

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061125\
 Data File : BM050295.D
 Acq On : 12 Jun 2025 05:10
 Operator : RC/JU
 Sample : PB168369TB
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168369TB

Quant Time: Jun 12 05:58:31 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 09 11:54:38 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.763	152	374014	20.000	ng	0.00
21) Naphthalene-d8	10.551	136	1396895	20.000	ng	0.00
39) Acenaphthene-d10	14.392	164	776202	20.000	ng	0.00
64) Phenanthrene-d10	17.127	188	1516518	20.000	ng	-0.01
76) Chrysene-d12	21.356	240	1557464	20.000	ng	-0.02
86) Perylene-d12	24.327	264	1641162	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	3376105	150.578	ng	0.00
7) Phenol-d6	6.934	99	4131453	139.977	ng	0.00
23) Nitrobenzene-d5	8.910	82	2436953	90.731	ng	0.00
42) 2,4,6-Tribromophenol	15.874	330	1349043	152.801	ng	0.00
45) 2-Fluorobiphenyl	13.015	172	5195965	90.628	ng	-0.01
79) Terphenyl-d14	19.756	244	7302911	88.857	ng	-0.01

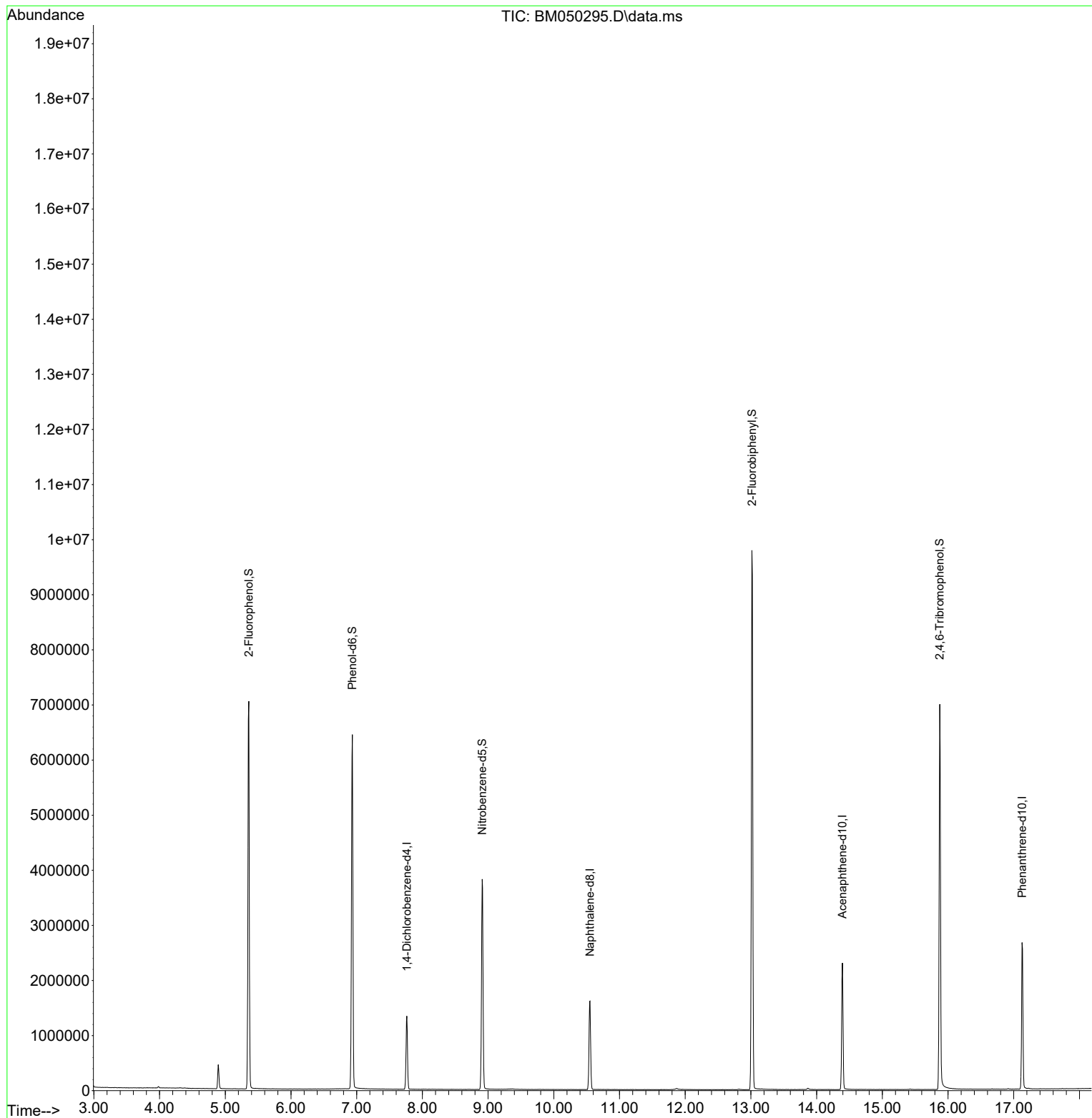
Target Compounds	Qvalue

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

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