

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
 Data File : BF088420.D  
 Acq On : 26 Jun 2016 16:03  
 Operator : UM/SJ  
 Sample : H3684-02  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 GROUNDWATER

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\_F\METHODS\8270-BF060716.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         | 1.690       | 7             | 9           | 25           | rVB      | 266994         | 500263        | 22.64%          | 2.628%        |
| 2         | 4.033       | 211           | 214         | 223          | rVB      | 78349          | 140974        | 6.38%           | 0.740%        |
| 3         | 4.661       | 267           | 269         | 282          | rBV      | 350758         | 525275        | 23.77%          | 2.759%        |
| 4         | 4.901       | 288           | 290         | 300          | rBV2     | 14076          | 36365         | 1.65%           | 0.191%        |
| 5         | 5.130       | 308           | 310         | 332          | rBV      | 623004         | 788942        | 35.71%          | 4.144%        |
| 6         | 5.530       | 344           | 345         | 358          | rVB      | 11294          | 26201         | 1.19%           | 0.138%        |
| 7         | 5.850       | 371           | 373         | 383          | rBV      | 55828          | 96652         | 4.37%           | 0.508%        |
| 8         | 6.216       | 402           | 405         | 412          | rBV      | 311658         | 450971        | 20.41%          | 2.369%        |
| 9         | 6.319       | 412           | 414         | 432          | rVB      | 1472932        | 1572641       | 71.18%          | 8.260%        |
| 10        | 6.547       | 432           | 434         | 440          | rBV      | 351312         | 386256        | 17.48%          | 2.029%        |
| 11        | 6.696       | 445           | 447         | 450          | rBV      | 1079478        | 1423783       | 64.44%          | 7.478%        |
| 12        | 7.130       | 482           | 485         | 487          | rBV      | 708912         | 1071182       | 48.48%          | 5.626%        |
| 13        | 7.759       | 538           | 540         | 544          | rBV2     | 30808          | 51348         | 2.32%           | 0.270%        |
| 14        | 7.827       | 544           | 546         | 549          | rBV      | 366465         | 542510        | 24.55%          | 2.849%        |
| 15        | 8.079       | 565           | 568         | 569          | rBV      | 853598         | 992108        | 44.90%          | 5.211%        |
| 16        | 8.102       | 569           | 570         | 576          | rVV      | 959107         | 948250        | 42.92%          | 4.981%        |
| 17        | 8.193       | 576           | 578         | 579          | rVV      | 56987          | 71467         | 3.23%           | 0.375%        |
| 18        | 8.216       | 579           | 580         | 587          | rVB      | 38410          | 56541         | 2.56%           | 0.297%        |
| 19        | 8.410       | 595           | 597         | 599          | rBV      | 941086         | 863887        | 39.10%          | 4.537%        |
| 20        | 8.627       | 615           | 616         | 620          | rBV2     | 44627          | 84347         | 3.82%           | 0.443%        |
| 21        | 8.913       | 638           | 641         | 643          | rBV      | 2197161        | 2209539       | 100.00%         | 11.605%       |
| 22        | 9.039       | 650           | 652         | 655          | rVB      | 33040          | 30557         | 1.38%           | 0.160%        |
| 23        | 9.313       | 674           | 676         | 680          | rVB      | 91024          | 85246         | 3.86%           | 0.448%        |
| 24        | 9.405       | 682           | 684         | 692          | rVB2     | 41773          | 67235         | 3.04%           | 0.353%        |
| 25        | 9.576       | 696           | 699         | 702          | rBV      | 708955         | 640064        | 28.97%          | 3.362%        |
| 26        | 9.976       | 731           | 734         | 736          | rBV      | 126314         | 174016        | 7.88%           | 0.914%        |
| 27        | 10.022      | 736           | 738         | 740          | rVB2     | 28506          | 28383         | 1.28%           | 0.149%        |
| 28        | 10.365      | 765           | 768         | 771          | rBV      | 1063336        | 1203377       | 54.46%          | 6.321%        |
| 29        | 10.491      | 778           | 779         | 783          | rVB      | 25701          | 28094         | 1.27%           | 0.148%        |
| 30        | 10.936      | 814           | 818         | 820          | rBV2     | 15976          | 24584         | 1.11%           | 0.129%        |
| 31        | 11.051      | 826           | 828         | 833          | rBV      | 620926         | 604561        | 27.36%          | 3.175%        |
| 32        | 11.222      | 841           | 843         | 846          | rBV      | 50307          | 53636         | 2.43%           | 0.282%        |
| 33        | 11.302      | 848           | 850         | 854          | rVB      | 30373          | 46251         | 2.09%           | 0.243%        |
| 34        | 11.622      | 877           | 878         | 882          | rVB      | 35959          | 38922         | 1.76%           | 0.204%        |

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Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

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:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |         |
|----|--------|------|------|------|------|---------|---------|--------|---------|
| 35 | 12.548 | 958  | 959  | 962  | rVB  | 35813   | 41346   | 1.87%  | 0.217%  |
| 36 | 12.639 | 964  | 967  | 969  | rVV  | 1772830 | 2066620 | 93.53% | 10.855% |
| 37 | 12.674 | 969  | 970  | 978  | rVB2 | 67891   | 80302   | 3.63%  | 0.422%  |
| 38 | 13.542 | 1043 | 1046 | 1050 | rBV2 | 35683   | 67023   | 3.03%  | 0.352%  |
| 39 | 13.679 | 1055 | 1058 | 1069 | rVV  | 425004  | 512953  | 23.22% | 2.694%  |
| 40 | 15.028 | 1174 | 1176 | 1185 | rBV  | 310295  | 406213  | 18.38% | 2.134%  |

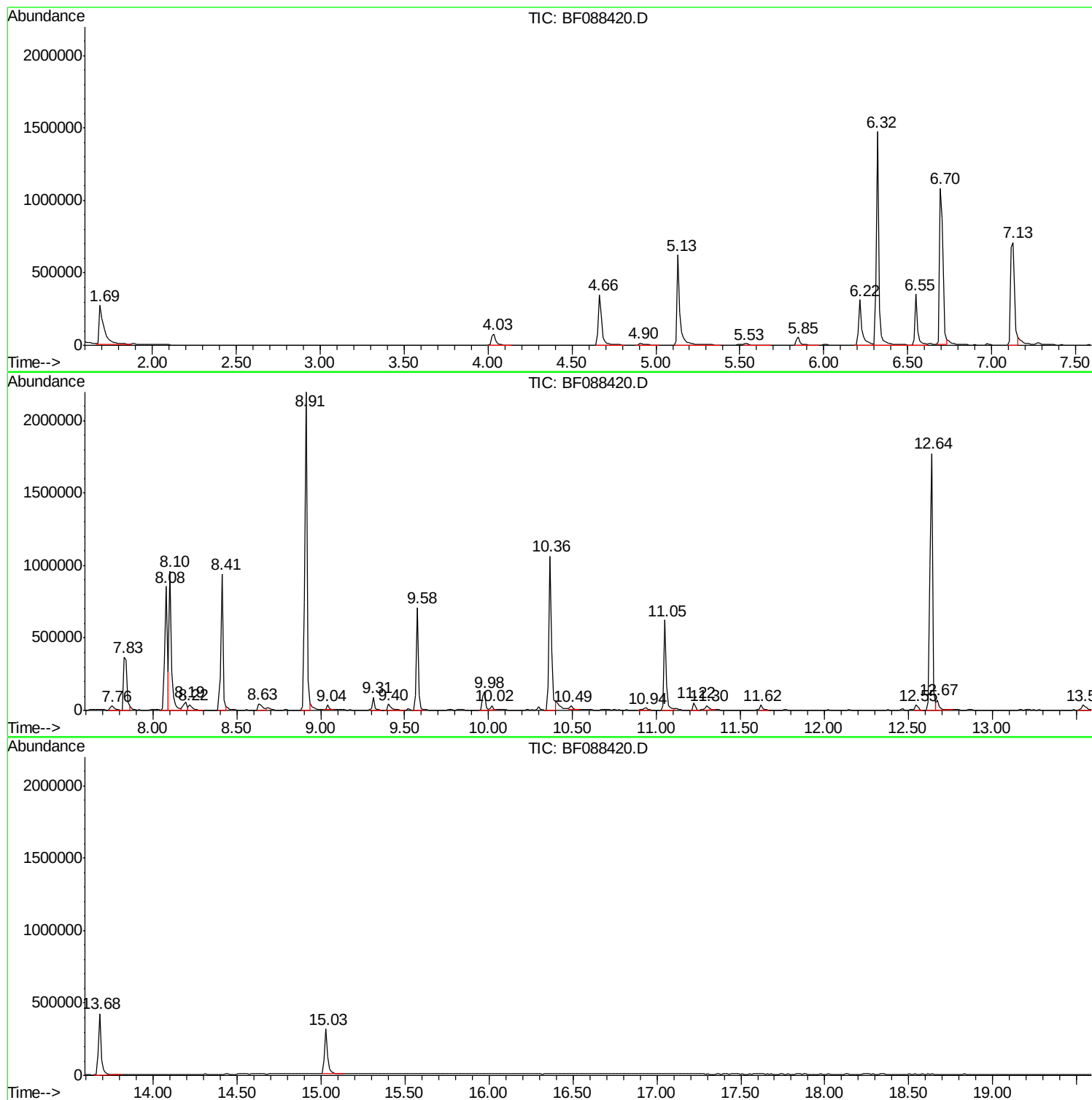
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : H3684-02  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

