

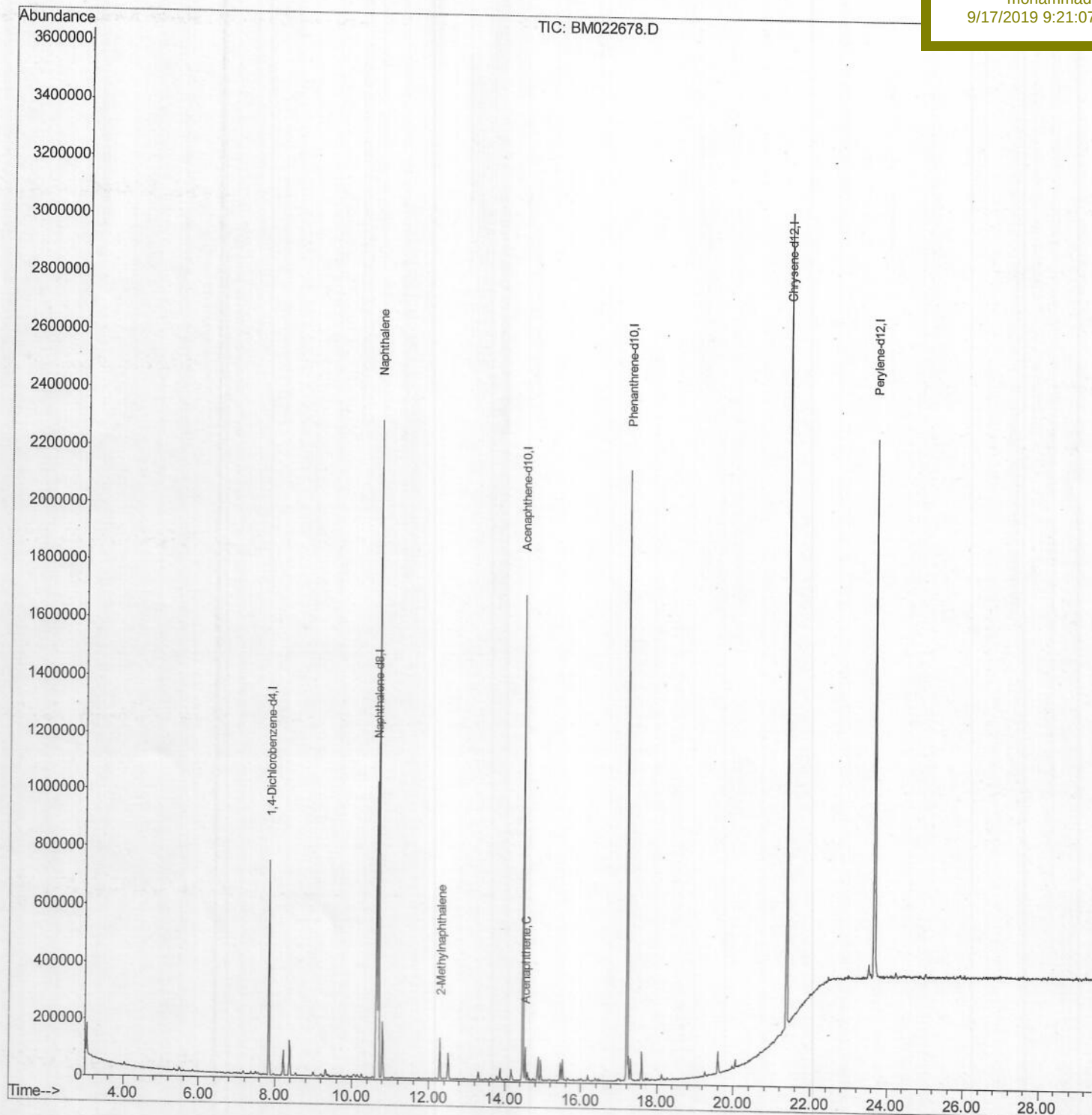
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\  
Data File : BM022678.D  
Acq On : 13 Sep 2019 15:58  
Operator : HP/JU  
Sample : K4706-07DL2 50X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Sep 14 01:21:33 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM090519MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Sep 09 07:45:00 2019  
Response via : Initial Calibration

