

Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
 Data File : BM005497.D  
 Acq On : 19 May 2016 07:49  
 Operator : UM/SJ  
 Sample : H2890-22RX  
 Misc :  
 ALS Vial : 37 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 D9X07RX

## Integration Parameters: LSCINT.P

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

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

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

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
 Title : SVOA CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.687       | 351           | 357         | 365          | rVB      | 23162          | 37697         | 3.26%           | 0.628%        |
| 2         | 7.169       | 774           | 779         | 788          | rBV      | 14007          | 22887         | 1.98%           | 0.381%        |
| 3         | 7.363       | 807           | 812         | 819          | rVB      | 14742          | 23149         | 2.00%           | 0.386%        |
| 4         | 7.834       | 887           | 892         | 899          | rVB      | 82981          | 133361        | 11.53%          | 2.222%        |
| 5         | 8.528       | 1003          | 1010        | 1015         | rVB3     | 17644          | 28144         | 2.43%           | 0.469%        |
| 6         | 8.986       | 1082          | 1088        | 1095         | rBV2     | 16609          | 28098         | 2.43%           | 0.468%        |
| 7         | 9.092       | 1100          | 1106        | 1112         | rVB3     | 15630          | 24657         | 2.13%           | 0.411%        |
| 8         | 9.704       | 1204          | 1210        | 1215         | rBV2     | 11933          | 18909         | 1.64%           | 0.315%        |
| 9         | 10.239      | 1295          | 1301        | 1306         | rBV2     | 23621          | 34680         | 3.00%           | 0.578%        |
| 10        | 10.622      | 1359          | 1366        | 1374         | rBV      | 136839         | 228254        | 19.74%          | 3.804%        |
| 11        | 11.139      | 1448          | 1454        | 1463         | rBV      | 42518          | 77409         | 6.69%           | 1.290%        |
| 12        | 13.874      | 1913          | 1919        | 1923         | rBV      | 50769          | 71579         | 6.19%           | 1.193%        |
| 13        | 13.921      | 1923          | 1927        | 1931         | rVB      | 69387          | 92904         | 8.03%           | 1.548%        |
| 14        | 14.157      | 1962          | 1967        | 1973         | rVB      | 45626          | 64618         | 5.59%           | 1.077%        |
| 15        | 14.257      | 1977          | 1984        | 1990         | rBV      | 24924          | 39070         | 3.38%           | 0.651%        |
| 16        | 14.368      | 1997          | 2003        | 2008         | rBV      | 20620          | 27296         | 2.36%           | 0.455%        |
| 17        | 14.463      | 2012          | 2019        | 2026         | rBV2     | 250151         | 379477        | 32.82%          | 6.324%        |
| 18        | 14.639      | 2045          | 2049        | 2056         | rBV2     | 9801           | 15550         | 1.34%           | 0.259%        |
| 19        | 15.321      | 2160          | 2165        | 2170         | rVB      | 38693          | 48901         | 4.23%           | 0.815%        |
| 20        | 15.457      | 2182          | 2188        | 2193         | rVB      | 65338          | 95723         | 8.28%           | 1.595%        |
| 21        | 15.563      | 2201          | 2206        | 2211         | rVB2     | 18589          | 25607         | 2.21%           | 0.427%        |
| 22        | 15.727      | 2224          | 2234        | 2238         | rBV3     | 10286          | 24003         | 2.08%           | 0.400%        |
| 23        | 16.180      | 2306          | 2311        | 2318         | rBV      | 53901          | 83224         | 7.20%           | 1.387%        |
| 24        | 16.968      | 2441          | 2445        | 2450         | rBV2     | 24143          | 32714         | 2.83%           | 0.545%        |
| 25        | 17.204      | 2478          | 2485        | 2490         | rBV      | 393098         | 558320        | 48.28%          | 9.304%        |
| 26        | 17.257      | 2490          | 2494        | 2497         | rVV2     | 18427          | 25546         | 2.21%           | 0.426%        |
| 27        | 17.298      | 2497          | 2501        | 2508         | rVB      | 72999          | 101925        | 8.81%           | 1.698%        |
| 28        | 17.574      | 2544          | 2548        | 2554         | rVB5     | 10667          | 15503         | 1.34%           | 0.258%        |
| 29        | 17.709      | 2566          | 2571        | 2576         | rBV      | 14271          | 19611         | 1.70%           | 0.327%        |
| 30        | 18.098      | 2632          | 2637        | 2644         | rBV      | 163525         | 220784        | 19.09%          | 3.679%        |
| 31        | 18.398      | 2684          | 2688        | 2694         | rBV3     | 10763          | 18987         | 1.64%           | 0.316%        |
| 32        | 18.951      | 2777          | 2782        | 2786         | rBV5     | 9113           | 16712         | 1.45%           | 0.278%        |
| 33        | 19.115      | 2806          | 2810        | 2815         | rBV      | 24052          | 31477         | 2.72%           | 0.525%        |
| 34        | 19.227      | 2826          | 2829        | 2832         | rVV4     | 18520          | 25011         | 2.16%           | 0.417%        |

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ClientSampleId :  
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## Integration Parameters: LSCINT.P

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Sampling : 1  
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Filtering: 5  
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Max Peaks: 100  
Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |         |         |         |
|----|--------|------|------|------|------|--------|---------|---------|---------|
| 35 | 19.256 | 2832 | 2834 | 2843 | rVB7 | 13002  | 23559   | 2.04%   | 0.393%  |
| 36 | 19.374 | 2849 | 2854 | 2864 | rBV2 | 199769 | 299452  | 25.89%  | 4.990%  |
| 37 | 19.451 | 2865 | 2867 | 2876 | rVV  | 17064  | 27215   | 2.35%   | 0.454%  |
| 38 | 19.586 | 2885 | 2890 | 2896 | rVV  | 105985 | 148921  | 12.88%  | 2.482%  |
| 39 | 19.645 | 2897 | 2900 | 2906 | rVB7 | 8763   | 13524   | 1.17%   | 0.225%  |
| 40 | 20.350 | 3016 | 3020 | 3028 | rBV  | 44001  | 84996   | 7.35%   | 1.416%  |
| 41 | 21.092 | 3142 | 3146 | 3156 | rBV  | 221537 | 298272  | 25.79%  | 4.970%  |
| 42 | 21.374 | 3189 | 3194 | 3200 | rBV  | 804833 | 993990  | 85.95%  | 16.564% |
| 43 | 23.509 | 3552 | 3557 | 3563 | rBV2 | 92555  | 170578  | 14.75%  | 2.842%  |
| 44 | 23.656 | 3575 | 3582 | 3592 | rVB  | 600633 | 1156411 | 100.00% | 19.270% |
| 45 | 24.309 | 3690 | 3693 | 3702 | rVB7 | 47549  | 92349   | 7.99%   | 1.539%  |

