

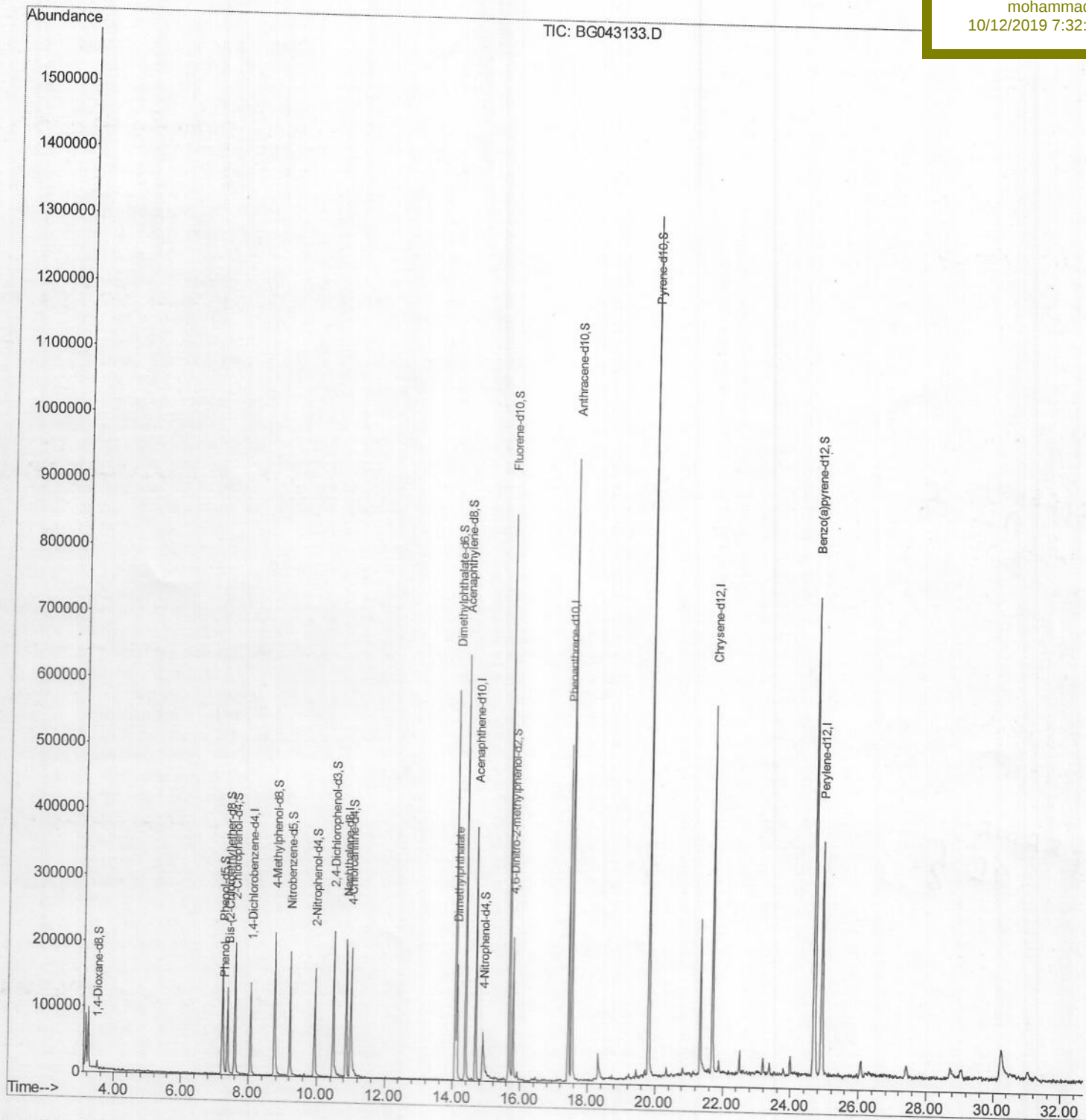
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\  
Data File : BG043133.D  
Acq On : 10 Oct 2019 12:15  
Operator : JU  
Sample : K5097-16  
Misc :  
ALS Vial : 30 Sample Multiplier: 1

