

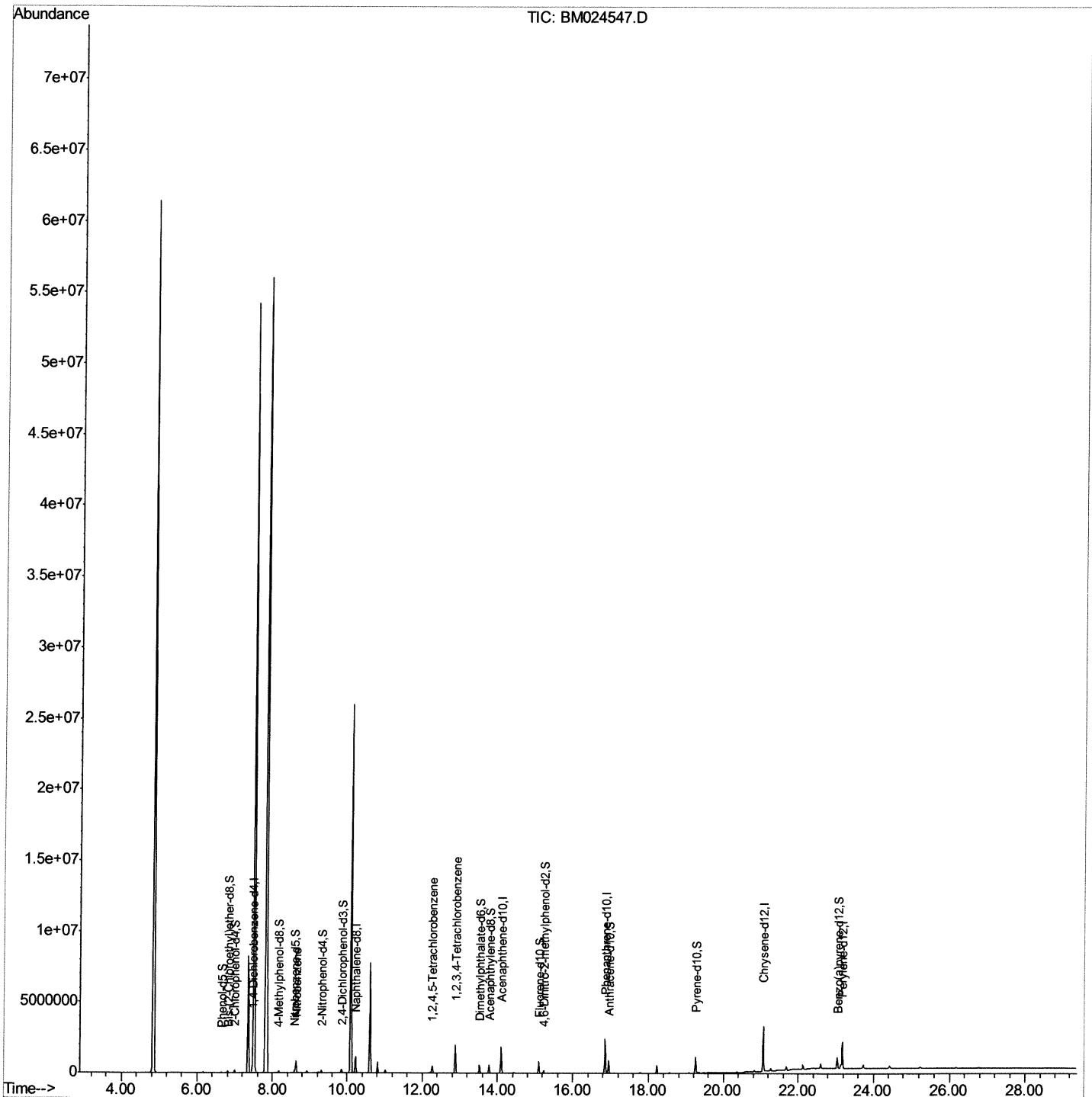
```
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012020\  
Data File : BM024547.D  
Acq On    : 20 Jan 2020  10:49  
Operator  : JU  
Sample    : L1139-05DL 5X  
Misc      :  
ALS Vial  : 4      Sample Multiplier: 1
```