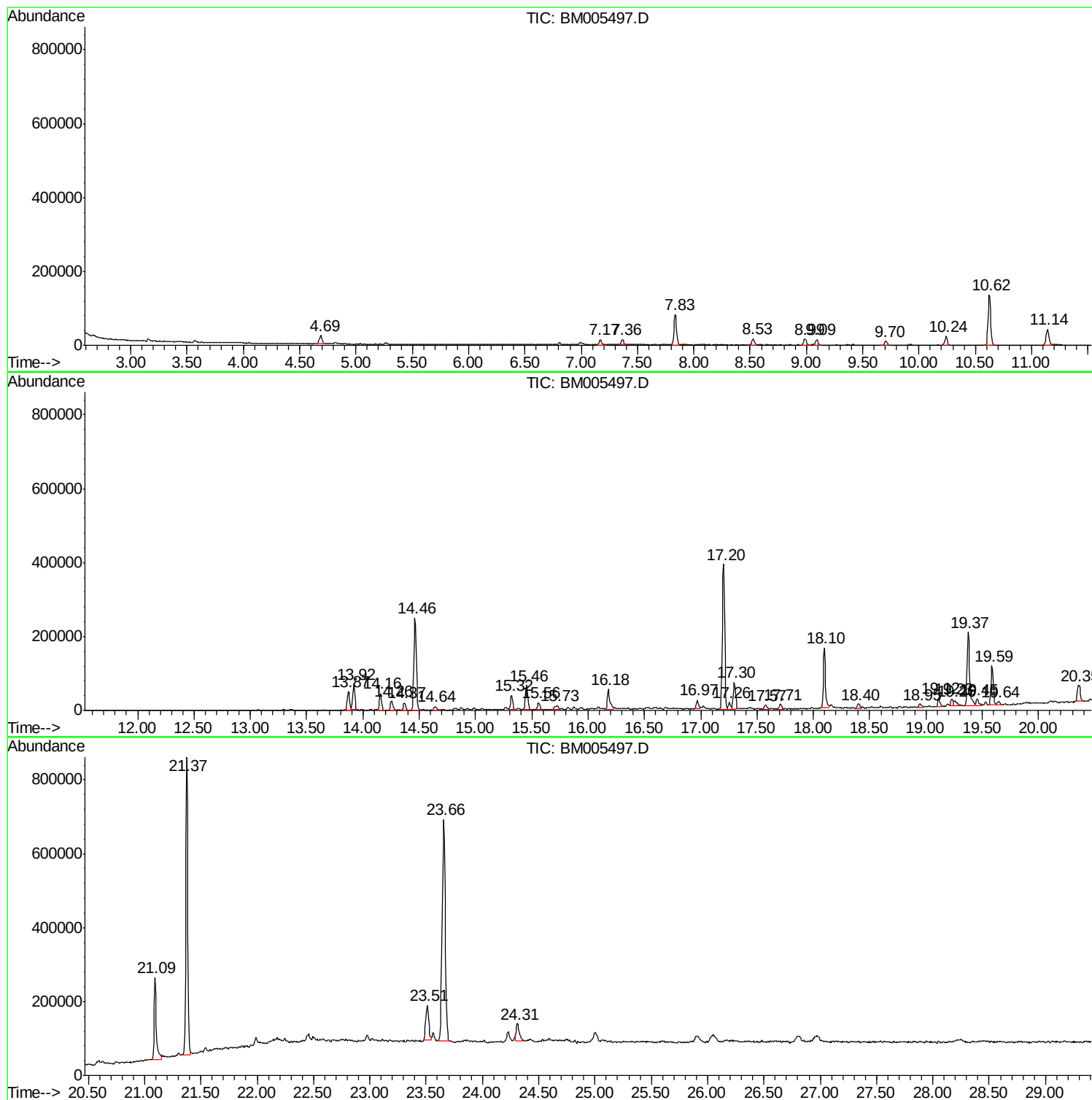
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D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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Sample : H2890-22RX  
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Instrument :  
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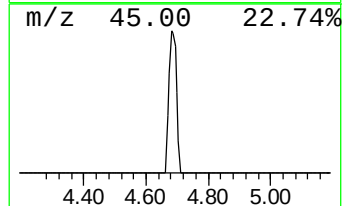
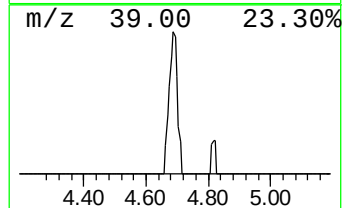
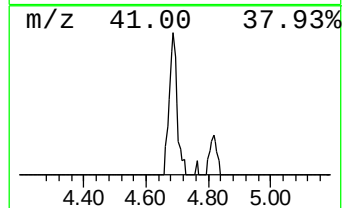
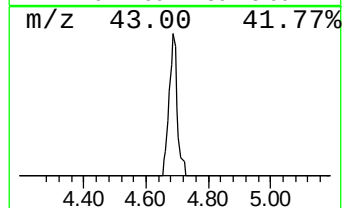
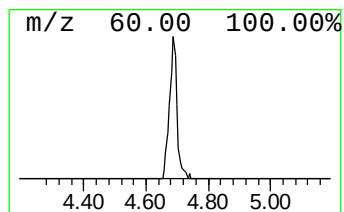
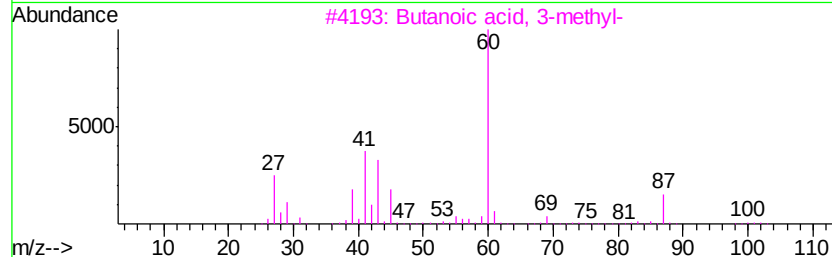
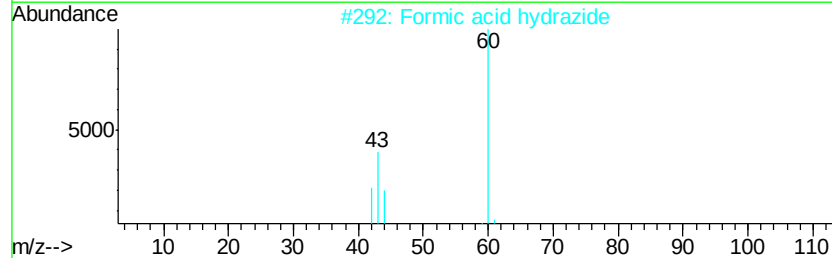
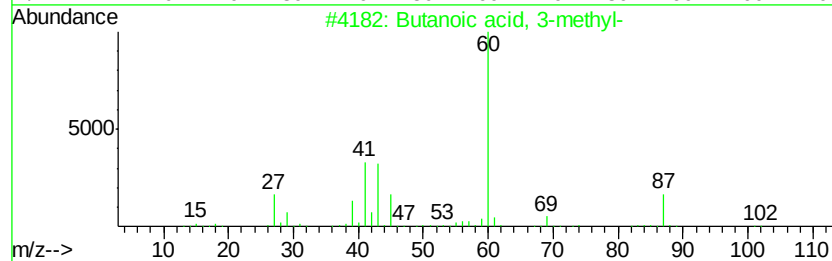
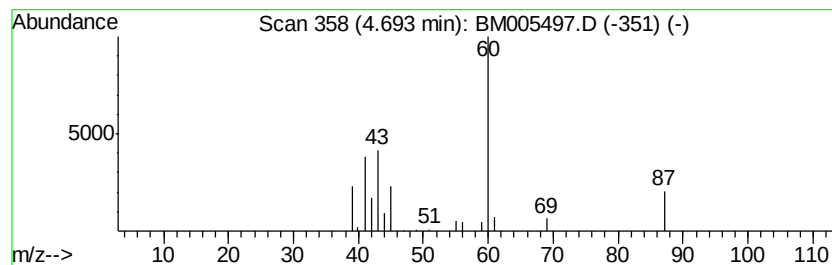
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butanoic acid, 3-methyl- Concentration Rank 5