Quant Time: Oct 10 14:16:58 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG100919MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Thu Oct 10 11:52:02 2019  
Response via : Initial Calibration

Instrument :  
BNA\_G  
ClientSampled :  
JLMF4

Manual Integrations  
APPROVED

mohammad  
10/12/2019 7:32:44 AM



# Quantitation Report (Qedit)

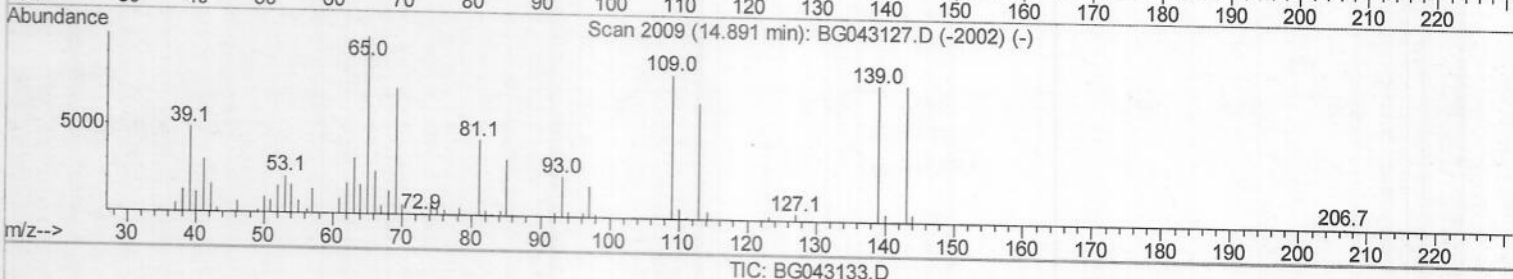
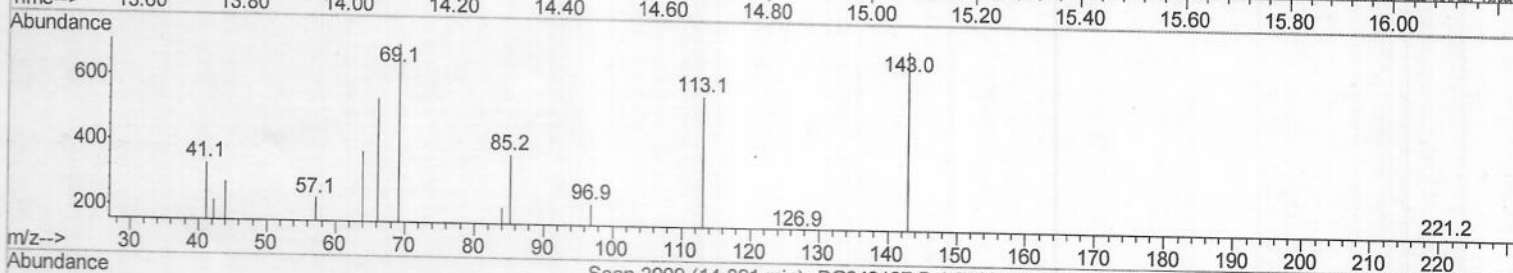
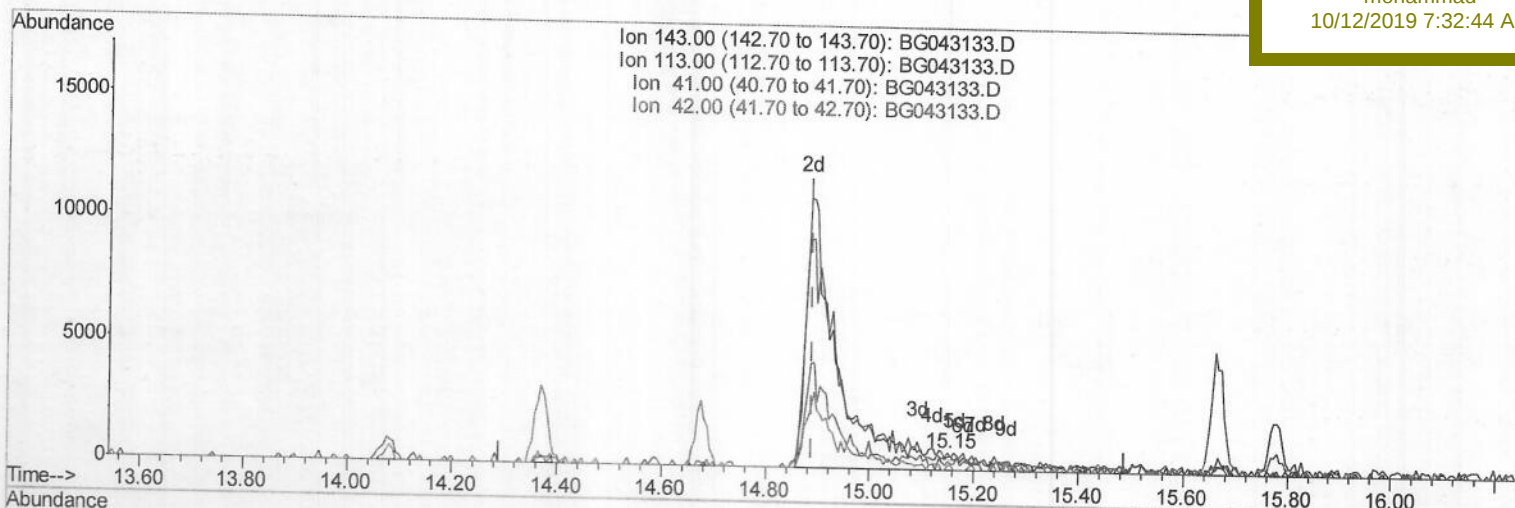
Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\  
 Data File : BG043133.D  
 Acq On : 10 Oct 2019 12:15  
 Operator : JU  
 Sample : K5097-16  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Oct 10 14:13:56 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG100919MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 10 11:52:02 2019  
 Response via : Initial Calibration

