----|------------|--------------|------|
| 1    | 16,22-Epoxy        | coprostan-3.beta.,26,... | 434 | C27H46O4   | 122337-91-1  | 25   |
| 2    | Cobaltocene,       | 1,1'-diacetyl-           | 273 | C14H14CoO2 | 073249-54-4  | 11   |
| 3    | 2-Naphthalen-      | 2-yl-isoindole-1,3-...   | 273 | C18H11NO2  | 1000310-57-1 | 11   |
| 4    | N-Methylpiperidino | [2,4-b,c]1,2,3...        | 273 | C18H27NO   | 1000128-54-7 | 11   |
| 5    | Octahydroxanthene- | 1,9-dione, 3,3,...       | 316 | C20H28O3   | 030038-63-2  | 11   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG042718\  
Data File : BG034060.D  
Acq On : 28 Apr 2018 5:48  
Operator : SJ/JU  
Sample : J2571-13  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

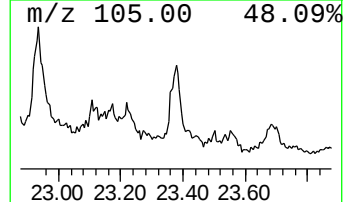
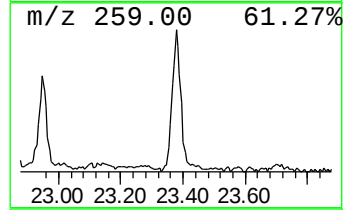
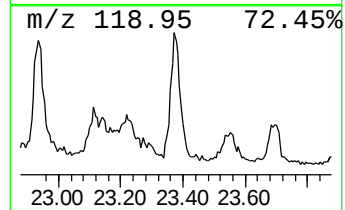
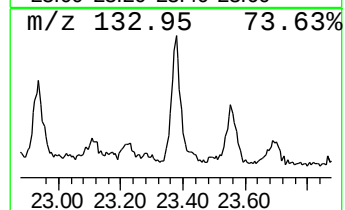
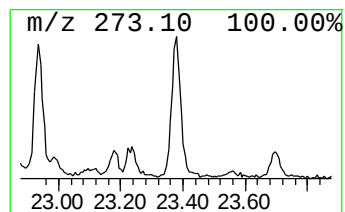
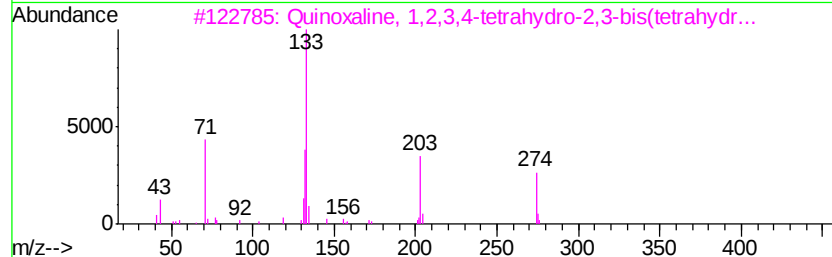
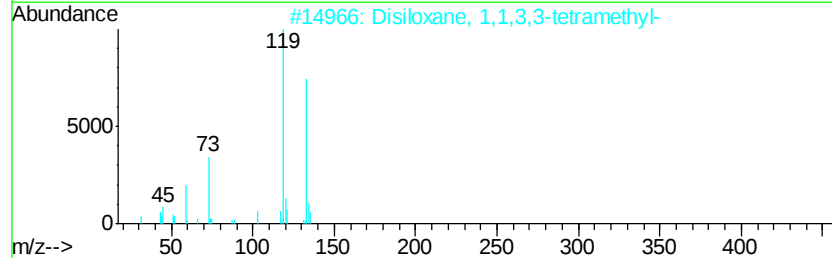
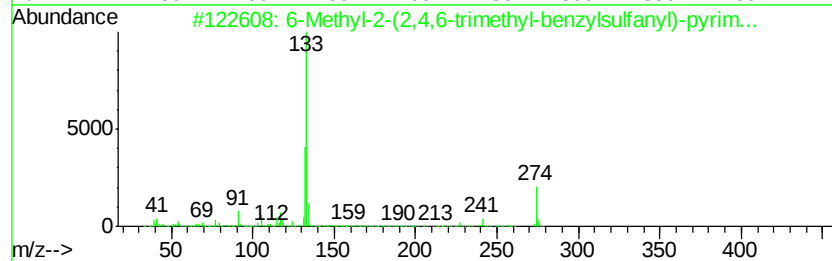
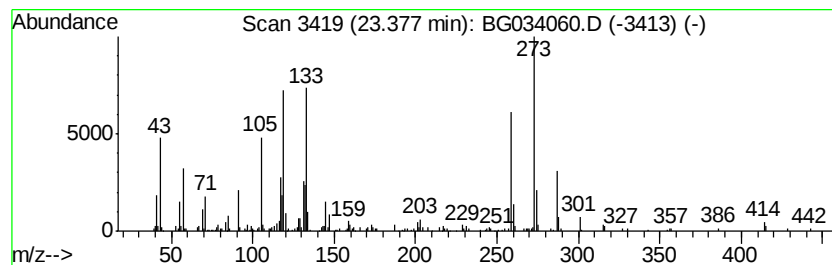
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 6-Methyl-2-(2,4,6-trimethyl... Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 23.38 | 6.28 ng | 419103 | Perylene-d12     | 24.39 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 6-Methyl-2-(2,4,6-trimethyl-benz... | 274 | C15H18N2OS | 1000294-32-1 | 22   |
| 2    |    | Disiloxane, 1,1,3,3-tetramethyl-    | 134 | C4H14OSi2  | 003277-26-7  | 22   |
| 3    |    | Quinoxaline, 1,2,3,4-tetrahydro...  | 274 | C16H22N2O2 | 018503-40-7  | 22   |
| 4    |    | N-Methylpiperidino[2,4-b,c]1,2,3... | 273 | C18H27NO   | 1000128-54-7 | 20   |
| 5    |    | 1H-Pyrazolo[3,4-b]quinoline, 3,6... | 273 | C18H15N3   | 1000302-05-0 | 18   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG042718\  
 Data File : BG034060.D  
 Acq On : 28 Apr 2018 5:48  
 Operator : SJ/JU  
 Sample : J2571-13  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 042418-JETB-E-D-M

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG041918.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 |
