

Data Path : Z:\svoasrv\HPCHEM1\BNA\_N\Data\BN051819\

Data File : BN005744.D

Acq On : 18 May 2019 09:20

Operator : JU/SJ

Sample : SSTD00573

Misc :

ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 20 03:39:40 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN051819MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat May 18 11:54:52 2019

Response via : Initial Calibration

Instrument :

BNA\_N

Client Sampled :

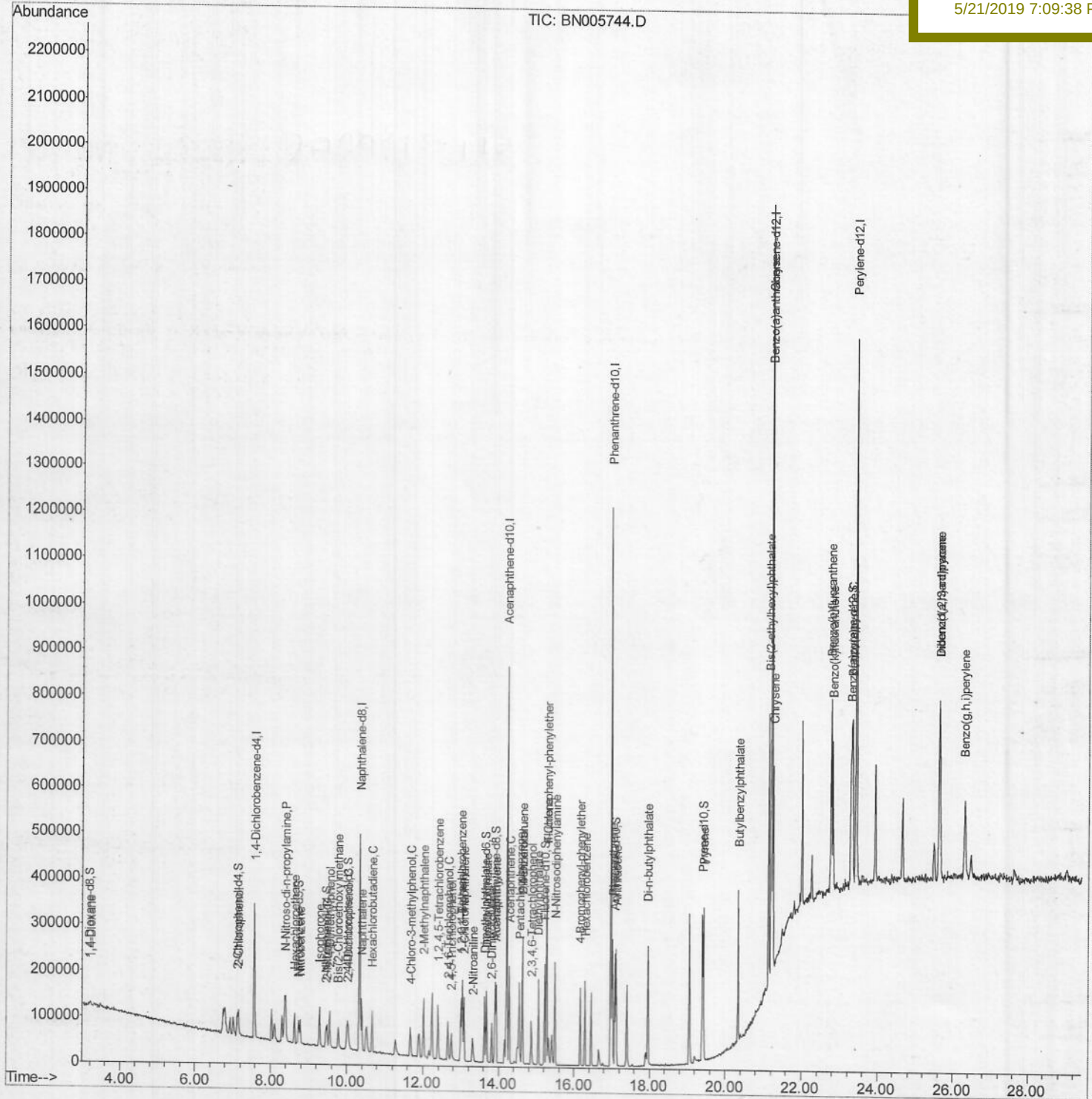
SSTD00573

Manual Integrations

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Acq On : 18 May 2019 09:20  
Operator : JU/SJ  
Sample : SSTD00573  
Misc :  
ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 22 09:51:58 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN051819MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat May 11 23:17:30 2019  
Response via : Initial Calibration

Instrument :

