

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
 Data File : BM016955.D  
 Acq On : 02 Oct 2018 11:28  
 Operator : MJ/SJ  
 Sample : J5126-01  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0B89

## Integration Parameters: LSCINT.P

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

Filtering: 5  
 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.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         | 3.187       | 31            | 34          | 39           | rVB      | 1181457        | 1242755       | 6.42%           | 0.573%        |
| 2         | 7.234       | 717           | 722         | 729          | rVB      | 25565          | 43089         | 0.22%           | 0.020%        |
| 3         | 7.710       | 795           | 803         | 814          | rVB      | 1376030        | 2192437       | 11.32%          | 1.011%        |
| 4         | 8.057       | 858           | 862         | 867          | rVB2     | 20198          | 32983         | 0.17%           | 0.015%        |
| 5         | 8.928       | 1006          | 1010        | 1016         | rVB2     | 14013          | 22127         | 0.11%           | 0.010%        |
| 6         | 10.351      | 1244          | 1252        | 1264         | rBV      | 2048798        | 3412802       | 17.63%          | 1.574%        |
| 7         | 10.486      | 1266          | 1275        | 1286         | rVV      | 1785074        | 3108357       | 16.06%          | 1.434%        |
| 8         | 10.651      | 1300          | 1303        | 1311         | rVB      | 26612          | 39902         | 0.21%           | 0.018%        |
| 9         | 10.869      | 1332          | 1340        | 1348         | rBV      | 398037         | 668695        | 3.45%           | 0.308%        |
| 10        | 11.404      | 1427          | 1431        | 1437         | rBV7     | 11830          | 24216         | 0.13%           | 0.011%        |
| 11        | 11.516      | 1445          | 1450        | 1455         | rBV      | 48507          | 84520         | 0.44%           | 0.039%        |
| 12        | 11.775      | 1488          | 1494        | 1496         | rBV6     | 21454          | 31307         | 0.16%           | 0.014%        |
| 13        | 11.922      | 1512          | 1519        | 1524         | rBV      | 64529          | 100268        | 0.52%           | 0.046%        |
| 14        | 12.139      | 1554          | 1556        | 1564         | rVB5     | 23853          | 39313         | 0.20%           | 0.018%        |
| 15        | 12.233      | 1566          | 1572        | 1578         | rBV8     | 17215          | 39158         | 0.20%           | 0.018%        |
| 16        | 12.363      | 1591          | 1594        | 1600         | rBV8     | 11611          | 21315         | 0.11%           | 0.010%        |
| 17        | 12.522      | 1613          | 1621        | 1631         | rBV      | 200465         | 369987        | 1.91%           | 0.171%        |
| 18        | 12.774      | 1661          | 1664        | 1670         | rVB5     | 36693          | 53447         | 0.28%           | 0.025%        |
| 19        | 12.833      | 1670          | 1674        | 1678         | rBV5     | 20597          | 35885         | 0.19%           | 0.017%        |
| 20        | 12.880      | 1679          | 1682        | 1687         | rVB      | 63003          | 87268         | 0.45%           | 0.040%        |
| 21        | 12.939      | 1688          | 1692        | 1696         | rBV4     | 24792          | 41606         | 0.21%           | 0.019%        |
| 22        | 13.139      | 1719          | 1726        | 1730         | rBV      | 416408         | 702847        | 3.63%           | 0.324%        |
| 23        | 13.174      | 1730          | 1732        | 1739         | rVB      | 65305          | 83696         | 0.43%           | 0.039%        |
| 24        | 13.274      | 1744          | 1749        | 1752         | rBV6     | 25028          | 42627         | 0.22%           | 0.020%        |
| 25        | 13.557      | 1794          | 1797        | 1802         | rBV6     | 15103          | 25104         | 0.13%           | 0.012%        |
| 26        | 13.663      | 1811          | 1815        | 1819         | rBV6     | 30906          | 51401         | 0.27%           | 0.024%        |
| 27        | 13.763      | 1828          | 1832        | 1837         | rVB      | 170875         | 228566        | 1.18%           | 0.105%        |
| 28        | 13.839      | 1841          | 1845        | 1848         | rVV      | 108650         | 143837        | 0.74%           | 0.066%        |
| 29        | 13.880      | 1849          | 1852        | 1860         | rVB4     | 63670          | 119755        | 0.62%           | 0.055%        |
| 30        | 14.086      | 1882          | 1887        | 1893         | rBV6     | 91779          | 185118        | 0.96%           | 0.085%        |
| 31        | 14.180      | 1898          | 1903        | 1910         | rVB      | 266667         | 430271        | 2.22%           | 0.198%        |
| 32        | 14.257      | 1911          | 1916        | 1920         | rBV2     | 119098         | 190633        | 0.98%           | 0.088%        |
| 33        | 14.351      | 1921          | 1932        | 1940         | rBV2     | 2338438        | 3932128       | 20.31%          | 1.814%        |
| 34        | 14.880      | 2018          | 2022        | 2029         | rBV4     | 171726         | 285077        | 1.47%           | 0.131%        |

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Instrument :  
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## Integration Parameters: LSCINT.P

Integrator: RTE  
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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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |         |        |
|----|--------|------|------|------|------|----------|----------|---------|--------|
| 35 | 15.068 | 2050 | 2054 | 2060 | rBV6 | 126256   | 251023   | 1.30%   | 0.116% |
| 36 | 15.139 | 2061 | 2066 | 2073 | rVV2 | 352819   | 600558   | 3.10%   | 0.277% |
| 37 | 15.621 | 2144 | 2148 | 2154 | rVB  | 385477   | 605937   | 3.13%   | 0.279% |
| 38 | 16.104 | 2222 | 2230 | 2235 | rBV2 | 1818416  | 3351073  | 17.31%  | 1.546% |
| 39 | 16.921 | 2366 | 2369 | 2376 | rVB  | 1308840  | 2055123  | 10.62%  | 0.948% |
| 40 | 17.098 | 2395 | 2399 | 2405 | rBV2 | 2563605  | 4167360  | 21.53%  | 1.922% |
| 41 | 19.027 | 2723 | 2727 | 2732 | rVB  | 4300567  | 5710897  | 29.50%  | 2.634% |
| 42 | 19.309 | 2771 | 2775 | 2780 | rVB  | 5849598  | 7277523  | 37.59%  | 3.357% |
| 43 | 19.739 | 2844 | 2848 | 2850 | rBV  | 2583476  | 3859452  | 19.94%  | 1.780% |
| 44 | 19.880 | 2868 | 2872 | 2876 | rVB  | 5565277  | 6402490  | 33.07%  | 2.953% |
| 45 | 19.927 | 2876 | 2880 | 2887 | rVB2 | 2153282  | 3974766  | 20.53%  | 1.833% |
| 46 | 20.021 | 2892 | 2896 | 2901 | rVB  | 13358447 | 16918962 | 87.39%  | 7.804% |
| 47 | 20.221 | 2927 | 2930 | 2938 | rVB2 | 2038956  | 2662936  | 13.75%  | 1.228% |
| 48 | 20.303 | 2939 | 2944 | 2948 | rVB  | 16015892 | 19069344 | 98.50%  | 8.796% |
| 49 | 20.356 | 2948 | 2953 | 2959 | rBV  | 4116363  | 5266591  | 27.20%  | 2.429% |
| 50 | 20.456 | 2966 | 2970 | 2978 | rVB2 | 5845643  | 9136773  | 47.19%  | 4.214% |
| 51 | 20.550 | 2983 | 2986 | 2989 | rVV2 | 1608043  | 1851565  | 9.56%   | 0.854% |
| 52 | 20.615 | 2990 | 2997 | 3003 | rVV  | 9218655  | 18450799 | 95.30%  | 8.511% |
| 53 | 20.668 | 3003 | 3006 | 3010 | rVB  | 1503750  | 1543317  | 7.97%   | 0.712% |
| 54 | 20.762 | 3018 | 3022 | 3028 | rVB  | 8264598  | 10227809 | 52.83%  | 4.718% |
| 55 | 20.827 | 3029 | 3033 | 3041 | rVB  | 4400951  | 5355831  | 27.66%  | 2.470% |
| 56 | 20.921 | 3046 | 3049 | 3053 | rBV2 | 998752   | 1037468  | 5.36%   | 0.479% |
| 57 | 20.962 | 3053 | 3056 | 3062 | rBV3 | 846311   | 1073249  | 5.54%   | 0.495% |
| 58 | 21.033 | 3064 | 3068 | 3073 | rVB  | 8000679  | 9305436  | 48.07%  | 4.292% |
| 59 | 21.092 | 3073 | 3078 | 3083 | rVB2 | 3969231  | 5500806  | 28.41%  | 2.537% |
| 60 | 21.145 | 3083 | 3087 | 3090 | rBV  | 1810717  | 2265284  | 11.70%  | 1.045% |
| 61 | 21.244 | 3101 | 3104 | 3107 | rVV  | 959017   | 1022397  | 5.28%   | 0.472% |
| 62 | 21.292 | 3107 | 3112 | 3114 | rVV  | 3010649  | 4515766  | 23.33%  | 2.083% |
| 63 | 21.321 | 3114 | 3117 | 3123 | rVB  | 15717579 | 19359851 | 100.00% | 8.930% |
| 64 | 21.633 | 3165 | 3170 | 3175 | rBV  | 6471484  | 8640345  | 44.63%  | 3.985% |
| 65 | 21.686 | 3176 | 3179 | 3186 | rVB  | 2890732  | 3866012  | 19.97%  | 1.783% |
| 66 | 21.756 | 3187 | 3191 | 3197 | rVB  | 3445520  | 4322039  | 22.32%  | 1.994% |
| 67 | 22.103 | 3246 | 3250 | 3255 | rVB  | 1082017  | 1405508  | 7.26%   | 0.648% |
| 68 | 22.321 | 3283 | 3287 | 3294 | rVB2 | 2479159  | 3459234  | 17.87%  | 1.596% |
| 69 | 23.521 | 3485 | 3491 | 3499 | rVB  | 2189079  | 4171991  | 21.55%  | 1.924% |

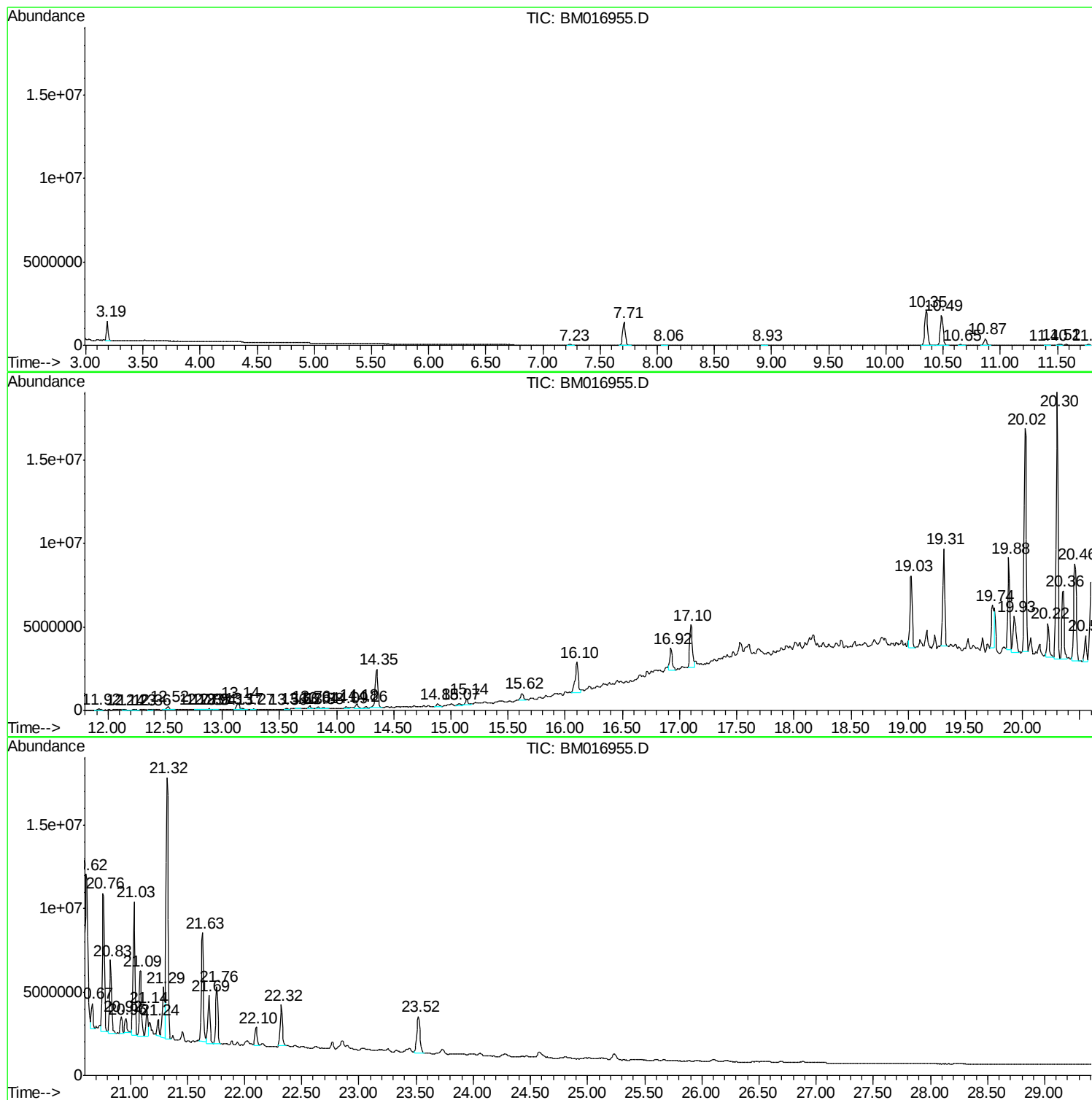
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Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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