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 4.69 | 5.65 ng/ul | 37697 | 1,4-Dichlorobenzene-d4 | 7.84 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Butanoic acid, 3-methyl- | 102 | C5H10O2 | 000503-74-2 | 78   |
| 2    |    |   | Formic acid hydrazide    | 60  | CH4N2O  | 000624-84-0 | 50   |
| 3    |    |   | Butanoic acid, 3-methyl- | 102 | C5H10O2 | 000503-74-2 | 43   |
| 4    |    |   | Hexanoic acid            | 116 | C6H12O2 | 000142-62-1 | 39   |
| 5    |    |   | 1-Pentanamine            | 87  | C5H13N  | 000110-58-7 | 12   |



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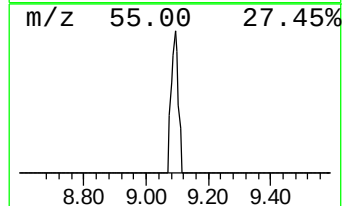
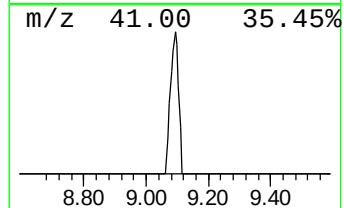
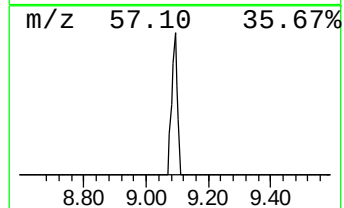
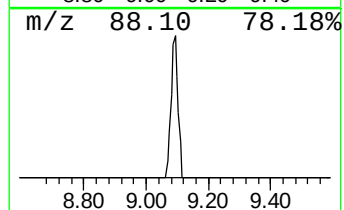
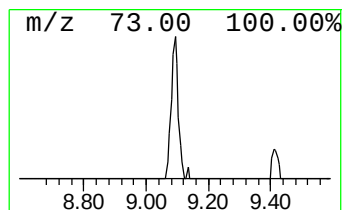
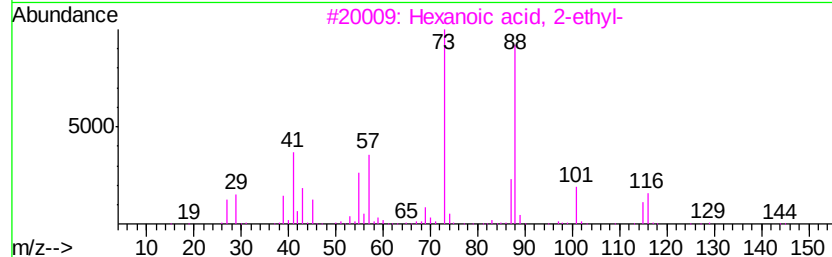
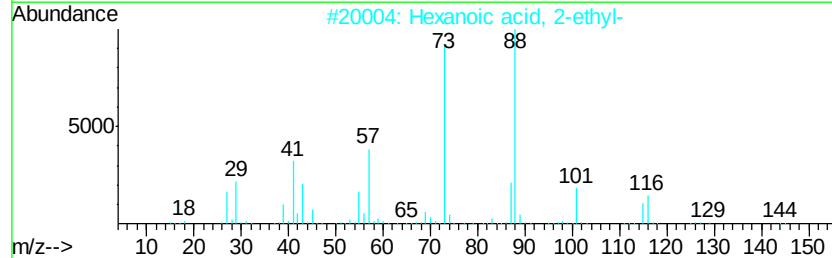
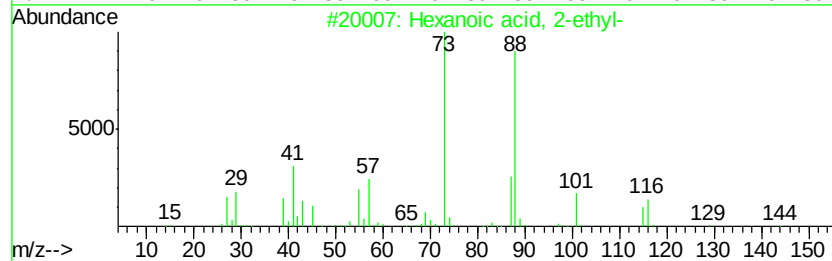
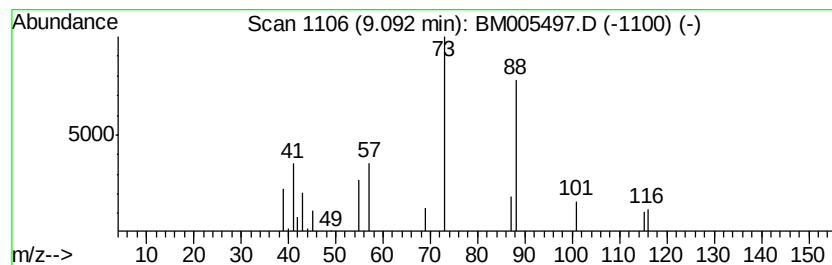
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Hexanoic acid, 2-ethyl- Concentration Rank 6

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 9.09 | 3.70 ng/ul | 24657 | 1,4-Dichlorobenzene-d4 | 7.84 |

| Hit# | of | Tentative ID                    | MW  | MolForm  | CAS#         | Qual |
|------|----|---------------------------------|-----|----------|--------------|------|
| 1    | 5  | Hexanoic acid, 2-ethyl-         | 144 | C8H16O2  | 000149-57-5  | 78   |
| 2    |    | Hexanoic acid, 2-ethyl-         | 144 | C8H16O2  | 000149-57-5  | 78   |
| 3    |    | Hexanoic acid, 2-ethyl-         | 144 | C8H16O2  | 000149-57-5  | 78   |
| 4    |    | Hexanoic acid, 2-ethyl-         | 144 | C8H16O2  | 000149-57-5  | 40   |
| 5    |    | 1-Methyl-1-silacyclopentan-1-ol | 116 | C5H12OSi | 1000216-84-4 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

