

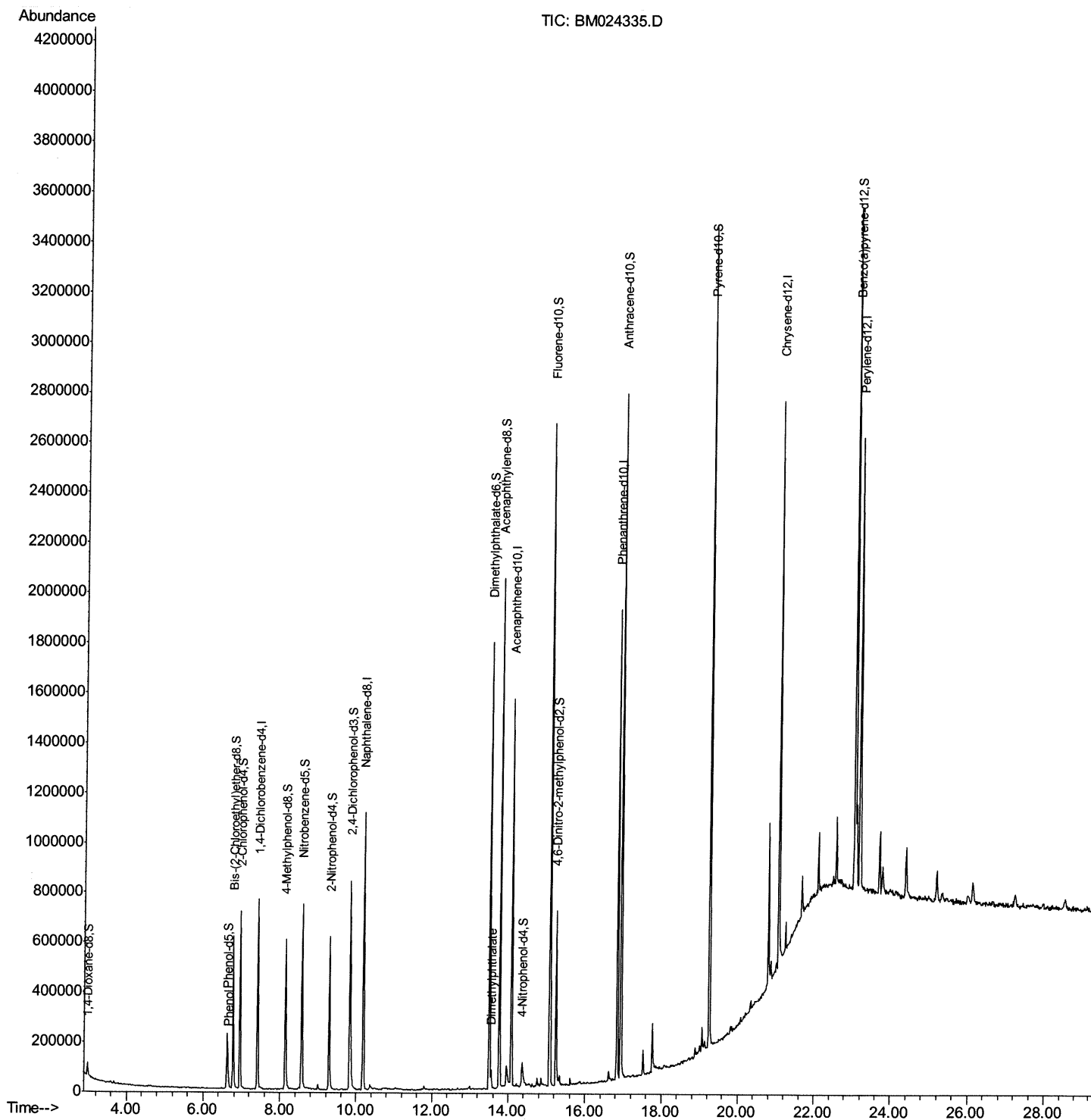
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122719\  
Data File : BM024335.D  
Acq On : 27 Dec 2019 20:19  
Operator : JU  
Sample : K6280-14  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFF36

Manual Integrations  
APPROVED

mohammad  
12/29/2019 5:29:22 PM

Quant Time: Dec 28 03:26:00 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121719MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 27 23:52:10 2019  
Response via : Initial Calibration



