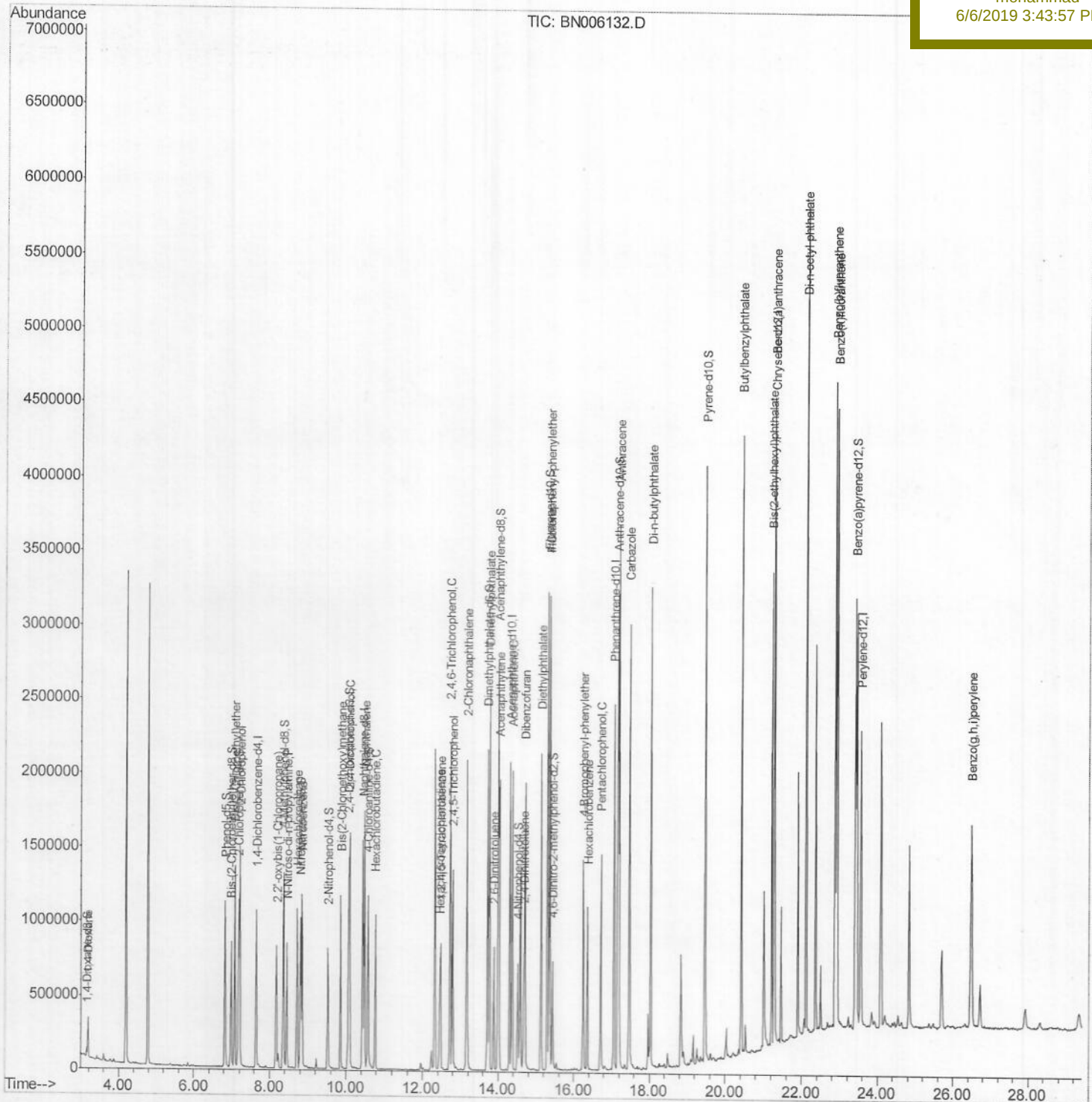


Quant Time: Jun 06 12:24:50 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 06 07:51:51 2019  
Response via : Initial Calibration