Instrument :  
 BNA\_G  
 ClientSampleId :  
 JLMF4

Manual Integrations  
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(51) 4-Nitrophenol-d4 (S)  
 15.151min (+0.262) 0.11ng/ul  
 response 167

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 143.00 | 100   | 100   |
| 113.00 | 84.80 | 78.87 |
| 41.00  | 42.80 | 46.81 |
| 42.00  | 28.10 | 30.64 |

# Quantitation Report (Qedit)

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\

Data File : BG043133.D

Acq On : 10 Oct 2019 12:15

Operator : JU

Sample : K5097-16

Misc :

ALS Vial : 30 Sample Multiplier: 1

Quant Time: Oct 10 14:13:56 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG100919MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Oct 10 11:52:02 2019

Response via : Initial Calibration

Instrument :

BNA\_G

ClientSampleId :

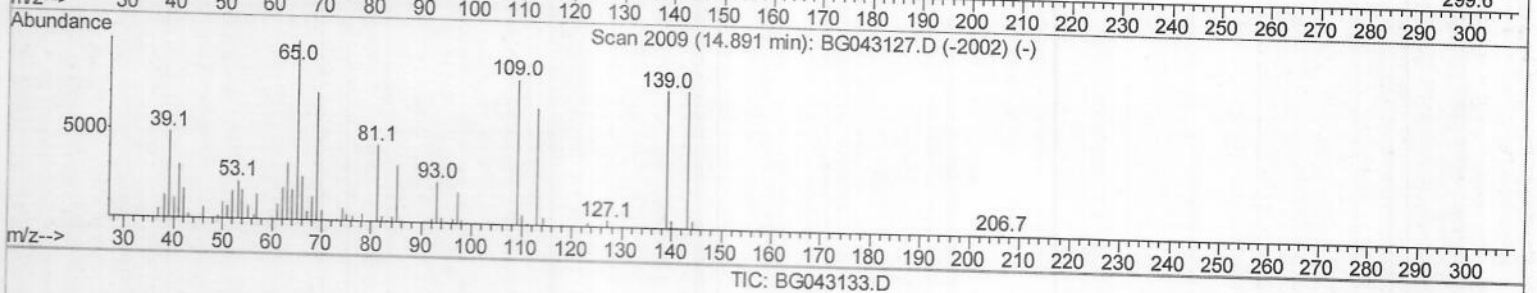
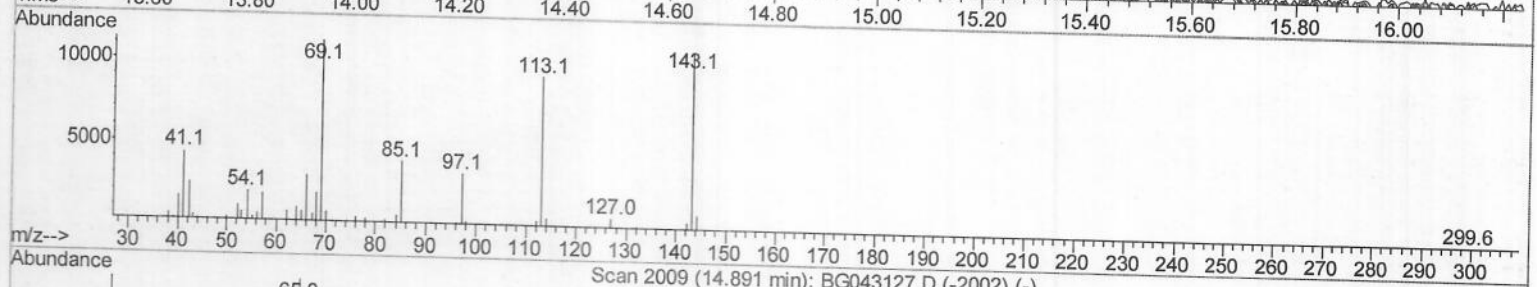
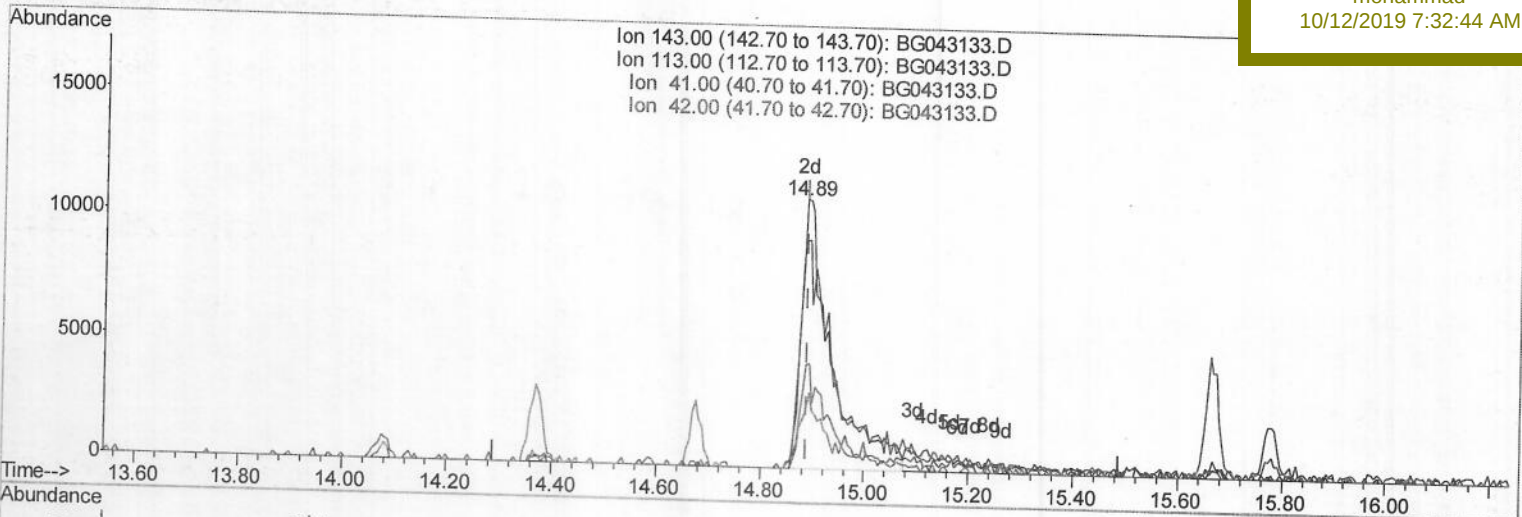
JLMF4

