

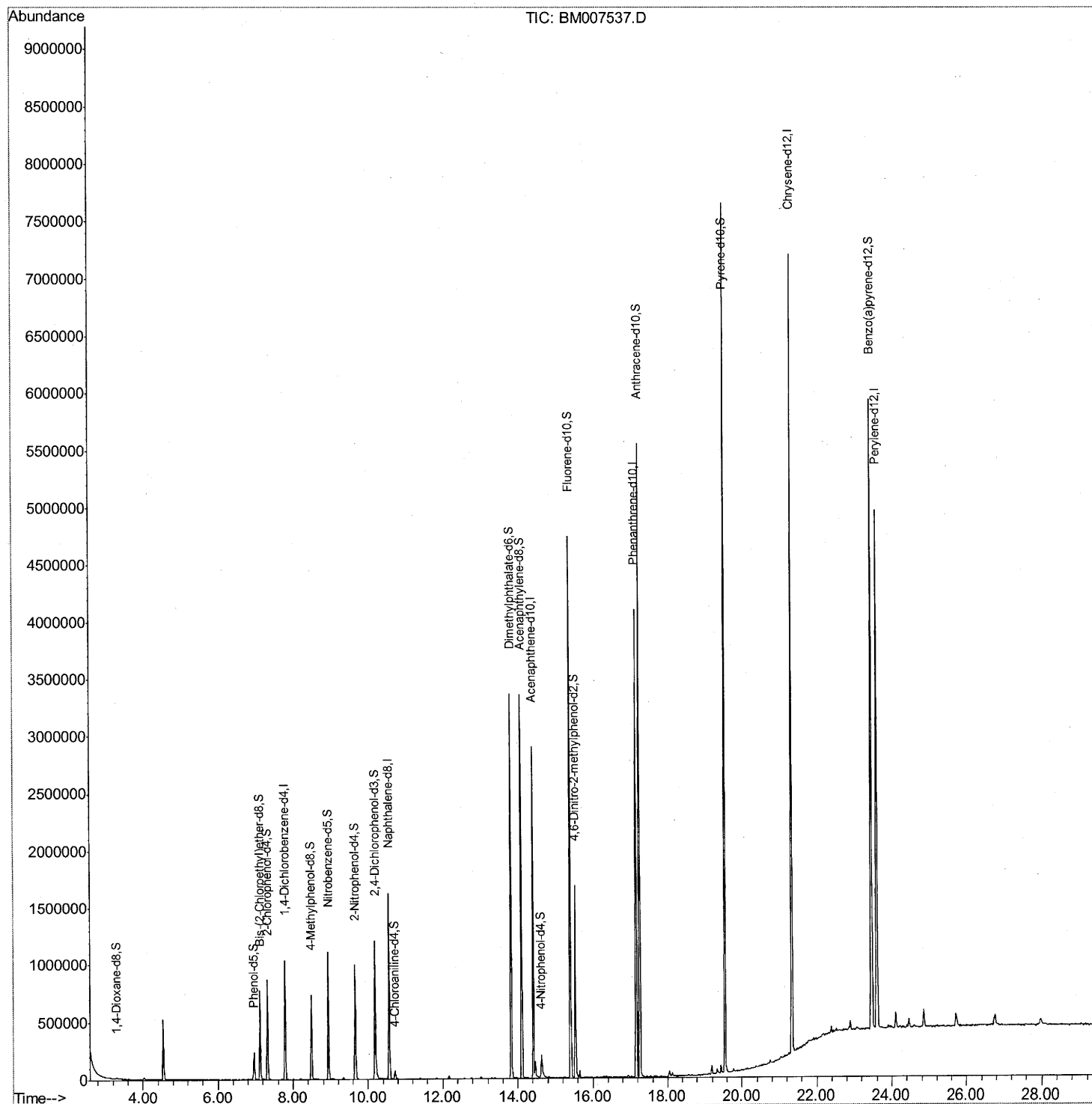
Data Path : Z:\HPCHEM1\BNA\_M\DATA\BM092316\  
Data File : BM007537.D  
Acq On : 23 Sep 2016 23:25  
Operator : UM/SJ  
Sample : H4855-07  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampled :  
H9012

Manual Integrations  
APPROVED

sohil  
9/26/2016 6:33:45 PM

Quant Time: Sep 24 02:47:11 2016  
Quant Method : Z:\HPCHEM1\BNA\_M\METHODS\SOM02.2-EPA-BM092316.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Sat Sep 24 02:24:23 2016  
Response via : Initial Calibration