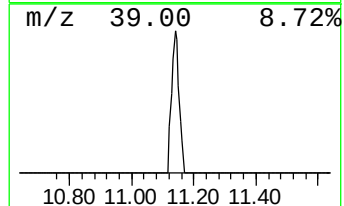
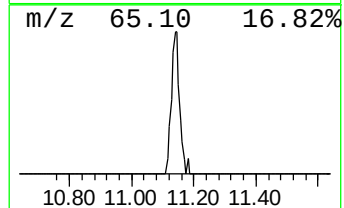
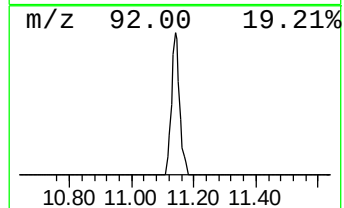
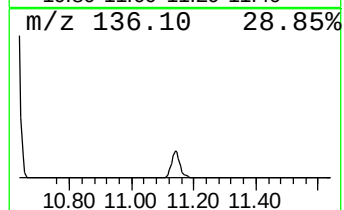
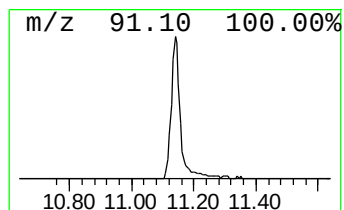
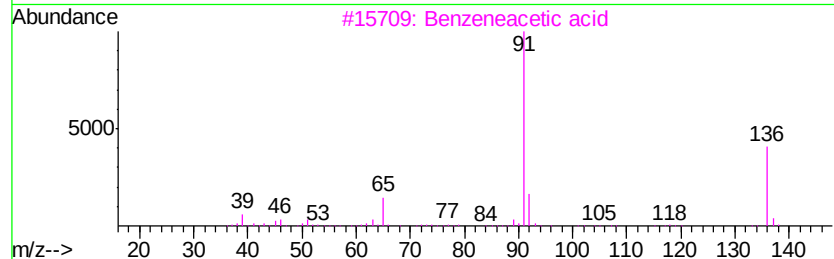
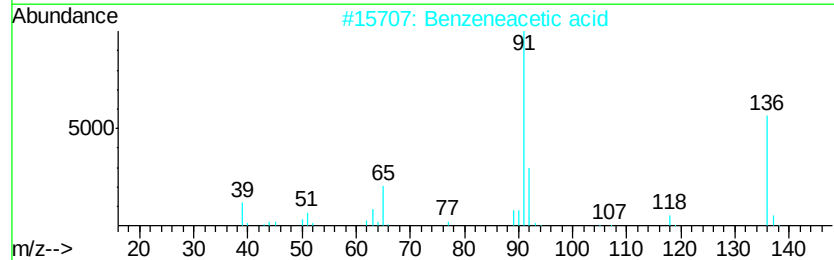
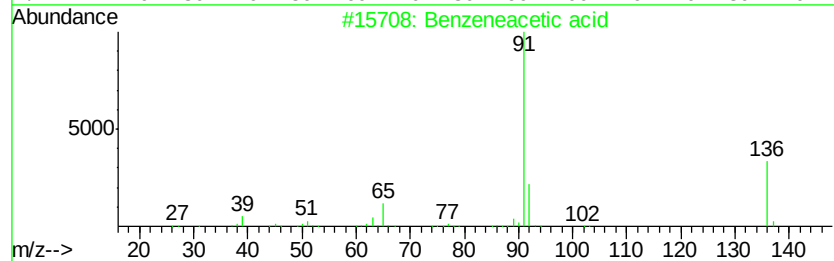
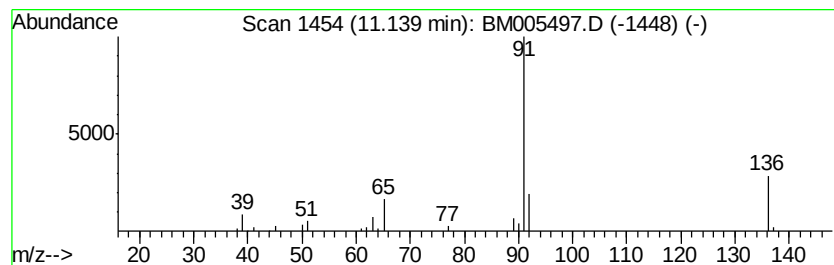
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 Benzeneacetic acid Concentration Rank 2

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 11.14 | 6.78 ng/ul | 77409 | Naphthalene-d8   | 10.62 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzeneacetic acid         | 136 | C8H8O2  | 000103-82-2 | 91   |
| 2    |    | Benzeneacetic acid         | 136 | C8H8O2  | 000103-82-2 | 86   |
| 3    |    | Benzeneacetic acid         | 136 | C8H8O2  | 000103-82-2 | 86   |
| 4    |    | Propanedioic acid, phenyl- | 180 | C9H8O4  | 002613-89-0 | 83   |
| 5    |    | Propanedioic acid, phenyl- | 180 | C9H8O4  | 002613-89-0 | 83   |



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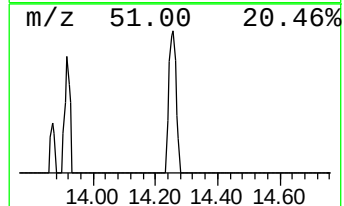
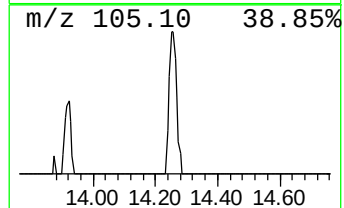
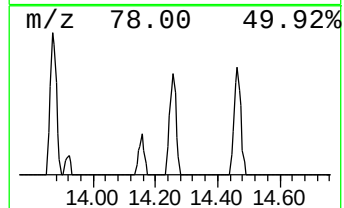
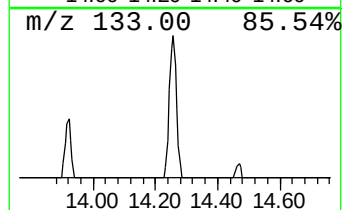
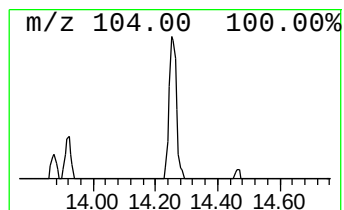
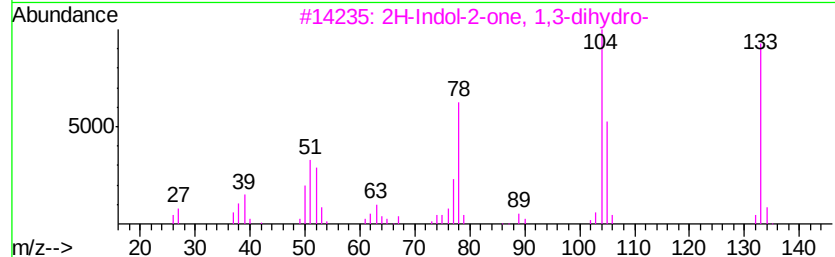
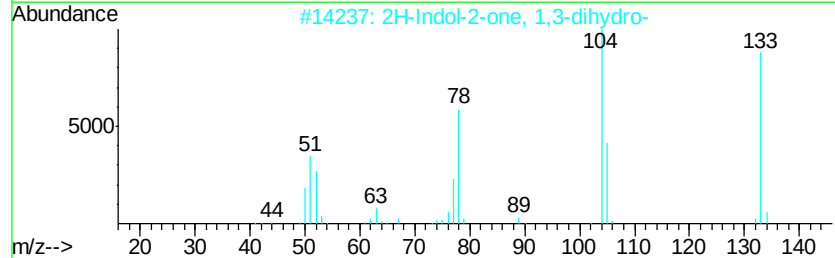
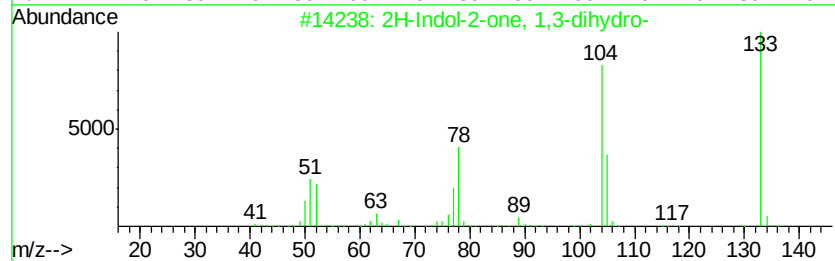
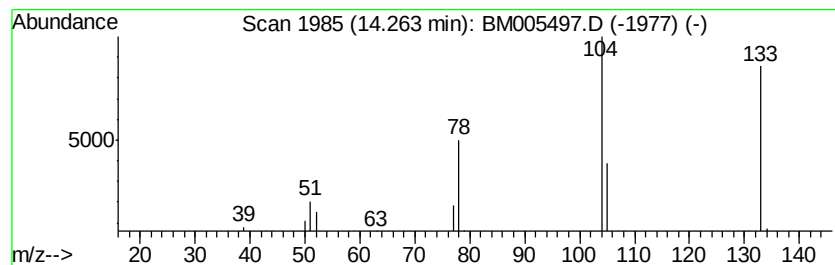
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 2H-Indol-2-one, 1,3-dihydro- Concentration Rank 9