Manual Integrations

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mohammad

10/12/2019 7:32:44 AM



(51) 4-Nitrophenol-d4 (S)

14.886min (-0.003) 28.80ng/ul m > JU 10/11/19

response 44277

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 143.00 | 100   | 100    |
| 113.00 | 84.80 | 84.58  |
| 41.00  | 42.80 | 39.18  |
| 42.00  | 28.10 | 22.04# |

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100919\  
 Data File : BG043133.D  
 Acq On : 10 Oct 2019 12:15  
 Operator : JU  
 Sample : K5097-16  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Quant Time: Oct 10 14:16:58 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_G\METHODS\SOM-EPA-BG100919MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Thu Oct 10 11:52:02 2019  
 Response via : Initial Calibration

Instrument :  
 BNA\_G  
 ClientSampleId :  
 JLMF4

Manual Integrations  
 APPROVED

mohammad  
 10/12/2019 7:32:44 AM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 8.06  | 152  | 46583    | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.86 | 136  | 186300   | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.67 | 164  | 144198   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.42 | 188  | 347631   | 20.00 | ng/ul | 0.00     |
| 77) Chrysene-d12          | 21.69 | 240  | 371840   | 20.00 | ng/ul | 0.00     |
| 85) Perylene-d12          | 24.91 | 264  | 436173   | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |       |       |      |
|--------------------------------|-------|-----|--------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.50  | 96  | 5799   | 6.02  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 7.22  | 99  | 112610 | 32.70 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl) ether-d | 7.38  | 67  | 67723  | 35.24 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.59  | 132 | 99061  | 35.89 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.76  | 113 | 98576  | 33.20 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 9.22  | 128 | 52426  | 39.97 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.94  | 143 | 59011  | 39.47 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.49 | 165 | 104887 | 33.53 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 11.00 | 131 | 131617 | 40.42 | ng/ul | 0.00 |
| 43) Dimethylphthalate-d6       | 14.08 | 166 | 442271 | 38.15 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.37 | 160 | 490801 | 38.92 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.89 | 143 | 44277m | 28.80 | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.67 | 176 | 386681 | 39.28 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.78 | 200 | 62086  | 27.84 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.52 | 188 | 646732 | 39.38 | ng/ul | 0.00 |
| 78) Pyrene-d10                 | 19.80 | 212 | 755170 | 41.19 | ng/ul | 0.00 |
| 89) Benzo(a)pyrene-d12         | 24.69 | 264 | 850989 | 39.48 | ng/ul | 0.00 |

## Target Compounds

|                       |       |     |        |        | Ovalue   |
|-----------------------|-------|-----|--------|--------|----------|
| 6) Phenol             | 7.25  | 94  | 13102  | 3.726  | ng/ul 91 |
| 44) Dimethylphthalate | 14.13 | 163 | 116072 | 10.366 | ng/ul 96 |

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