# Quantitation Report (Qedit)

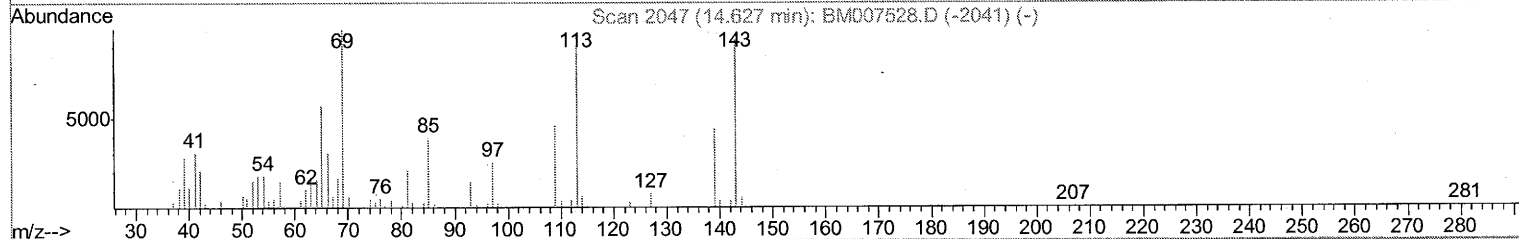
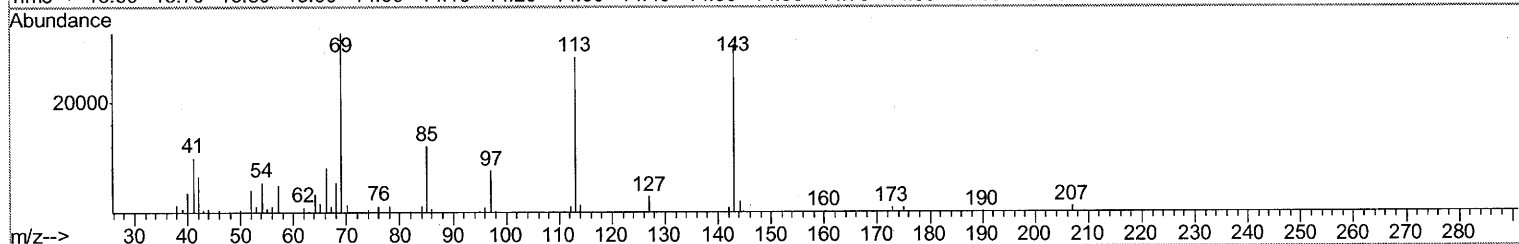
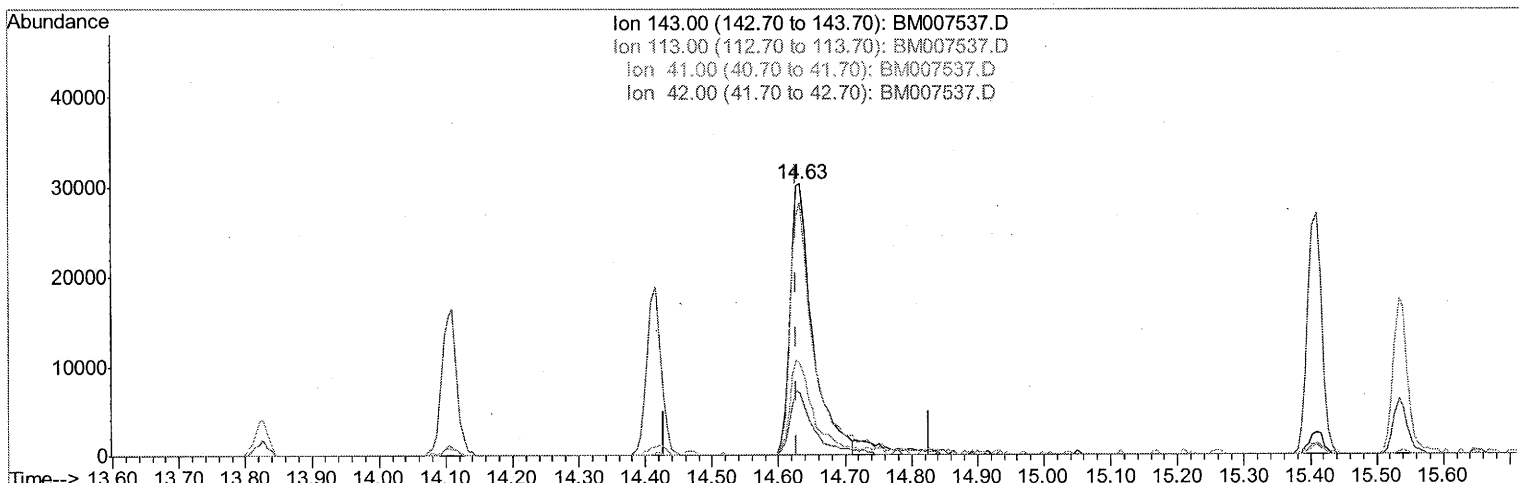
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TIC: BM007537.D