BNA\_N

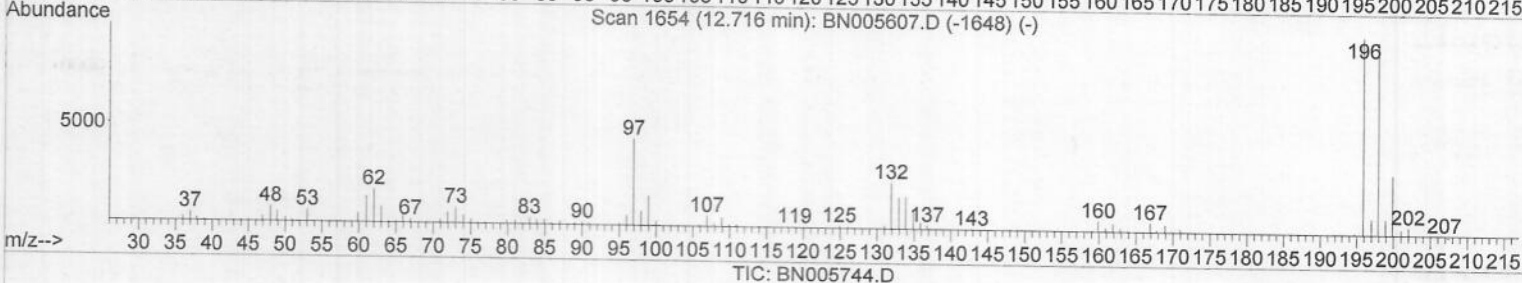
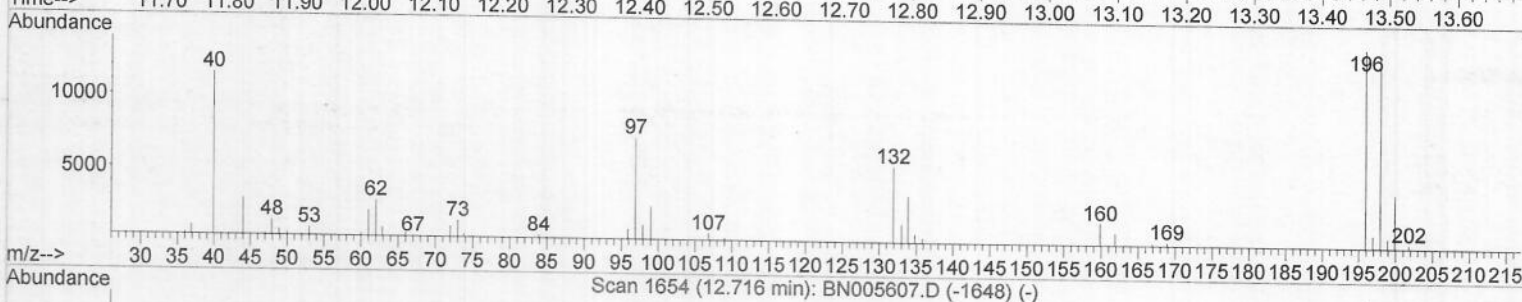
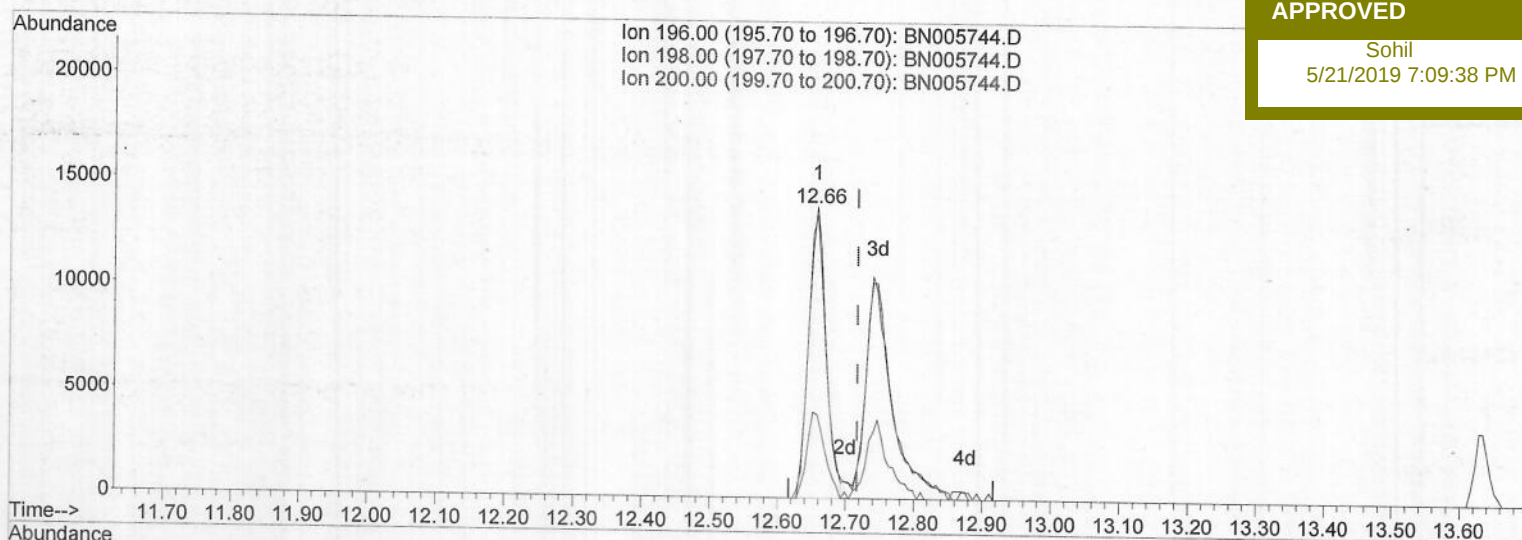
Client Sampled :