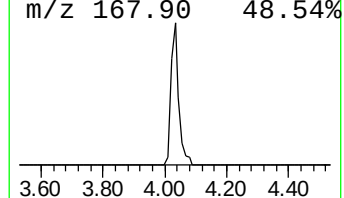
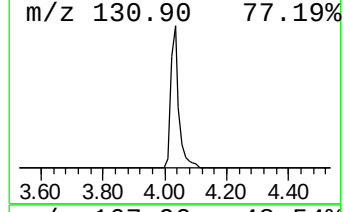
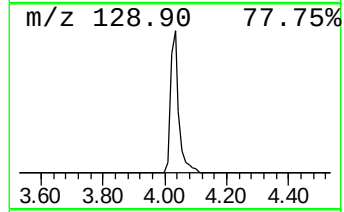
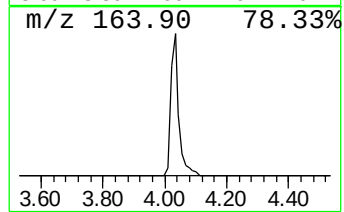
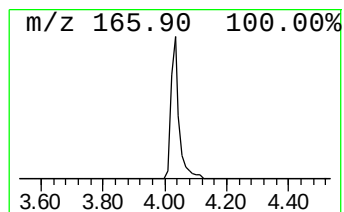
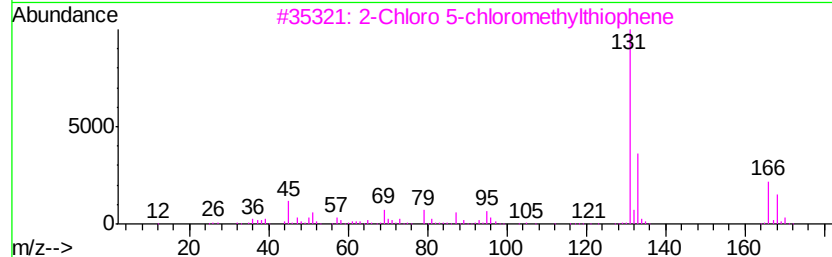
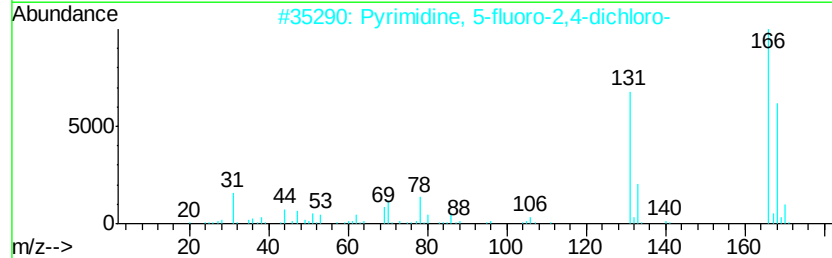
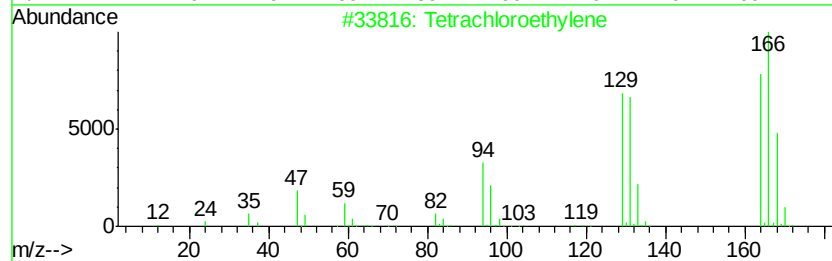
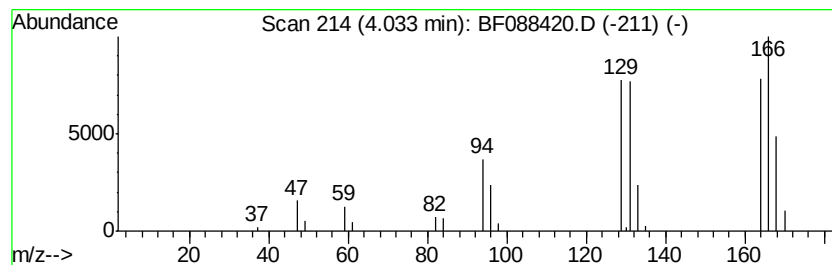
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Tetrachloroethylene Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.03 | 7.30 ng | 140974 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                       | MW  | MolForm   | CAS#        | Qual |
|------|----|------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Tetrachloroethylene                | 164 | C2Cl4     | 000127-18-4 | 98   |
| 2    |    | Pyrimidine, 5-fluoro-2,4-dichloro- | 166 | C4HCl2FN2 | 002927-71-1 | 43   |
| 3    |    | 2-Chloro 5-chloromethylthiophene   | 166 | C5H4Cl2S  | 023784-96-5 | 12   |
| 4    |    | 2-Chloroquinoxaline                | 164 | C8H5ClN2  | 001448-87-9 | 9    |
| 5    |    | 4,6-Dichloro-2-fluoropyrimidine    | 166 | C4HCl2FN2 | 003824-45-1 | 9    |



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ALS Vial : 10 Sample Multiplier: 1

Instrument :  
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GROUNDWATER

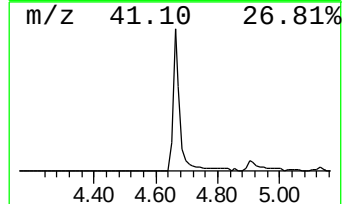
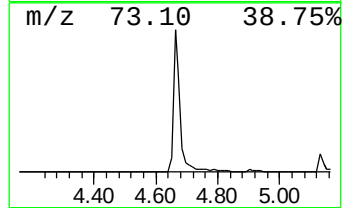
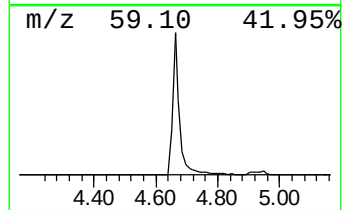
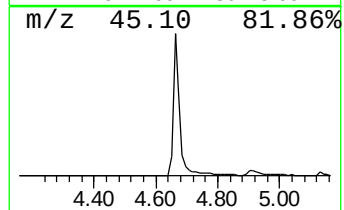
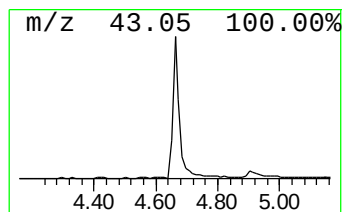
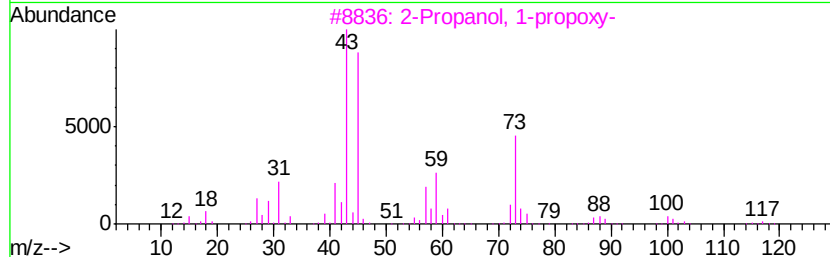
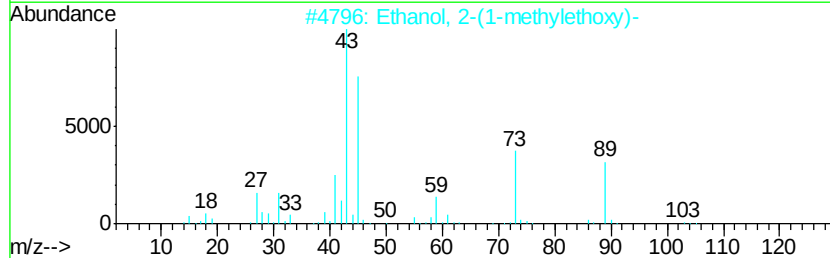
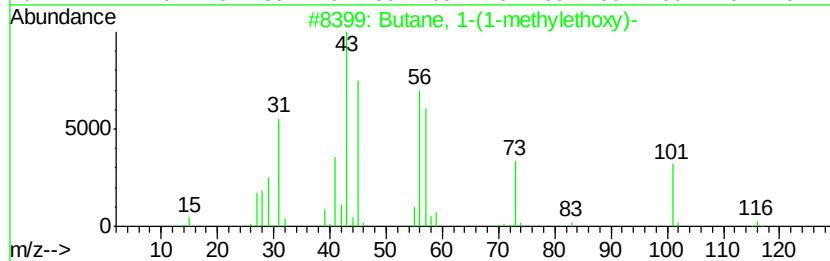
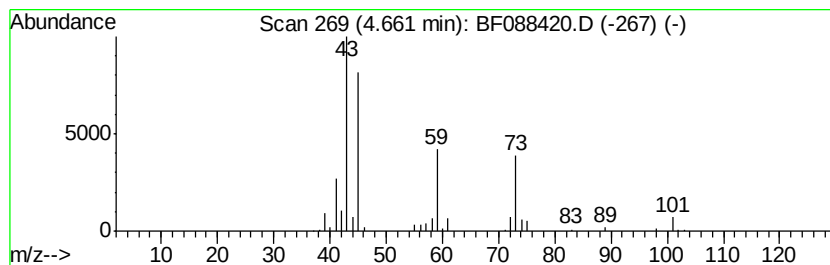
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Butane, 1-(1-methylethoxy)- Concentration Rank 6

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.66 | 27.20 ng | 525275 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 1-(1-methylethoxy)-     | 116 | C7H16O  | 001860-27-1 | 64   |
| 2    |    | Ethanol, 2-(1-methylethoxy)-    | 104 | C5H12O2 | 000109-59-1 | 59   |
| 3    |    | 2-Propanol, 1-propoxy-          | 118 | C6H14O2 | 001569-01-3 | 56   |
| 4    |    | 2-Propanol, 1-(1-methylethoxy)- | 118 | C6H14O2 | 003944-36-3 | 56   |
| 5    |    | Butane, 2-ethoxy-               | 102 | C6H14O  | 002679-87-0 | 47   |