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Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

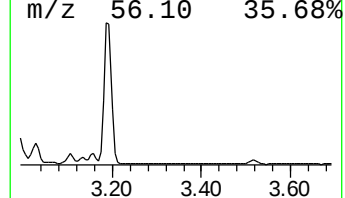
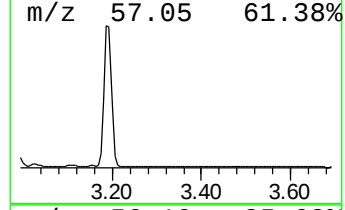
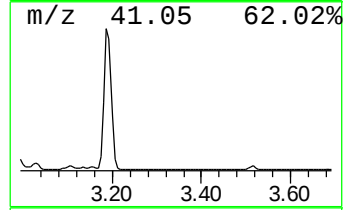
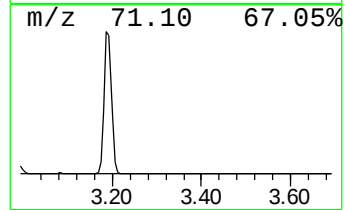
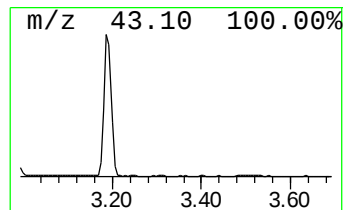
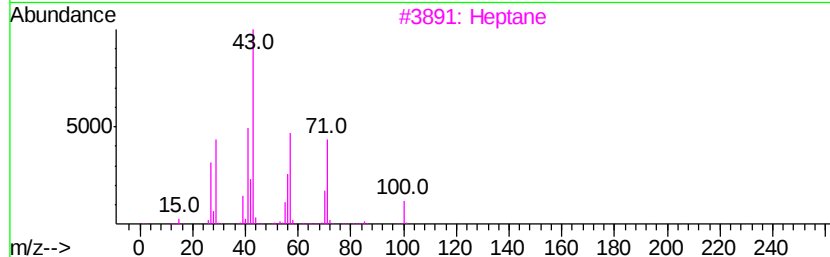
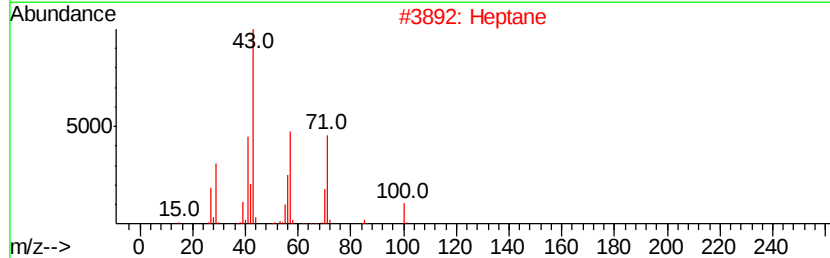
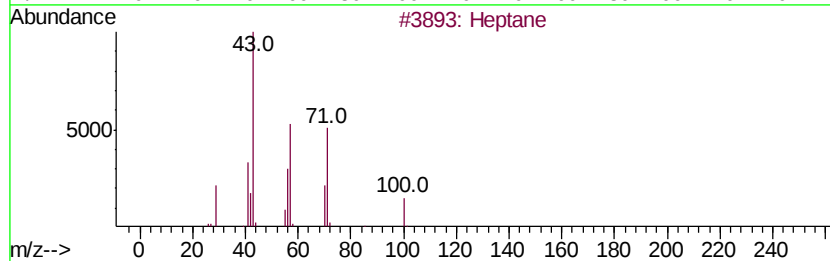
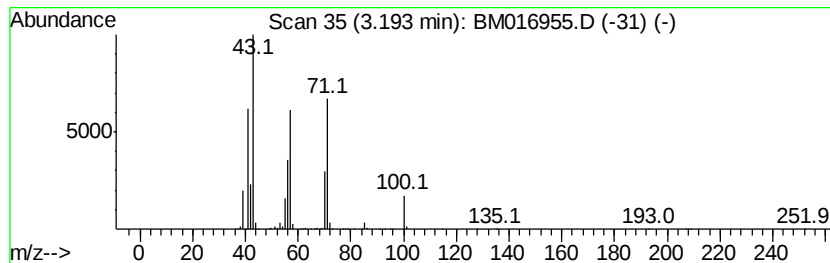
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 3.19 | 11.34 ng/ul | 1242760 | 1,4-Dichlorobenzene-d4 | 7.71 |

| Hit# | of | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------|-----|---------|-------------|------|
| 1    | 5  | Heptane           | 100 | C7H16   | 000142-82-5 | 91   |
| 2    |    | Heptane           | 100 | C7H16   | 000142-82-5 | 91   |
| 3    |    | Heptane           | 100 | C7H16   | 000142-82-5 | 91   |
| 4    |    | Heptane           | 100 | C7H16   | 000142-82-5 | 90   |
| 5    |    | Hexane, 3-methyl- | 100 | C7H16   | 000589-34-4 | 40   |



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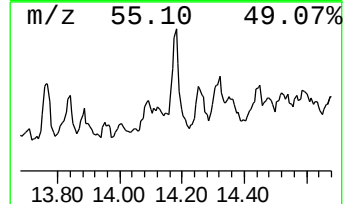
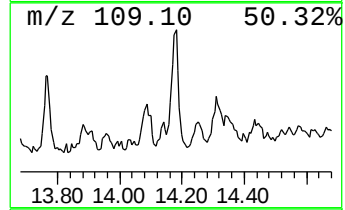
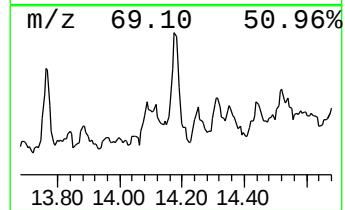
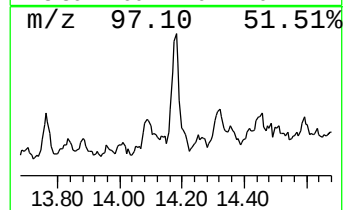
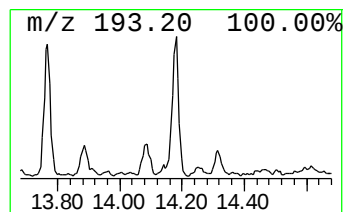
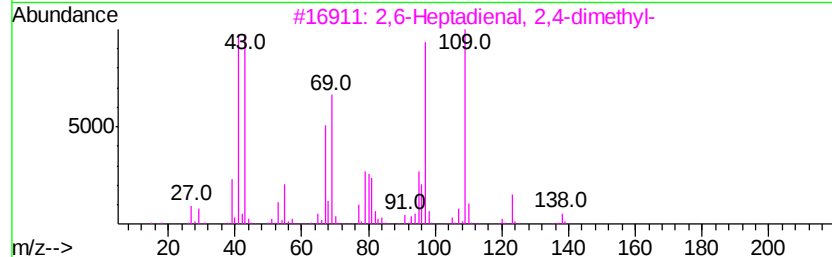
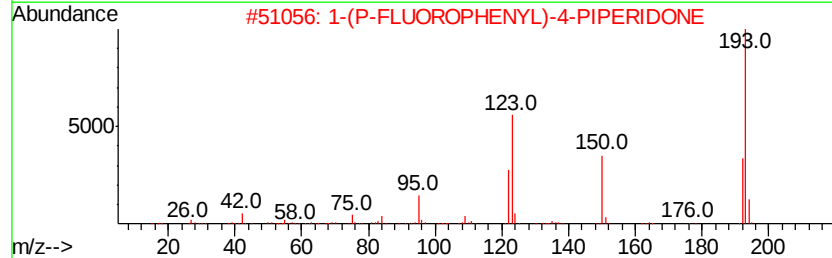
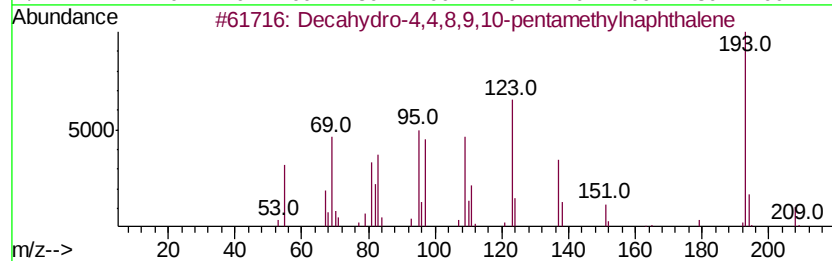
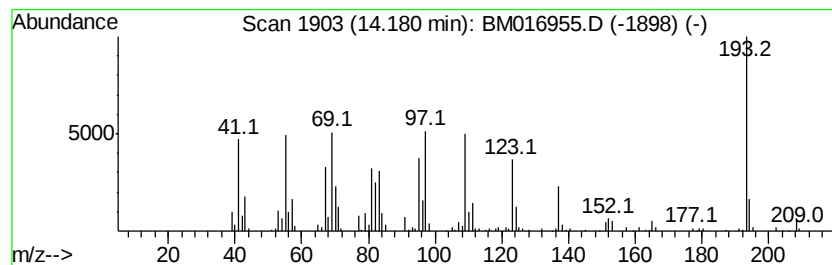
Instrument :  
BNA\_M  
ClientSampled :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Decahydro-4,4,8,9,10-pentam... Concentration Rank 33

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.            |
|---------|------------|-------------------------------------|------------------|-----------------|
| 14.18   | 2.19 ng/ul | 430271                              | Acenaphthene-d10 | 14.35           |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |            | Decahydro-4,4,8,9,10-pentamethyl... | 208 C15H28       | 080655-44-3 64  |
| 2       |            | 1-(P-FLUOROPHENYL)-4-PIPERIDONE     | 193 C11H12FNO    | 1000238-56-7 43 |
| 3       |            | 2,6-Heptadienal, 2,4-dimethyl-      | 138 C9H14O       | 085136-08-9 30  |
| 4       |            | 1H-Indole, 2-phenyl-                | 193 C14H11N      | 000948-65-2 25  |
| 5       |            | 2-Anthracenamine                    | 193 C14H11N      | 000613-13-8 22  |



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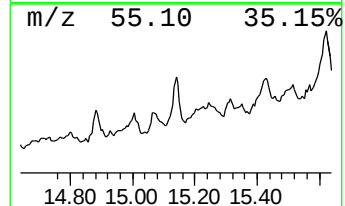
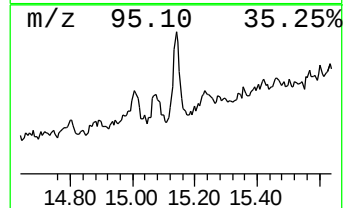
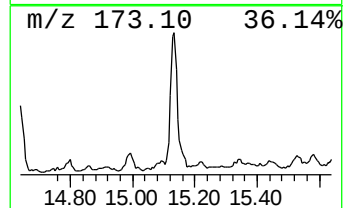
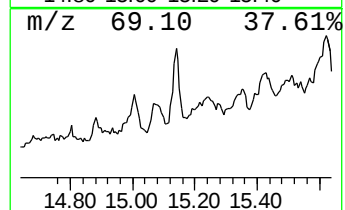
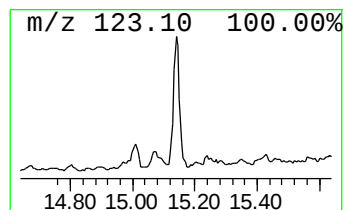
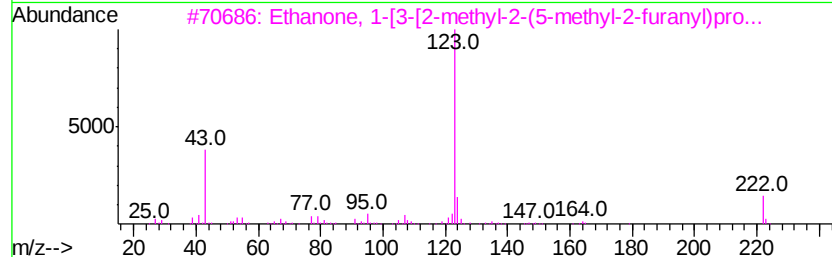
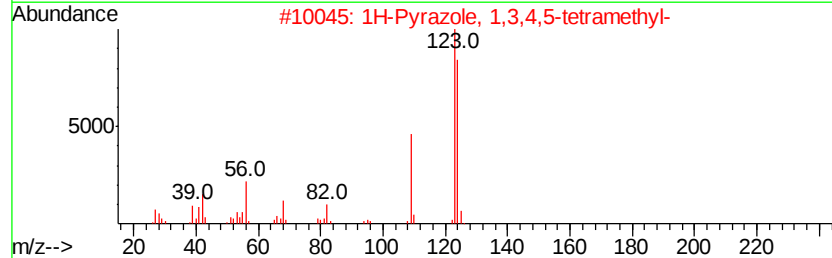
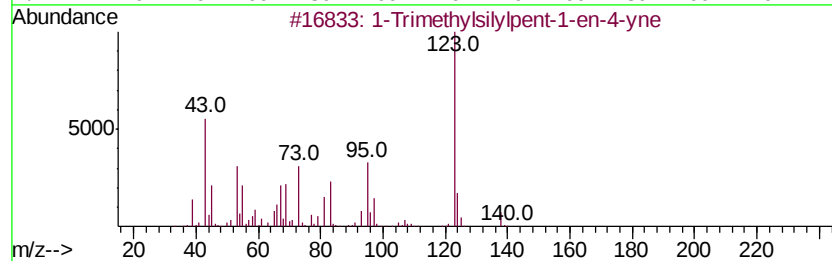
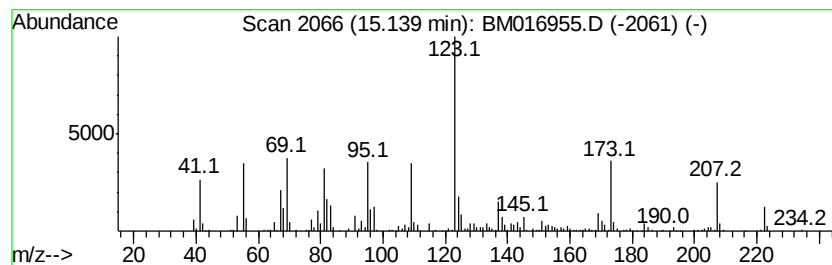
Instrument :  
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Quant Title : SVOA CALIBRATION