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(39) 2,4,5-Trichlorophenol

12.657min (-0.059) 4.34ng/ul

response 24591

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 196.00 | 100   | 100   |
| 198.00 | 93.70 | 97.06 |
| 200.00 | 30.30 | 28.80 |
| 0.00   | 0.00  | 0.00  |

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Acq On : 18 May 2019 09:20

Operator : JU/SJ

Sample : SSTD00573

Misc :

ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 18 12:04:39 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN051819MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Sat May 18 11:54:52 2019

Response via : Initial Calibration

Instrument :

BNA\_N

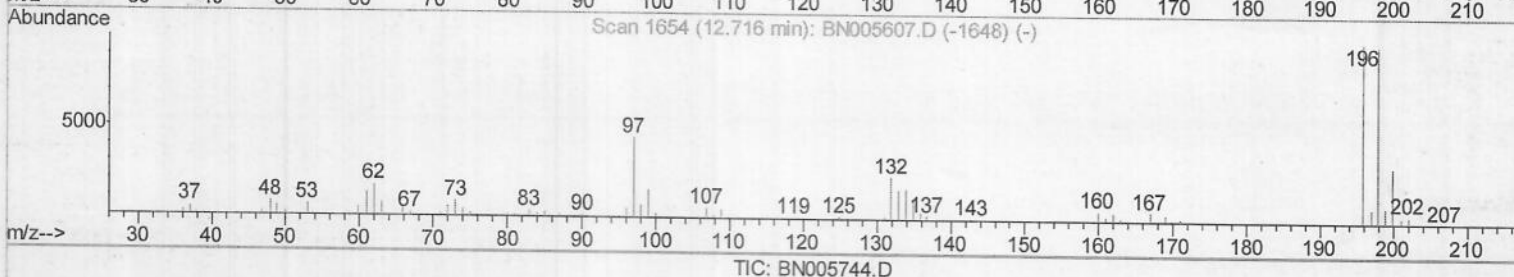
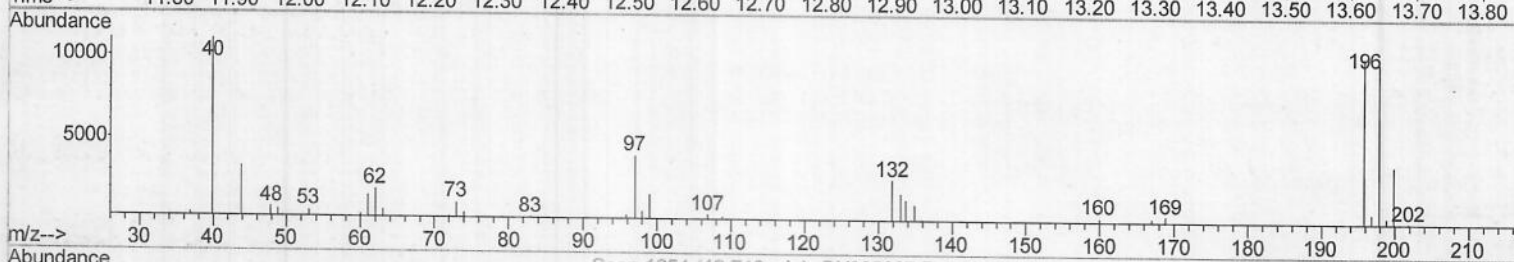
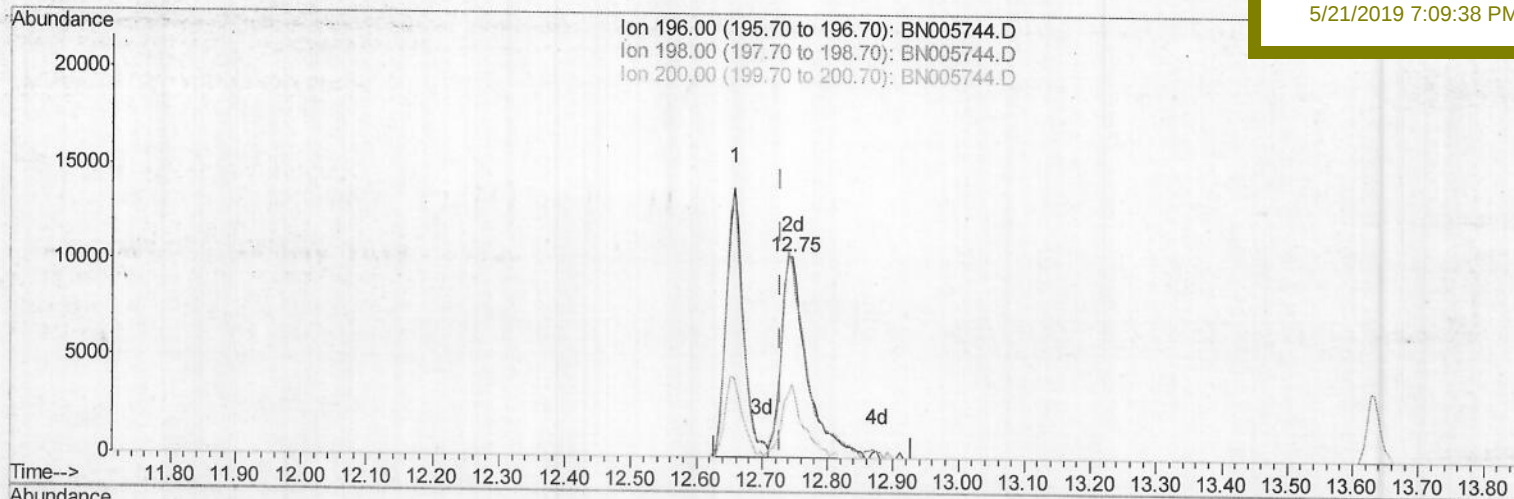
ClientSampleId :