(51) 4-Nitrophenol-d4 (S)

14.633min (+0.006) 5.83ng/ul

response 70641

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 143.00 | 100   | 100   |
| 113.00 | 93.70 | 93.02 |
| 41.00  | 36.40 | 33.46 |
| 42.00  | 22.60 | 22.54 |

# Quantitation Report (Qedit)

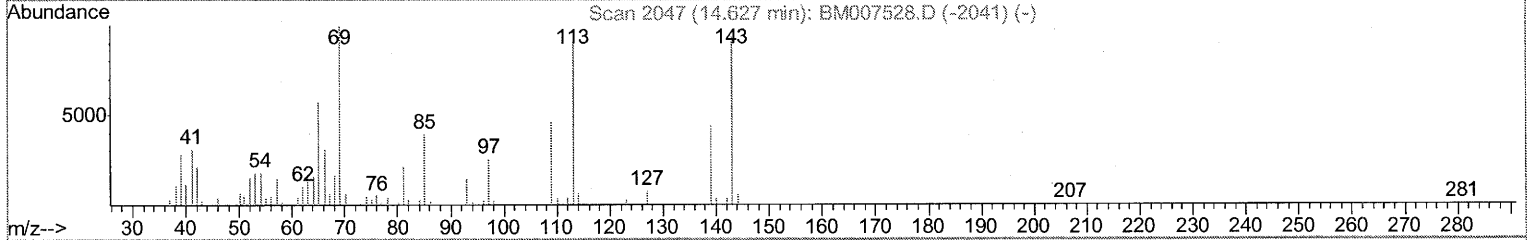
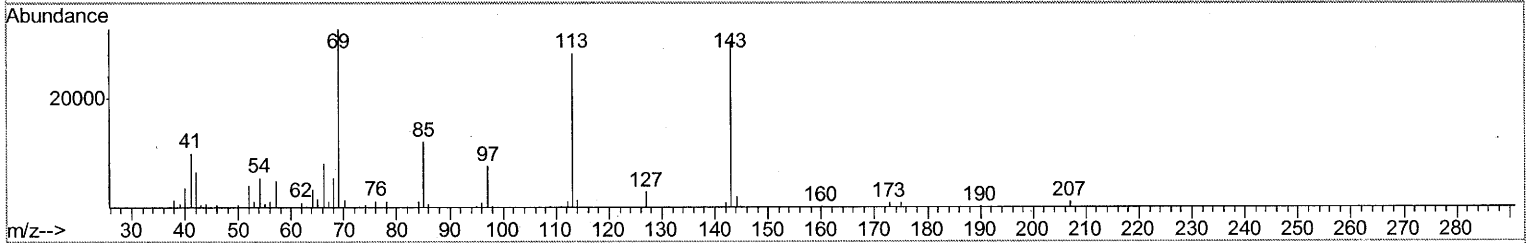
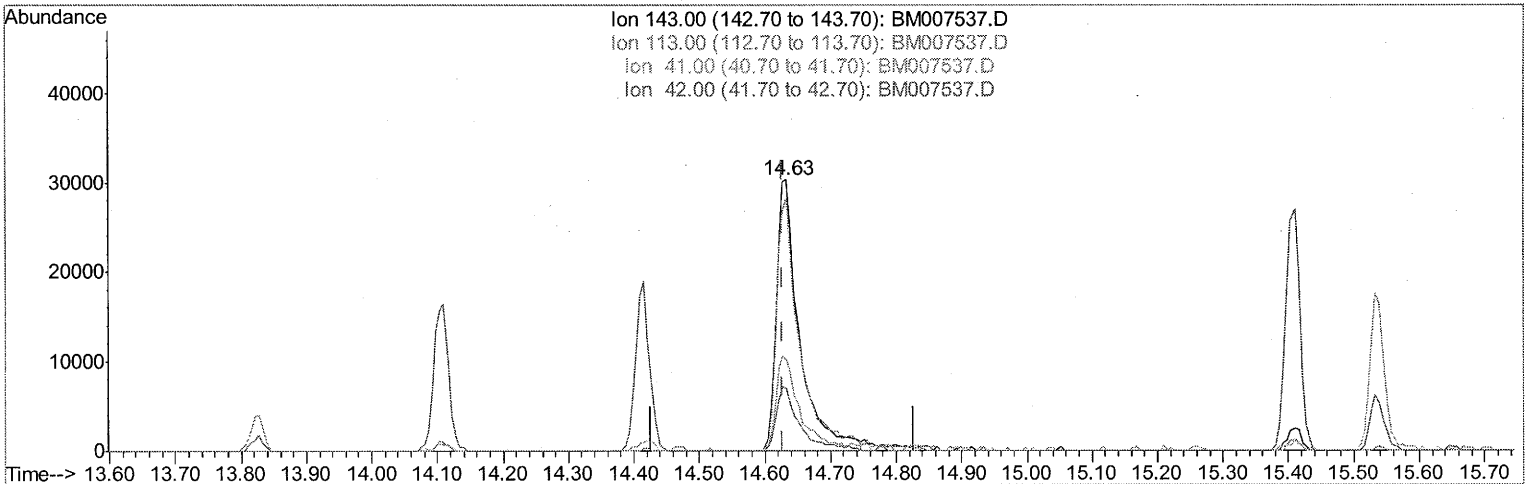
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 APPROVED