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

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Peak Number 5 unknown-01 Concentration Rank 32

| R.T.      | EstConc                                               | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------------------------|--------|------------------|--------------|------|
| 15.14     | 3.05 ng/ul                                            | 600558 | Acenaphthene-d10 | 14.35        |      |
| Hit# of 5 | Tentative ID                                          | MW     | MolForm          | CAS#         | Qual |
| 1         | 1-Trimethylsilylpent-1-en-4-yne                       | 138    | C8H14Si          | 079516-25-9  | 43   |
| 2         | 1H-Pyrazole, 1,3,4,5-tetramethyl-                     | 124    | C7H12N2          | 001073-20-7  | 38   |
| 3         | Ethanone, 1-[3-[2-methyl-2-(5-methyl-2-furanyl)pro... | 222    | C13H18O3         | 080114-28-9  | 38   |
| 4         | 3-Fluorobenzoic acid, 2-pentadec...                   | 350    | C22H35FO2        | 1000280-60-7 | 38   |
| 5         | 3-Fluorobenzoic acid, 2-ethylcyc...                   | 250    | C15H19FO2        | 1000279-04-9 | 38   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

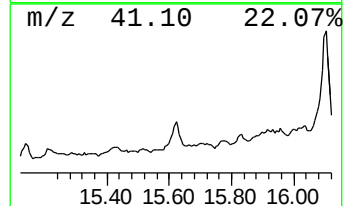
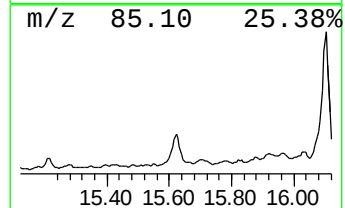
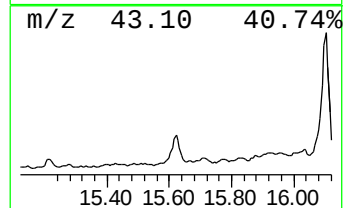
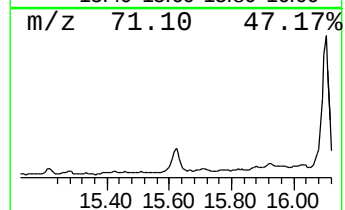
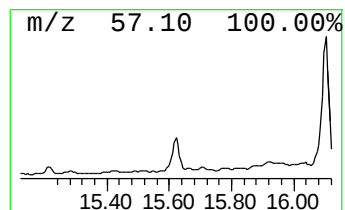
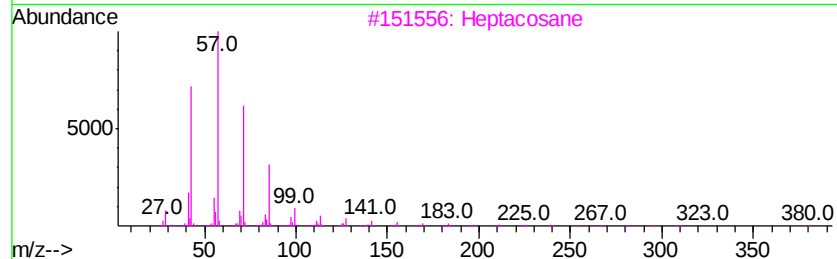
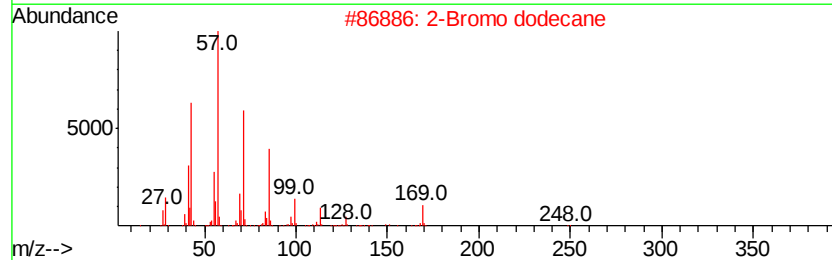
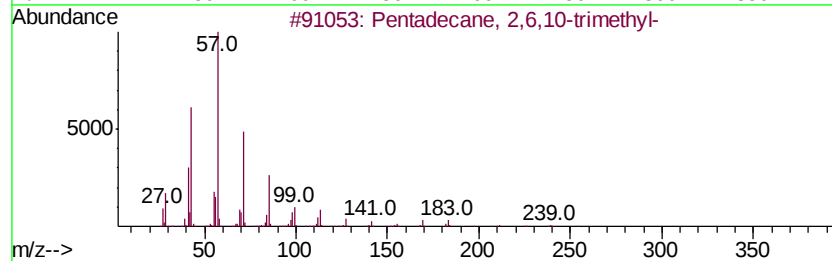
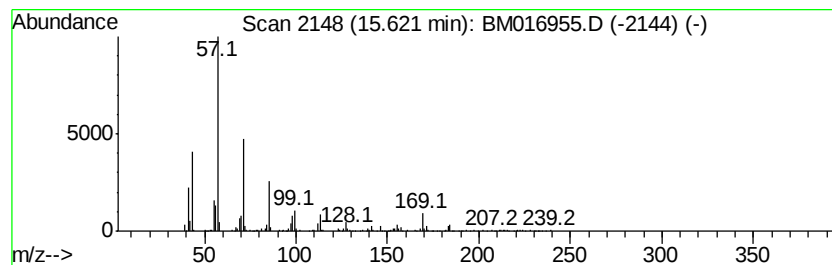
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 31

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.62 | 3.08 ng/ul | 605937 | Acenaphthene-d10 | 14.35 |

| Hit# | of | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------|-----|----------|-------------|------|
| 1    | 5  | Pentadecane, 2,6,10-trimethyl- | 254 | C18H38   | 003892-00-0 | 80   |
| 2    |    | 2-Bromo dodecane               | 248 | C12H25Br | 013187-99-0 | 80   |
| 3    |    | Heptacosane                    | 380 | C27H56   | 000593-49-7 | 80   |
| 4    |    | Dodecane, 1-iodo-              | 296 | C12H25I  | 004292-19-7 | 64   |
| 5    |    | Decane, 2,6,8-trimethyl-       | 184 | C13H28   | 062108-26-3 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

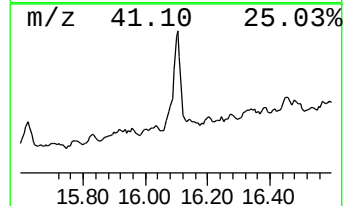
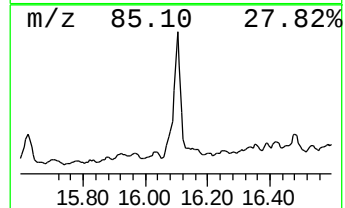
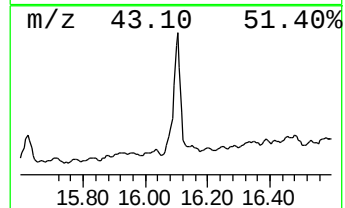
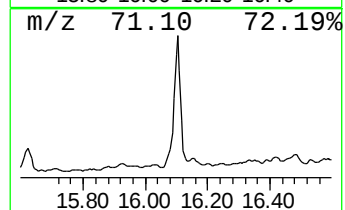
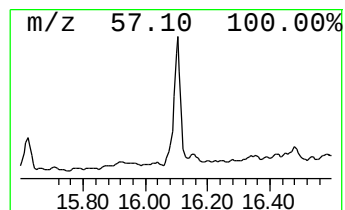
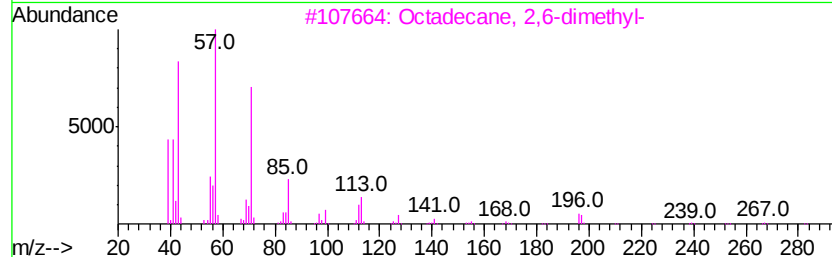
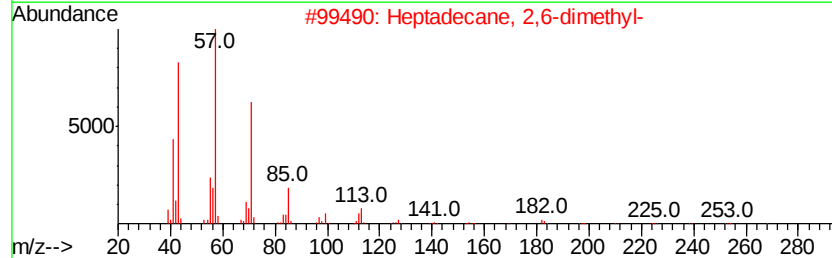
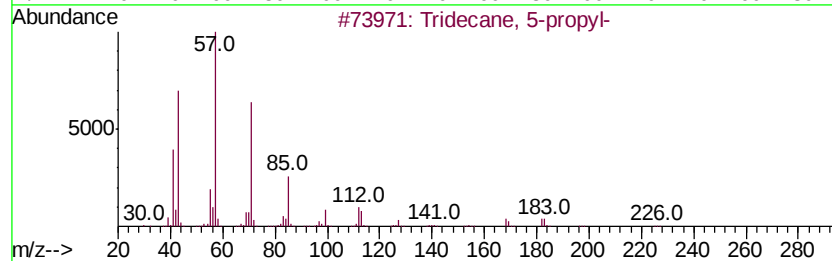
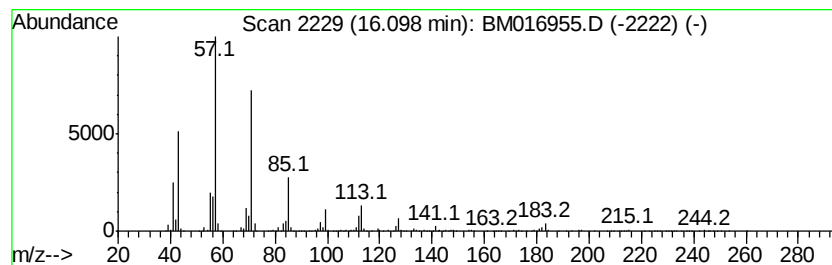
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 16.10 | 16.08 ng/ul | 3351070 | Phenanthrene-d10 | 17.10 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 5-propyl-       | 226 | C16H34  | 055045-11-9 | 94   |
| 2    |    | Heptadecane, 2,6-dimethyl- | 268 | C19H40  | 054105-67-8 | 87   |
| 3    |    | Octadecane, 2,6-dimethyl-  | 282 | C20H42  | 075163-97-2 | 78   |
| 4    |    | Heptane, 2,6-dimethyl-     | 128 | C9H20   | 001072-05-5 | 74   |
| 5    |    | Tridecane, 7-hexyl-        | 268 | C19H40  | 007225-66-3 | 74   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

