

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071425\
Data File : BM050432.D
Acq On : 14 Jul 2025 18:54
Operator : RC/JU
Sample : PB168820BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
PB168820BL

Quant Time: Jul 15 02:02:22 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	365994	20.000	ng	0.00
21) Naphthalene-d8	10.657	136	1297145	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	778529	20.000	ng	0.00
64) Phenanthrene-d10	17.221	188	1483211	20.000	ng	0.00
76) Chrysene-d12	21.439	240	1497440	20.000	ng	-0.02
86) Perylene-d12	24.468	264	1585798	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3429626	161.300	ng	0.00
7) Phenol-d6	7.022	99	4208201	157.001	ng	0.00
23) Nitrobenzene-d5	9.016	82	2423333	95.312	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	1554980	164.388	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	6014177	95.399	ng	0.00
79) Terphenyl-d14	19.833	244	9795135	115.091	ng	-0.01

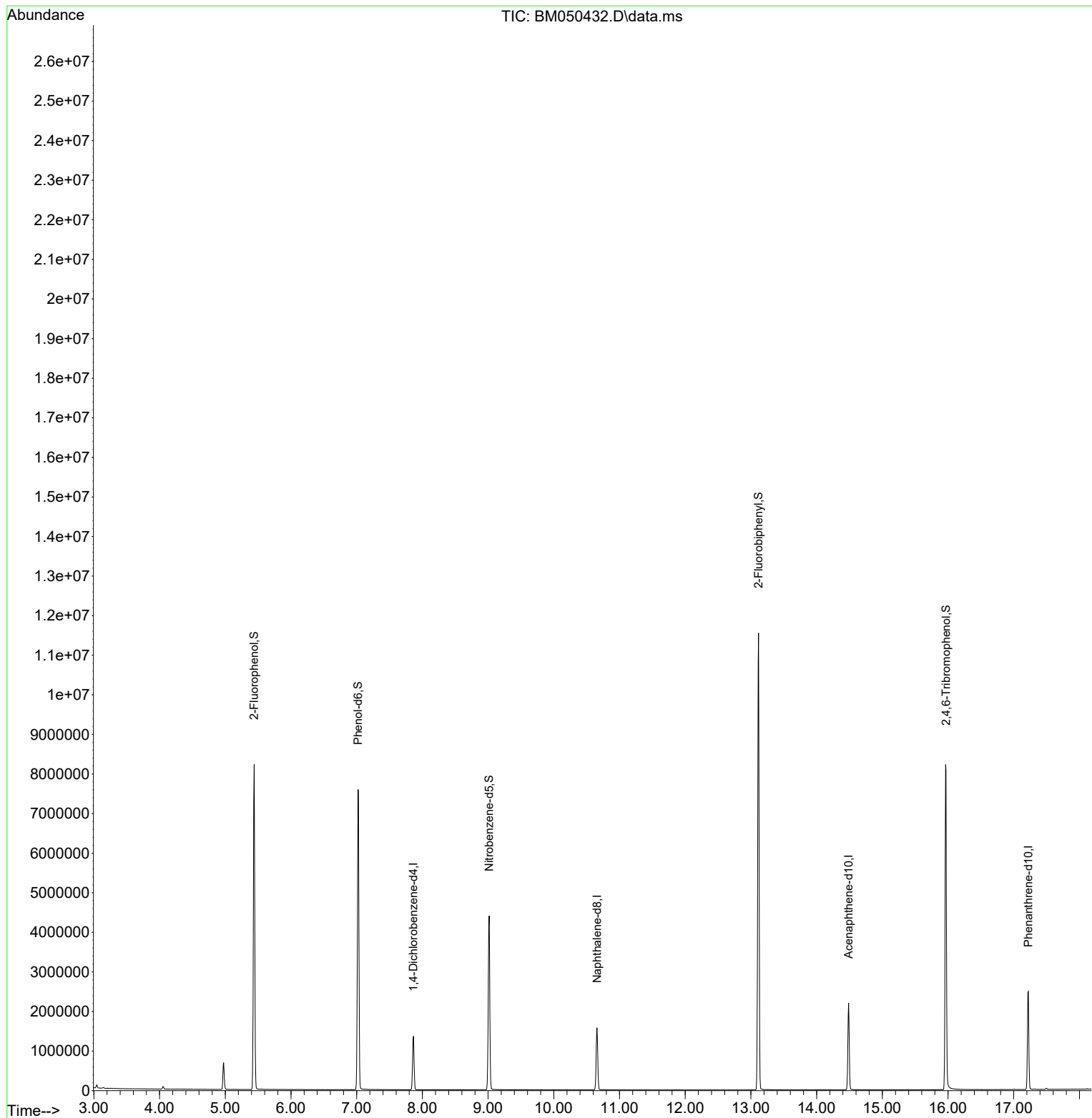
Target Compounds	Qvalue

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

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