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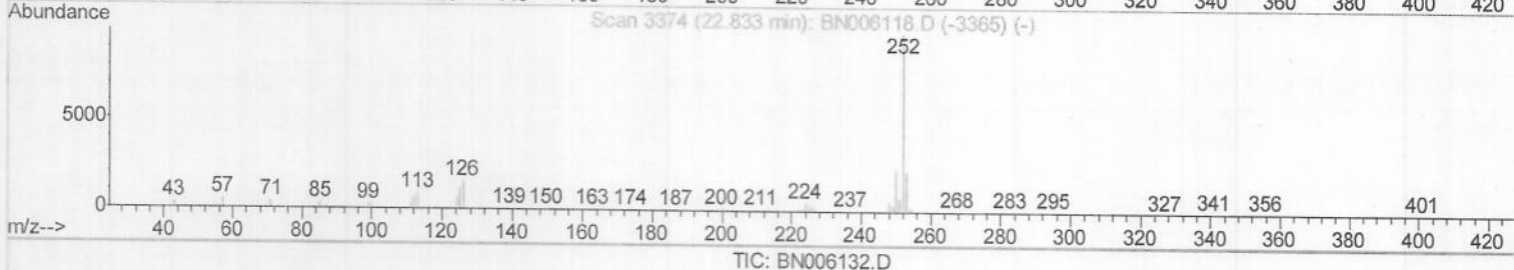
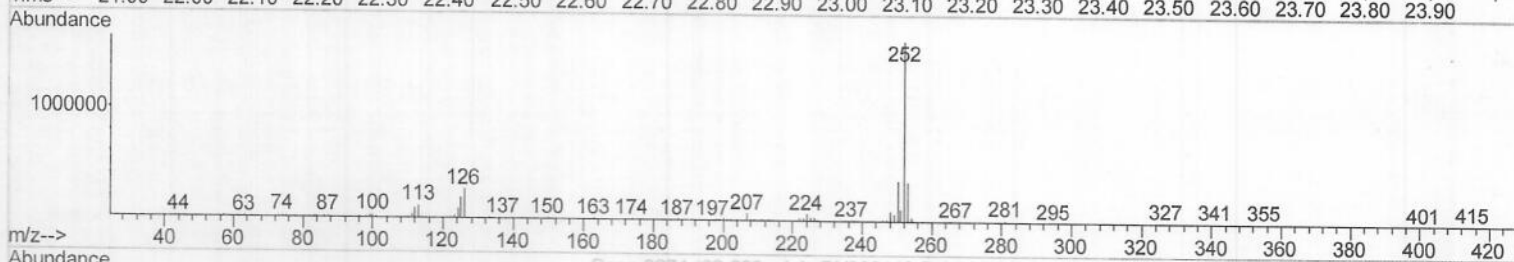
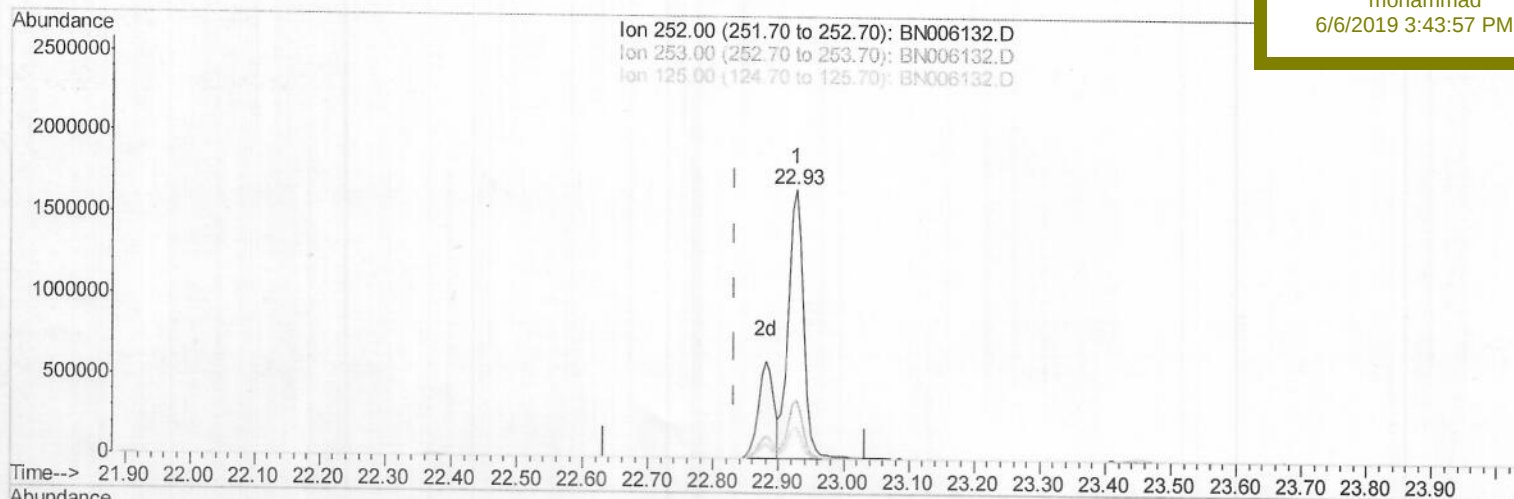
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060519\  
Data File : BN006132.D  
Acq On : 06 Jun 2019 11:48  
Operator : JU/SJ  
Sample : K3016-26  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Quant Time: Jun 06 12:21:43 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN060419MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 06 07:51:51 2019  
Response via : Initial Calibration

