

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022121\  
Data File : BN013716.D  
Acq On : 21 Feb 2021 09:05  
Operator : CG/JU  
Sample : M1429-09  
Misc :  
ALS Vial : 25 Sample Multiplier: 1

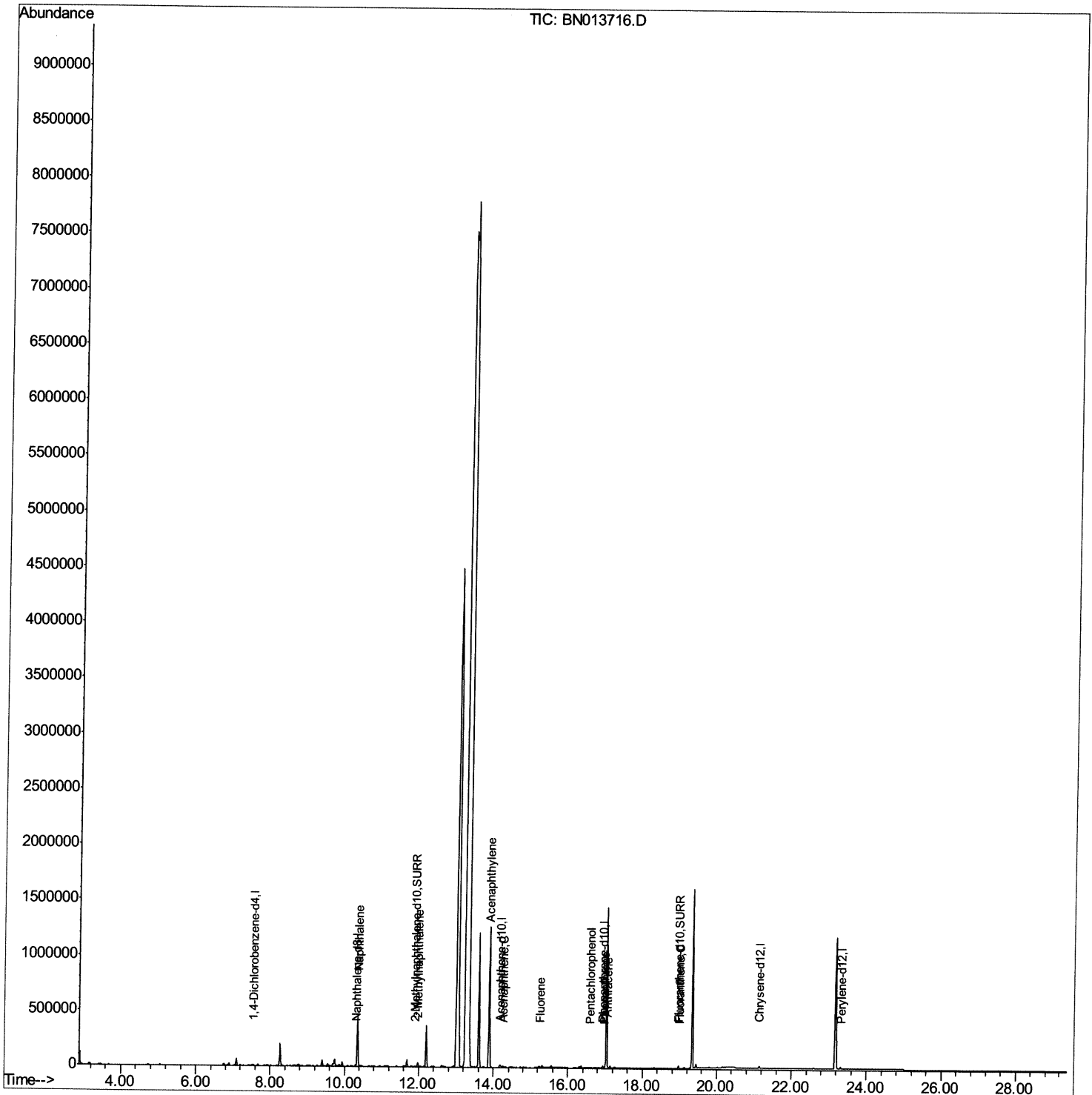
Instrument :  
BNA\_N  
ClientSampleId :  
DBGS1

Manual Integrations  
APPROVED

mohammad

2/22/2021 2:14:27 PM

Quant Time: Feb 22 01:40:04 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN022121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Feb 22 01:25:38 2021  
Response via : Initial Calibration



## Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN022121\  
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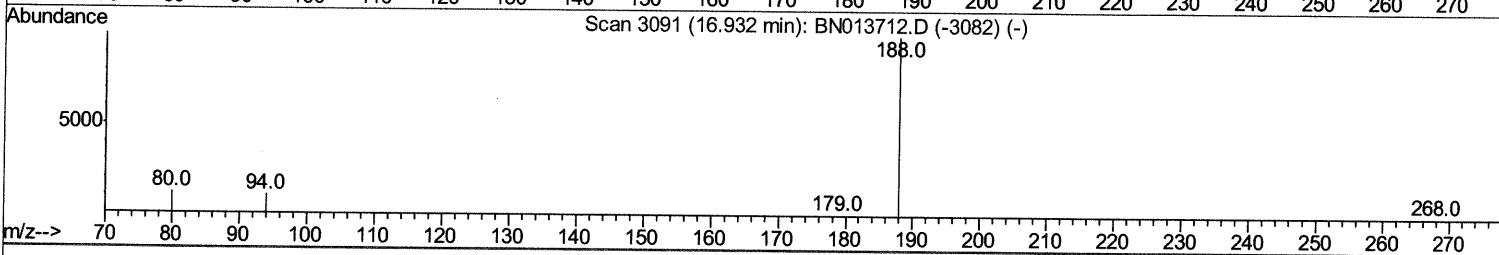
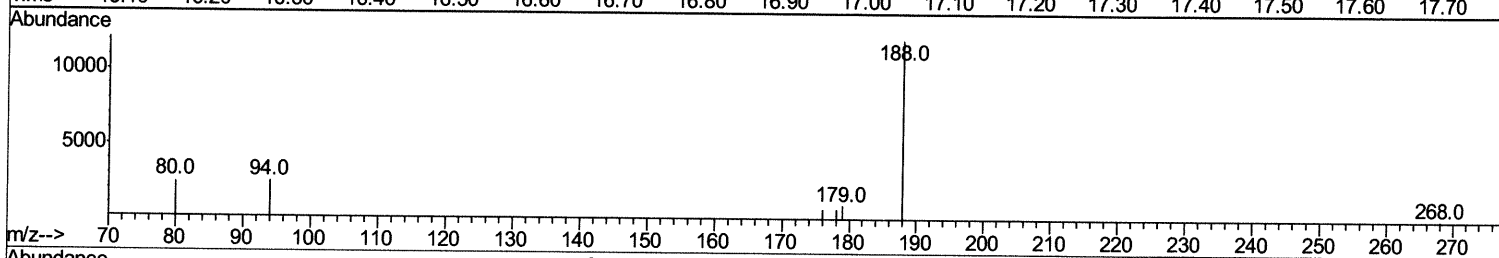
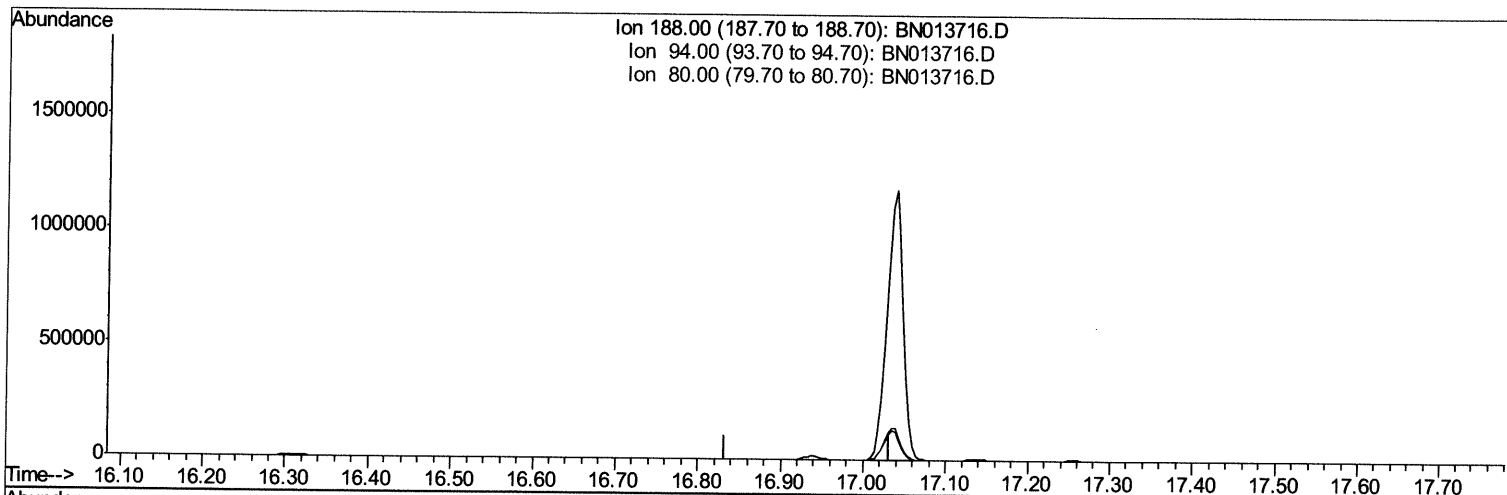
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Quant Time: Feb 22 01:38:12 2021  
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TIC: BN013716.D