| 2-Pentanone, 4-hy... | 4.89  | 7.0 ng  |       | 107115   | 1                     | 7.77  | 305010  | 20.0 |
| unknown7.30          | 7.30  | 60.2 ng |       | 918664   | 1                     | 7.77  | 305010  | 20.0 |
| Benzene, (1-methy... | 17.06 | 4.3 ng  |       | 243141   | 4                     | 17.14 | 1136090 | 20.0 |
| Benzene, (1-methy... | 17.82 | 11.0 ng |       | 625401   | 4                     | 17.14 | 1136090 | 20.0 |
| 9,12-Octadecadien... | 18.98 | 4.4 ng  |       | 251349   | 4                     | 17.14 | 1136090 | 20.0 |
| unknown20.84         | 20.84 | 3.8 ng  |       | 345634   | 5                     | 21.39 | 1845630 | 20.0 |
| 3-Acetyl-2-(4-hyd... | 20.92 | 3.9 ng  |       | 363190   | 5                     | 21.39 | 1845630 | 20.0 |
| 1,3,5-Triazin-2-a... | 21.02 | 5.0 ng  |       | 466466   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown21.12         | 21.12 | 6.1 ng  |       | 560588   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown21.62         | 21.62 | 9.4 ng  |       | 864375   | 5                     | 21.39 | 1845630 | 20.0 |
| Pregnan-18-ol, (5... | 21.71 | 10.0 ng |       | 926051   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown21.94         | 21.94 | 8.5 ng  |       | 784284   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.04         | 22.04 | 10.7 ng |       | 988804   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.18         | 22.18 | 6.7 ng  |       | 616025   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.28         | 22.28 | 8.9 ng  |       | 817336   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.54         | 22.54 | 5.1 ng  |       | 472206   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.66         | 22.66 | 7.6 ng  |       | 698481   | 5                     | 21.39 | 1845630 | 20.0 |
| unknown22.94         | 22.94 | 5.8 ng  |       | 389075   | 6                     | 24.39 | 1335040 | 20.0 |
| 6-Methyl-2-(2,4,6... | 23.38 | 6.3 ng  |       | 419103   | 6                     | 24.39 | 1335040 | 20.0 |