Instrument :  
BNA\_N  
ClientSampleId :  
A51H2

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(87) Benzo(b)fluoranthene

22.927min (+0.094) 36.21ng/ul

response 2776280

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 21.80 | 21.81 |
| 125.00 | 10.20 | 11.63 |
| 0.00   | 0.00  | 0.00  |

## Quantitation Report (Qedit)

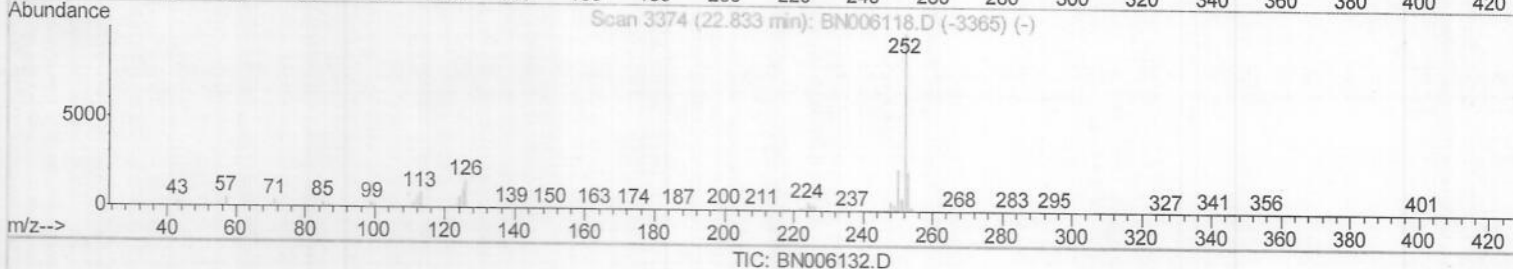
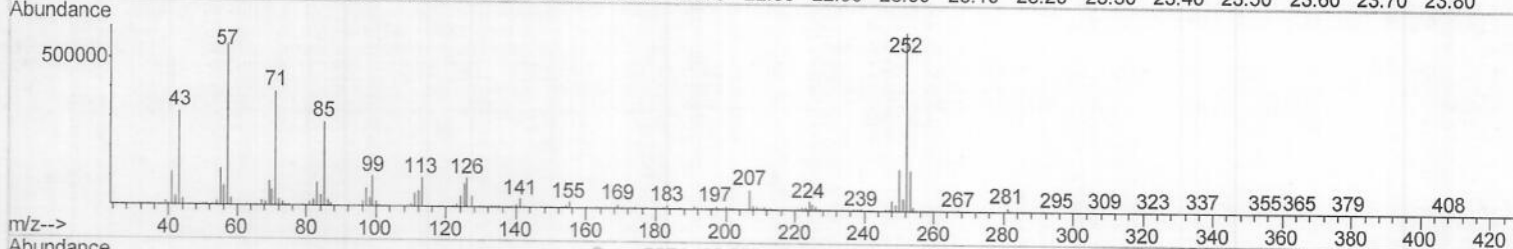
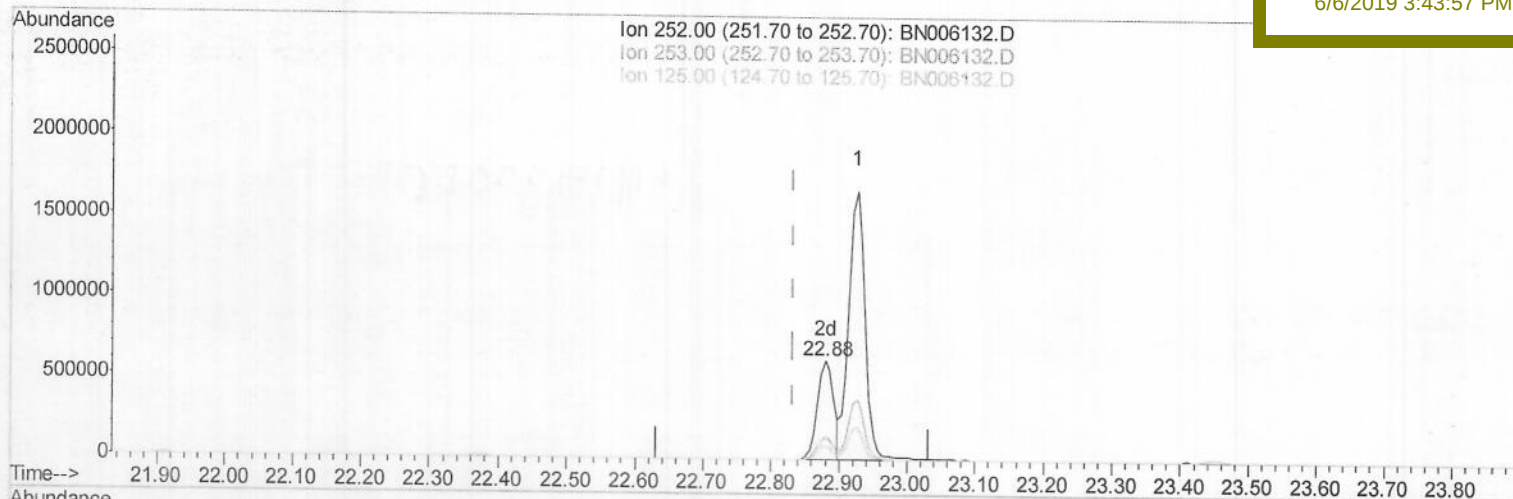
Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060519\  
Data File : BN006132.D  
Acq On : 06 Jun 2019 11:48  
Operator : JU/SJ  
Sample : K3016-26  
Misc :  
ALS Vial : 41 Sample Multiplier: 1