**Instrument :**  
BNA\_M  
**ClientSampleId :**  
C0D51DL

## Manual Integrations

mohammad  
1/21/2020 2:54:05 PM

Quant Time: Jan 20 15:58:19 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jan 20 15:42:34 2020  
Response via : Initial Calibration



## Quantitation Report (Qedit)

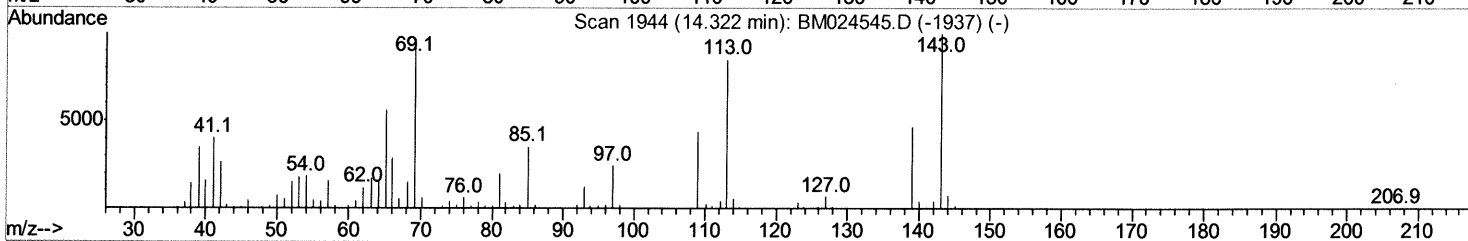
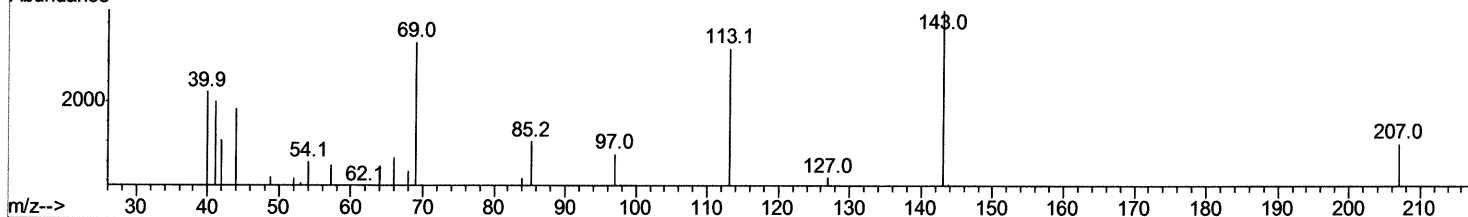
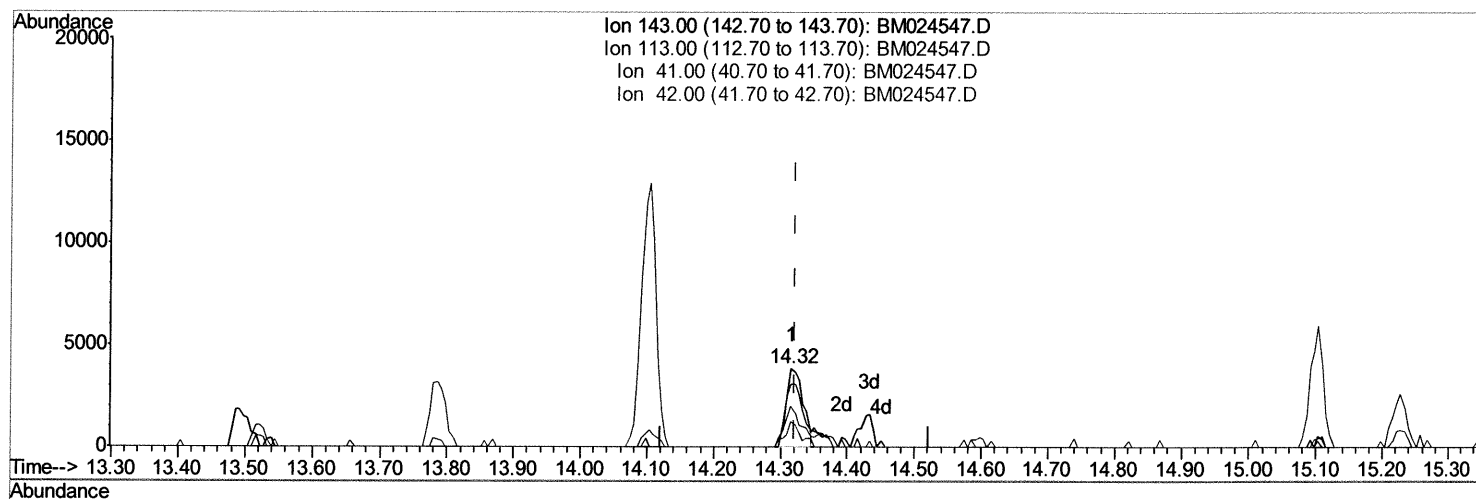
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012020\  
Data File : BM024547.D  
Acq On : 20 Jan 2020 10:49  
Operator : JU  
Sample : L1139-05DL 5X  
Misc :  
ALS Vial : 4 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
C0D51DL