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 14.26 | 2.06 ng/ul | 39070 | Acenaphthene-d10 | 14.46 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2H-Indol-2-one, 1,3-dihydro-     | 133 | C8H7NO  | 000059-48-3 | 94   |
| 2    |    | 2H-Indol-2-one, 1,3-dihydro-     | 133 | C8H7NO  | 000059-48-3 | 91   |
| 3    |    | 2H-Indol-2-one, 1,3-dihydro-     | 133 | C8H7NO  | 000059-48-3 | 91   |
| 4    |    | 1H-Indole, 2,3-dihydro-4-methyl- | 133 | C9H11N  | 062108-16-1 | 91   |
| 5    |    | 1H-Benzotriazole, 4-methyl-      | 133 | C7H7N3  | 029878-31-7 | 90   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

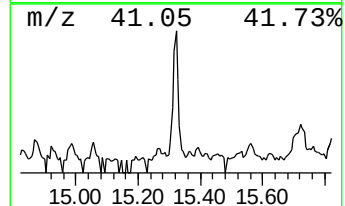
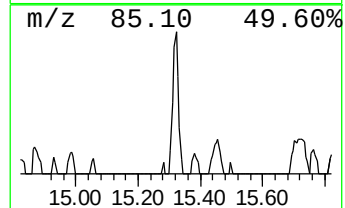
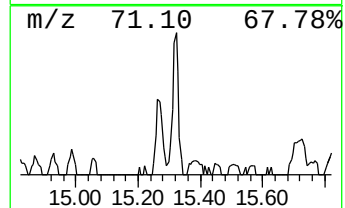
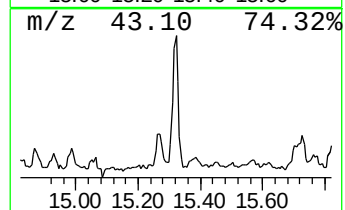
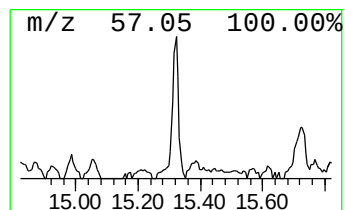
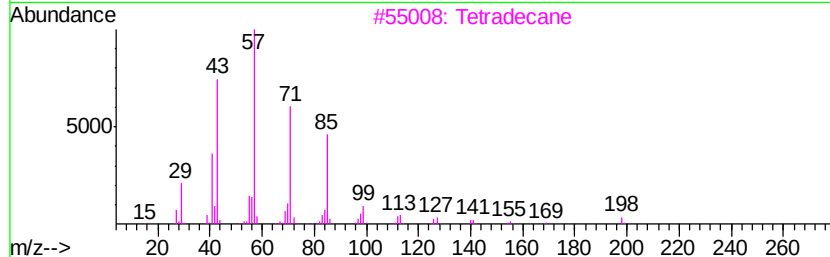
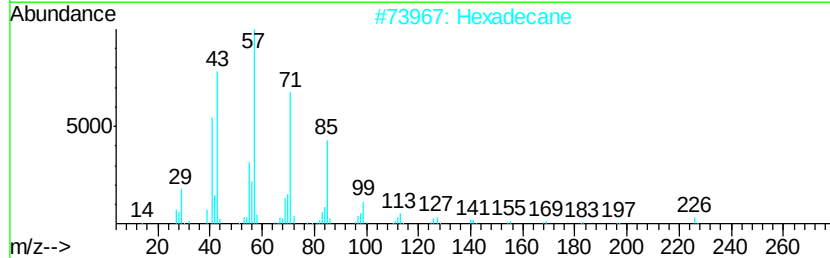
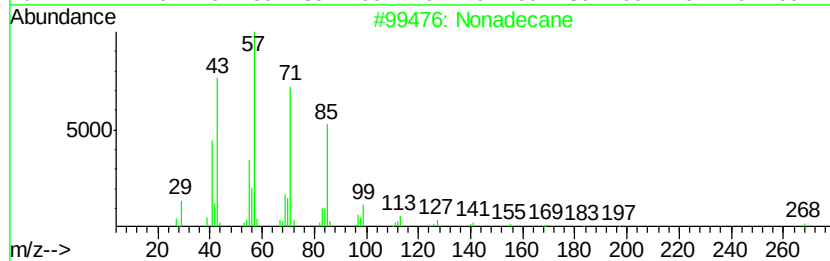
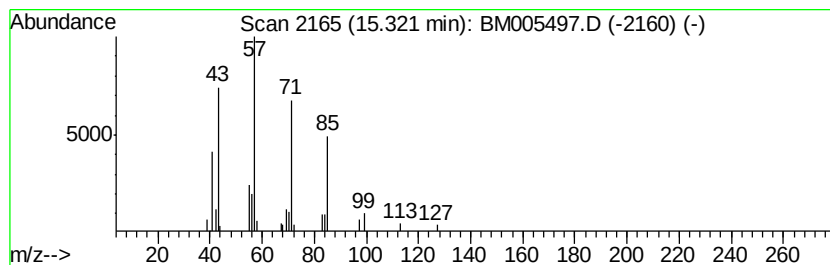
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 15.32 | 2.58 ng/ul | 48901 | Acenaphthene-d10 | 14.46 |

| Hit# of 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|-----------|--------------|-----|---------|-------------|------|
| 1         | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |
| 2         | Hexadecane   | 226 | C16H34  | 000544-76-3 | 86   |
| 3         | Tetradecane  | 198 | C14H30  | 000629-59-4 | 86   |
| 4         | Pentadecane  | 212 | C15H32  | 000629-62-9 | 78   |
| 5         | Pentadecane  | 212 | C15H32  | 000629-62-9 | 72   |





Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

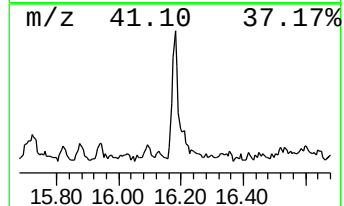
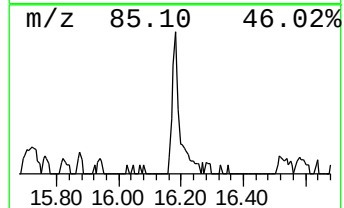
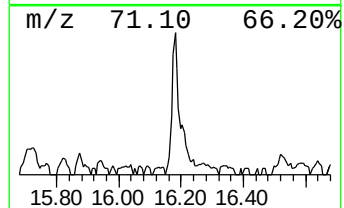
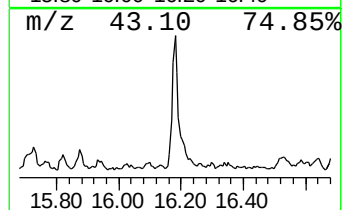
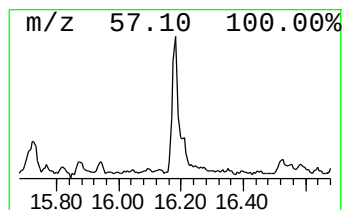
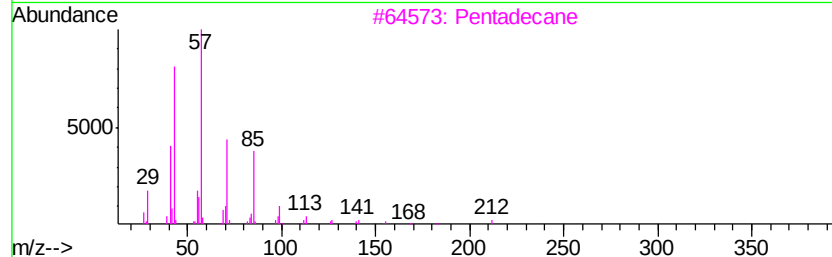
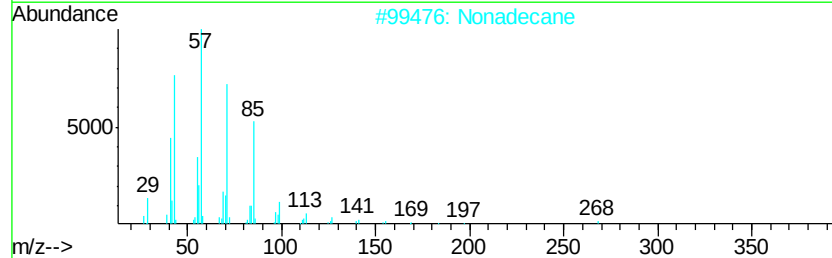
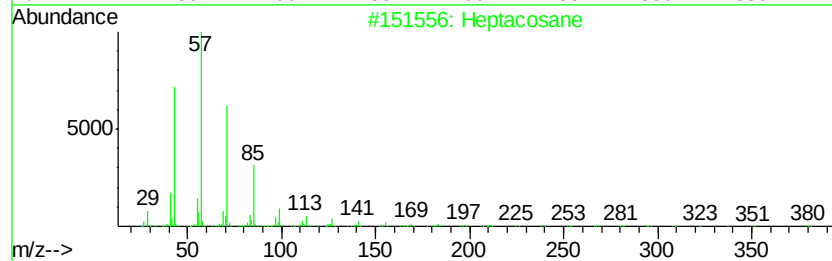
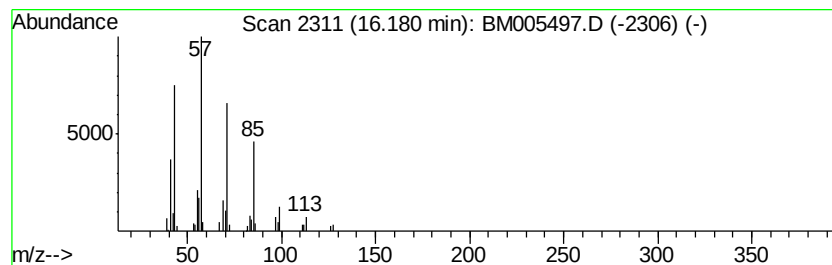
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 16.18 | 2.98 ng/ul | 83224 | Phenanthrene-d10 | 17.20 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Heptacosane  | 380 | C27H56  | 000593-49-7 | 90   |
| 2    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 87   |
| 3    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 86   |
| 4    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 86   |
| 5    |    |   | Pentadecane  | 212 | C15H32  | 000629-62-9 | 86   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

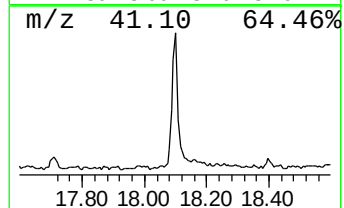
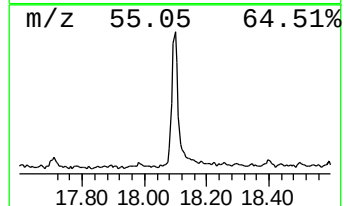
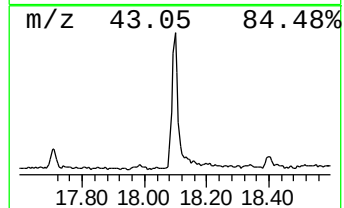
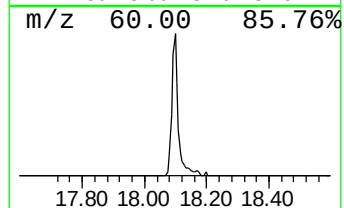
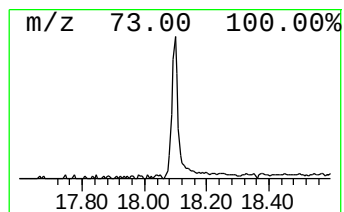
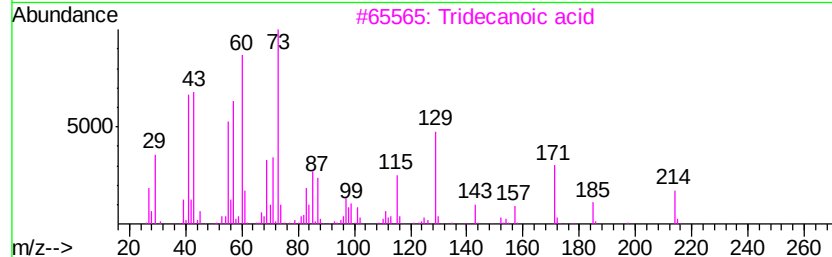
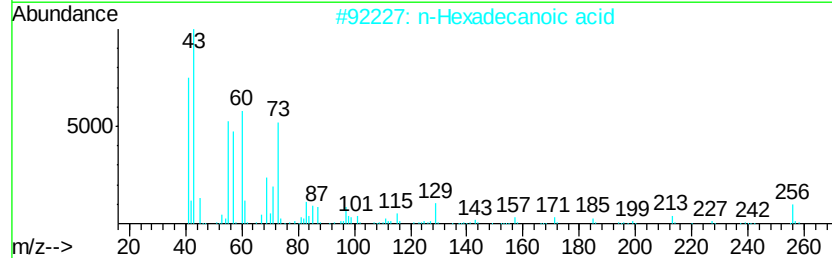
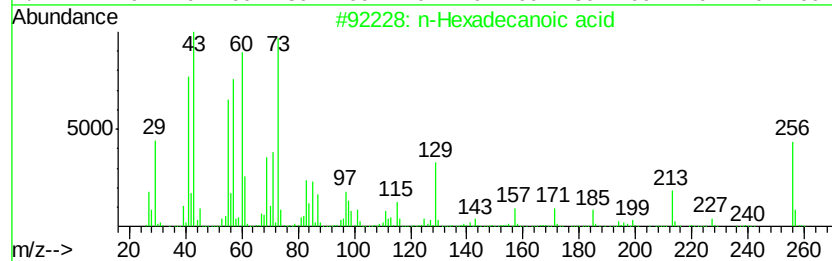
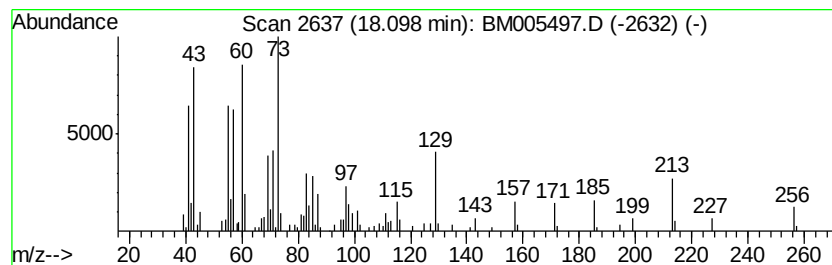
Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