Instrument :  
BNA\_M  
ClientSampleId :  
BF573DL2

Manual Integrations  
APPROVED

mohammad  
9/17/2019 9:21:07 AM



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\  
Data File : BM022678.D  
Acq On : 13 Sep 2019 15:58  
Operator : HP/JU  
Sample : K4706-07DL2 50X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Sep 14 00:42:18 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM090519MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Sep 09 07:45:00 2019  
Response via : Initial Calibration

Instrument :

BNA\_M

ClientSampleId :

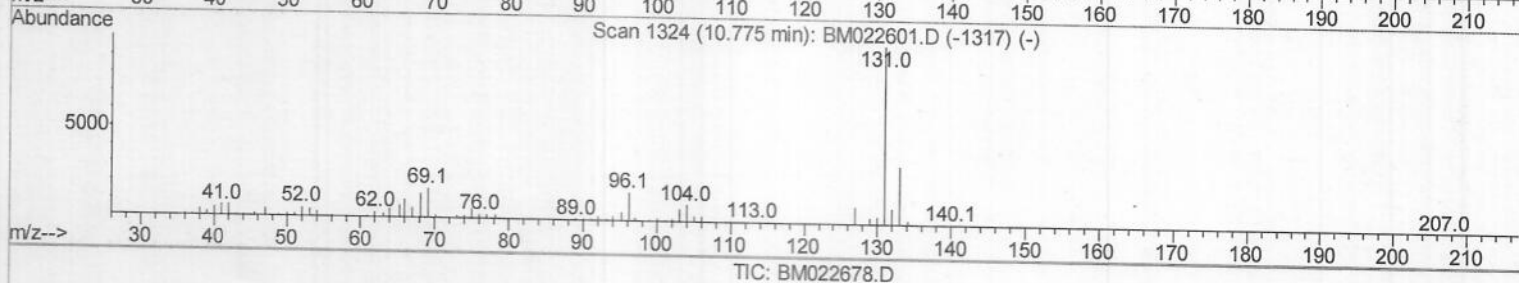
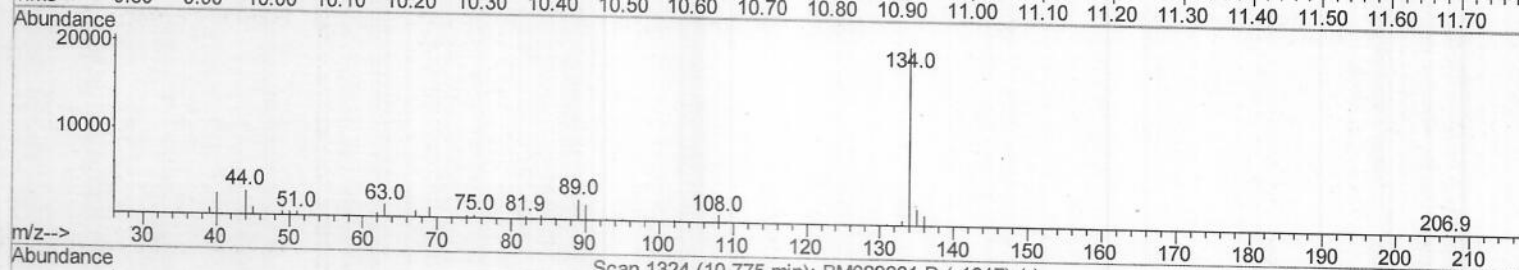
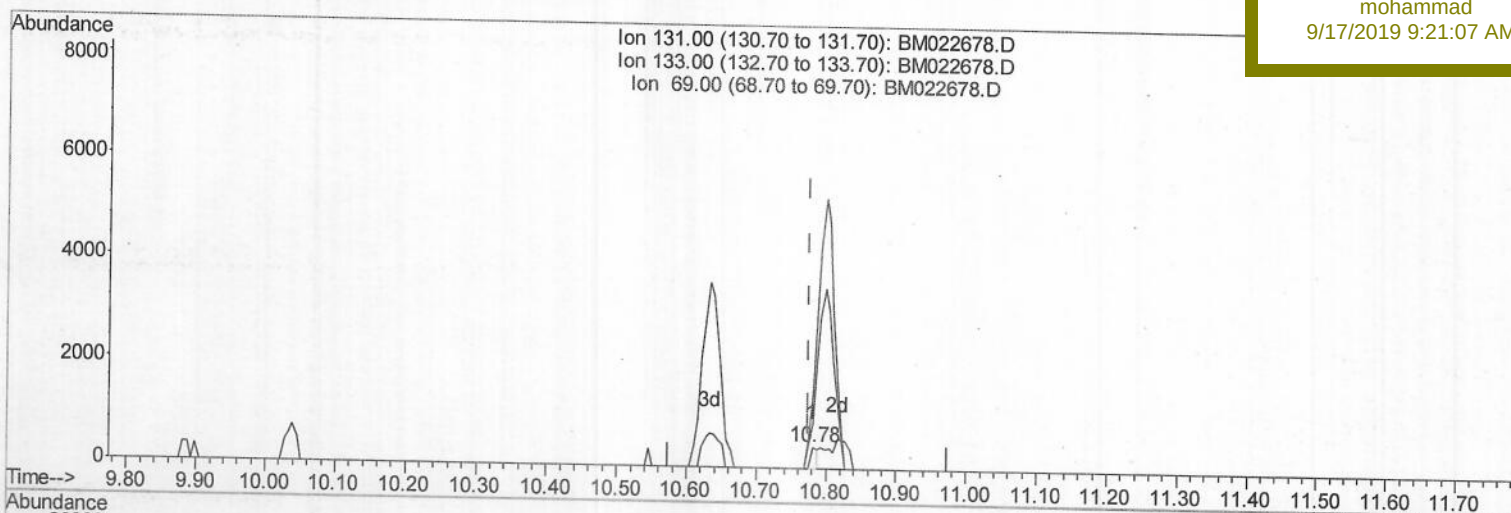
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Manual Integrations