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Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

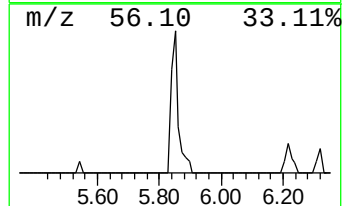
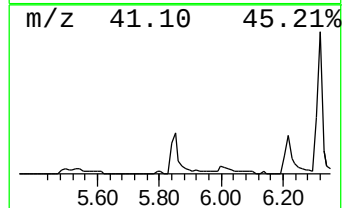
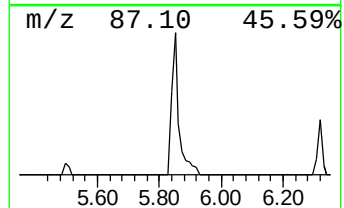
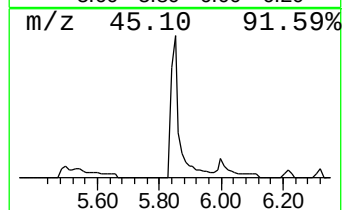
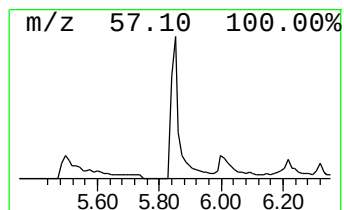
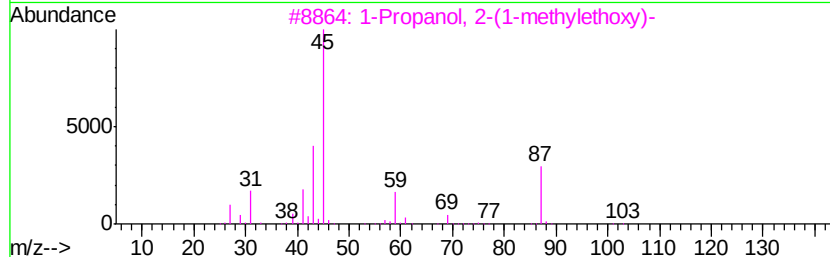
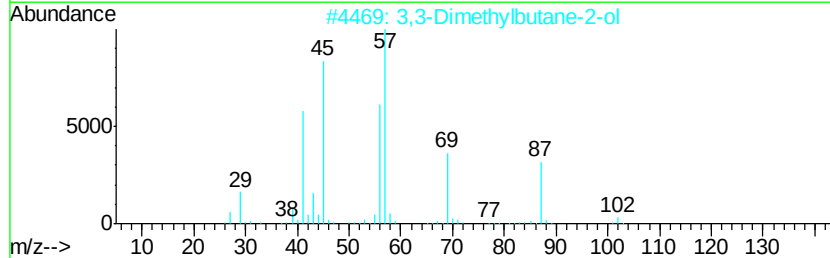
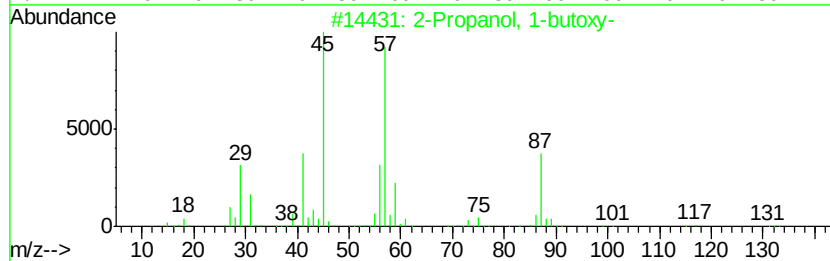
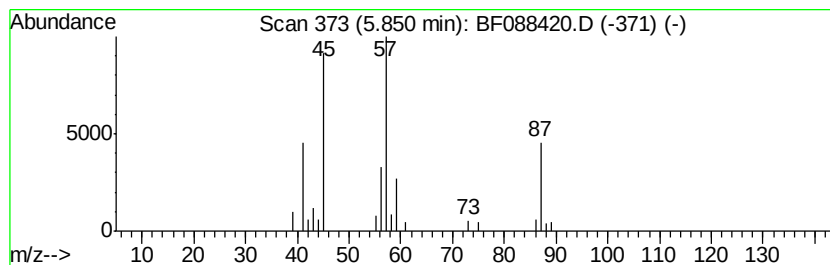
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 2-Propanol, 1-butoxy- Concentration Rank 10

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 5.85 | 5.00 ng | 96652 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2-Propanol, 1-butoxy-             | 132 | C7H16O2 | 005131-66-8 | 90   |
| 2    |    |   | 3,3-Dimethylbutane-2-ol           | 102 | C6H14O  | 000464-07-3 | 42   |
| 3    |    |   | 1-Propanol, 2-(1-methylethoxy)-   | 118 | C6H14O2 | 003944-37-4 | 38   |
| 4    |    |   | 4-Heptanol, 3,5-dimethyl-         | 144 | C9H20O  | 019549-79-2 | 33   |
| 5    |    |   | Formic acid, 1-methylpropyl ester | 102 | C5H10O2 | 000589-40-2 | 33   |



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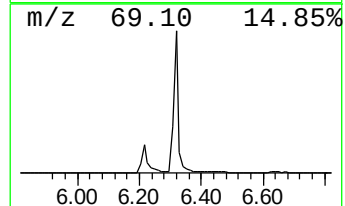
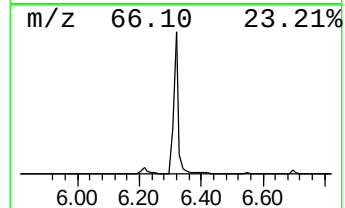
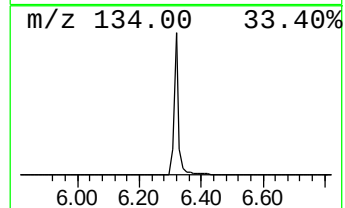
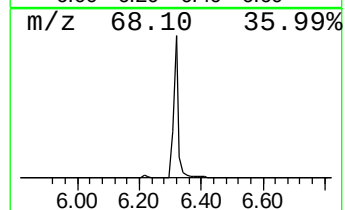
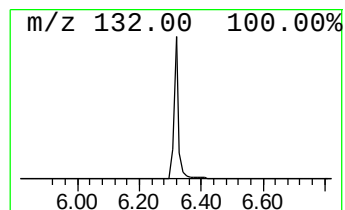
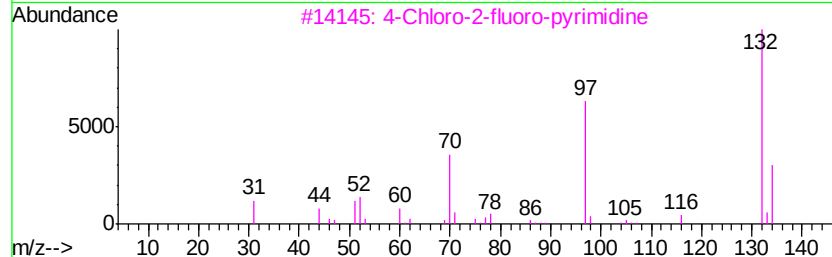
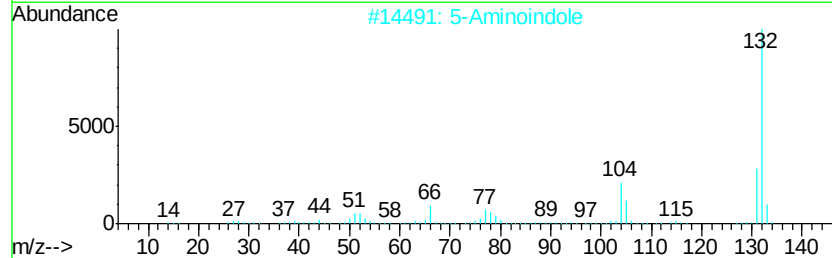
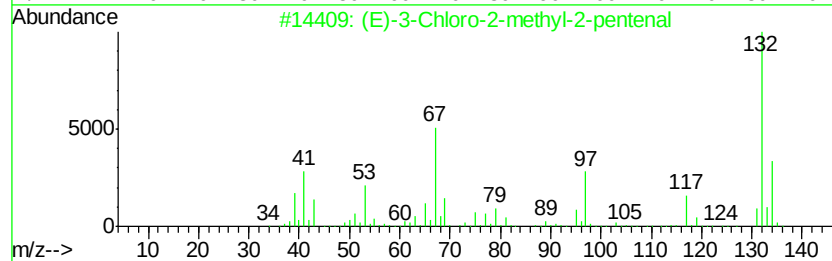
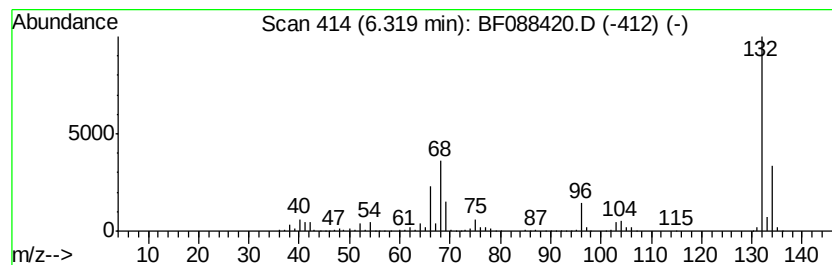
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown6.32 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.32 | 81.43 ng | 1572640 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of                               | 5   | Tentative ID | MW          | MolForm | CAS# | Qual |
|------|----------------------------------|-----|--------------|-------------|---------|------|------|
| 1    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO      | 031357-76-3 | 17      |      |      |
| 2    | 5-Aminoindole                    | 132 | C8H8N2       | 005192-03-0 | 12      |      |      |
| 3    | 4-Chloro-2-fluoro-pyrimidine     | 132 | C4H2ClFN2    | 051422-00-5 | 9       |      |      |
| 4    | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2    | 051422-01-6 | 9       |      |      |
| 5    | 1-Methylpyrrolo[1,2-a]pyrazine   | 132 | C8H8N2       | 064608-59-9 | 9       |      |      |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