## Quantitation Report (Qedit)

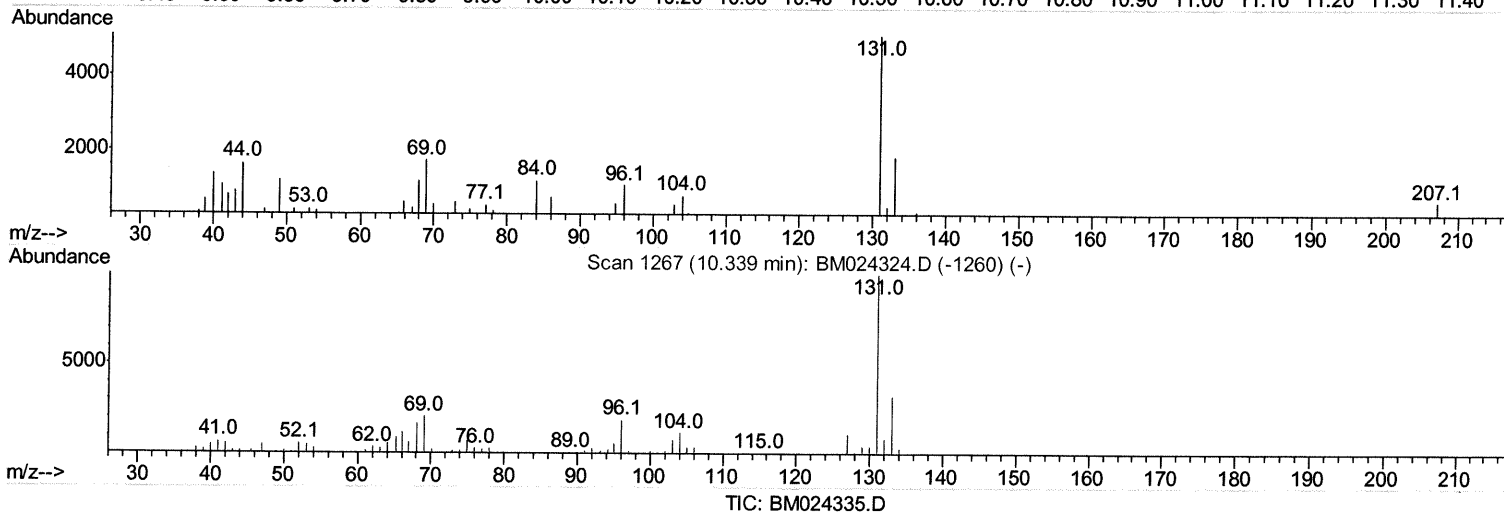
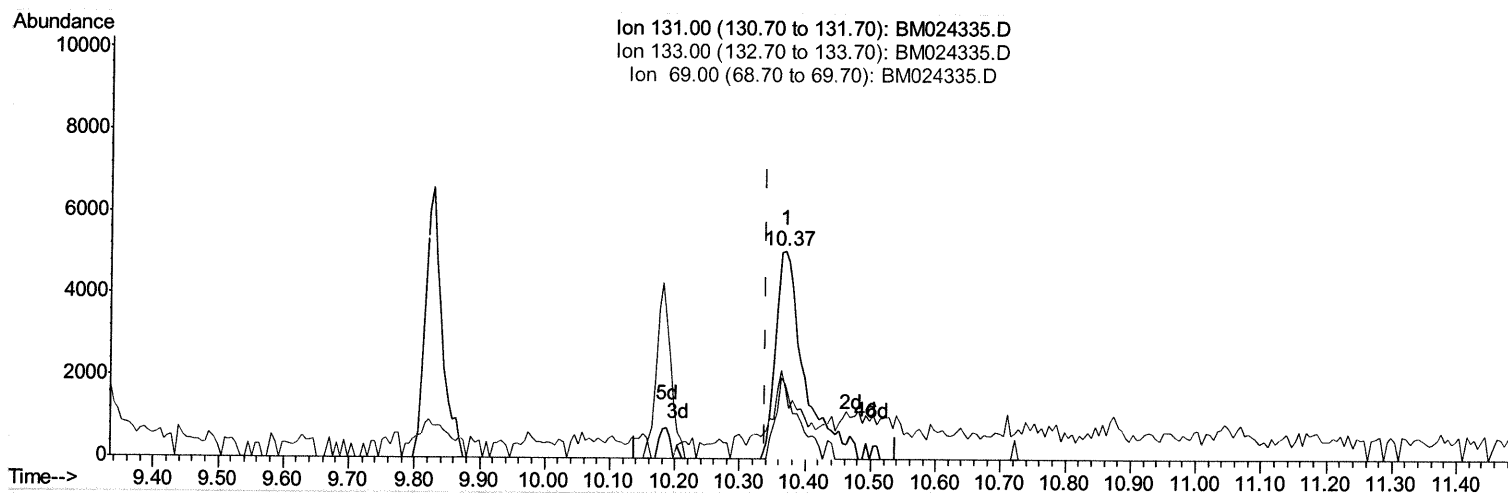
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122719\  
Data File : BM024335.D  
Acq On : 27 Dec 2019 20:19  
Operator : JU  
Sample : K6280-14  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFF36