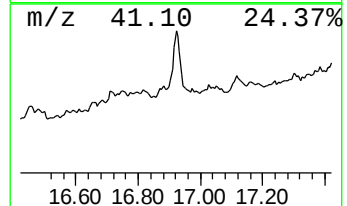
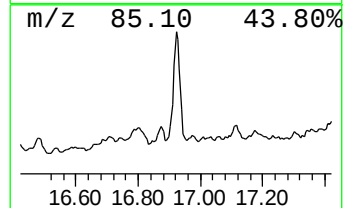
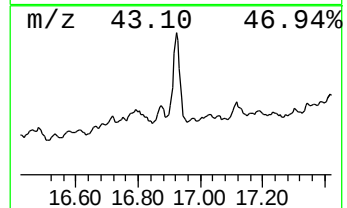
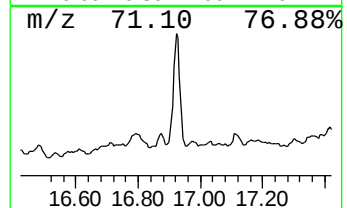
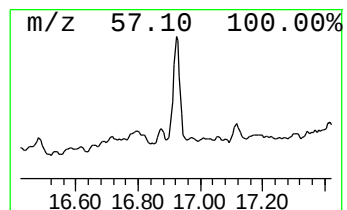
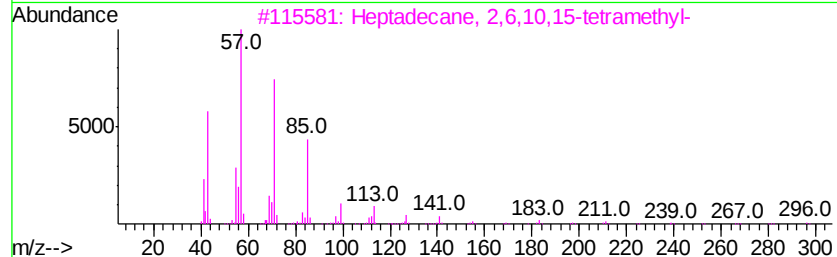
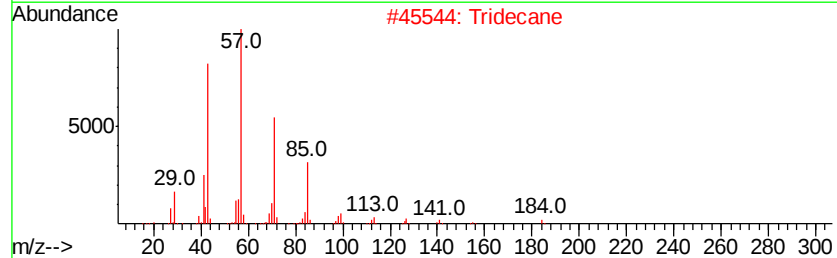
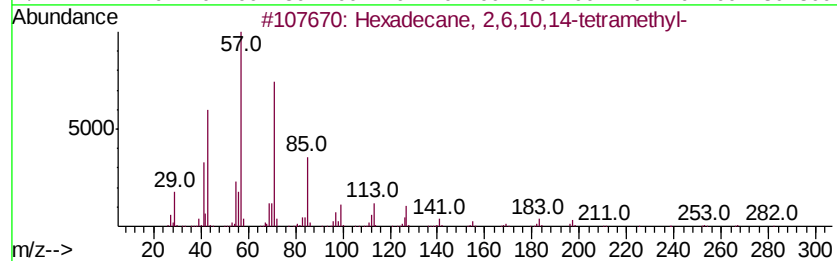
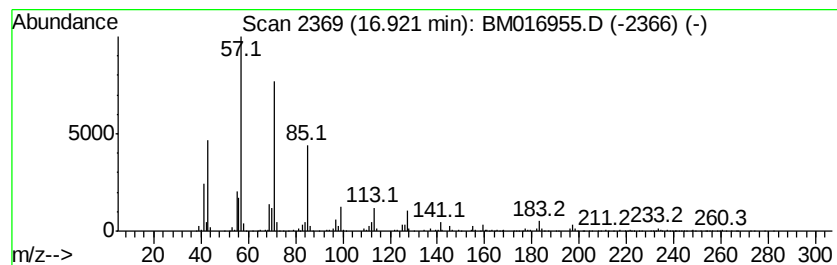
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 24

| R.T.  | EstConc    | Area    | Relative to ISTD | R.T.  |
|-------|------------|---------|------------------|-------|
| 16.92 | 9.86 ng/ul | 2055120 | Phenanthrene-d10 | 17.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8  | 98   |
| 2    |    | Tridecane                           | 184 | C13H28  | 000629-50-5  | 87   |
| 3    |    | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6  | 87   |
| 4    |    | Tridecanol, 2-ethyl-2-methyl-       | 242 | C16H34O | 1000115-66-1 | 86   |
| 5    |    | Octadecane, 1-iodo-                 | 380 | C18H37I | 000629-93-6  | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

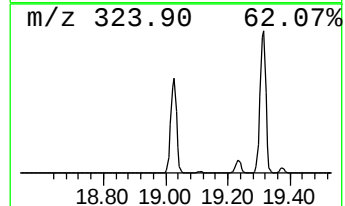
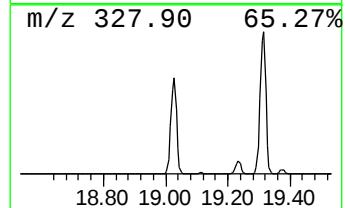
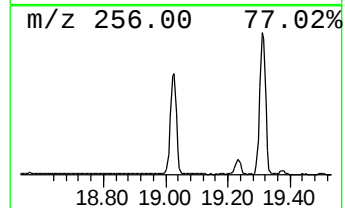
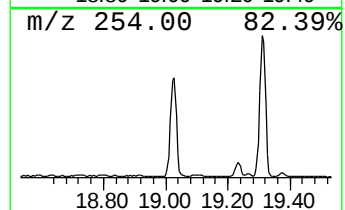
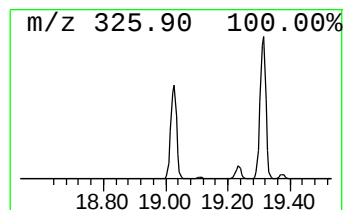
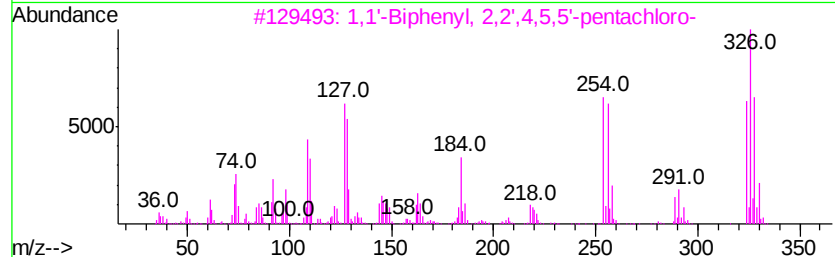
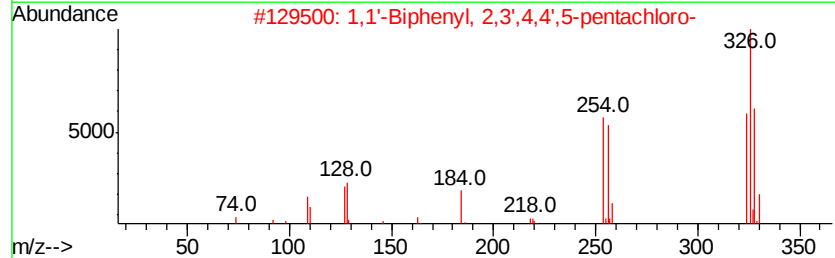
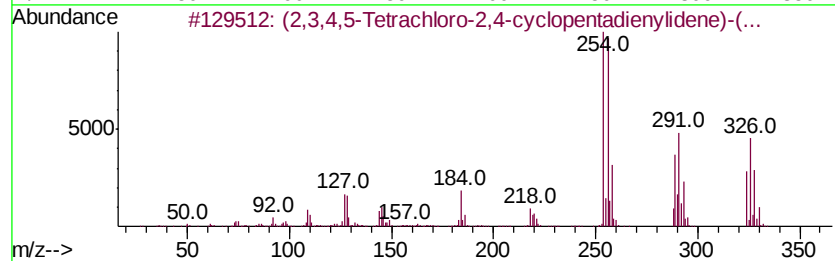
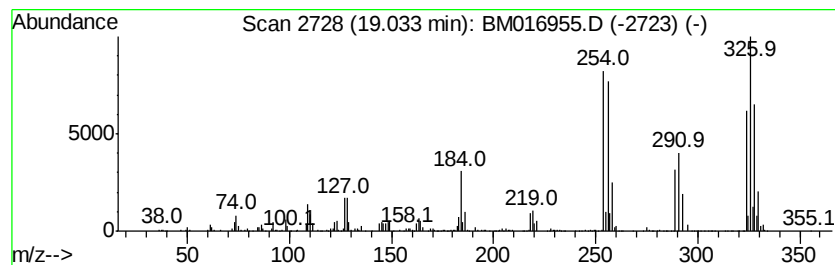
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 (2,3,4,5-Tetrachloro-2,4-cy... Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.03 | 27.41 ng/ul | 5710900 | Phenanthrene-d10 | 17.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | (2,3,4,5-Tetrachloro-2,4-cyclope... | 324 | C12H5Cl5 | 029887-33-0 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,3',4,4',5-penta... | 324 | C12H5Cl5 | 031508-00-6 | 97   |
| 3    |    | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 97   |
| 4    |    | 1,1'-Biphenyl, 2,2',3,4,6-Pentac... | 324 | C12H5Cl5 | 055215-17-3 | 96   |
| 5    |    | 1,1'-Biphenyl, 2,2',3,3',4-penta... | 324 | C12H5Cl5 | 052663-62-4 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

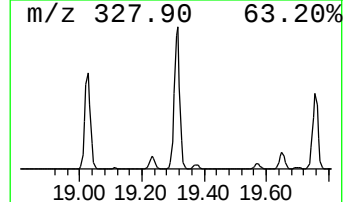
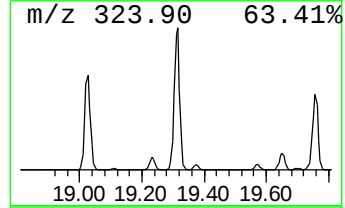
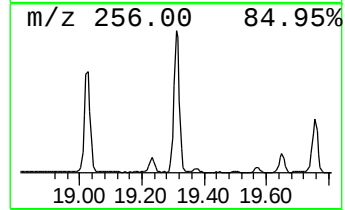
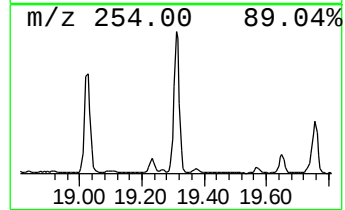
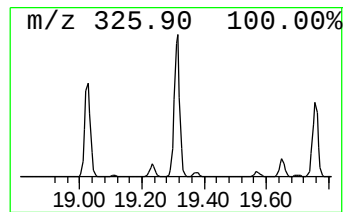
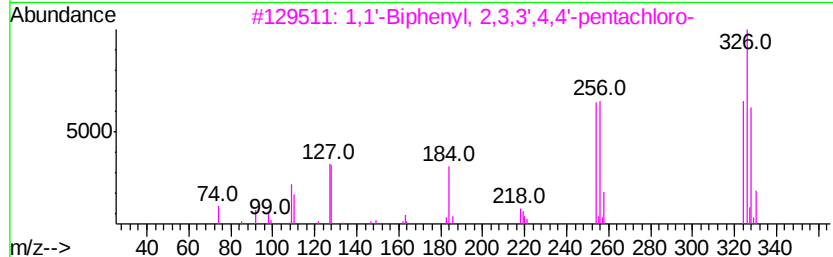
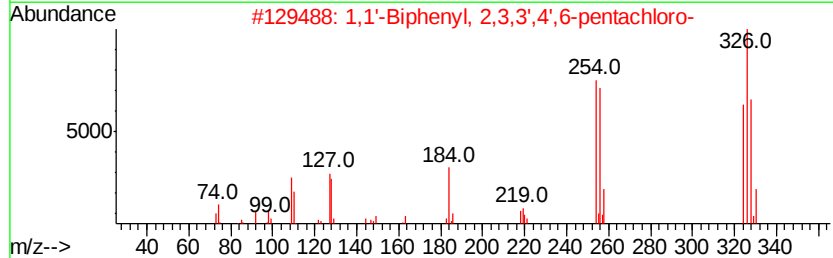
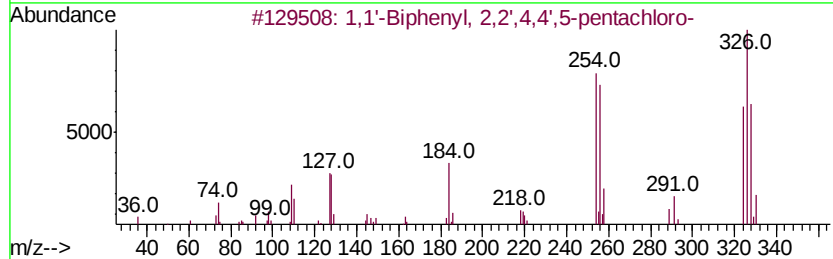
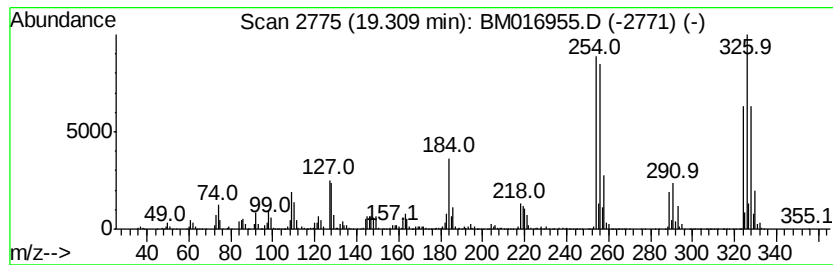
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 1,1'-Biphenyl, 2,2',4,4',5-... Concentration Rank 8

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 19.31   | 32.23 ng/ul                         | 7277520      | Chrysene-d12     | 21.29     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 C12H5Cl5 | 038380-01-7      | 99        |
| 2       | 1,1'-Biphenyl, 2,3,3',4',6-penta... | 324 C12H5Cl5 | 038380-03-9      | 99        |
| 3       | 1,1'-Biphenyl, 2,3,3',4,4'-penta... | 324 C12H5Cl5 | 032598-14-4      | 99        |
| 4       | 1,1'-Biphenyl, 2,3',4,4',5-penta... | 324 C12H5Cl5 | 031508-00-6      | 99        |
| 5       | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 C12H5Cl5 | 038380-01-7      | 99        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