| R.T.    | EstConc    | Area                | Relative to ISTD |          | R.T.        |      |
|---------|------------|---------------------|------------------|----------|-------------|------|
| 18.10   | 7.91 ng/ul | 220784              | Phenanthrene-d10 |          | 17.20       |      |
| Hit# of | 5          | Tentative ID        | MW               | MolForm  | CAS#        | Qual |
| 1       |            | n-Hexadecanoic acid | 256              | C16H32O2 | 000057-10-3 | 98   |
| 2       |            | n-Hexadecanoic acid | 256              | C16H32O2 | 000057-10-3 | 98   |
| 3       |            | Tridecanoic acid    | 214              | C13H26O2 | 000638-53-9 | 91   |
| 4       |            | Tridecanoic acid    | 214              | C13H26O2 | 000638-53-9 | 89   |
| 5       |            | n-Decanoic acid     | 172              | C10H20O2 | 000334-48-5 | 76   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

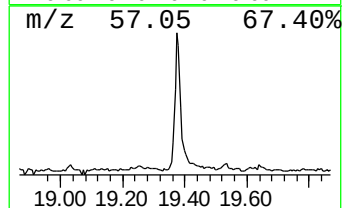
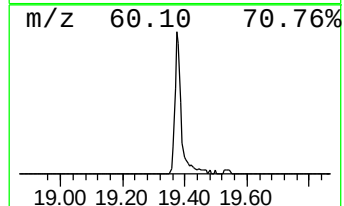
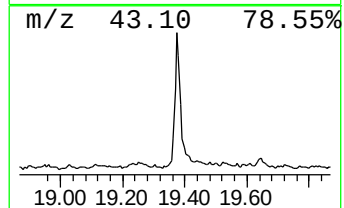
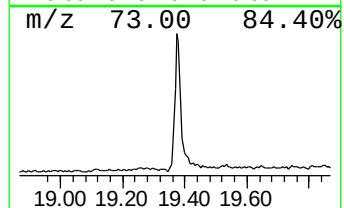
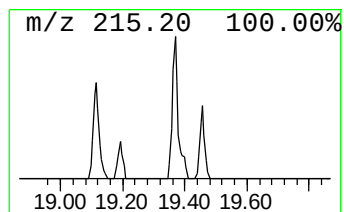
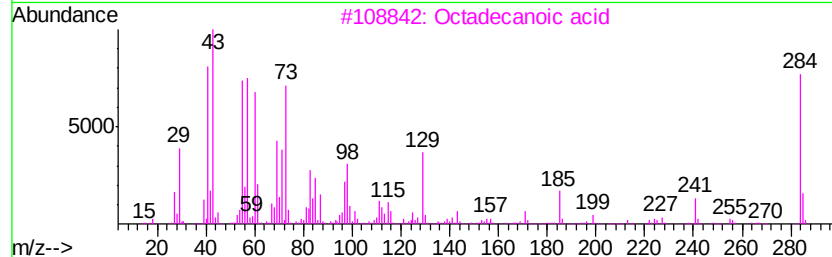
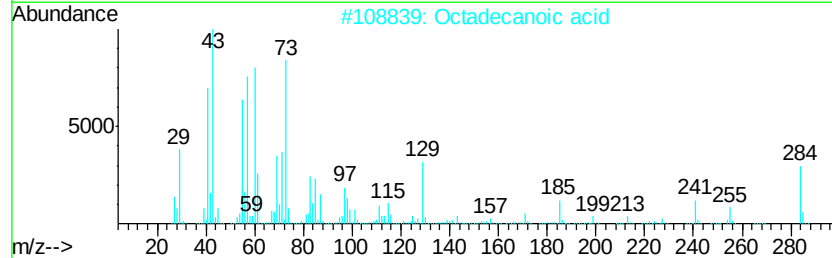
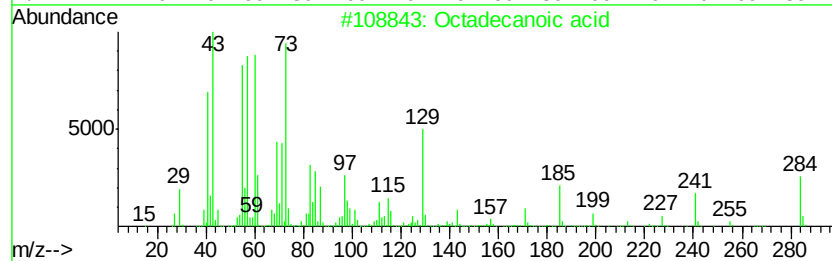
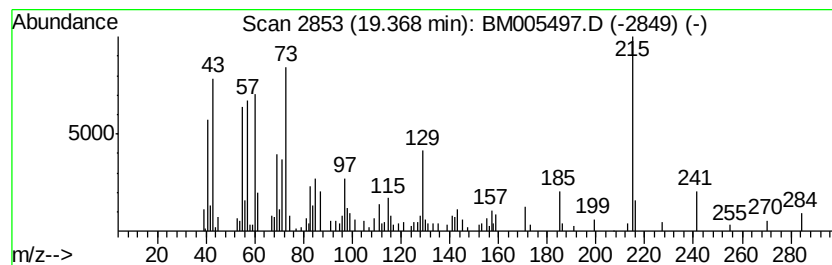
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Octadecanoic acid Concentration Rank 3