(10) Phenanthrene-d10 (I)

16.933min (-16.933) 0.00ng/ul

response 0

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 188.00 | 100   | 0.00  |
| 94.00  | 10.70 | 0.00# |
| 80.00  | 11.60 | 0.00# |
| 0.00   | 0.00  | 0.00  |

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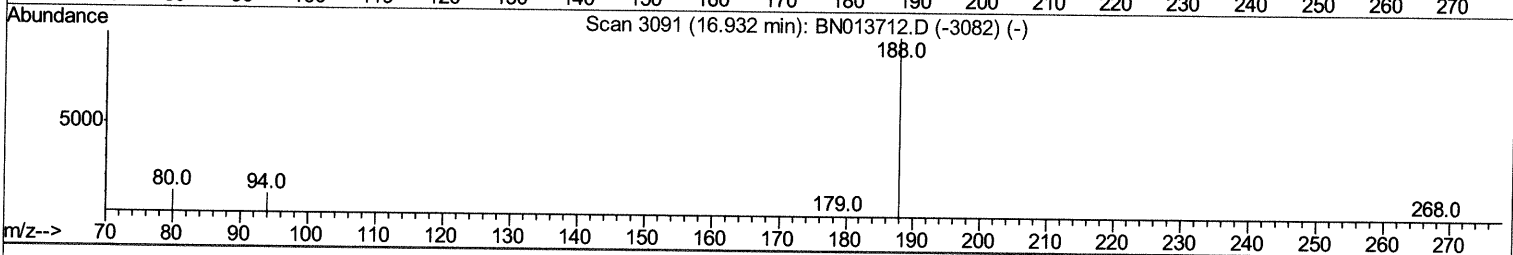
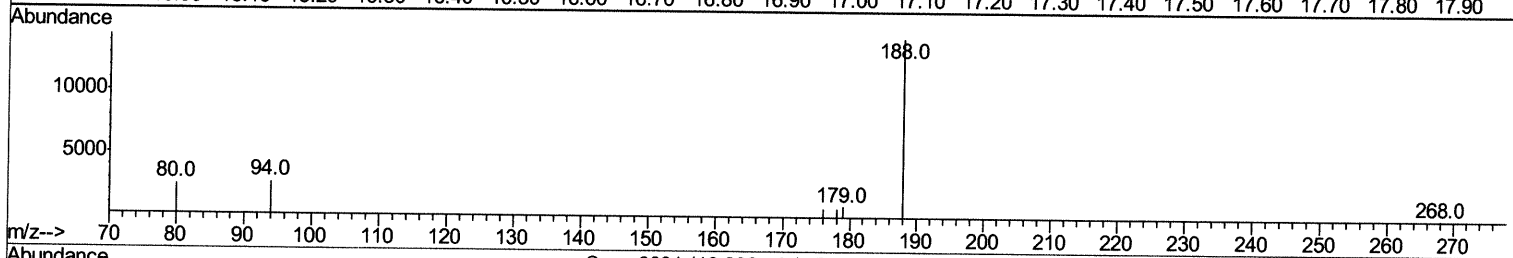
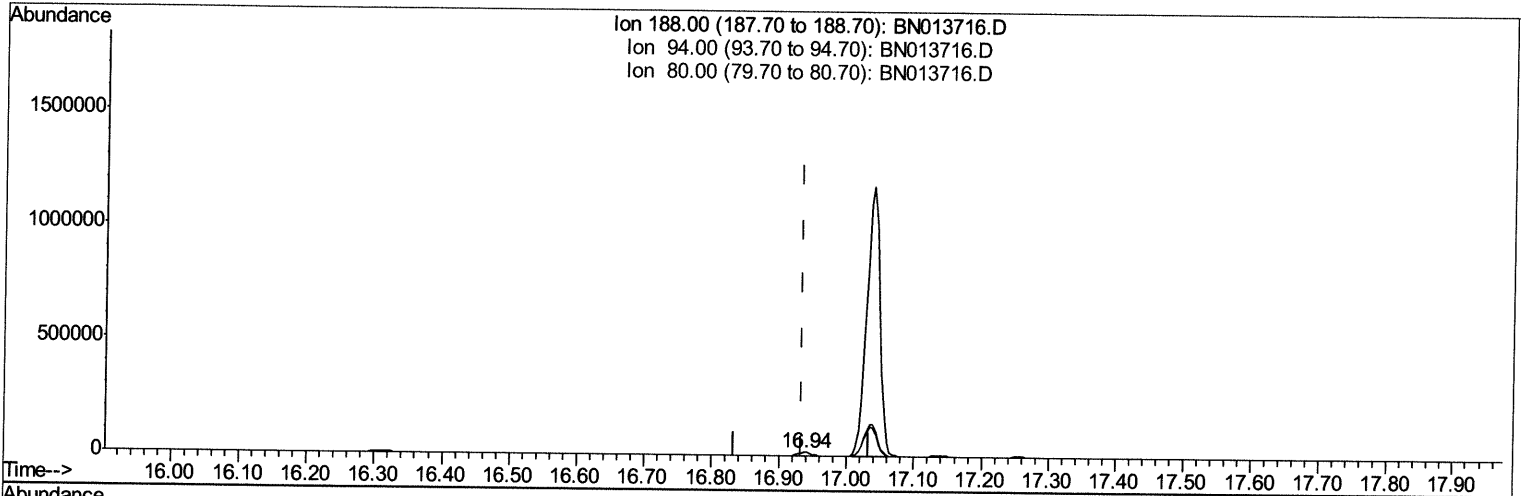
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TIC: BN013716.D