Manual Integrations  
APPROVED

mohammad  
12/29/2019 5:29:22 PM

Quant Time: Dec 27 23:54:55 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121719MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 27 23:52:10 2019  
Response via : Initial Calibration



(29) 4-Chloroaniline-d4 (S)

10.369min (+0.030) 0.98ng/ul m *JU 12/28/19*

response 15458

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 131.00 | 100   | 100    |
| 133.00 | 32.50 | 36.10  |
| 69.00  | 19.00 | 34.18# |
| 0.00   | 0.00  | 0.00   |

## Quantitation Report (Qedit)

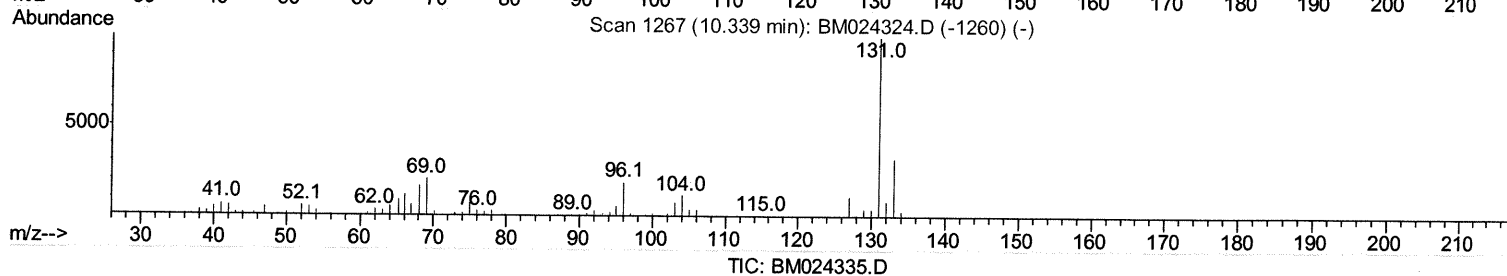
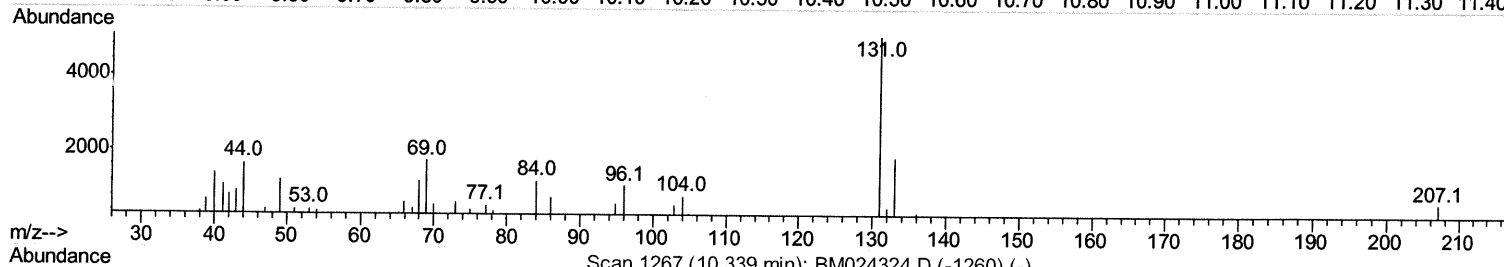
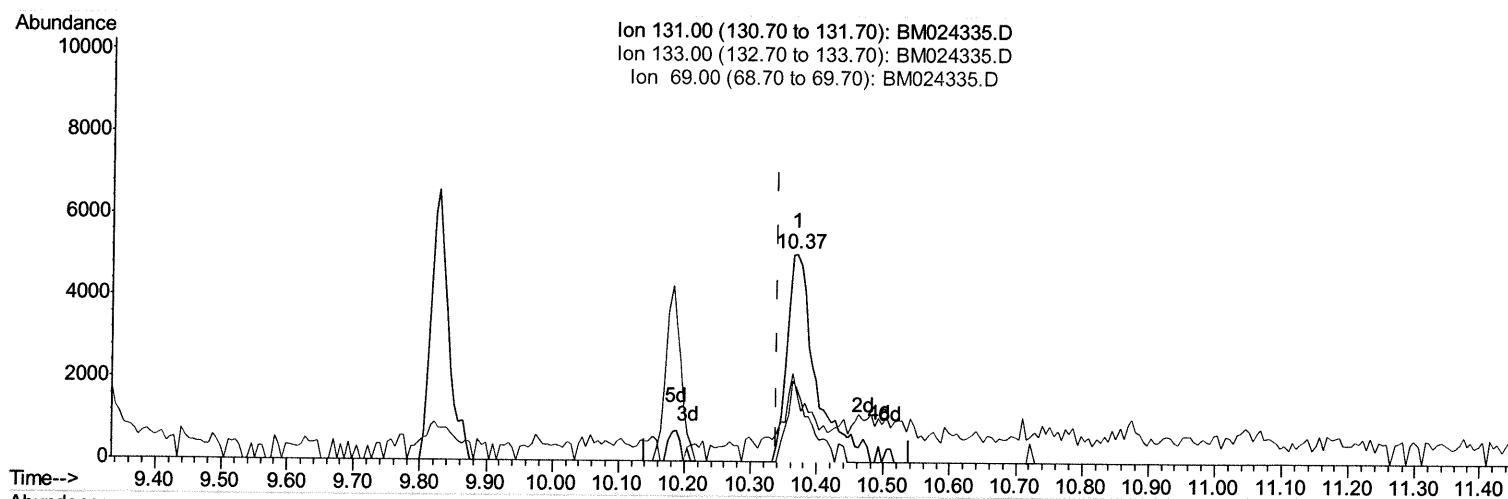
Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122719\  
Data File : BM024335.D  
Acq On : 27 Dec 2019 20:19  
Operator : JU  
Sample : K6280-14  
Misc :  
ALS Vial : 13 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
BFF36