Manual Integrations  
APPROVED

mohammad  
1/21/2020 2:54:05 PM

Quant Time: Jan 20 15:49:23 2020  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
Quant Title : SVOA CALIBRATION  
QLast Update : Mon Jan 20 15:42:34 2020  
Response via : Initial Calibration



TIC: BM024547.D

(52) 4-Nitrophenol-d4 (S)

14.316min (-0.006) 0.81ng/ul

response 6682

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 143.00 | 100   | 100   |
| 113.00 | 90.90 | 80.01 |
| 41.00  | 43.40 | 52.00 |
| 42.00  | 30.10 | 31.74 |

# Quantitation Report (Qedit)

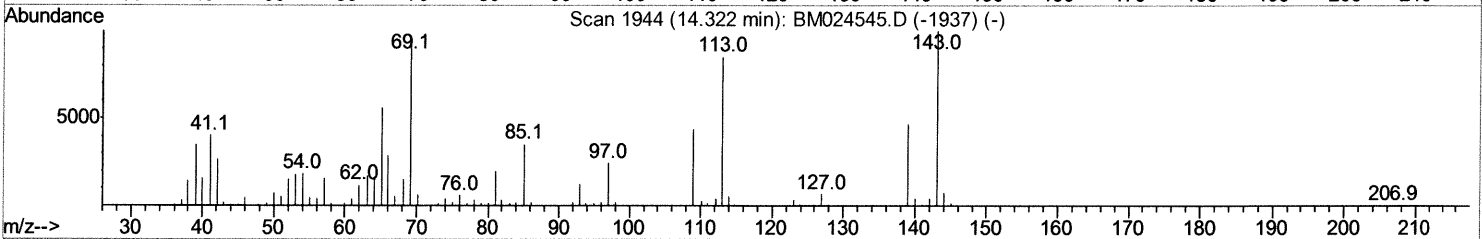
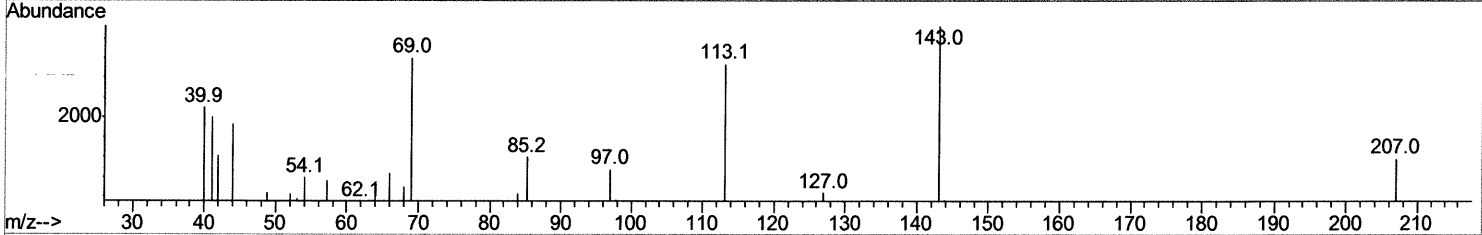
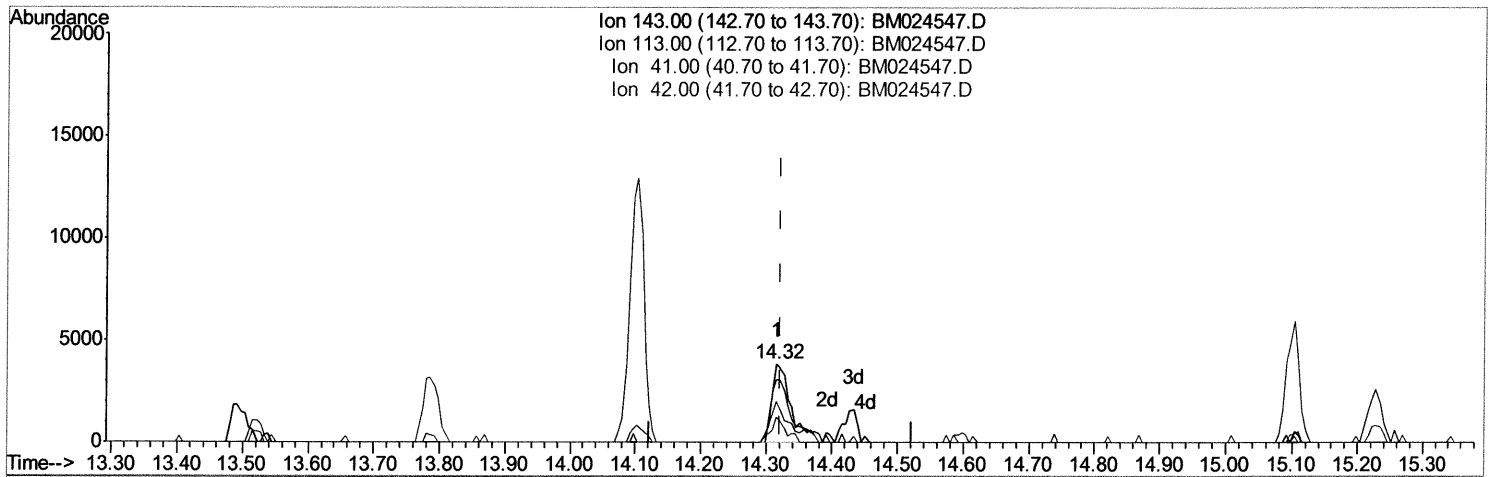
Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012020\  
 Data File : BM024547.D  
 Acq On : 20 Jan 2020 10:49  
 Operator : JU  
 Sample : L1139-05DL 5X  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0D51DL

Manual Integrations  
 APPROVED

mohammad  
 1/21/2020 2:54:05 PM

Quant Time: Jan 20 15:49:23 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 20 15:42:34 2020  
 Response via : Initial Calibration



TIC: BM024547.D

(52) 4-Nitrophenol-d4 (S)

14.316min (-0.006) 0.94ng/ul m> 57 1/21/20

response 7748

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 143.00 | 100   | 100   |
| 113.00 | 90.90 | 80.01 |
| 41.00  | 43.40 | 52.00 |
| 42.00  | 30.10 | 31.74 |

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM012020\  
 Data File : BM024547.D  
 Acq On : 20 Jan 2020 10:49  
 Operator : JU  
 Sample : L1139-05DL 5X  
 Misc :  
 ALS Vial : 4 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 C0D51DL