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.37 | 6.03 ng/ul | 299452 | Chrysene-d12     | 21.37 |

| Hit# | of | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------|-----|----------|-------------|------|
| 1    | 5  | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 94   |
| 3    |    | Octadecanoic acid  | 284 | C18H36O2 | 000057-11-4 | 94   |
| 4    |    | Heptadecanoic acid | 270 | C17H34O2 | 000506-12-7 | 92   |
| 5    |    | Heptadecanoic acid | 270 | C17H34O2 | 000506-12-7 | 86   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

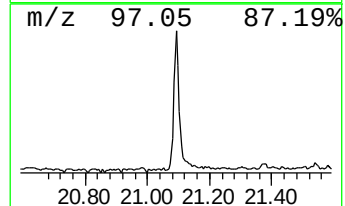
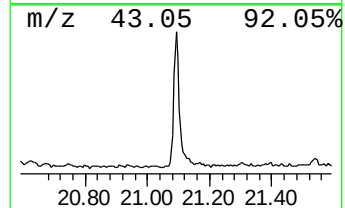
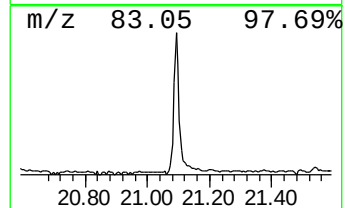
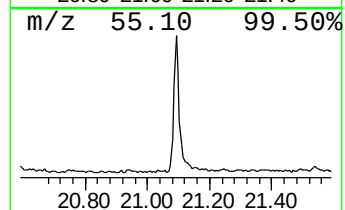
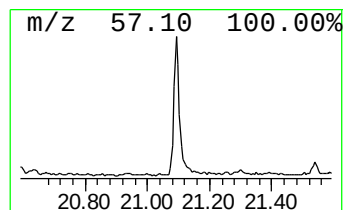
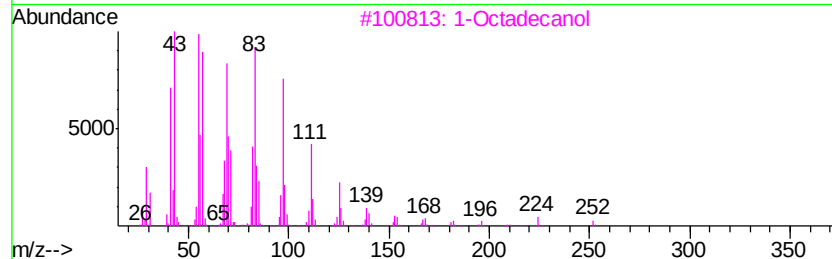
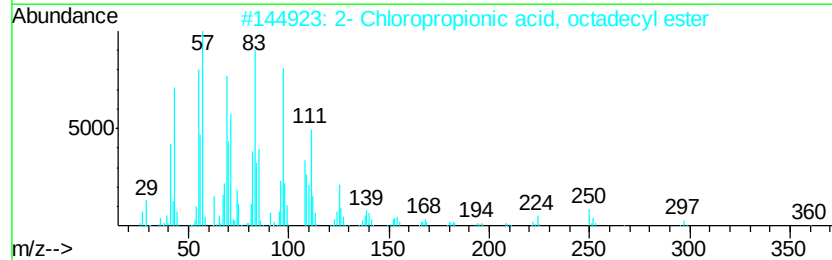
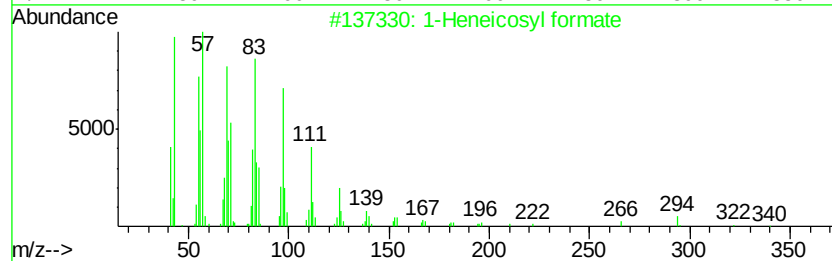
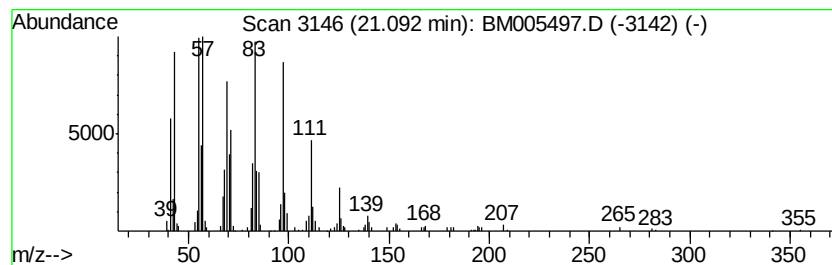
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 1-Heneicosyl formate Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.09 | 6.00 ng/ul | 298272 | Chrysene-d12     | 21.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | 1-Heneicosyl formate                | 340 | C22H44O2    | 077899-03-7 | 94   |
| 2    |    | 2- Chloropropionic acid, octadec... | 360 | C21H41ClO2  | 088104-31-8 | 91   |
| 3    |    | 1-Octadecanol                       | 270 | C18H38O     | 000112-92-5 | 91   |
| 4    |    | Trichloroacetic acid, tetradecyl... | 358 | C16H29Cl3O2 | 074339-52-9 | 87   |
| 5    |    | 1-Nonadecene                        | 266 | C19H38      | 018435-45-5 | 87   |



Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM051816\  
Data File : BM005497.D  
Acq On : 19 May 2016 07:49  
Operator : UM/SJ  
Sample : H2890-22RX  
Misc :  
ALS Vial : 37 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
D9X07RX

Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM051816.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Butanoic acid, 3-... | 4.69  | 5.7     | ng/ul | 37697    | 1                     | 7.84  | 133361 | 20.0 |
| Hexanoic acid, 2-... | 9.09  | 3.7     | ng/ul | 24657    | 1                     | 7.84  | 133361 | 20.0 |
| Benzeneacetic acid   | 11.14 | 6.8     | ng/ul | 77409    | 2                     | 10.62 | 228254 | 20.0 |
| 2H-Indol-2-one, 1... | 14.26 | 2.1     | ng/ul | 39070    | 3                     | 14.46 | 379477 | 20.0 |
| (DEL) Alkane: Str... | 15.32 | 2.6     | ng/ul | 48901    | 3                     | 14.46 | 379477 | 20.0 |
| (DEL) Alkane: Str... | 16.18 | 3.0     | ng/ul | 83224    | 4                     | 17.20 | 558320 | 20.0 |
| n-Hexadecanoic acid  | 18.10 | 7.9     | ng/ul | 220784   | 4                     | 17.20 | 558320 | 20.0 |
| Octadecanoic acid    | 19.37 | 6.0     | ng/ul | 299452   | 5                     | 21.37 | 993990 | 20.0 |
| 1-Heneicosyl formate | 21.09 | 6.0     | ng/ul | 298272   | 5                     | 21.37 | 993990 | 20.0 |