SSTD00573

Manual Integrations  
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TIC: BN005744.D

(39) 2,4,5-Trichlorophenol

12.746min (+0.018) 5.01ng/ul m

&gt; JU/05/22/19

response 28341

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 196.00 | 100   | 100    |
| 198.00 | 93.70 | 96.35  |
| 200.00 | 30.30 | 36.47# |
| 0.00   | 0.00  | 0.00   |

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Misc :

ALS Vial : 2 Sample Multiplier: 1

Quant Time: May 18 12:04:39 2019

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Quant Title : SVOA CALIBRATION

QLast Update : Sat May 18 11:54:52 2019

Response via : Initial Calibration

Instrument :

BNA\_N

ClientSampleId :

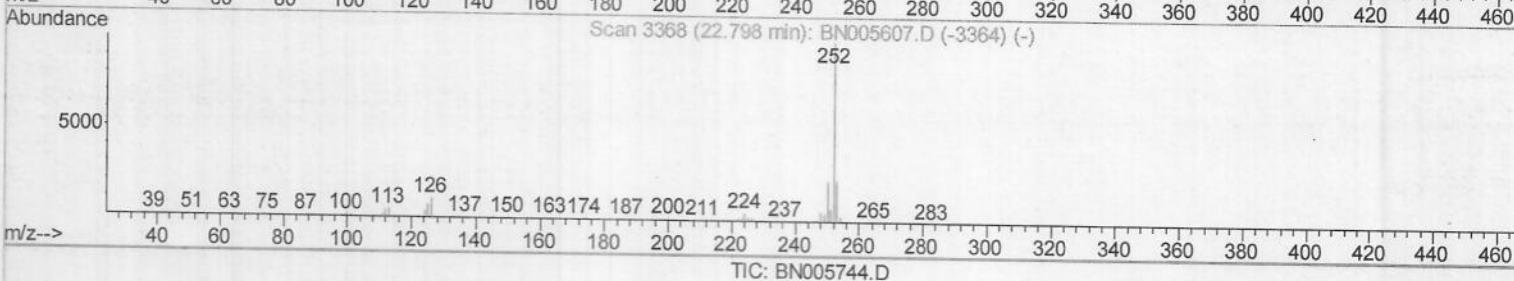
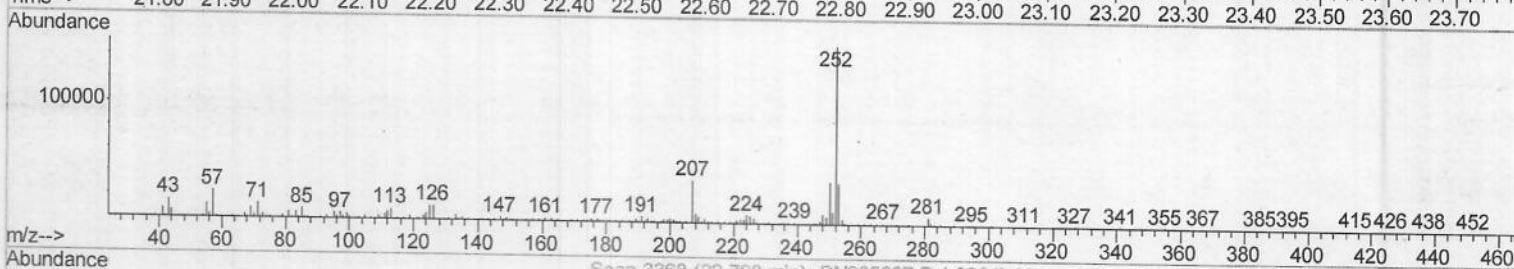
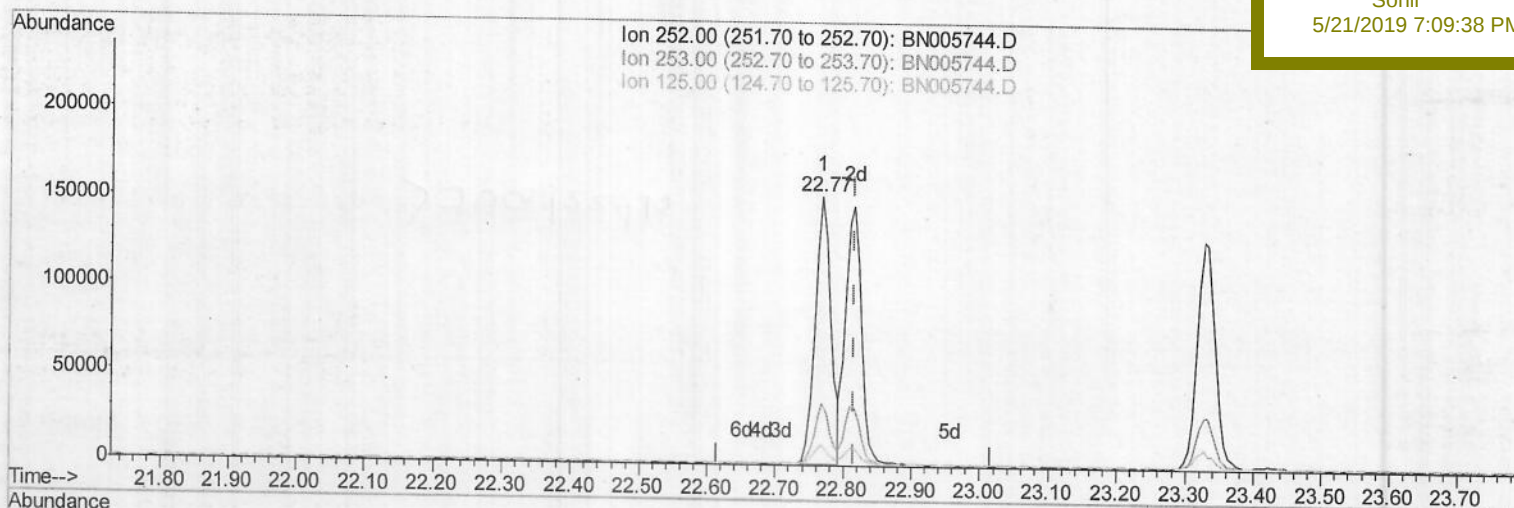
SSTD00573

Manual Integrations

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(88) Benzo(k)fluoranthene

22.769min (-0.047) 4.87ng/ul

response 243944

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 21.70 | 22.66 |
| 125.00 | 9.40  | 7.70  |
| 0.00   | 0.00  | 0.00  |

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Sample : SST00573

Misc :

ALS Vial : 2 Sample Multiplier: 1

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QLast Update : Sat May 18 11:54:52 2019

Response via : Initial Calibration

Instrument :