Manual Integrations  
 APPROVED

mohammad  
 1/21/2020 2:54:05 PM

Quant Time: Jan 20 15:58:19 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM011520MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jan 20 15:42:34 2020  
 Response via : Initial Calibration

| Internal Standards        | R.T.  | QIon | Response | Conc  | Units | Dev(Min) |
|---------------------------|-------|------|----------|-------|-------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.46  | 152  | 217679   | 20.00 | ng/ul | 0.00     |
| 18) Naphthalene-d8        | 10.22 | 136  | 884424   | 20.00 | ng/ul | 0.00     |
| 36) Acenaphthene-d10      | 14.10 | 164  | 627127   | 20.00 | ng/ul | 0.00     |
| 62) Phenanthrene-d10      | 16.85 | 188  | 1457784  | 20.00 | ng/ul | 0.00     |
| 78) Chrysene-d12          | 21.06 | 240  | 1436602  | 20.00 | ng/ul | 0.00     |
| 86) Perylene-d12          | 23.18 | 264  | 1392044  | 20.00 | ng/ul | 0.00     |

## System Monitoring Compounds

|                                |       |     |        |      |       |      |
|--------------------------------|-------|-----|--------|------|-------|------|
| 3) 1,4-Dioxane-d8              | 0.00  | 96  | 0d     | 0.00 | ng/uL |      |
| 5) Phenol-d5                   | 6.65  | 99  | 18552  | 1.15 | ng/ul | 0.00 |
| 7) Bis-(2-Chloroethyl)ether-d  | 6.80  | 67  | 54157  | 6.04 | ng/ul | 0.00 |
| 9) 2-Chlorophenol-d4           | 7.00  | 132 | 69286  | 5.03 | ng/ul | 0.00 |
| 13) 4-Methylphenol-d8          | 8.16  | 113 | 40922  | 2.75 | ng/ul | 0.00 |
| 19) Nitrobenzene-d5            | 8.59  | 128 | 45286  | 7.29 | ng/ul | 0.00 |
| 22) 2-Nitrophenol-d4           | 9.31  | 143 | 49602  | 7.37 | ng/ul | 0.00 |
| 26) 2,4-Dichlorophenol-d3      | 9.85  | 165 | 86260  | 5.69 | ng/ul | 0.00 |
| 29) 4-Chloroaniline-d4         | 10.36 | 131 | 5527   | 0.33 | ng/ul | 0.00 |
| 44) Dimethylphthalate-d6       | 13.52 | 166 | 364186 | 7.41 | ng/ul | 0.00 |
| 47) Acenaphthylene-d8          | 13.79 | 160 | 409143 | 7.23 | ng/ul | 0.00 |
| 52) 4-Nitrophenol-d4           | 14.32 | 143 | 7748m  | 0.94 | ng/ul | 0.00 |
| 58) Fluorene-d10               | 15.10 | 176 | 314136 | 7.26 | ng/ul | 0.00 |
| 63) 4,6-Dinitro-2-methylphenol | 15.23 | 200 | 42076  | 5.06 | ng/ul | 0.00 |
| 71) Anthracene-d10             | 16.95 | 188 | 513241 | 7.57 | ng/ul | 0.00 |
| 79) Pyrene-d10                 | 19.26 | 212 | 612737 | 7.98 | ng/ul | 0.00 |
| 90) Benzo(a)pyrene-d12         | 23.04 | 264 | 542141 | 7.51 | ng/ul | 0.00 |

SD 1/21/20

## Target Compounds

|                                |       |     |        | Ovalue |          |
|--------------------------------|-------|-----|--------|--------|----------|
| 20) Nitrobenzene               | 8.63  | 77  | 418491 | 27.827 | ng/ul 96 |
| 37) 1,2,4,5-Tetrachlorobenzene | 12.27 | 216 | 168633 | 8.173  | ng/ul 99 |
| 73) 1,2,3,4-Tetrachlorobenzene | 12.89 | 216 | 611068 | 29.478 | ng/uL 99 |

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