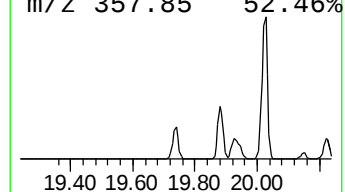
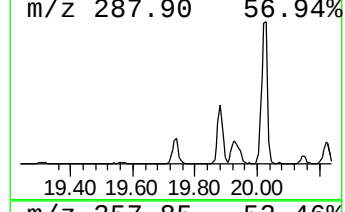
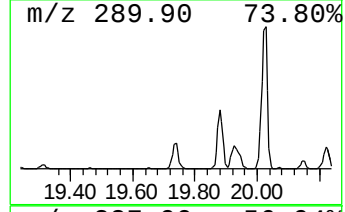
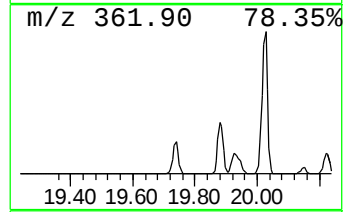
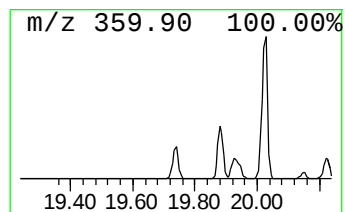
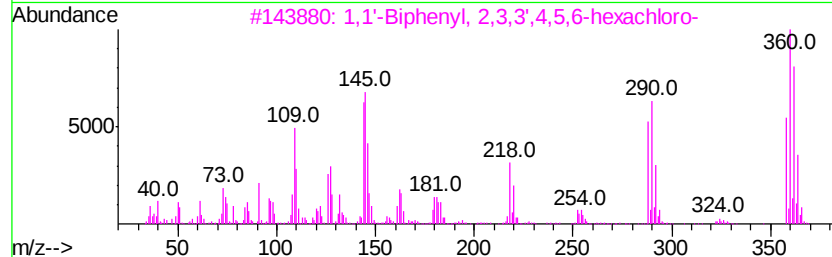
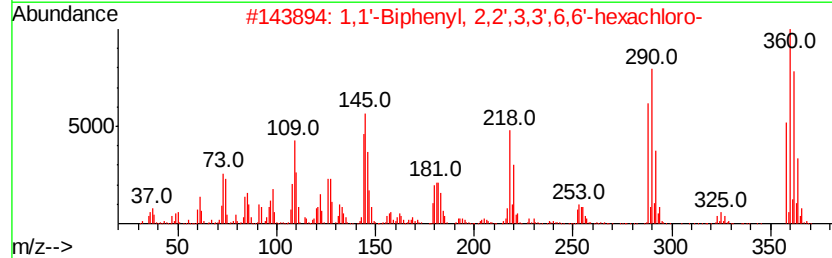
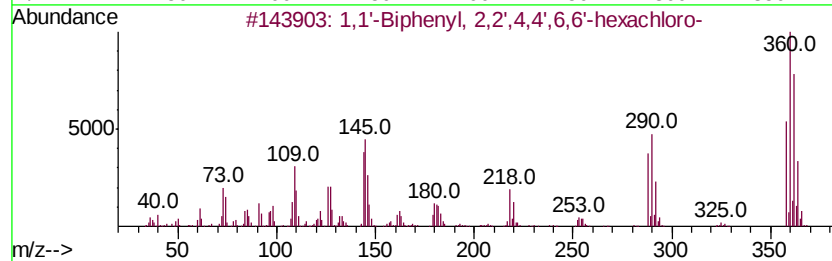
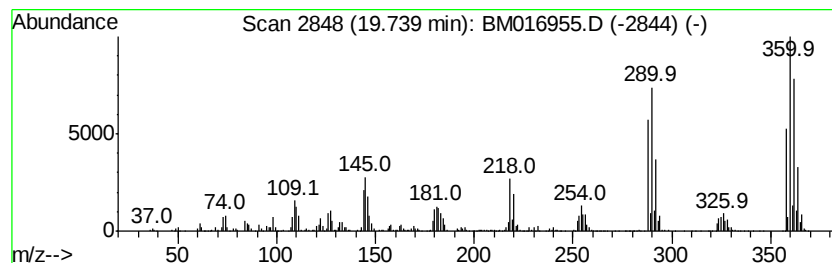
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 1,1'-Biphenyl, 2,2',4,4',6,... Concentration Rank 18

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.74 | 17.09 ng/ul | 3859450 | Chrysene-d12     | 21.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',4,4',6,6'-he... | 358 | C12H4Cl6 | 033979-03-2 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,2',3,3',6,6'-he... | 358 | C12H4Cl6 | 038411-22-2 | 99   |
| 3    |    | 1,1'-Biphenyl, 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 4    |    | 1,1'-Biphenyl, 2,2',3,4',5,5'-he... | 358 | C12H4Cl6 | 051908-16-8 | 96   |
| 5    |    | 1,1'-Biphenyl, 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 96   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

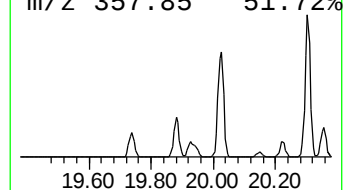
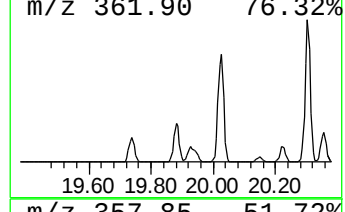
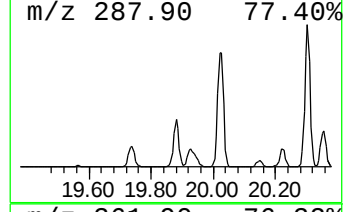
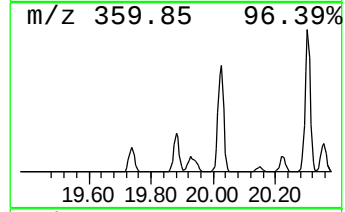
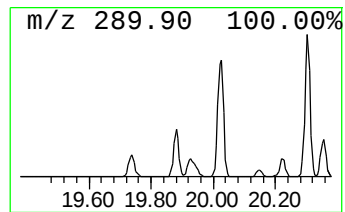
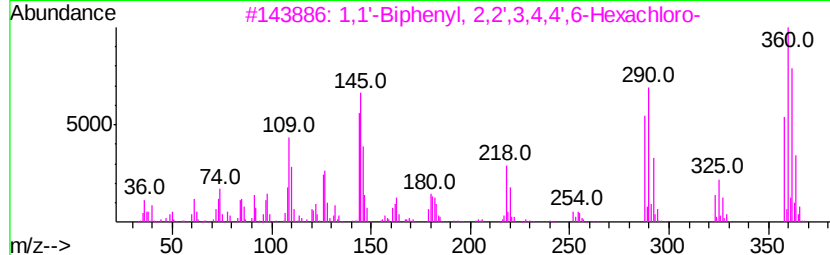
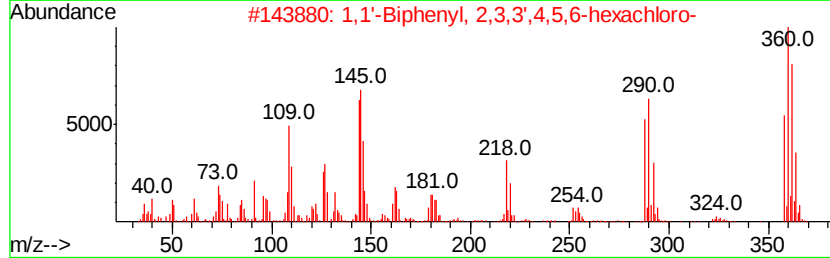
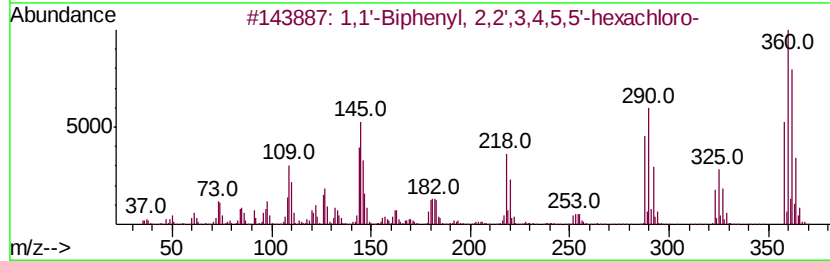
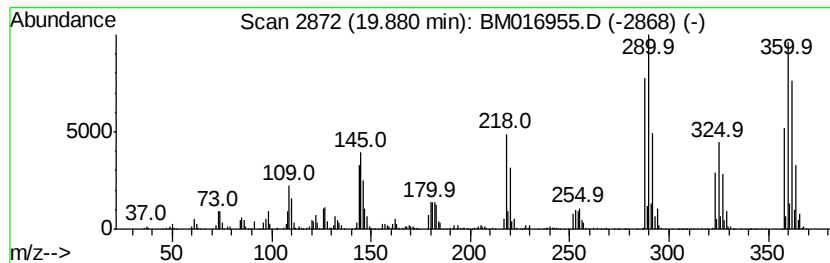
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 1,1'-Biphenyl, 2,2',3,4,5,5... Concentration Rank 9

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.88 | 28.36 ng/ul | 6402490 | Chrysene-d12     | 21.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',3,4,5,5'-hex... | 358 | C12H4Cl6 | 052712-04-6 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 3    |    | 1,1'-Biphenyl, 2,2',3,4,4',6-Hex... | 358 | C12H4Cl6 | 056030-56-9 | 99   |
| 4    |    | 1,1'-Biphenyl, 2,2',3,3',4,6'-he... | 358 | C12H4Cl6 | 038380-05-1 | 99   |
| 5    |    | 1,1'-Biphenyl, 2,2',3,3',4,4'-he... | 358 | C12H4Cl6 | 038380-07-3 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

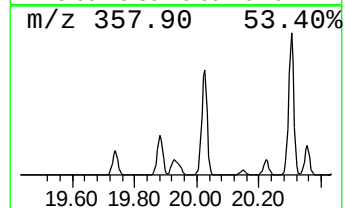
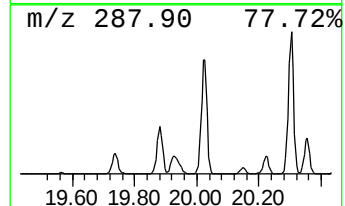
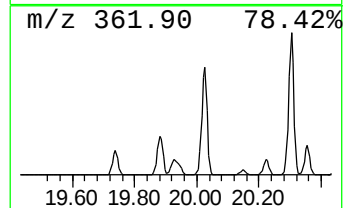
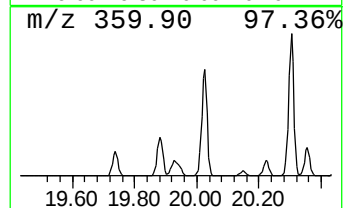
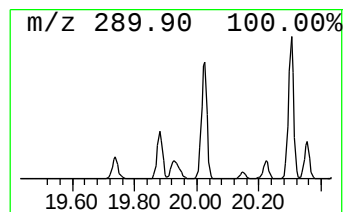
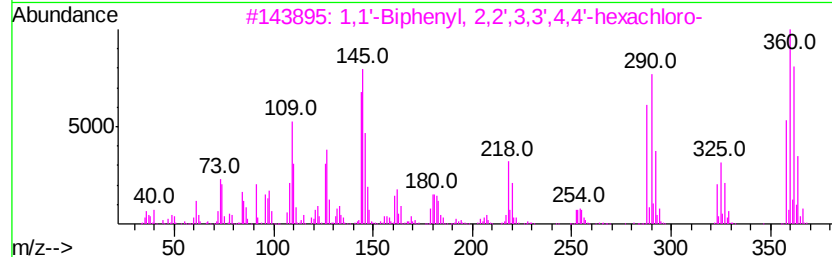
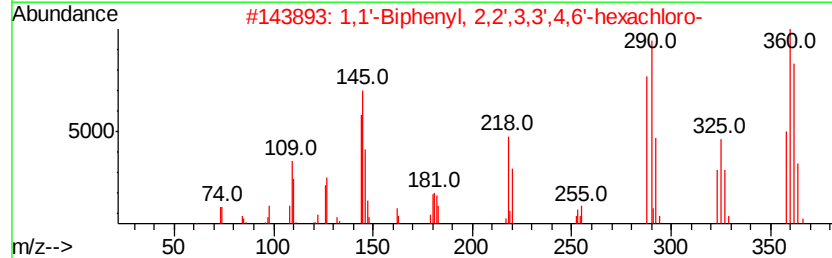
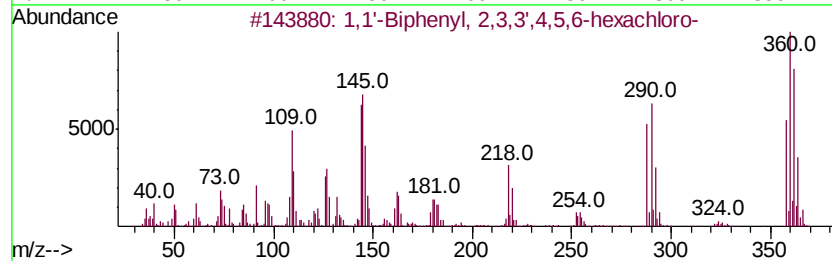
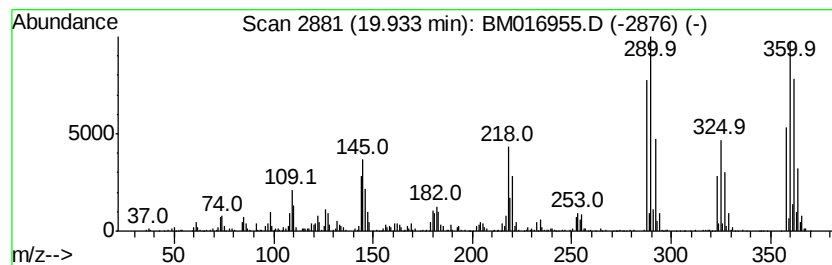
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 1,1'-Biphenyl, 2,3,3',4,5,6... Concentration Rank 16