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(29) 4-Chloroaniline-d4 (S)

10.780min (+0.006) 0.01ng/ul

response 273

| Ion    | Exp%  | Act%    |
|--------|-------|---------|
| 131.00 | 100   | 100     |
| 133.00 | 32.10 | 206.22# |
| 69.00  | 16.90 | 341.87# |
| 0.00   | 0.00  | 0.00    |

# Quantitation Report (Qedit)

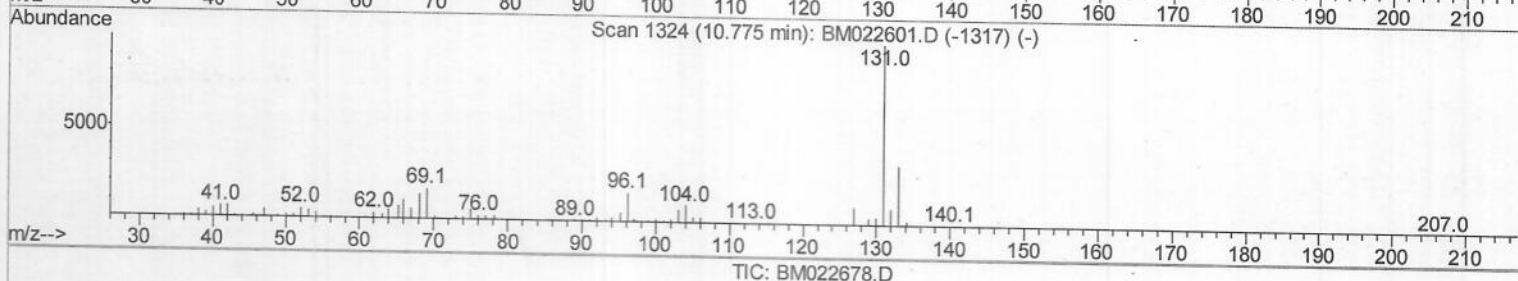
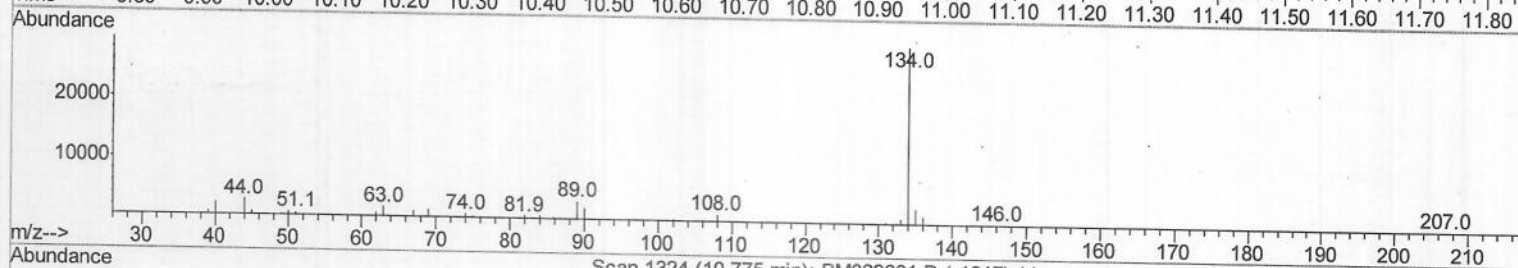
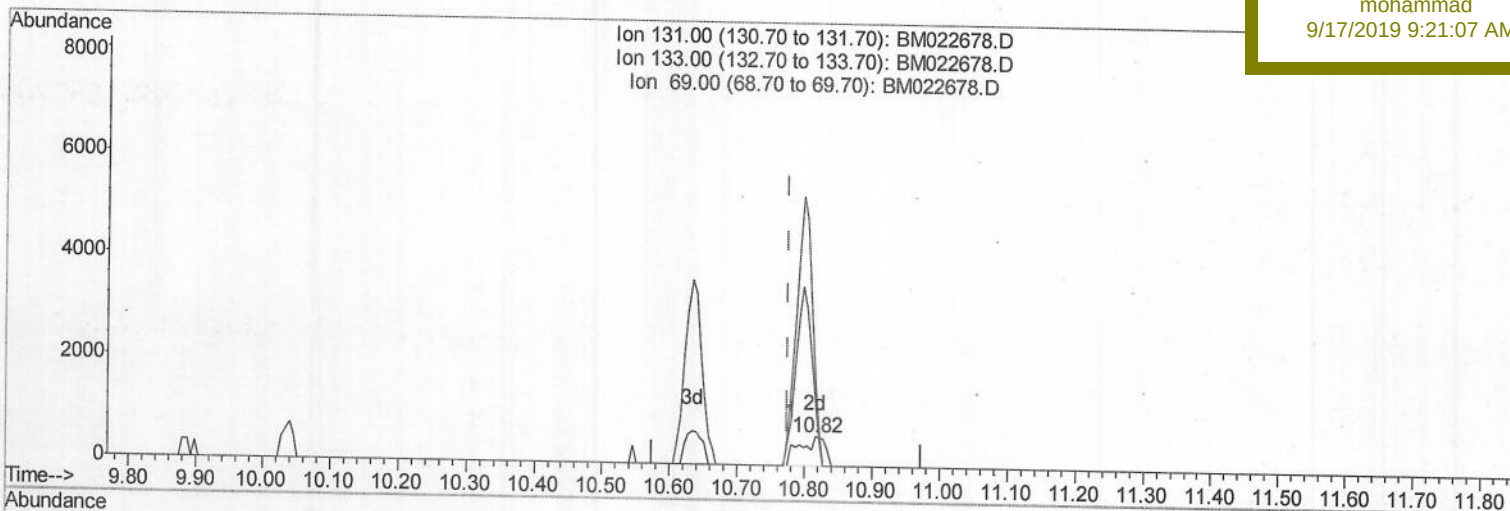
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 ALS Vial : 4 Sample Multiplier: 1

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 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Sep 09 07:45:00 2019  
 Response via : Initial Calibration

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BF573DL2

Manual Integrations  
 APPROVED

mohammad  
 9/17/2019 9:21:07 AM



(29) 4-Chloroaniline-d4 (S)