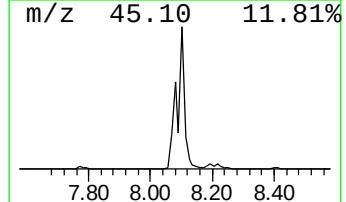
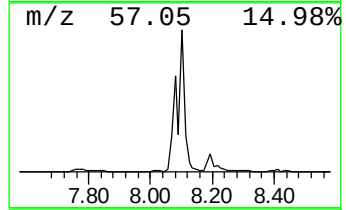
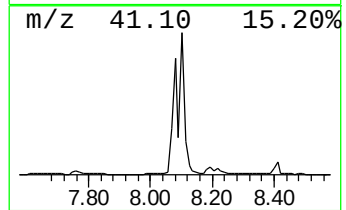
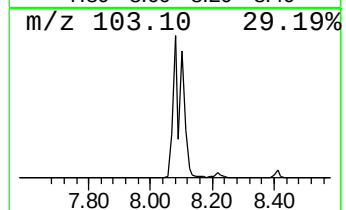
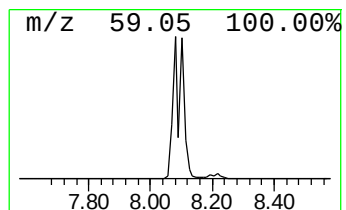
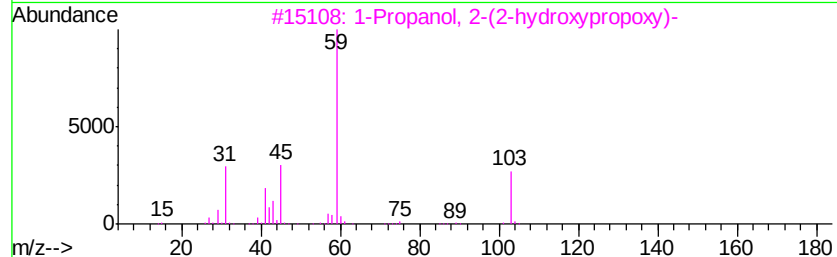
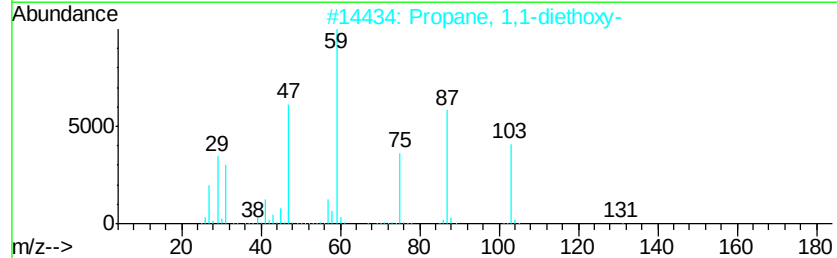
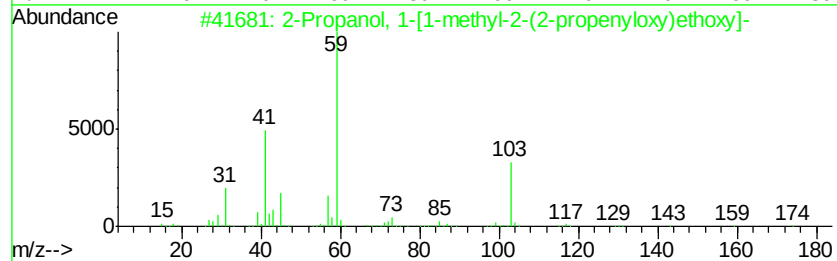
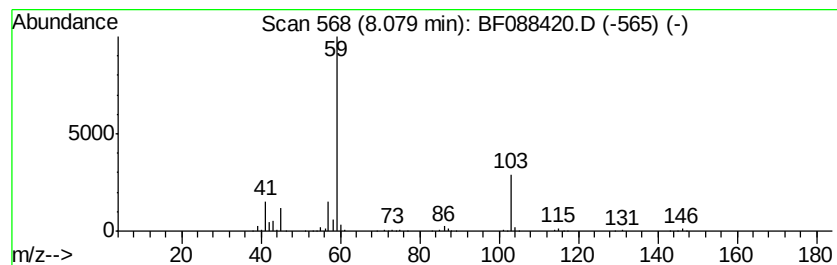
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 2-Propanol, 1-[1-methyl-2-(... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.08 | 36.57 ng | 992108 | Napthalene-d8    | 7.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2-Propanol, 1-[1-methyl-2-(2-pro... | 174 | C9H18O3  | 055956-25-7  | 83   |
| 2    |    | Propane, 1,1-diethoxy-              | 132 | C7H16O2  | 004744-08-5  | 74   |
| 3    |    | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134 | C6H14O3  | 000106-62-7  | 64   |
| 4    |    | 1-(1-Methoxypropan-2-yloxy)propa... | 232 | C12H24O4 | 1000367-13-4 | 59   |
| 5    |    | 2-Pentanol, 2,3-dimethyl-           | 116 | C7H16O   | 004911-70-0  | 59   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

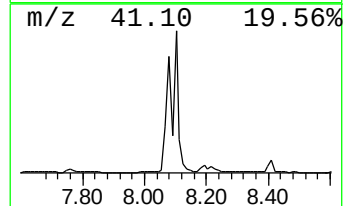
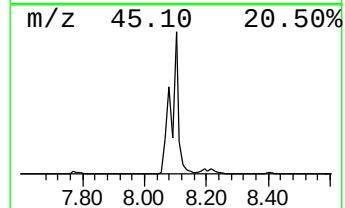
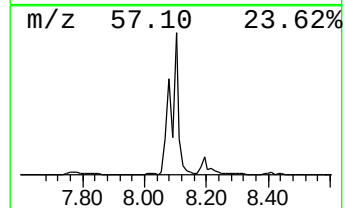
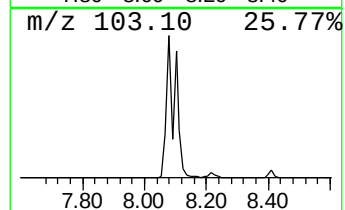
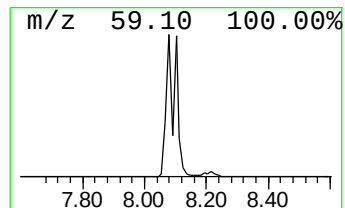
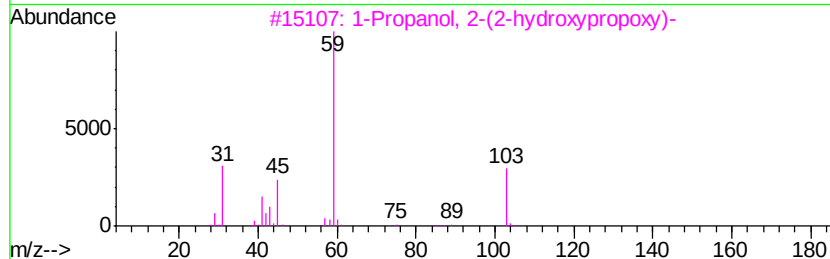
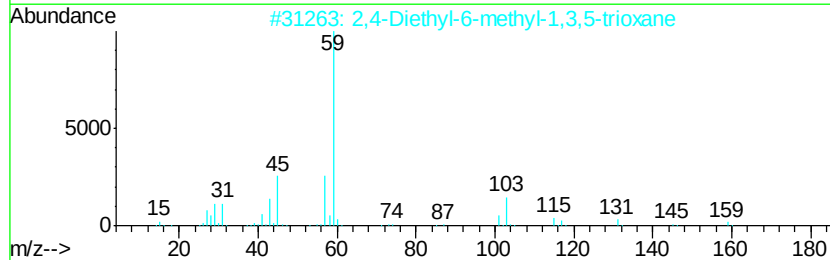
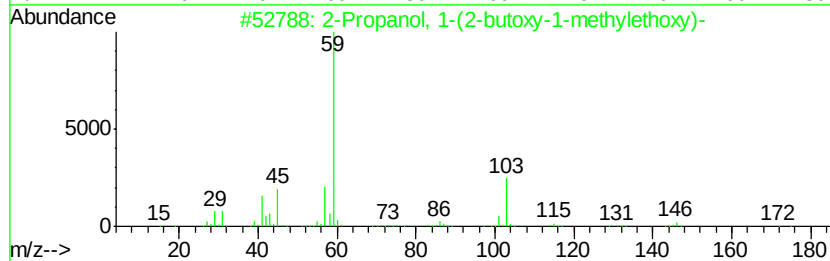
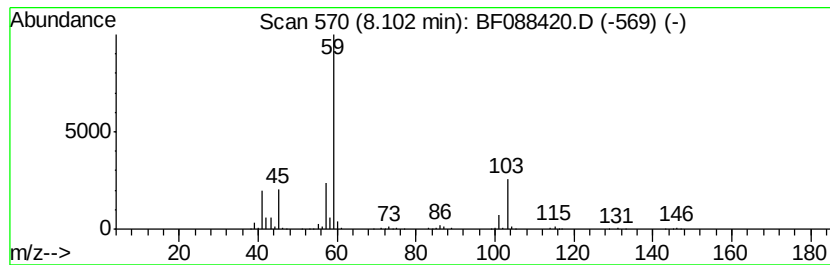
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 2-Propanol, 1-(2-butoxy-1-m... Concentration Rank 4