Quant Time: Jun 06 12:21:43 2019  
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Quant Title : SVOA CALIBRATION  
QLast Update : Thu Jun 06 07:51:51 2019  
Response via : Initial Calibration

Instrument :  
BNA\_N  
ClientSampleId :  
A51H2

Manual Integrations  
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TIC: BN006132.D

(87) Benzo(b)fluoranthene

22.880min (+0.047) 13.11ng/ul m

3006107119

response 1005175

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 252.00 | 100   | 100    |
| 253.00 | 21.80 | 22.46  |
| 125.00 | 10.20 | 13.26# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN060519\  
 Data File : BN006132.D  
 Acq On : 06 Jun 2019 11:48  
 Operator : JU/SJ  
 Sample : K3016-26  
 Misc :  
 ALS Vial : 41 Sample Multiplier: 1

Quant Time: Jun 06 12:24:50 2019  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 06 07:51:51 2019  
 Response via : Initial Calibration

Instrument :  
 BNA\_N  
 Client Sampled :  
 A51H2

Manual Integrations  
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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.65  | 152  | 282097   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.43 | 136  | 1253010  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.30 | 164  | 673477   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.06 | 188  | 1397086  | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.29 | 240  | 1266024  | 20.00 | ng/ul | 0.02     |
| 85) Perylene-d12          | 23.55 | 264  | 1438379  | 20.00 | ng/ul | 0.05     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.15  | 96  | 31029   | 4.77  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.82  | 99  | 673849  | 29.90 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl) ether-d | 6.98  | 67  | 386797  | 30.01 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.18  | 132 | 510908  | 28.21 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.36  | 113 | 509954  | 29.20 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.80  | 128 | 247161  | 27.64 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.52  | 143 | 266773  | 26.88 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.06 | 165 | 509939  | 28.43 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.57 | 131 | 645928  | 30.15 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 13.71 | 166 | 1435045 | 30.04 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.99 | 160 | 1820069 | 29.43 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.52 | 143 | 200921  | 23.81 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.30 | 176 | 1199006 | 29.70 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.43 | 200 | 138107  | 17.67 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.16 | 188 | 1822899 | 30.58 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.47 | 212 | 1999192 | 32.68 | ng/ul | 0.01 |
| 89) Benzo(a)pyrene-d12         | 23.41 | 264 | 2005177 | 31.31 | ng/ul | 0.05 |