Manual Integrations  
APPROVED

mohammad  
12/29/2019 5:29:22 PM

Quant Time: Dec 28 04:20:27 2019  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121719MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Fri Dec 27 23:52:10 2019  
Response via : Initial Calibration



(29) 4-Chloroaniline-d4 (S)

10.369min (+0.030) 0.92ng/ul

response 14619

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 131.00 | 100   | 100    |
| 133.00 | 32.50 | 36.10  |
| 69.00  | 19.00 | 34.18# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122719\  
 Data File : BM024335.D  
 Acq On : 27 Dec 2019 20:19  
 Operator : JU  
 Sample : K6280-14  
 Misc :  
 ALS Vial : 13 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 BFF36

Manual Integrations  
 APPROVED

mohammad  
 12/29/2019 5:29:22 PM

Quant Time: Dec 28 03:26:00 2019  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM121719MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Fri Dec 27 23:52:10 2019  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.42  | 152  | 201608   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.18 | 136  | 849285   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.08 | 164  | 492918   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.84 | 188  | 1006426  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.05 | 240  | 957619   | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.17 | 264  | 1182054  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 2.98  | 96  | 31811   | 6.97  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.62  | 99  | 144303  | 7.56  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl) ether-d | 6.77  | 67  | 305215  | 26.66 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 6.96  | 132 | 318779  | 23.41 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.14  | 113 | 238654  | 15.29 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.57  | 128 | 170968  | 28.01 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.29  | 143 | 192541  | 28.70 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.83  | 165 | 345931  | 26.52 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.37 | 131 | 15458m  | 0.98  | ng/ul | 0.03 |
| 44) Dimethylphthalate-d6       | 13.50 | 166 | 1130734 | 31.02 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.77 | 160 | 1350700 | 30.93 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.36 | 143 | 40157   | 6.14  | ng/ul | 0.02 |
| 58) Fluorene-d10               | 15.08 | 176 | 985953  | 30.86 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.24 | 200 | 138124  | 22.50 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.94 | 188 | 1435694 | 31.35 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.24 | 212 | 1581736 | 34.68 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.04 | 264 | 1882749 | 32.07 | ng/ul | 0.00 |

Ju12/28/19

## Target Compounds

|                       |       |     |       |       | Ovalue   |
|-----------------------|-------|-----|-------|-------|----------|
| 6) Phenol             | 6.65  | 94  | 20793 | 1.055 | ng/ul 93 |
| 45) Dimethylphthalate | 13.55 | 163 | 41517 | 1.096 | ng/ul 98 |

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