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.10 | 34.96 ng | 948250 | Napthalene-d8    | 7.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Propanol, 1-(2-butoxy-1-methyl... | 190 | C10H22O3 | 029911-28-2 | 90   |
| 2    |    | 2,4-Diethyl-6-methyl-1,3,5-trioxane | 160 | C8H16O3  | 117888-04-7 | 78   |
| 3    |    | 1-Propanol, 2-(2-hydroxypropoxy)-   | 134 | C6H14O3  | 000106-62-7 | 78   |
| 4    |    | 2-Propanol, 1-[1-methyl-2-(2-pro... | 174 | C9H18O3  | 055956-25-7 | 64   |
| 5    |    | Hexaethylene glycol dimethyl ether  | 310 | C14H30O7 | 001072-40-8 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

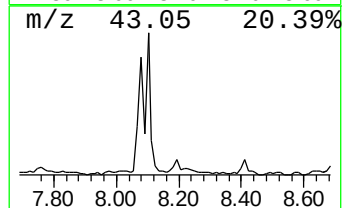
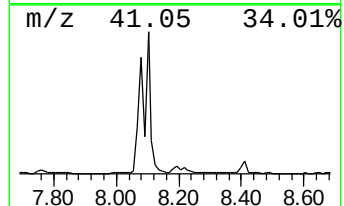
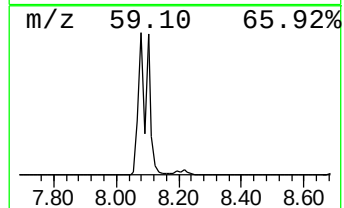
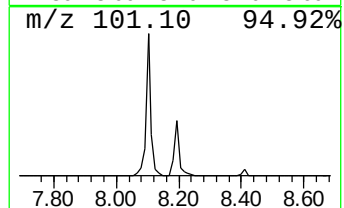
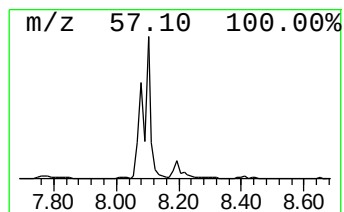
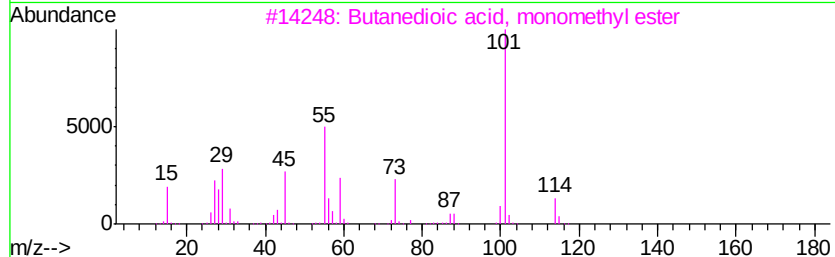
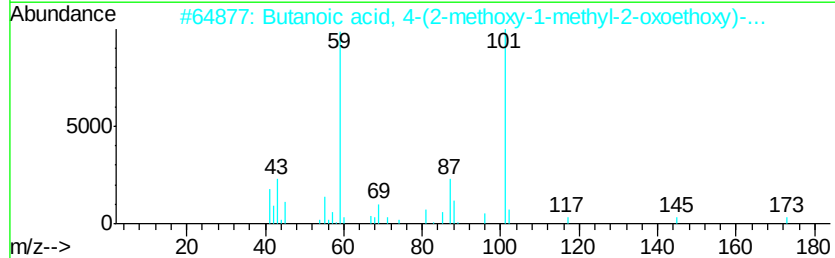
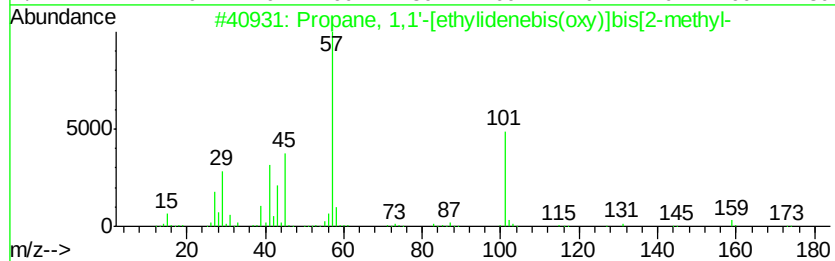
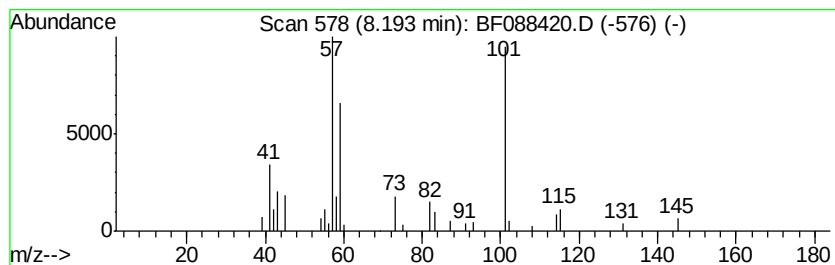
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown8.19 Concentration Rank 12

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 8.19 | 2.63 ng | 71467 | Naphthalene-d8   | 7.84 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Propane, 1,1'-[ethylidenebis(oxv... | 174 | C10H22O2 | 005669-09-0 | 47   |
| 2    |    |   | Butanoic acid, 4-(2-methoxy-1-me... | 204 | C9H16O5  | 054966-46-0 | 42   |
| 3    |    |   | Butanedioic acid, monomethyl ester  | 132 | C5H8O4   | 003878-55-5 | 38   |
| 4    |    |   | Butanoic acid, 4-[(tetrahydro-2H... | 202 | C10H18O4 | 093691-87-3 | 38   |
| 5    |    |   | Butyric acid, 4-isopropoxy-, met... | 160 | C8H16O3  | 029006-05-1 | 32   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

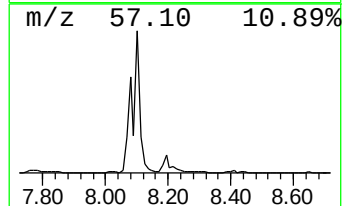
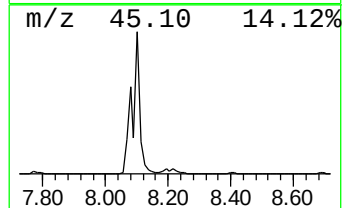
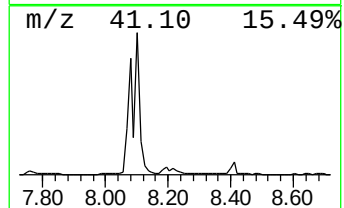
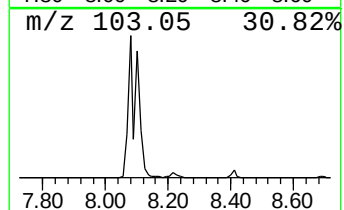
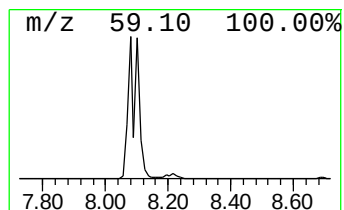
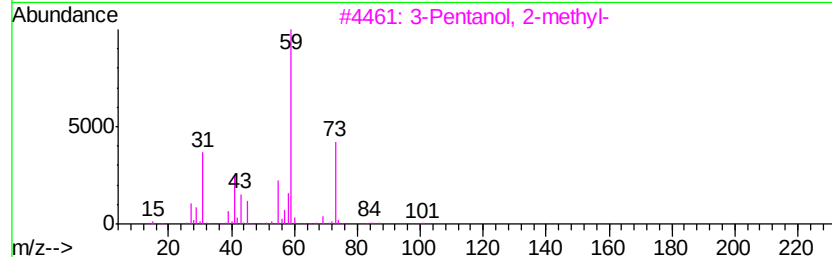
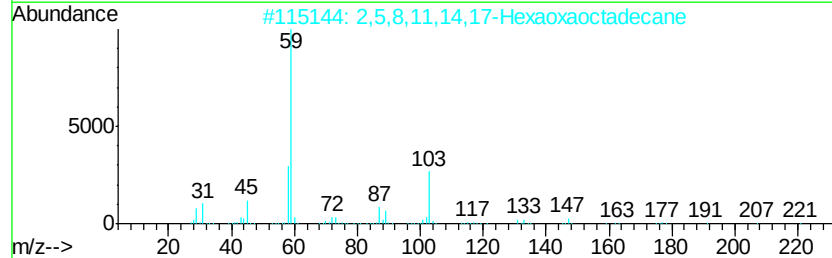
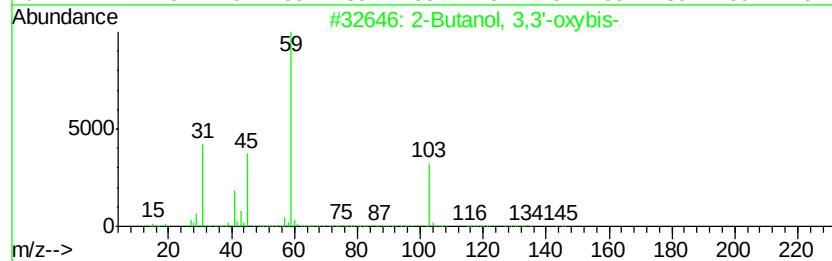
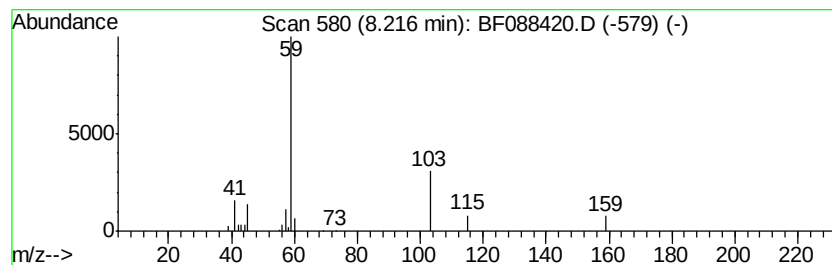
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown8.22 Concentration Rank 15