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TIC: BM007537.D

(51) 4-Nitrophenol-d4 (S)

14.633min (+0.006) 6.07ng/ul m

response 73565

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 143.00 | 100   | 100   |
| 113.00 | 93.70 | 93.02 |
| 41.00  | 36.40 | 33.46 |
| 42.00  | 22.60 | 22.54 |

UM

09/26/16

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 ClientSampled :  
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Manual Integrations  
 APPROVED

sohil  
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Quant Time: Sep 24 02:47:11 2016  
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 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.78  | 152  | 277712   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.56 | 136  | 1319628  | 20.00 | ng/ul | 0.00     |
| 35) Acenaphthene-d10      | 14.42 | 164  | 944013   | 20.00 | ng/ul | 0.00     |
| 61) Phenanthrene-d10      | 17.16 | 188  | 2471505  | 20.00 | ng/ul | 0.00     |
| 75) Chrysene-d12          | 21.34 | 240  | 3453563  | 20.00 | ng/ul | 0.00     |
| 83) Perylene-d12          | 23.61 | 264  | 3470205  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |         |       |       |      |
|--------------------------------|-------|-----|---------|-------|-------|------|
| 3) 1,4-Dioxane-d8              | 3.30  | 96  | 7174    | 1.27  | ng/uL | 0.00 |
| 5) Phenol-d5                   | 6.96  | 99  | 138175  | 5.61  | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 7.12  | 67  | 350723  | 26.46 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.32  | 132 | 381524  | 21.31 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.49  | 113 | 274909  | 13.07 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.94  | 128 | 257471  | 27.29 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.66  | 143 | 304722  | 28.43 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 10.19 | 165 | 496688  | 23.26 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.72 | 131 | 43715   | 1.95  | ng/ul | 0.01 |
| 43) Dimethylphthalate-d6       | 13.83 | 166 | 2158605 | 31.44 | ng/ul | 0.00 |
| 46) Acenaphthylene-d8          | 14.10 | 160 | 2414073 | 30.30 | ng/ul | 0.00 |
| 51) 4-Nitrophenol-d4           | 14.63 | 143 | 73565m  | 6.07  | ng/ul | 0.00 |
| 57) Fluorene-d10               | 15.41 | 176 | 1861685 | 30.91 | ng/ul | 0.00 |
| 62) 4,6-Dinitro-2-methylphenol | 15.54 | 200 | 400569  | 26.56 | ng/ul | 0.00 |
| 70) Anthracene-d10             | 17.26 | 188 | 3149460 | 27.73 | ng/ul | 0.00 |
| 76) Pyrene-d10                 | 19.55 | 212 | 4024374 | 29.36 | ng/ul | 0.00 |
| 87) Benzo(a)pyrene-d12         | 23.47 | 264 | 4289259 | 27.93 | ng/ul | 0.00 |

UM  
 09/26/16

Target Compounds Qvalue

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