BNA\_N

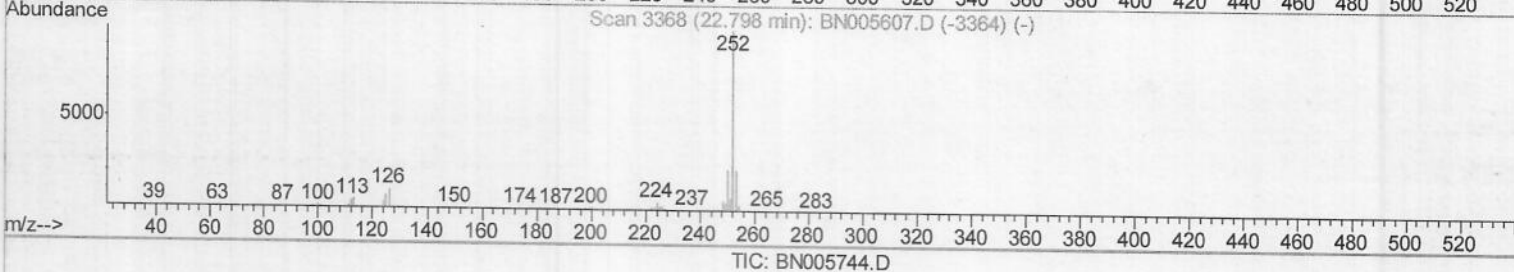
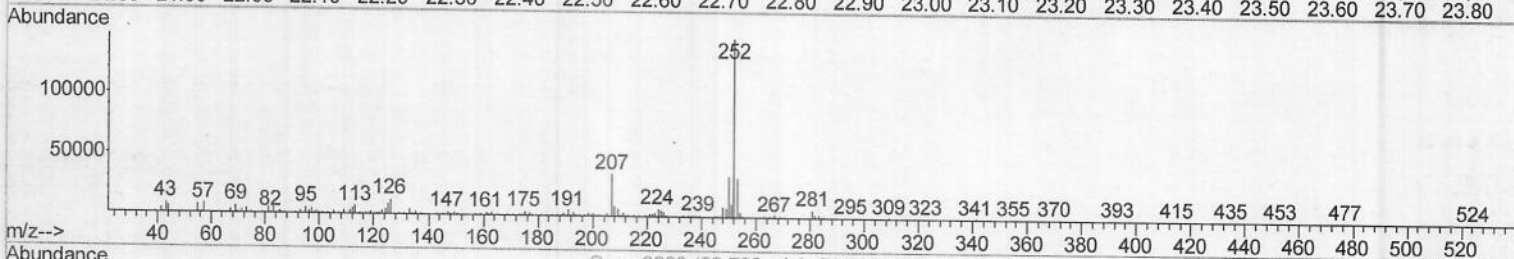
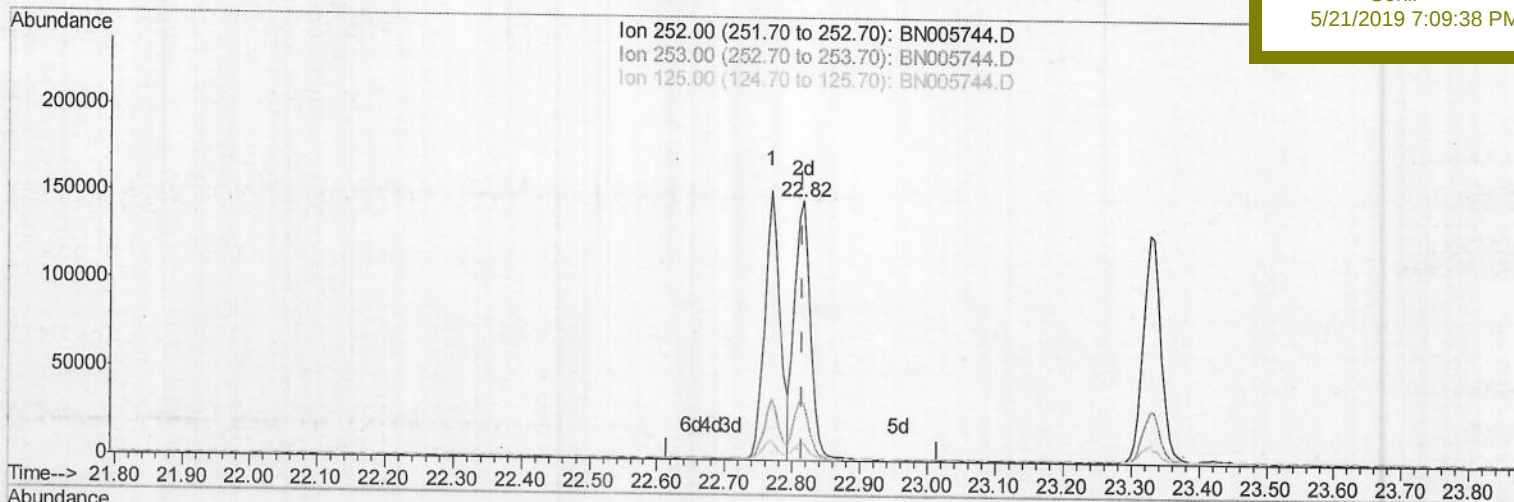
Client Sample Id :