10.816min (+0.041) 0.08ng/ul m > JU 09/17/19

response 1494

| Ion    | Exp%  | Act%    |
|--------|-------|---------|
| 131.00 | 100   | 100     |
| 133.00 | 32.10 | 208.60# |
| 69.00  | 16.90 | 283.51# |
| 0.00   | 0.00  | 0.00    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\

Data File : BM022678.D

Acq On : 13 Sep 2019 15:58

Operator : HP/JU

Sample : K4706-07DL2 50X

Misc :

ALS Vial : 4 Sample Multiplier: 1

Quant Time: Sep 14 01:21:33 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM090519MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Mon Sep 09 07:45:00 2019

Response via : Initial Calibration

Instrument :

BNA\_M

ClientSampleId :

BF573DL2

Manual Integrations

APPROVED

mohammad

9/17/2019 9:21:07 AM

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.84  | 152  | 197849   | 20.00 | ng/ul | -0.02    |
| 18) Naphthalene-d8        | 10.63 | 136  | 862901   | 20.00 | ng/ul | -0.02    |
| 36) Acenaphthene-d10      | 14.47 | 164  | 554072   | 20.00 | ng/ul | -0.02    |
| 62) Phenanthrene-d10      | 17.21 | 188  | 1195441  | 20.00 | ng/ul | -0.01    |
| 78) Chrysene-d12          | 21.39 | 240  | 1184854  | 20.00 | ng/ul | -0.01    |
| 86) Perylene-d12          | 23.66 | 264  | 1229741  | 20.00 | ng/ul | -0.02    |

## System Monitoring Compounds

|                                |       |     |        |      |       |       |
|--------------------------------|-------|-----|--------|------|-------|-------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0      | 0.00 | ng/ul |       |
| 5) Phenol-d5                   | 7.00  | 99  | 2221   | 0.12 | ng/ul | -0.01 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.17  | 67  | 6255   | 0.60 | ng/ul | 0.00  |
| 9) 2-Chlorophenol-d4           | 7.37  | 132 | 5869   | 0.41 | ng/ul | -0.01 |
| 13) 4-Methylphenol-d8          | 8.54  | 113 | 3585   | 0.24 | ng/ul | -0.02 |
| 19) Nitrobenzene-d5            | 8.99  | 128 | 3580   | 0.58 | ng/ul | -0.01 |
| 22) 2-Nitrophenol-d4           | 9.71  | 143 | 2424   | 0.42 | ng/ul | -0.02 |
| 26) 2,4-Dichlorophenol-d3      | 10.25 | 165 | 5433   | 0.37 | ng/ul | -0.02 |
| 29) 4-Chloroaniline-d4         | 10.82 | 131 | 1494m) | 0.08 | ng/ul | 0.04  |
| 44) Dimethylphthalate-d6       | 13.88 | 166 | 26844  | 0.60 | ng/ul | -0.01 |
| 47) Acenaphthylene-d8          | 14.16 | 160 | 26640  | 0.50 | ng/ul | -0.02 |
| 52) 4-Nitrophenol-d4           | 14.65 | 143 | 498    | 0.06 | ng/ul | -0.01 |
| 58) Fluorene-d10               | 15.46 | 176 | 24733  | 0.66 | ng/ul | -0.01 |
| 63) 4,6-Dinitro-2-methylphenol | 15.57 | 200 | 1332   | 0.25 | ng/ul | -0.02 |
| 71) Anthracene-d10             | 17.30 | 188 | 37969  | 0.62 | ng/ul | -0.02 |
| 79) Pyrene-d10                 | 19.60 | 212 | 39384  | 0.60 | ng/ul | -0.01 |
| 90) Benzo(a)pyrene-d12         | 23.51 | 264 | 32694  | 0.49 | ng/ul | -0.02 |

## Target Compounds

|                         |       |     |         |        |       | Ovalue |
|-------------------------|-------|-----|---------|--------|-------|--------|
| 28) Naphthalene         | 10.68 | 128 | 1918128 | 40.093 | ng/ul | 99     |
| 34) 2-Methylnaphthalene | 12.29 | 142 | 67772   | 1.880  | ng/ul | 98     |
| 50) Acenaphthene        | 14.53 | 153 | 45528   | 1.178  | ng/ul | 97     |

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