

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
 Data File : BM050430.D
 Acq On : 14 Jul 2025 17:34
 Operator : RC/JU
 Sample : PB168788BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB168788BL

Quant Time: Jul 14 18:42:25 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 18:32:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	390165	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1399531	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	845767	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	1660365	20.000	ng	0.00
76) Chrysene-d12	21.439	240	1715059	20.000	ng	-0.02
86) Perylene-d12	24.474	264	1805974	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3420172	150.890	ng	0.00
7) Phenol-d6	7.022	99	4207432	147.248	ng	0.00
23) Nitrobenzene-d5	9.016	82	2410969	87.889	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1583921	154.136	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6032236	88.079	ng	0.00
79) Terphenyl-d14	19.833	244	10134767	103.971	ng	-0.01

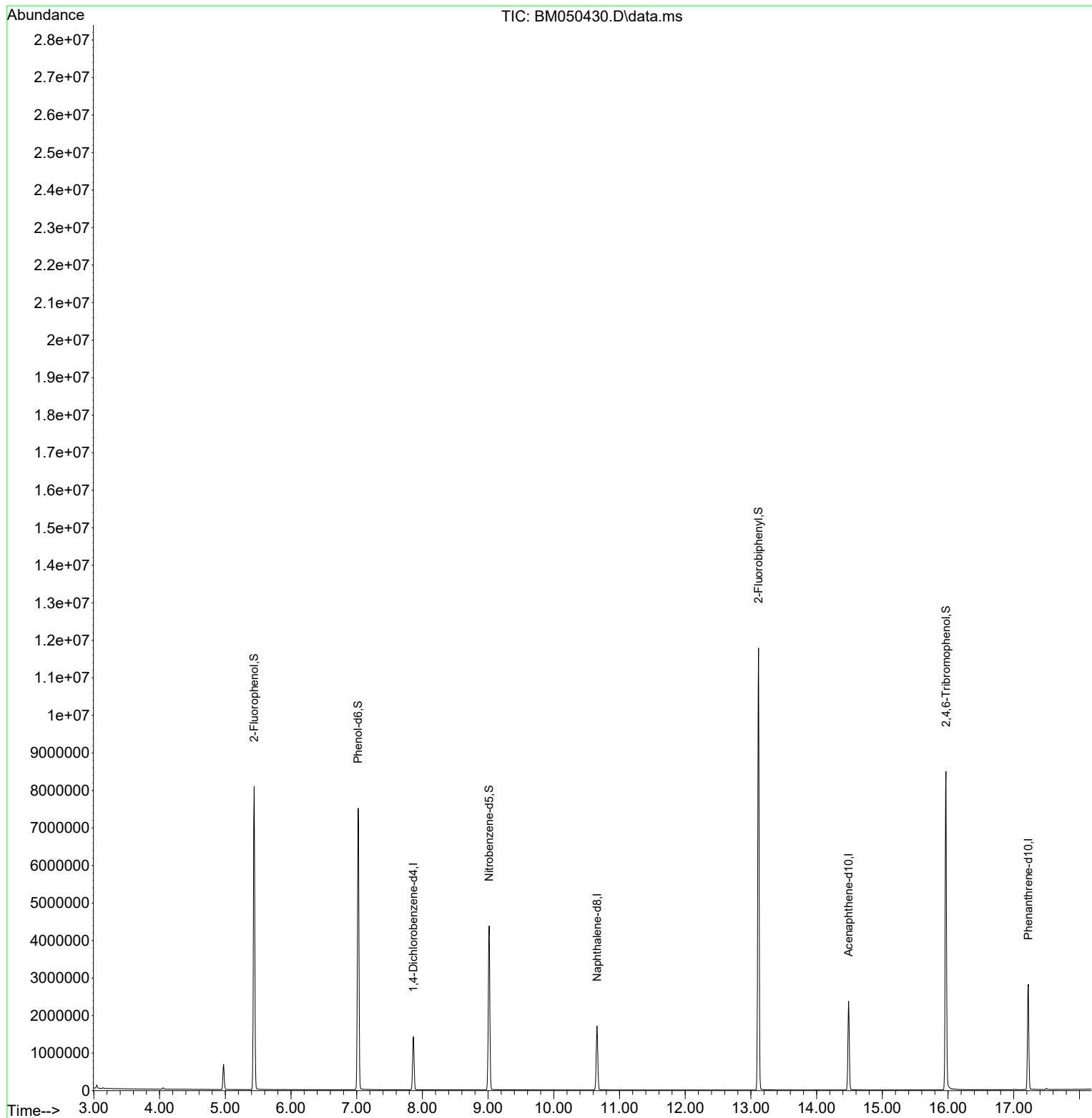
Target Compounds	Qvalue

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

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