SSTD00573

Manual Integrations  
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(88) Benzo(k)fluoranthene

22.816min (+0.000) 4.82ng/ul m

response 241310

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 252.00 | 100   | 100   |
| 253.00 | 21.70 | 22.19 |
| 125.00 | 9.40  | 6.54# |
| 0.00   | 0.00  | 0.00  |

&gt; JU/05/22/19

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 Sample : SST00573  
 Misc :  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 Client Sampled :  
 SST00573

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Manual Integrations  
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| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.58  | 152  | 87166    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.34 | 136  | 398112   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.22 | 164  | 282274   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 16.97 | 188  | 729205   | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.21 | 240  | 782316   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 23.43 | 264  | 932719   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.19  | 96  | 3237   | 1.79 | ng/uL | 0.00 |
| 5) Phenol-d5                   | 0.00  | 99  | 0d     | 0.00 | ng/ul |      |
| 7) Bis-(2-Chloroethyl) ether-d | 0.00  | 67  | 0d     | 0.00 | ng/ul |      |
| 9) 2-Chlorophenol-d4           | 7.12  | 132 | 24694  | 4.71 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 0.00  | 113 | 0d     | 0.00 | ng/ul |      |
| 19) Nitrobenzene-d5            | 8.73  | 128 | 13347  | 5.45 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.45  | 143 | 14249  | 5.74 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.99  | 165 | 28502  | 5.01 | ng/ul | 0.02 |
| 29) 4-Chloroaniline-d4         | 0.00  | 131 | 0d     | 0.00 | ng/ul |      |
| 43) Dimethylphthalate-d6       | 13.63 | 166 | 104674 | 5.11 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 13.90 | 160 | 125328 | 4.98 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 0.00  | 143 | 0d     | 0.00 | ng/ul |      |
| 57) Fluorene-d10               | 15.22 | 176 | 90132  | 5.24 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 0.00  | 200 | 0d     | 0.00 | ng/ul |      |
| 70) Anthracene-d10             | 17.08 | 188 | 148570 | 4.86 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.39 | 212 | 177808 | 4.44 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 23.29 | 264 | 199589 | 4.90 | ng/ul | 0.00 |

## Target Compounds

|                                |       |     |        |       | Qvalue |     |
|--------------------------------|-------|-----|--------|-------|--------|-----|
| 2) 1,4-Dioxane                 | 3.22  | 88  | 3416   | 1.759 | ng/uL# | 90  |
| 10) 2-Chlorophenol             | 7.15  | 128 | 25313  | 4.673 | ng/ul  | 92  |
| 15) N-Nitroso-di-n-propylamine | 8.38  | 70  | 21178  | 4.623 | ng/ul# | 89  |
| 17) Hexachloroethane           | 8.63  | 117 | 11845  | 5.036 | ng/ul  | 89  |
| 20) Nitrobenzene               | 8.78  | 77  | 31378  | 4.653 | ng/ul  | 93  |
| 21) Isophorone                 | 9.28  | 82  | 62769  | 4.701 | ng/ul  | 95  |
| 23) 2-Nitrophenol              | 9.48  | 139 | 15266  | 5.423 | ng/ul  | 95  |
| 24) 2,4-Dimethylphenol         | 9.55  | 107 | 33320  | 4.727 | ng/ul  | 98  |
| 25) Bis(2-Chloroethoxy)methane | 9.78  | 93  | 35680  | 4.252 | ng/ul  | 97  |
| 27) 2,4-Dichlorophenol         | 10.02 | 162 | 26824  | 4.773 | ng/ul  | 95  |
| 28) Naphthalene                | 10.39 | 128 | 91419  | 4.940 | ng/ul  | 96  |
| 31) Hexachlorobutadiene        | 10.66 | 225 | 21451  | 5.405 | ng/ul  | 94  |
| 33) 4-Chloro-3-methylphenol    | 11.66 | 107 | 32195  | 5.088 | ng/ul  | 93  |
| 34) 2-Methylnaphthalene        | 12.02 | 142 | 69887  | 5.157 | ng/ul  | 97  |
| 36) 1,2,4,5-Tetrachlorobenzene | 12.40 | 216 | 41959  | 4.803 | ng/ul  | 97  |
| 38) 2,4,6-Trichlorophenol      | 12.66 | 196 | 24591  | 4.668 | ng/ul  | 97  |
| 39) 2,4,5-Trichlorophenol      | 12.75 | 196 | 28341m | 5.007 | ng/ul  | 97  |
| 40) 1,1'-Biphenyl              | 13.04 | 154 | 97740  | 4.647 | ng/ul# | 97  |
| 41) 2-Chloronaphthalene        | 13.09 | 162 | 76007  | 4.699 | ng/ul  | 99  |
| 42) 2-Nitroaniline             | 13.31 | 65  | 19777  | 4.610 | ng/ul  | 96  |
| 44) Dimethylphthalate          | 13.68 | 163 | 103943 | 5.091 | ng/ul# | 99  |
| 45) 2,6-Dinitrotoluene         | 13.82 | 165 | 19704  | 5.589 | ng/ul  | 93  |
| 47) Acenaphthylene             | 13.93 | 152 | 121106 | 4.897 | ng/ul  | 100 |