(10) Phenanthrene-d10 (I)

16.937min (+0.004) 0.40ng/ul m

response 19276

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 188.00 | 100   | 100    |
| 94.00  | 10.70 | 18.42# |
| 80.00  | 11.60 | 17.33# |
| 0.00   | 0.00  | 0.00   |

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 ALS Vial : 25 Sample Multiplier: 1

Instrument :  
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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 | 7.54  | 152  | 4222     | 0.40 | ng/ul | 0.00     |
| 2) Naphthalene-d8         | 10.31 | 136  | 16240    | 0.40 | ng/ul | 0.00     |
| 6) Acenaphthene-d10       | 14.19 | 164  | 11530    | 0.40 | ng/ul | 0.01     |
| 10) Phenanthrene-d10      | 16.94 | 188  | 19276m   | 0.40 | ng/ul | 0.00     |
| 16) Chrysene-d12          | 21.14 | 240  | 16986    | 0.40 | ng/ul | 0.00     |
| 20) Perylene-d12          | 23.31 | 264  | 15810    | 0.40 | ng/ul | 0.01     |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 4) 2-Methylnaphthalene-d10  | 11.91 | 152  | 10601    | 0.45 | ng/ul | 0.00     |
| 14) Fluoranthene-d10        | 18.98 | 212  | 21565    | 0.42 | ng/ul | 0.00     |

| Target Compounds       | R.T.  | QIon | Response | Conc   | Units  | Ovalue |
|------------------------|-------|------|----------|--------|--------|--------|
| 3) Naphthalene         | 10.36 | 128  | 604355   | 14.112 | ng/ul  | 99     |
| 5) 2-Methylnaphthalene | 11.98 | 142  | 29385    | 1.009  | ng/ul  | 98     |
| 7) Acenaphthylene      | 13.91 | 152  | 1523     | 0.034  | ng/ul# | 1      |
| 8) Acenaphthene        | 14.26 | 153  | 7093     | 0.183  | ng/ul  | 98     |
| 9) Fluorene            | 15.25 | 166  | 7472     | 0.167  | ng/ul# | 95     |
| 11) Pentachlorophenol  | 16.60 | 266  | 136      | 0.030  | ng/ul  | 92     |
| 12) Phenanthrene       | 16.98 | 178  | 3213     | 0.055  | ng/ul# | 34     |
| 13) Anthracene         | 17.07 | 178  | 1030     | 0.020  | ng/ul# | 1      |
| 15) Fluoranthene       | 19.01 | 202  | 1453     | 0.022  | ng/ul# | 2      |

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