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 8.22 | 2.08 ng | 56541 | Napthalene-d8    | 7.84 |

| Hit# | of                               | 5 | Tentative ID | MW  | MolForm  | CAS#        | Qual |
|------|----------------------------------|---|--------------|-----|----------|-------------|------|
| 1    | 2-Butanol, 3,3'-oxybis-          |   |              | 162 | C8H18O3  | 054305-61-2 | 39   |
| 2    | 2,5,8,11,14,17-Hexaoxaoctadecane |   |              | 266 | C12H26O6 | 001191-87-3 | 39   |
| 3    | 3-Pentanol, 2-methyl-            |   |              | 102 | C6H14O   | 000565-67-3 | 9    |
| 4    | 3-Heptanol, 4-methyl-            |   |              | 130 | C8H18O   | 014979-39-6 | 9    |
| 5    | 2-Pentanol, 2,3-dimethyl-        |   |              | 116 | C7H16O   | 004911-70-0 | 9    |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

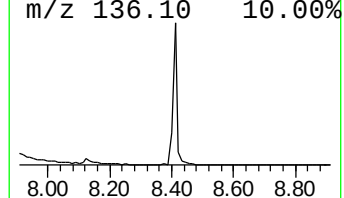
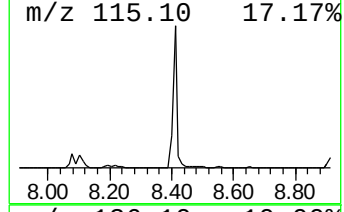
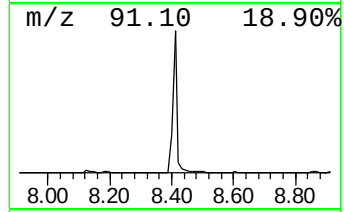
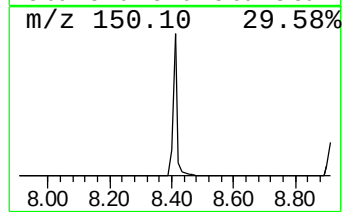
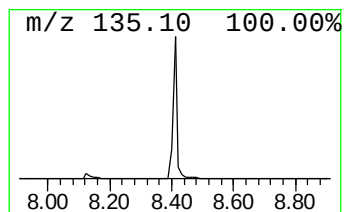
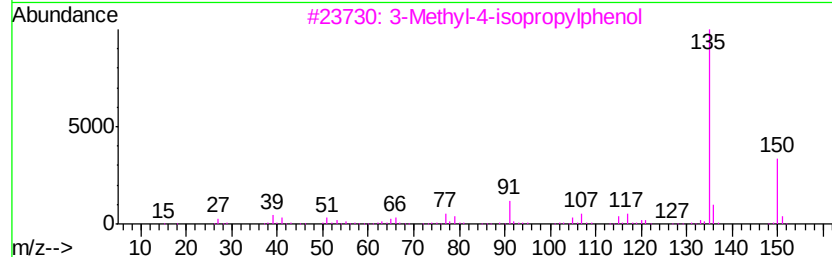
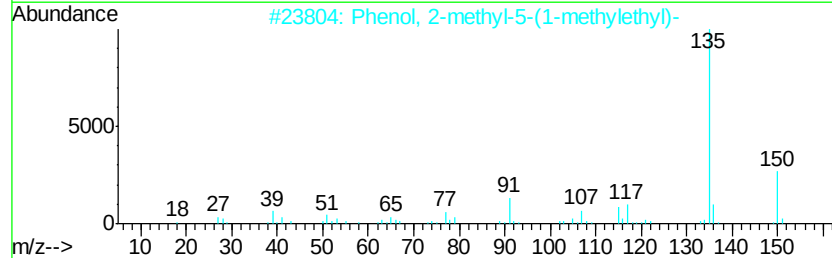
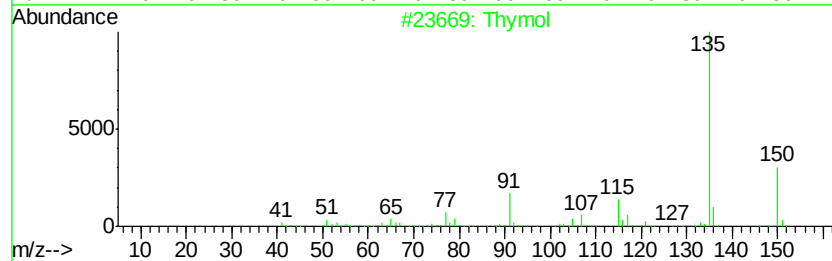
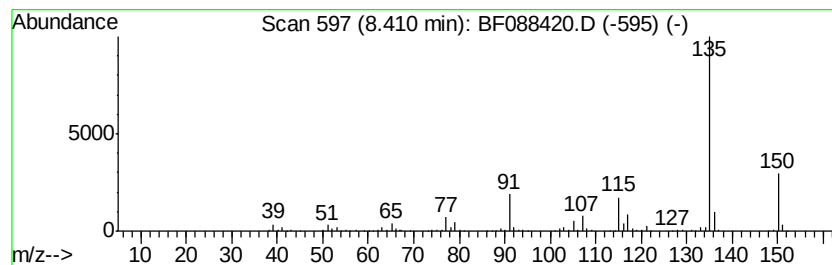
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Thymol Concentration Rank 5

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 8.41 | 31.85 ng | 863887 | Napthalene-d8    | 7.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Thymol                              | 150 | C10H14O | 000089-83-8 | 95   |
| 2    |    | Phenol, 2-methyl-5-(1-methylethyl)- | 150 | C10H14O | 000499-75-2 | 90   |
| 3    |    | 3-Methyl-4-isopropylphenol          | 150 | C10H14O | 003228-02-2 | 90   |
| 4    |    | Phenol, 2-(1,1-dimethylethyl)-      | 150 | C10H14O | 000088-18-6 | 80   |
| 5    |    | Phenol, 2,3,4,6-tetramethyl-        | 150 | C10H14O | 003238-38-8 | 80   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

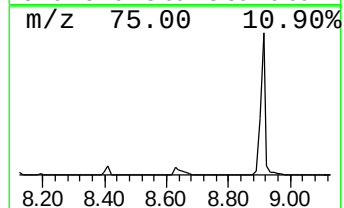
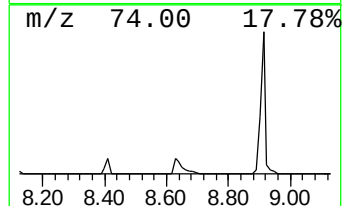
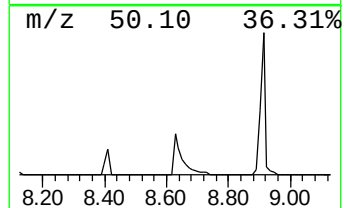
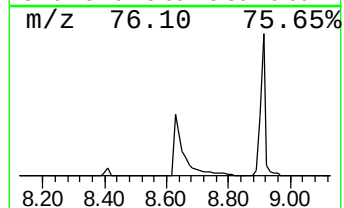
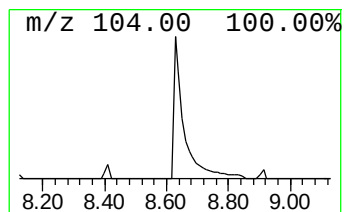
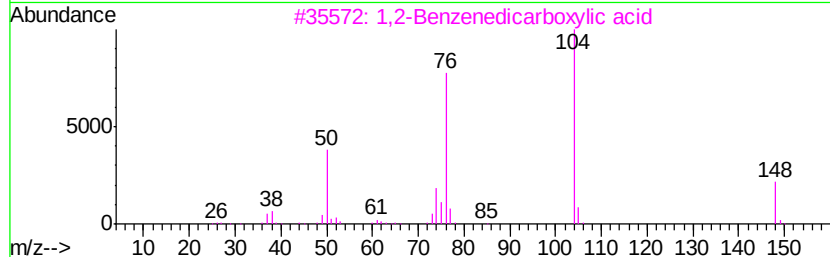
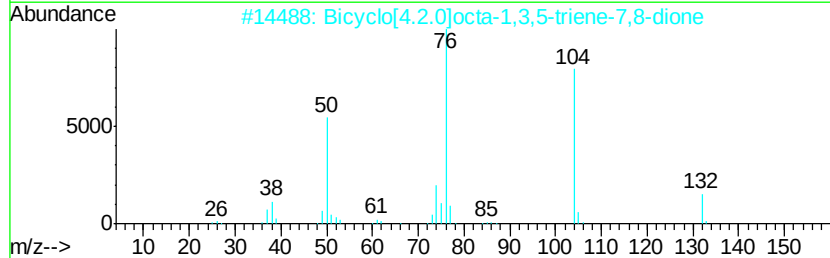
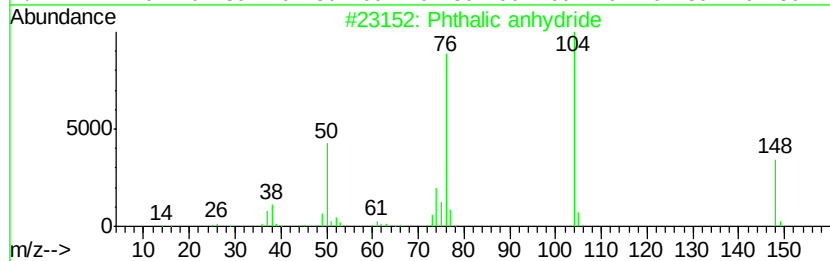
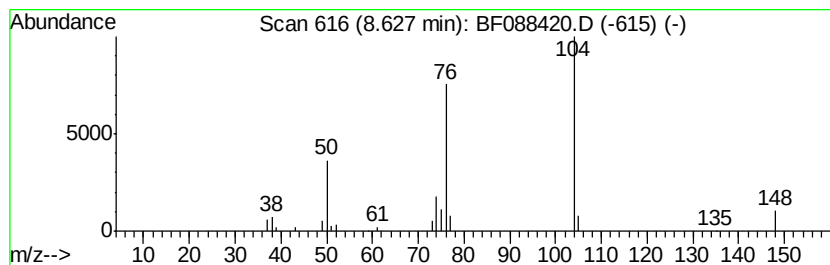
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 Phthalic anhydride Concentration Rank 11