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 19.93 | 17.60 ng/ul | 3974770 | Chrysene-d12     | 21.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,2',3,3',4,6'-he... | 358 | C12H4Cl6 | 038380-05-1 | 99   |
| 3    |    | 1,1'-Biphenyl, 2,2',3,3',4,4'-he... | 358 | C12H4Cl6 | 038380-07-3 | 99   |
| 4    |    | 1,1'-Biphenyl, 2,2',4,4',6,6'-he... | 358 | C12H4Cl6 | 033979-03-2 | 99   |
| 5    |    | 1,1'-Biphenyl, 2,2',4,4',5',6-He... | 358 | C12H4Cl6 | 060145-22-4 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

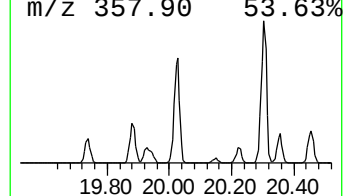
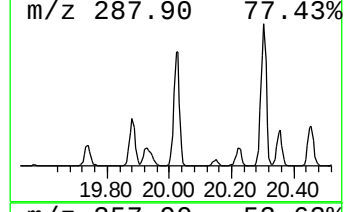
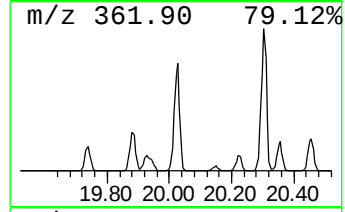
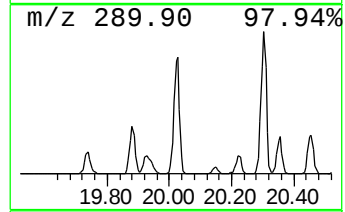
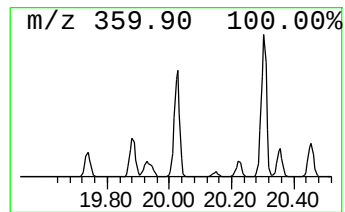
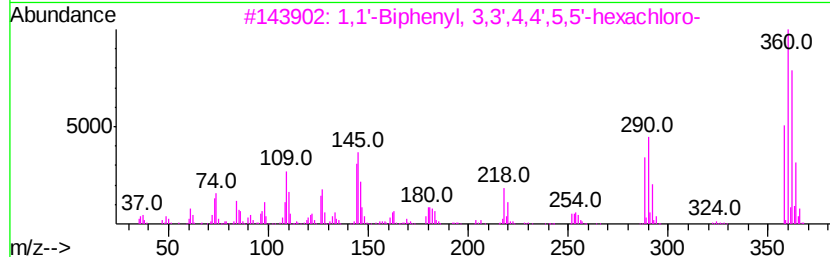
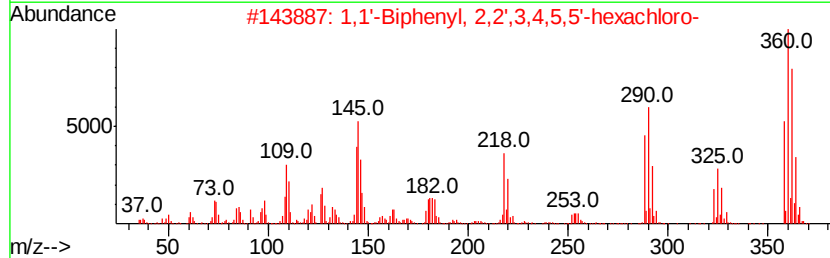
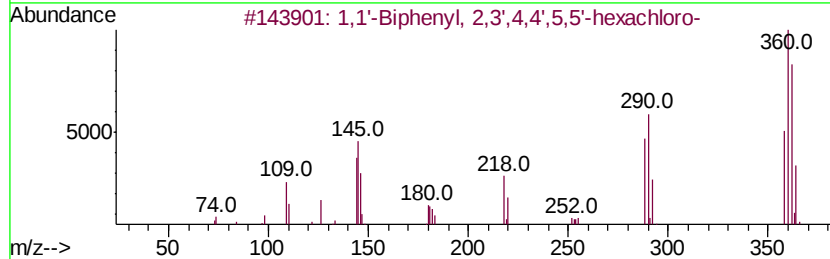
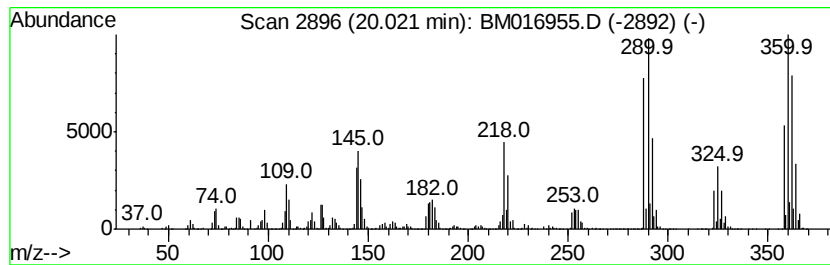
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 1,1'-Biphenyl, 2,3',4,4',5,... Concentration Rank 3

| R.T.  | EstConc     | Area       | Relative to ISTD     |     | R.T.     |             |      |
|-------|-------------|------------|----------------------|-----|----------|-------------|------|
| 20.02 | 74.93 ng/ul | 16919000   | Chrysene-d12         |     | 21.29    |             |      |
| Hit#  | of          | 5          | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
| 1     | 1,1'        | -Biphenyl, | 2,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 052663-72-6 | 99   |
| 2     | 1,1'        | -Biphenyl, | 2,2',3,4,5,5'-hex... | 358 | C12H4Cl6 | 052712-04-6 | 99   |
| 3     | 1,1'        | -Biphenyl, | 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99   |
| 4     | 1,1'        | -Biphenyl, | 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 5     | 1,1'        | -Biphenyl, | 2,2',3,3',6,6'-he... | 358 | C12H4Cl6 | 038411-22-2 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

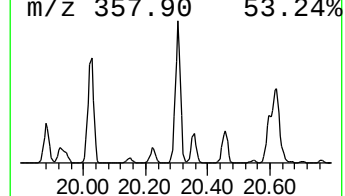
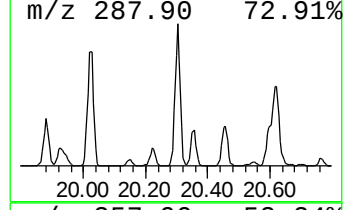
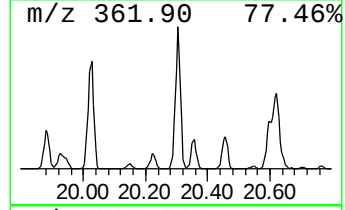
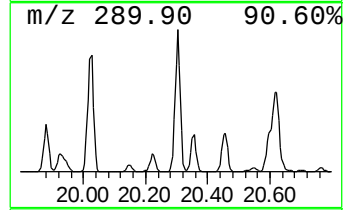
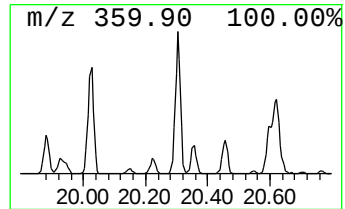
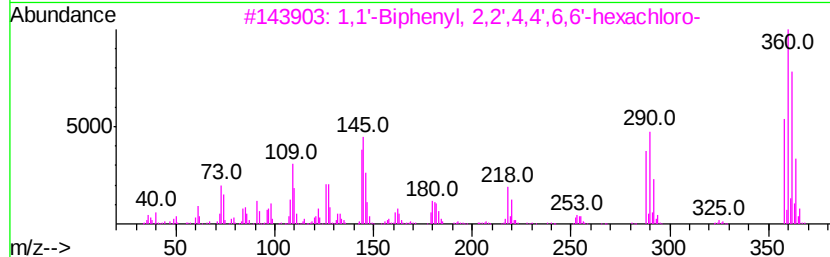
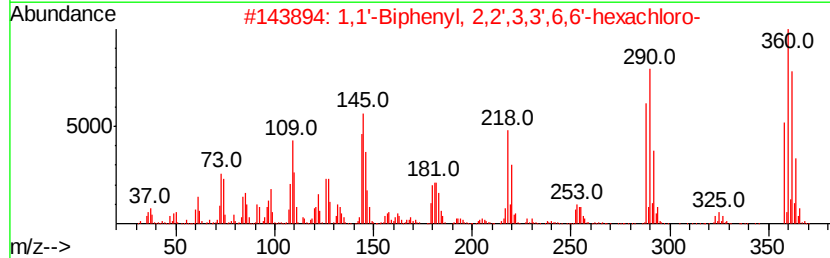
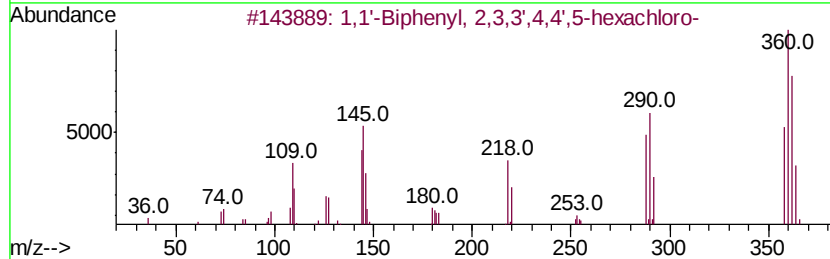
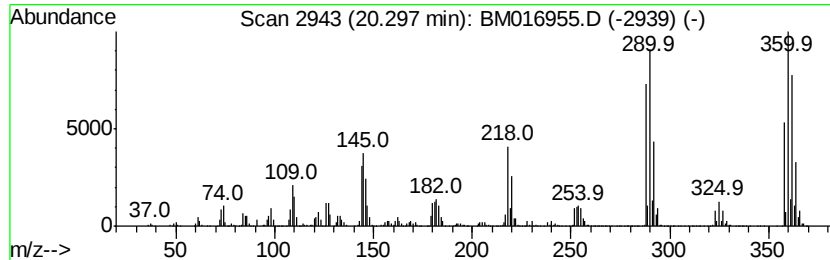
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 1,1'-Biphenyl, 2,3,3',4,4',... Concentration Rank 1

| R.T.  | EstConc     | Area       | Relative to ISTD     |     | R.T.     |             |      |
|-------|-------------|------------|----------------------|-----|----------|-------------|------|
| 20.30 | 84.46 ng/ul | 19069300   | Chrysene-d12         |     | 21.29    |             |      |
| Hit#  | of          | 5          | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
| 1     | 1,1'        | -Biphenyl, | 2,3,3',4,4',5-hex... | 358 | C12H4Cl6 | 038380-08-4 | 99   |
| 2     | 1,1'        | -Biphenyl, | 2,2',3,3',6,6'-he... | 358 | C12H4Cl6 | 038411-22-2 | 99   |
| 3     | 1,1'        | -Biphenyl, | 2,2',4,4',6,6'-he... | 358 | C12H4Cl6 | 033979-03-2 | 99   |
| 4     | 1,1'        | -Biphenyl, | 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 5     | 1,1'        | -Biphenyl, | 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

