

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP100520\  
 Data File : BP003518.D  
 Acq On : 06 Oct 2020 09:16  
 Operator : CG/JU  
 Sample : L4190-18  
 Misc :  
 ALS Vial : 30 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 PAV-B18-092920-DP

Quant Time: Oct 06 10:46:13 2020  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP091520.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Oct 05 14:33:14 2020  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min) |
|-----------------------------|--------|------|----------|--------|-------|----------|
| -----                       |        |      |          |        |       |          |
| Internal Standards          |        |      |          |        |       |          |
| 1) 1,4-Dichlorobenzene-d4   | 7.499  | 152  | 92536    | 20.00  | ng    | # 0.00   |
| 21) Naphthalene-d8          | 10.275 | 136  | 367706   | 20.00  | ng    | # 0.00   |
| 39) Acenaphthene-d10        | 14.157 | 164  | 225722   | 20.00  | ng    | 0.00     |
| 64) Phenanthrene-d10        | 16.916 | 188  | 510788   | 20.00  | ng    | # 0.00   |
| 76) Chrysene-d12            | 21.022 | 240  | 538796   | 20.00  | ng    | 0.00     |
| 86) Perylene-d12            | 23.198 | 264  | 487358   | 20.00  | ng    | 0.00     |
| System Monitoring Compounds |        |      |          |        |       |          |
| 5) 2-Fluorophenol           | 5.117  | 112  | 736547   | 138.99 | ng    | 0.00     |
| 7) Phenol-d6                | 6.711  | 99   | 849276   | 120.23 | ng    | 0.00     |
| 23) Nitrobenzene-d5         | 8.652  | 82   | 700013   | 111.84 | ng    | 0.00     |
| 42) 2,4,6-Tribromophenol    | 15.663 | 330  | 400001   | 116.19 | ng    | 0.00     |
| 45) 2-Fluorobiphenyl        | 12.775 | 172  | 1478455  | 95.79  | ng    | 0.00     |
| 79) Terphenyl-d14           | 19.504 | 244  | 2383345  | 92.17  | ng    | 0.00     |

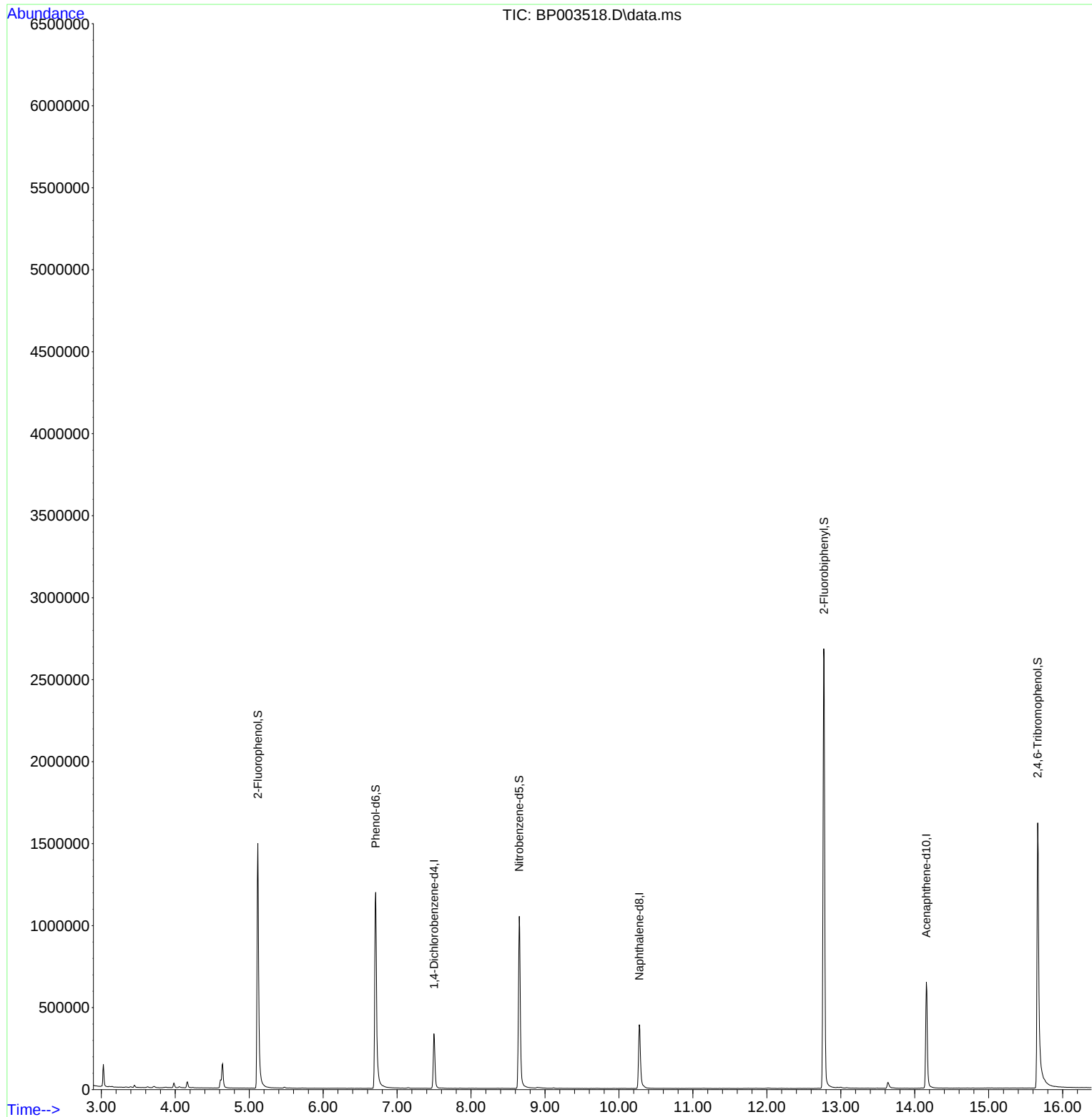
|                  |        |
|------------------|--------|
| Target Compounds | Qvalue |
| -----            |        |

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

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