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 8.63 | 3.11 ng | 84347 | Naphthalene-d8   | 7.84 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm     | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Phthalic anhydride                   | 148 | C8H4O3      | 000085-44-9 | 91   |
| 2    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene-...  | 132 | C8H4O2      | 006383-11-5 | 86   |
| 3    |    |   | 1,2-Benzenedicarboxylic acid         | 166 | C8H6O4      | 000088-99-3 | 83   |
| 4    |    |   | N-[2-Acetamidoethylthiol]phthalimide | 264 | C12H12N2O3S | 025158-14-9 | 78   |
| 5    |    |   | 1H-Isoindole-1,3(2H)-dione, 2-(h...  | 177 | C9H7NO3     | 000118-29-6 | 78   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

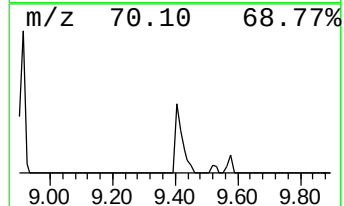
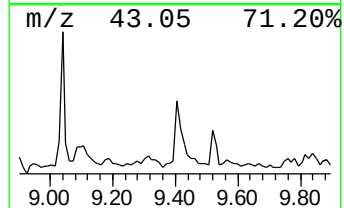
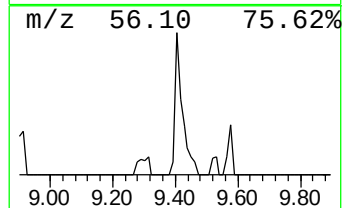
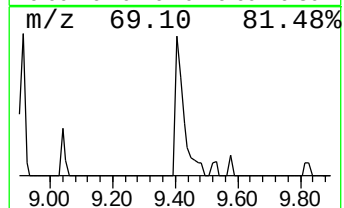
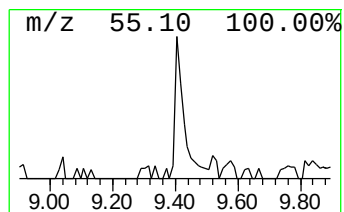
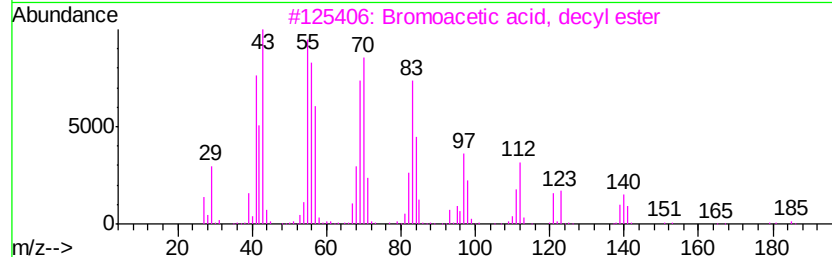
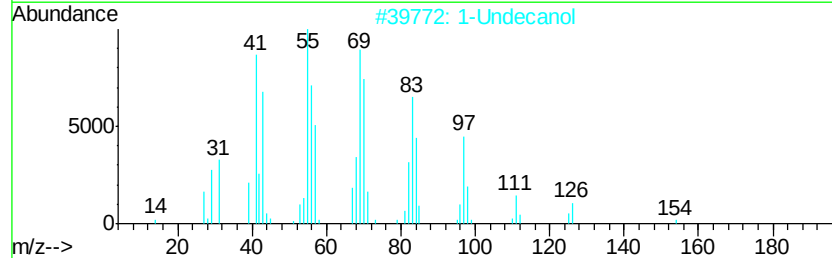
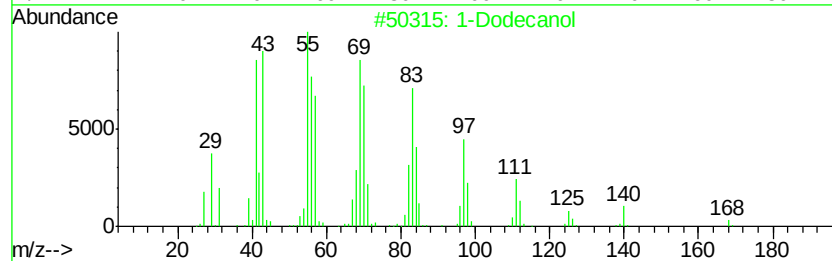
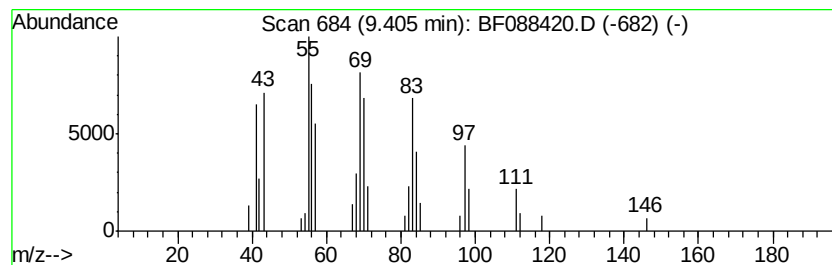
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 1-Dodecanol Concentration Rank 14

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 9.40 | 2.10 ng | 67235 | Acenaphthene-d10 | 9.58 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------|-----|------------|-------------|------|
| 1    |    |   | 1-Dodecanol                   | 186 | C12H26O    | 000112-53-8 | 91   |
| 2    |    |   | 1-Undecanol                   | 172 | C11H24O    | 000112-42-5 | 87   |
| 3    |    |   | Bromoacetic acid, decyl ester | 278 | C12H23BrO2 | 005436-93-1 | 86   |
| 4    |    |   | 1-Decanol                     | 158 | C10H22O    | 000112-30-1 | 83   |
| 5    |    |   | 1-Dodecene                    | 168 | C12H24     | 000112-41-4 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

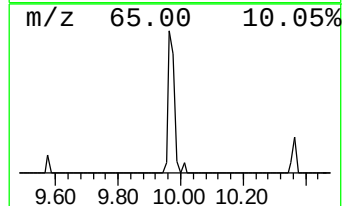
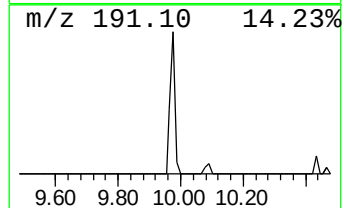
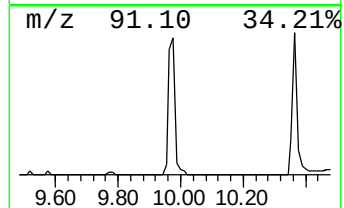
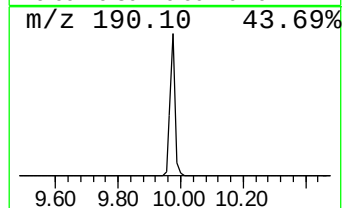
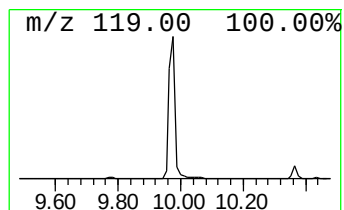
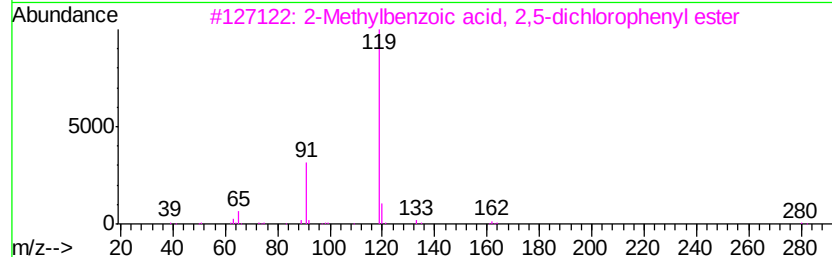
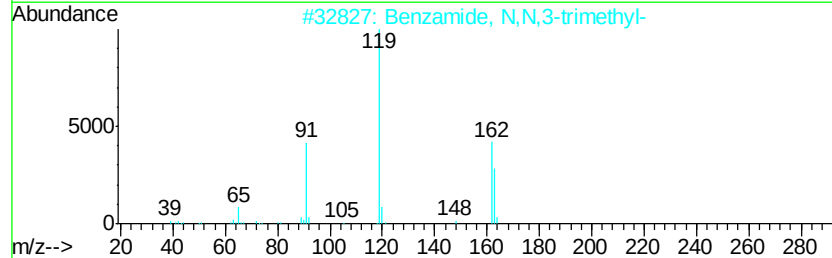
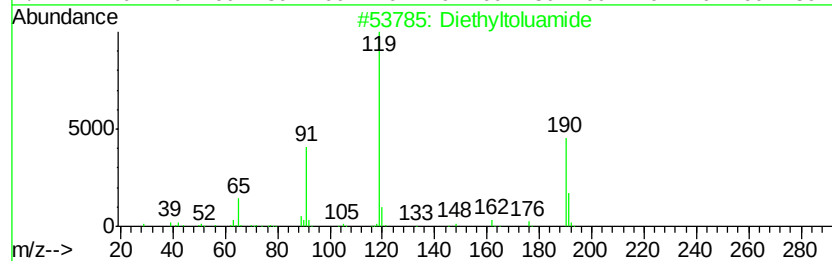
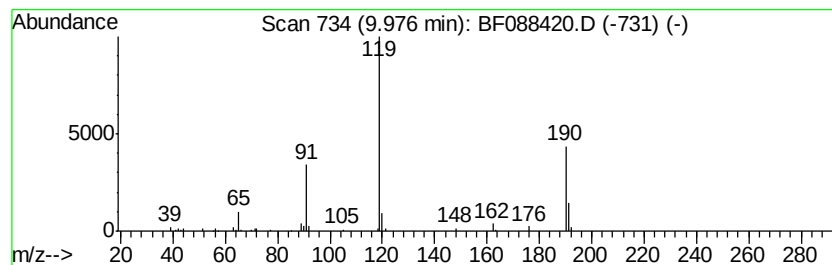
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Diethyltoluamide Concentration Rank 9