## Target Compounds

|                                 |       |     |         |        |       | Qvalue |
|---------------------------------|-------|-----|---------|--------|-------|--------|
| 2) 1,4-Dioxane                  | 3.18  | 88  | 119070  | 17.692 | ng/uL | 95     |
| 8) Bis(2-Chloroethyl) ether     | 7.08  | 93  | 735980  | 41.167 | ng/ul | 98     |
| 10) 2-Chlorophenol              | 7.21  | 128 | 617433  | 33.391 | ng/ul | 97     |
| 12) 2,2'-oxybis(1-Chloropropan  | 8.18  | 45  | 814402  | 34.088 | ng/ul | 98     |
| 15) N-Nitroso-di-n-propylamine  | 8.45  | 70  | 296963  | 21.210 | ng/ul | 99     |
| 17) Hexachloroethane            | 8.72  | 117 | 210640  | 27.088 | ng/ul | 98     |
| 20) Nitrobenzene                | 8.84  | 77  | 649159  | 30.688 | ng/ul | 100    |
| 25) Bis(2-Chloroethoxy) methane | 9.85  | 93  | 682759  | 27.209 | ng/ul | 98     |
| 27) 2,4-Dichlorophenol          | 10.08 | 162 | 518852  | 29.686 | ng/ul | 99     |
| 28) Naphthalene                 | 10.48 | 128 | 1382305 | 23.392 | ng/ul | 99     |
| 31) Hexachlorobutadiene         | 10.76 | 225 | 209204  | 18.723 | ng/ul | 98     |
| 36) 1,2,4,5-Tetrachlorobenzene  | 12.47 | 216 | 244919  | 12.149 | ng/ul | 96     |
| 37) Hexachlorocyclopentadiene   | 12.45 | 237 | 172756  | 15.005 | ng/ul | 97     |
| 38) 2,4,6-Trichlorophenol       | 12.72 | 196 | 502732  | 38.972 | ng/ul | 99     |
| 39) 2,4,5-Trichlorophenol       | 12.79 | 196 | 352999  | 25.658 | ng/ul | 97     |
| 41) 2-Chloronaphthalene         | 13.17 | 162 | 993117  | 25.391 | ng/ul | 99     |
| 44) Dimethylphthalate           | 13.76 | 163 | 1642832 | 34.667 | ng/ul | 100    |
| 45) 2,6-Dinitrotoluene          | 13.89 | 165 | 156296  | 16.243 | ng/ul | 97     |
| 47) Acenaphthylene              | 14.02 | 152 | 1303345 | 21.463 | ng/ul | 99     |
| 49) Acenaphthene                | 14.36 | 153 | 789844  | 19.330 | ng/ul | 100    |
| 53) Dibenzofuran                | 14.70 | 168 | 1218171 | 21.126 | ng/ul | 98     |
| 54) 2,4-Dinitrotoluene          | 14.68 | 165 | 181601  | 13.444 | ng/ul | 98     |
| 56) Diethylphthalate            | 15.13 | 149 | 1113786 | 23.469 | ng/ul | 100    |

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 ALS Vial : 41 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 Client Sampled :  
 A51H2

Manual Integrations  
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 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Jun 06 07:51:51 2019  
 Response via : Initial Calibration

| Internal Standards             | R.T.  | QIon | Response  | Conc   | Units | Dev (Min) |
|--------------------------------|-------|------|-----------|--------|-------|-----------|
| 58) Fluorene                   | 15.36 | 166  | 1013322   | 22.447 | ng/ul | 99        |
| 59) 4-Chlorophenyl-phenylether | 15.35 | 204  | 293433    | 13.039 | ng/ul | 98        |
| 65) 4-Bromophenyl-phenylether  | 16.25 | 248  | 263452    | 19.200 | ng/ul | 99        |
| 66) Hexachlorobenzene          | 16.36 | 284  | 208253    | 13.819 | ng/ul | 97        |
| 68) Pentachlorophenol          | 16.71 | 266  | 240199    | 28.029 | ng/ul | 96        |
| 71) Anthracene                 | 17.19 | 178  | 2322976   | 33.177 | ng/ul | 100       |
| 74) Carbazole                  | 17.46 | 167  | 1992419   | 31.856 | ng/ul | 100       |
| 75) Di-n-butylphthalate        | 18.03 | 149  | 2213533   | 27.823 | ng/ul | 99        |
| 80) Butylbenzylphthalate       | 20.42 | 149  | 1137589   | 31.230 | ng/ul | 97        |
| 82) Benzo(a)anthracene         | 21.28 | 228  | 1971869   | 25.334 | ng/ul | 100       |
| 83) Bis(2-ethylhexyl)phthalate | 21.22 | 149  | 1051466   | 19.992 | ng/ul | 99        |
| 86) Di-n-octyl phthalate       | 22.12 | 149  | 3438233   | 39.617 | ng/ul | 100       |
| 87) Benzo(b)fluoranthene       | 22.88 | 252  | 1005175m) | 13.109 | ng/ul |           |
| 88) Benzo(k)fluoranthene       | 22.93 | 252  | 2776280   | 36.988 | ng/ul | 98        |
| 93) Benzo(g,h,i)perylene       | 26.48 | 276  | 1664326   | 21.664 | ng/ul | 97        |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

) JU06107119