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Manual Integrations  
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| Internal Standards             | R.T.  | QIon | Response | Conc  | Units  | Dev(Min) |
|--------------------------------|-------|------|----------|-------|--------|----------|
| 49) Acenaphthene               | 14.28 | 153  | 82947    | 4.913 | ng/ul  | 96       |
| 53) Dibenzofuran               | 14.63 | 168  | 123291   | 5.072 | ng/ul  | 97       |
| 54) 2,4-Dinitrotoluene         | 14.63 | 165  | 29889    | 5.881 | ng/ul# | 74       |
| 55) 2,3,4,6-Tetrachlorophenol  | 14.87 | 232  | 24448    | 5.192 | ng/ul# | 91       |
| 56) Diethylphthalate           | 15.05 | 149  | 104989   | 5.128 | ng/ul  | 98       |
| 58) Fluorene                   | 15.27 | 166  | 100881   | 5.318 | ng/ul  | 98       |
| 59) 4-Chlorophenyl-phenylether | 15.27 | 204  | 55112    | 5.477 | ng/ul  | 96       |
| 64) N-Nitrosodiphenylamine     | 15.49 | 169  | 87668    | 4.393 | ng/ul  | 96       |
| 65) 4-Bromophenyl-phenylether  | 16.17 | 248  | 35593    | 4.845 | ng/ul  | 91       |
| 66) Hexachlorobenzene          | 16.29 | 284  | 40881    | 5.226 | ng/ul  | 94       |
| 69) Phenanthrene               | 17.02 | 178  | 176311   | 4.965 | ng/ul  | 98       |
| 71) Anthracene                 | 17.12 | 178  | 172187   | 4.760 | ng/ul  | 100      |
| 72) 1,2,3,4-Tetrachlorobenzene | 13.01 | 216  | 44266    | 4.156 | ng/uL  | 98       |
| 73) Pentachlorobenzene         | 14.55 | 250  | 45419    | 4.738 | ng/uL  | 95       |
| 75) Di-n-butylphthalate        | 17.96 | 149  | 186068   | 4.772 | ng/ul  | 98       |
| 79) Pyrene                     | 19.42 | 202  | 226740   | 4.501 | ng/ul# | 90       |
| 80) Butylbenzylphthalate       | 20.34 | 149  | 91145    | 4.570 | ng/ul  | 95       |
| 82) Benzo(a)anthracene         | 21.20 | 228  | 237873   | 4.888 | ng/ul  | 98       |
| 83) Bis(2-ethylhexyl)phthalate | 21.13 | 149  | 140822   | 4.586 | ng/ul# | 97       |
| 84) Chrysene                   | 21.25 | 228  | 226784   | 5.021 | ng/ul  | 98       |
| 87) Benzo(b)fluoranthene       | 22.77 | 252  | 243944   | 4.653 | ng/ul# | 97       |
| 88) Benzo(k)fluoranthene       | 22.82 | 252  | 241310m  | 4.819 | ng/ul  | 97       |
| 90) Benzo(a)pyrene             | 23.33 | 252  | 236413   | 4.797 | ng/ul# | 97       |
| 91) Indeno(1,2,3-cd)pyrene     | 25.63 | 276  | 299121   | 5.490 | ng/ul# | 92       |
| 92) Dibenzo(a,h)anthracene     | 25.63 | 278  | 252918   | 5.470 | ng/ul# | 94       |
| 93) Benzo(g,h,i)perylene       | 26.31 | 276  | 248138   | 5.508 | ng/ul# | 91       |

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