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 9.98 | 5.44 ng | 174016 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Diethyltoluamide                    | 191 | C12H17NO    | 000134-62-3  | 95   |
| 2    |    | Benzamide, N,N,3-trimethyl-         | 163 | C10H13NO    | 006935-65-5  | 50   |
| 3    |    | 2-Methylbenzoic acid, 2,5-dichlo... | 280 | C14H10Cl2O2 | 1000325-60-9 | 50   |
| 4    |    | p-Toluric acid                      | 193 | C10H11NO3   | 027115-50-0  | 50   |
| 5    |    | 2-Isopropyl-5-methyl-[1,4]benzoq... | 297 | C18H19NO3   | 1000301-09-7 | 50   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062616\  
Data File : BF088420.D  
Acq On : 26 Jun 2016 16:03  
Operator : UM/SJ  
Sample : H3684-02  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

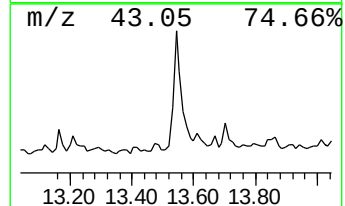
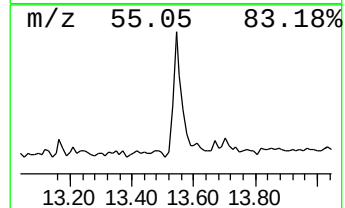
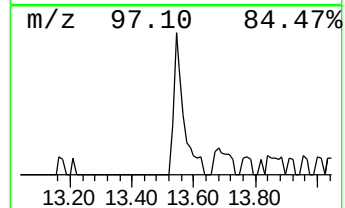
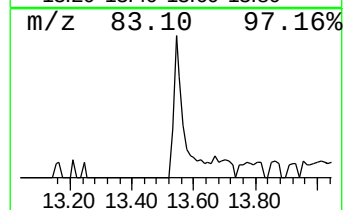
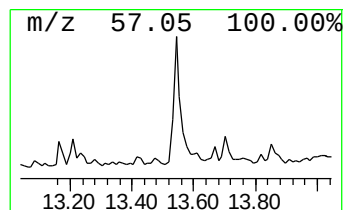
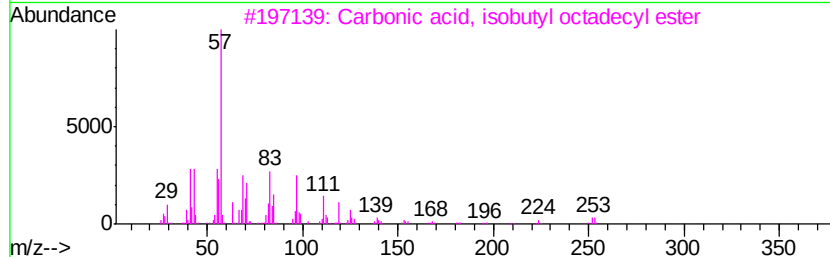
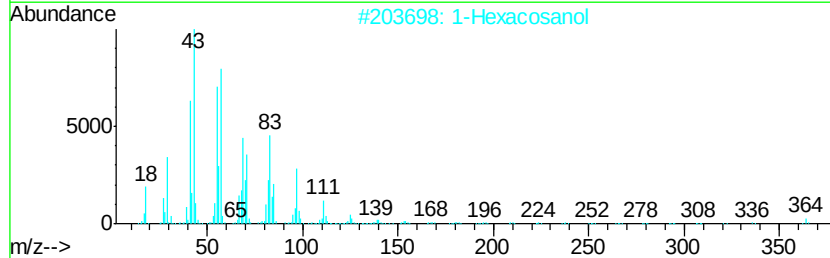
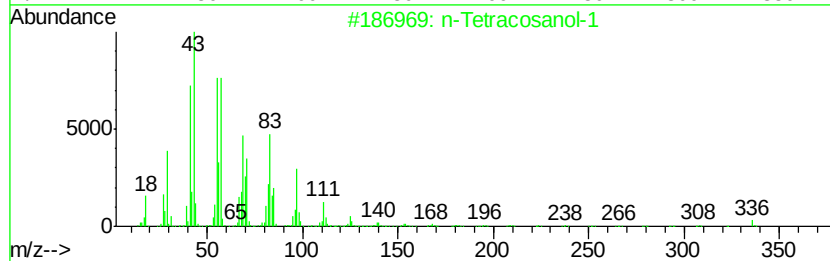
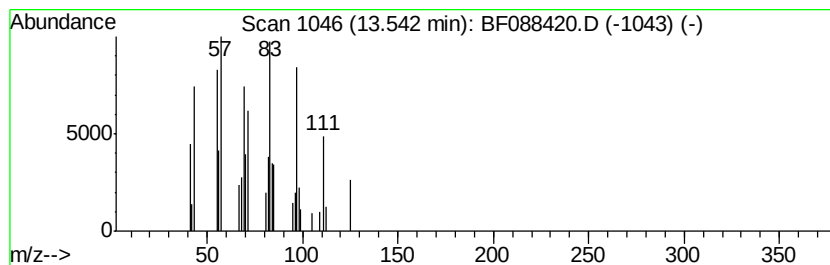
Instrument :  
BNA\_F  
ClientSampleId :  
GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.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 15 n-Tetracosanol-1 Concentration Rank 13

| R.T.      | EstConc                              | Area  | Relative to ISTD | R.T.         |      |
|-----------|--------------------------------------|-------|------------------|--------------|------|
| 13.54     | 2.61 ng                              | 67023 | Chrysene-d12     | 13.68        |      |
| Hit# of 5 | Tentative ID                         | MW    | MolForm          | CAS#         | Qual |
| 1         | n-Tetracosanol-1                     | 354   | C24H50O          | 000506-51-4  | 90   |
| 2         | 1-Hexacosanol                        | 382   | C26H54O          | 000506-52-5  | 81   |
| 3         | Carbonic acid, isobutyl octadecyl... | 370   | C23H46O3         | 1000314-61-5 | 59   |
| 4         | Isobutyl tetradecyl carbonate        | 314   | C19H38O3         | 959275-58-2  | 35   |
| 5         | 2(5H)-Furanone, 5,5-dimethyl-        | 112   | C6H8O2           | 020019-64-1  | 27   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062616\  
 Data File : BF088420.D  
 Acq On : 26 Jun 2016 16:03  
 Operator : UM/SJ  
 Sample : H3684-02  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 GROUNDWATER

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.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 |
| Tetrachloroethylene  | 4.03  | 7.3     | ng    | 140974   | 1                     | 6.55  | 386256 | 20.0 |
| Butane, 1-(1-meth... | 4.66  | 27.2    | ng    | 525275   | 1                     | 6.55  | 386256 | 20.0 |
| 2-Propanol, 1-but... | 5.85  | 5.0     | ng    | 96652    | 1                     | 6.55  | 386256 | 20.0 |
| unknown6.32          | 6.32  | 81.4    | ng    | 1572640  | 1                     | 6.55  | 386256 | 20.0 |
| 2-Propanol, 1-[1-... | 8.08  | 36.6    | ng    | 992108   | 2                     | 7.84  | 542510 | 20.0 |
| 2-Propanol, 1-(2-... | 8.10  | 35.0    | ng    | 948250   | 2                     | 7.84  | 542510 | 20.0 |
| unknown8.19          | 8.19  | 2.6     | ng    | 71467    | 2                     | 7.84  | 542510 | 20.0 |
| unknown8.22          | 8.22  | 2.1     | ng    | 56541    | 2                     | 7.84  | 542510 | 20.0 |
| Thymol               | 8.41  | 31.9    | ng    | 863887   | 2                     | 7.84  | 542510 | 20.0 |
| Phthalic anhydride   | 8.63  | 3.1     | ng    | 84347    | 2                     | 7.84  | 542510 | 20.0 |
| 1-Dodecanol          | 9.40  | 2.1     | ng    | 67235    | 3                     | 9.58  | 640064 | 20.0 |
| Diethyltoluamide     | 9.98  | 5.4     | ng    | 174016   | 3                     | 9.58  | 640064 | 20.0 |
| n-Tetracosanol-1     | 13.54 | 2.6     | ng    | 67023    | 5                     | 13.68 | 512953 | 20.0 |