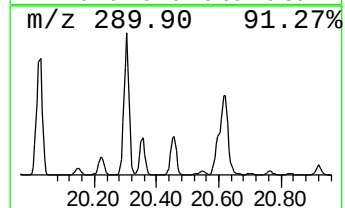
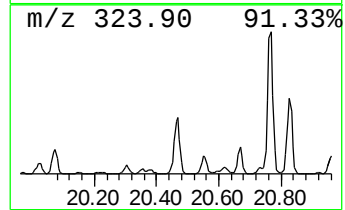
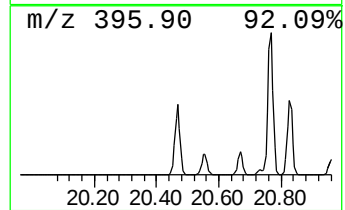
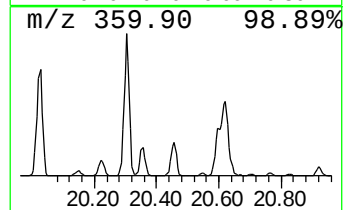
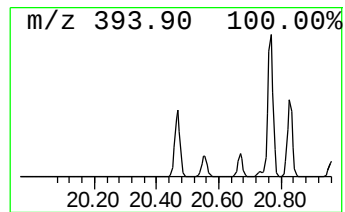
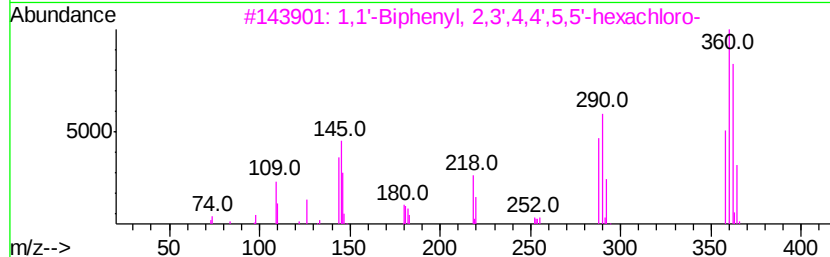
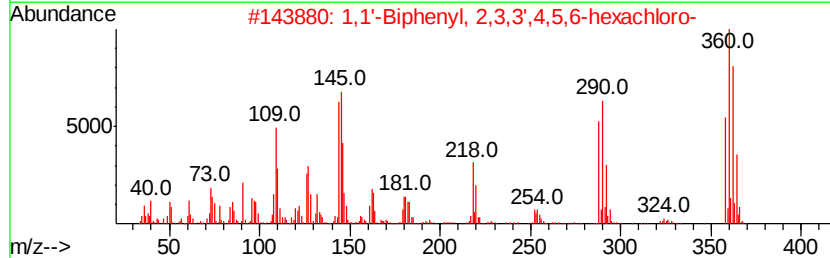
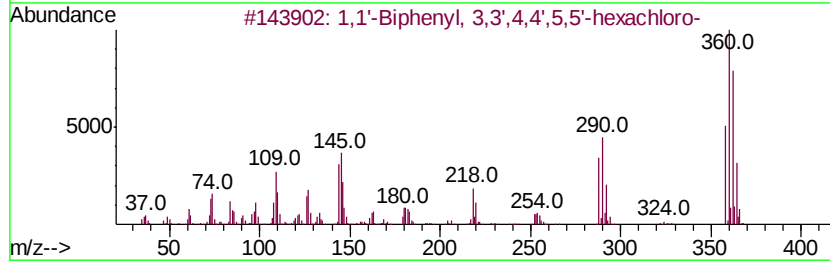
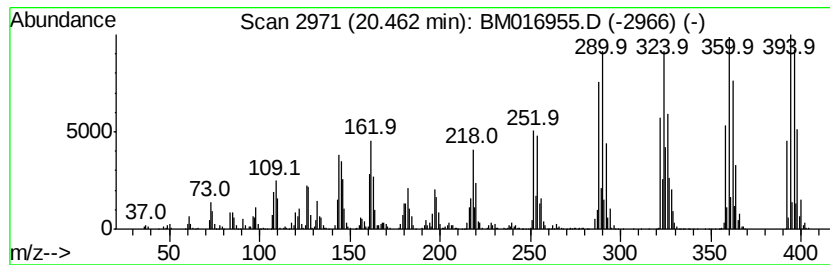
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 1,1'-Biphenyl, 3,3',4,4',5,... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 20.46 | 40.47 ng/ul | 9136770 | Chrysene-d12     | 21.29 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 99   |
| 2    |    | 1,1'-Biphenyl, 2,3,3',4,5,6-hexa... | 358 | C12H4Cl6 | 041411-62-5 | 99   |
| 3    |    | 1,1'-Biphenyl, 2,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 052663-72-6 | 99   |
| 4    |    | 1,1'-Biphenyl, 2,2',4,4',6,6'-he... | 358 | C12H4Cl6 | 033979-03-2 | 99   |
| 5    |    | 1,1'-Biphenyl, 2,2',3,4,5,5'-hex... | 358 | C12H4Cl6 | 052712-04-6 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

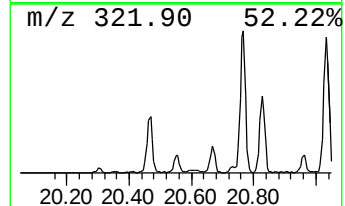
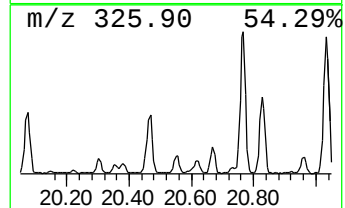
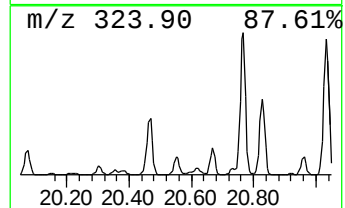
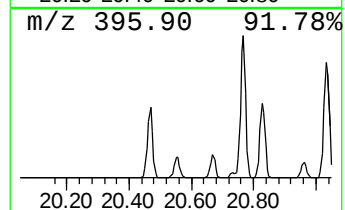
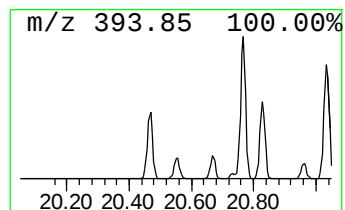
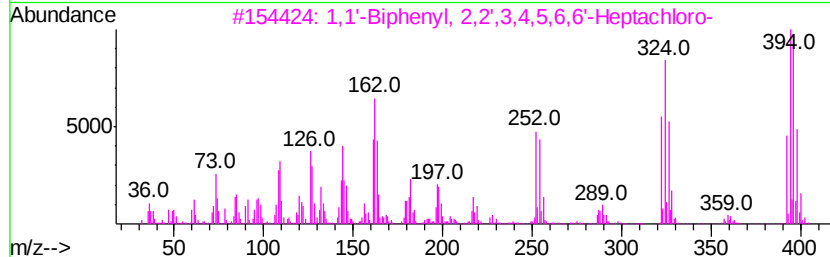
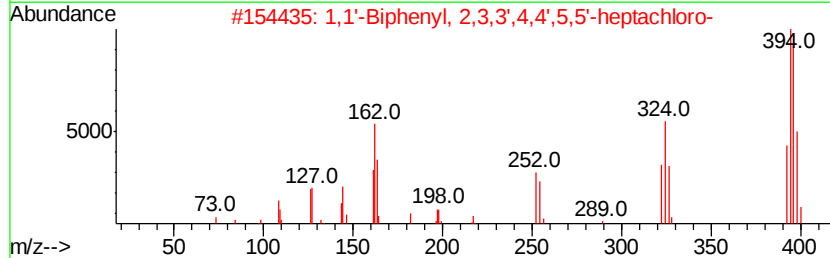
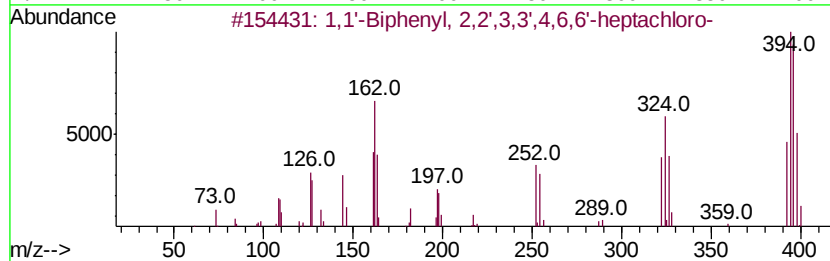
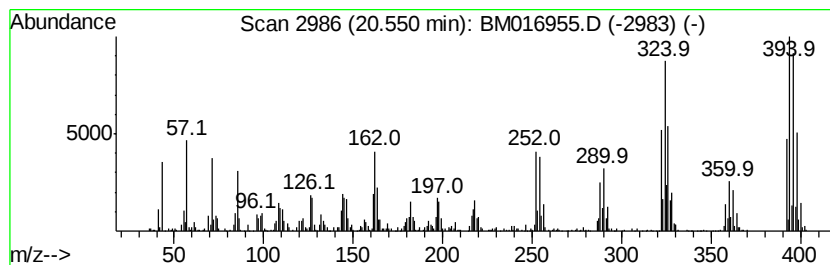
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BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 19 1,1'-Biphenyl, 2,2',3,3',4,... Concentration Rank 25

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.        |      |      |
|---------|-------------------------------------|--------------|------------------|-------------|------|------|
| 20.55   | 8.20 ng/ul                          | 1851570      | Chrysene-d12     | 21.29       |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1       | 1,1'-Biphenyl, 2,2',3,3',4,6,6'-... | 392          | C12H3Cl7         | 052663-65-7 | 99   |      |
| 2       | 1,1'-Biphenyl, 2,3,3',4,4',5,5'-... | 392          | C12H3Cl7         | 039635-31-9 | 99   |      |
| 3       | 1,1'-Biphenyl, 2,2',3,4,5,6,6'-H... | 392          | C12H3Cl7         | 074472-49-4 | 99   |      |
| 4       | 1,1'-Biphenyl, 2,3,3',4,5,5',6-h... | 392          | C12H3Cl7         | 074472-51-8 | 99   |      |
| 5       | 1,1'-Biphenyl, 2,2',3,4,4',5',6-... | 392          | C12H3Cl7         | 052663-69-1 | 96   |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

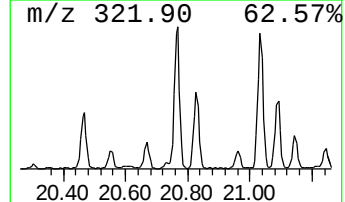
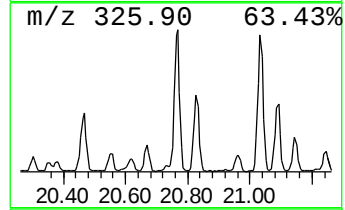
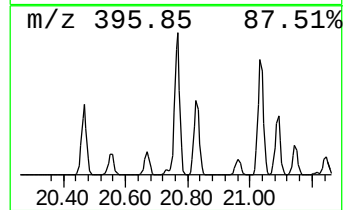
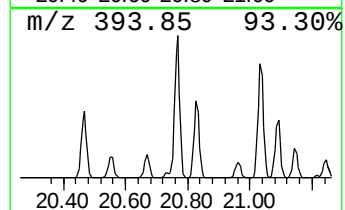
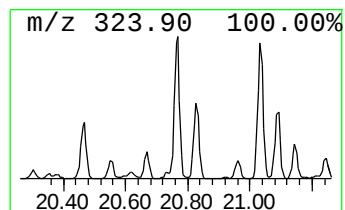
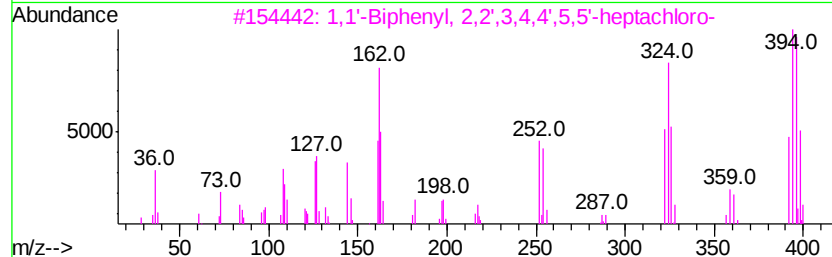
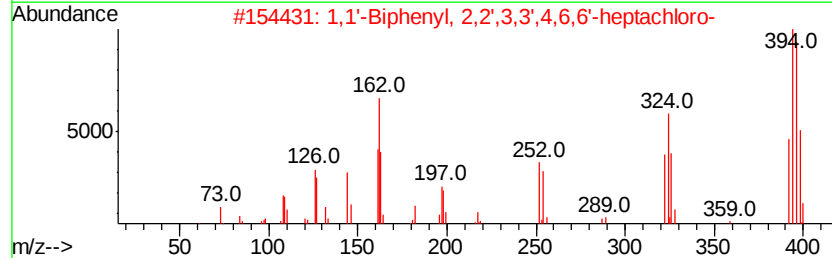
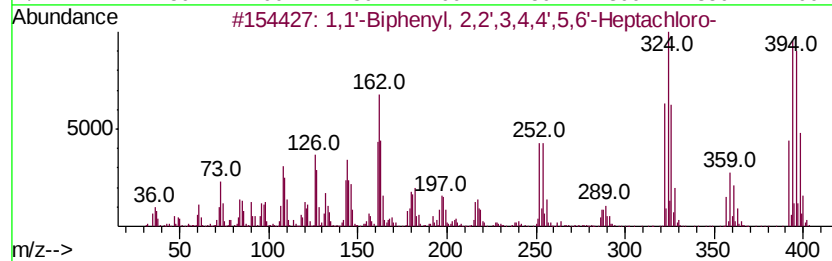
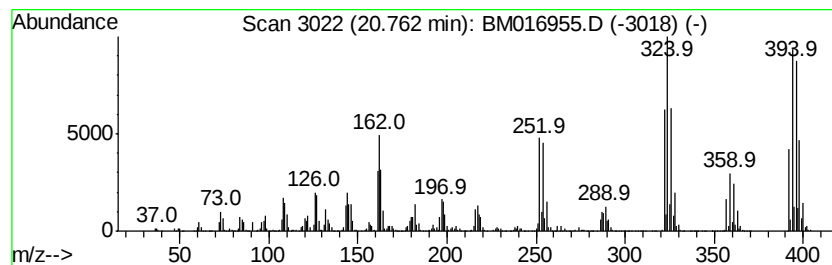
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 22 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 4

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.      |
|---------|-------------------------------------|--------------|------------------|-----------|
| 20.76   | 45.30 ng/ul                         | 10227800     | Chrysene-d12     | 21.29     |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | 1,1'-Biphenyl, 2,2',3,4,4',5,6'-... | 392 C12H3Cl7 | 060145-23-5      | 99        |
| 2       | 1,1'-Biphenyl, 2,2',3,3',4,6,6'-... | 392 C12H3Cl7 | 052663-65-7      | 99        |
| 3       | 1,1'-Biphenyl, 2,2',3,4,4',5,5'-... | 392 C12H3Cl7 | 035065-29-3      | 99        |
| 4       | 2,2',3,3',4,5',6-Heptachlorobiph... | 392 C12H3Cl7 | 040186-70-7      | 99        |
| 5       | 1,1'-Biphenyl, 2,2',3,3',5,5',6-... | 392 C12H3Cl7 | 052663-67-9      | 99        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

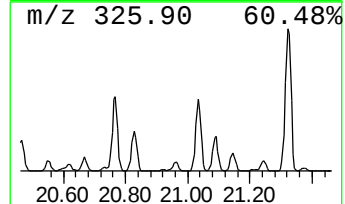
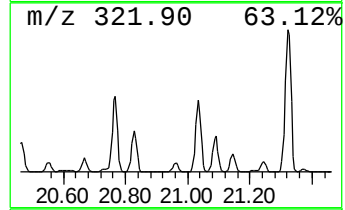
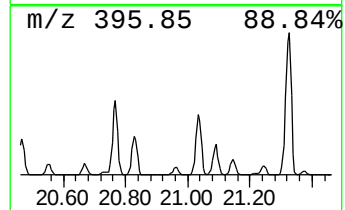
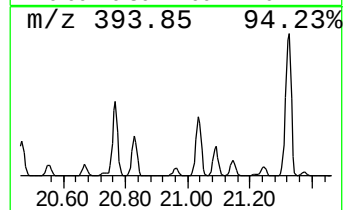
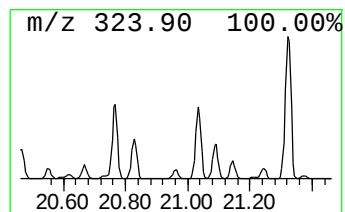
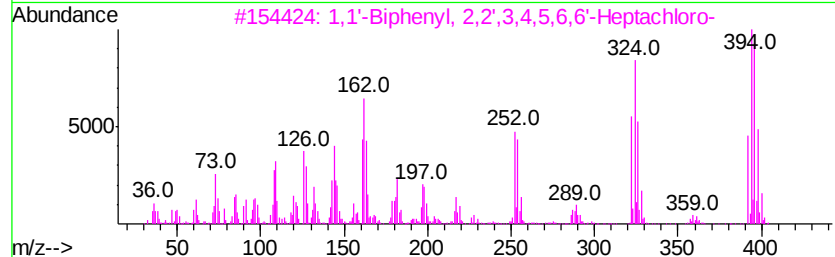
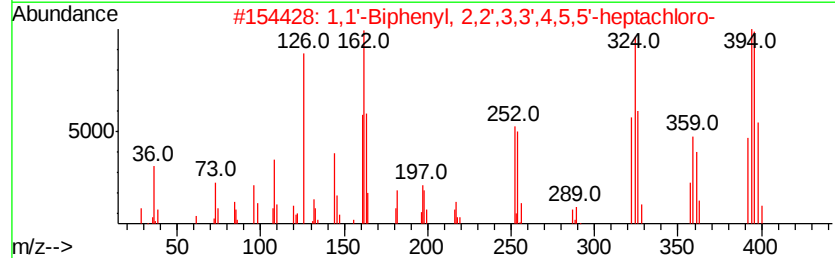
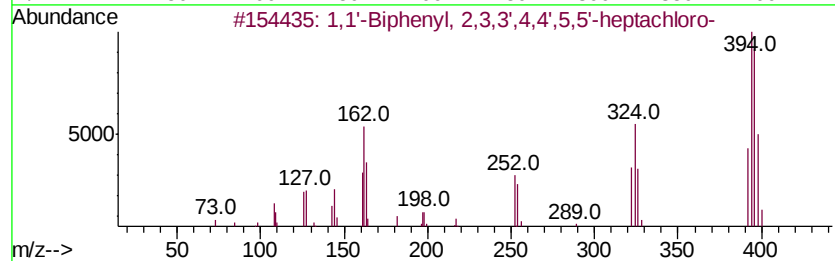
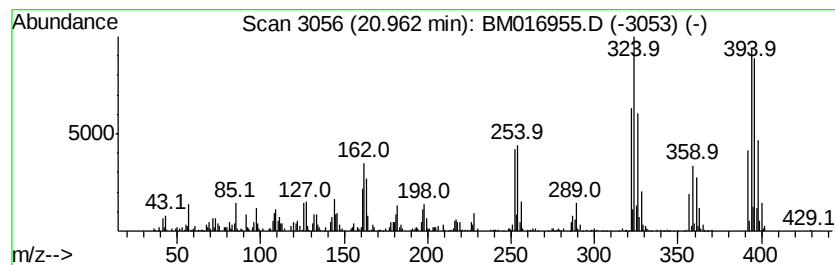
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 25 1,1'-Biphenyl, 2,3,3',4,4',... Concentration Rank 28

| R.T.    | EstConc        | Area                 | Relative to ISTD | R.T.     |             |      |
|---------|----------------|----------------------|------------------|----------|-------------|------|
| 20.96   | 4.75 ng/ul     | 1073250              | Chrysene-d12     | 21.29    |             |      |
| Hit# of | 5              | Tentative ID         | MW               | MolForm  | CAS#        | Qual |
| 1       | 1,1'-Biphenyl, | 2,3,3',4,4',5,5'-... | 392              | C12H3Cl7 | 039635-31-9 | 99   |
| 2       | 1,1'-Biphenyl, | 2,2',3,3',4,5,5'-... | 392              | C12H3Cl7 | 052663-74-8 | 99   |
| 3       | 1,1'-Biphenyl, | 2,2',3,4,5,6,6'-H... | 392              | C12H3Cl7 | 074472-49-4 | 99   |
| 4       | 1,1'-Biphenyl, | 2,2',3,4,4',5,6-H... | 392              | C12H3Cl7 | 074472-47-2 | 99   |
| 5       | 1,1'-Biphenyl, | 2,2',3,4,4',5',6-... | 392              | C12H3Cl7 | 052663-69-1 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM100118\  
Data File : BM016955.D  
Acq On : 02 Oct 2018 11:28  
Operator : MJ/SJ  
Sample : J5126-01  
Misc :  
ALS Vial : 40 Sample Multiplier: 1

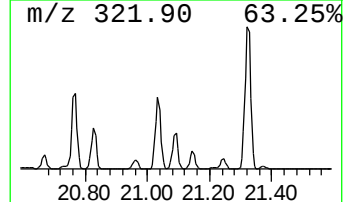
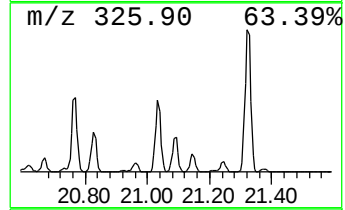
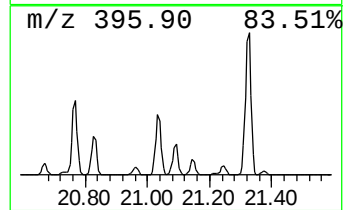
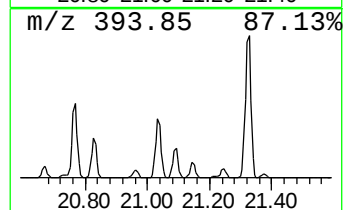
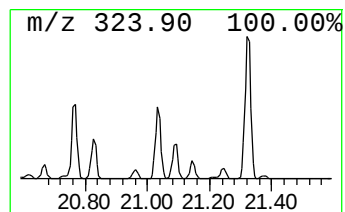
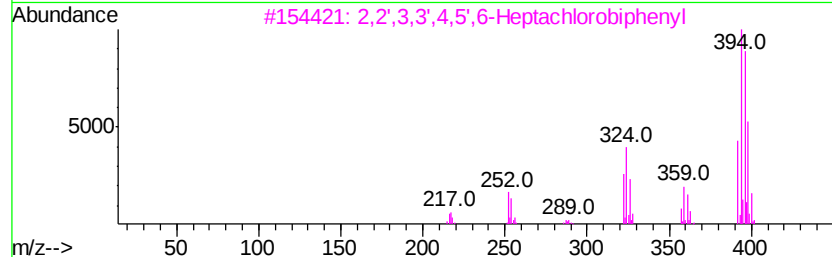
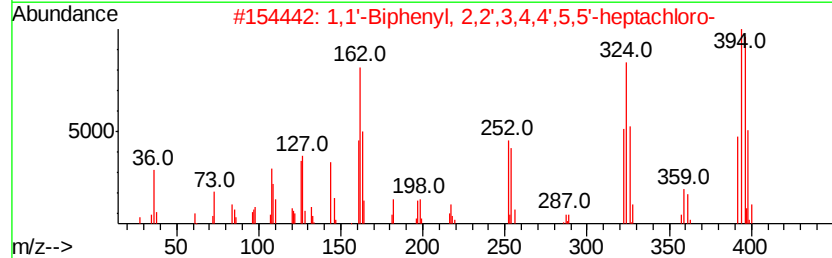
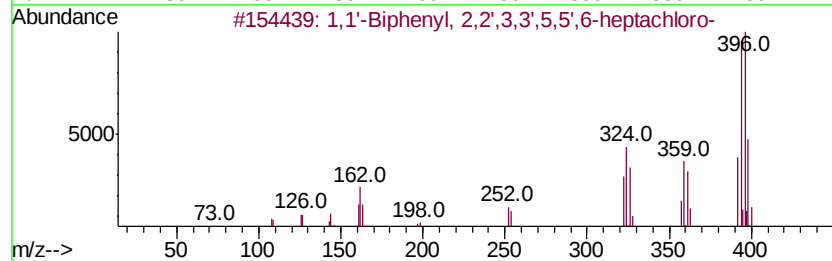
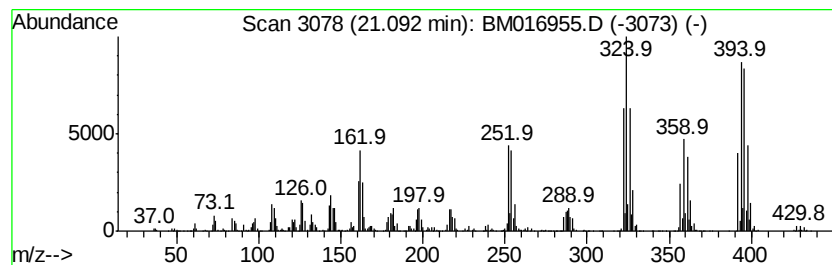
Instrument :  
BNA\_M  
ClientSampleId :  
C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 27 1,1'-Biphenyl, 2,2',3,3',5,... Concentration Rank 11

| R.T.    | EstConc                             | Area                            | Relative to ISTD | R.T.     |             |      |
|---------|-------------------------------------|---------------------------------|------------------|----------|-------------|------|
| 21.09   | 24.36 ng/ul                         | 5500810                         | Chrysene-d12     | 21.29    |             |      |
| Hit# of | 5                                   | Tentative ID                    | MW               | MolForm  | CAS#        | Qual |
| 1       | 1,1'                                | -Biphenyl, 2,2',3,3',5,5',6-... | 392              | C12H3Cl7 | 052663-67-9 | 99   |
| 2       | 1,1'                                | -Biphenyl, 2,2',3,4,4',5,5'-... | 392              | C12H3Cl7 | 035065-29-3 | 99   |
| 3       | 2,2',3,3',4,5',6-Heptachlorobiph... |                                 | 392              | C12H3Cl7 | 040186-70-7 | 99   |
| 4       | 1,1'                                | -Biphenyl, 2,2',3,3',4,5,5'-... | 392              | C12H3Cl7 | 052663-74-8 | 99   |
| 5       | 1,1'                                | -Biphenyl, 2,2',3,4,4',5',6-... | 392              | C12H3Cl7 | 052663-69-1 | 99   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM100118\  
 Data File : BM016955.D  
 Acq On : 02 Oct 2018 11:28  
 Operator : MJ/SJ  
 Sample : J5126-01  
 Misc :  
 ALS Vial : 40 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0B89

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM091018MA.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 |
| (DEL) Alkane: Str... | 3.19  | 11.3    | ng/ul | 1242760  | 1                     | 7.71  | 2192440 | 20.0 |
| Decahydro-4,4,8,9... | 14.18 | 2.2     | ng/ul | 430271   | 3                     | 14.35 | 3932130 | 20.0 |
| unknown-01           | 15.14 | 3.0     | ng/ul | 600558   | 3                     | 14.35 | 3932130 | 20.0 |
| (DEL) Alkane: Str... | 15.62 | 3.1     | ng/ul | 605937   | 3                     | 14.35 | 3932130 | 20.0 |
| (DEL) Alkane: Str... | 16.10 | 16.1    | ng/ul | 3351070  | 4                     | 17.10 | 4167360 | 20.0 |
| (DEL) Alkane: Str... | 16.92 | 9.9     | ng/ul | 2055120  | 4                     | 17.10 | 4167360 | 20.0 |
| (2,3,4,5-Tetrachl... | 19.03 | 27.4    | ng/ul | 5710900  | 4                     | 17.10 | 4167360 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.31 | 32.2    | ng/ul | 7277520  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.74 | 17.1    | ng/ul | 3859450  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.88 | 28.4    | ng/ul | 6402490  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.93 | 17.6    | ng/ul | 3974770  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.02 | 74.9    | ng/ul | 16919000 | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.30 | 84.5    | ng/ul | 19069300 | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 3,... | 20.46 | 40.5    | ng/ul | 9136770  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.55 | 8.2     | ng/ul | 1851570  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.76 | 45.3    | ng/ul | 10227800 | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.96 | 4.8     | ng/ul | 1073250  | 5                     | 21.29 | 4515770 | 20.0 |
| 1,1'-Biphenyl, 2,... | 21.09 | 24.4    | ng/ul | 5500810  | 5                     | 21.29 | 4515770 | 20.0 |