

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082624\
 Data File : BM047353.D
 Acq On : 26 Aug 2024 19:03
 Operator : RC/JU
 Sample : PB162917BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162917BL

Quant Time: Aug 27 04:29:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 26 18:11:40 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.351	152	266712	20.000	ng	0.00
21) Naphthalene-d8	10.104	136	957975	20.000	ng	0.00
39) Acenaphthene-d10	14.010	164	646074	20.000	ng	0.00
64) Phenanthrene-d10	16.774	188	1332896	20.000	ng	0.00
76) Chrysene-d12	21.015	240	930328	20.000	ng	0.00
86) Perylene-d12	23.715	264	873449	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	2114704	152.118	ng	0.00
7) Phenol-d6	6.569	99	2704541	142.812	ng	0.00
23) Nitrobenzene-d5	8.498	82	1750820	101.407	ng	0.00
42) 2,4,6-Tribromophenol	15.521	330	1172283	150.393	ng	0.00
45) 2-Fluorobiphenyl	12.616	172	4151531	98.732	ng	0.00
79) Terphenyl-d14	19.445	244	5503082	105.288	ng	0.00

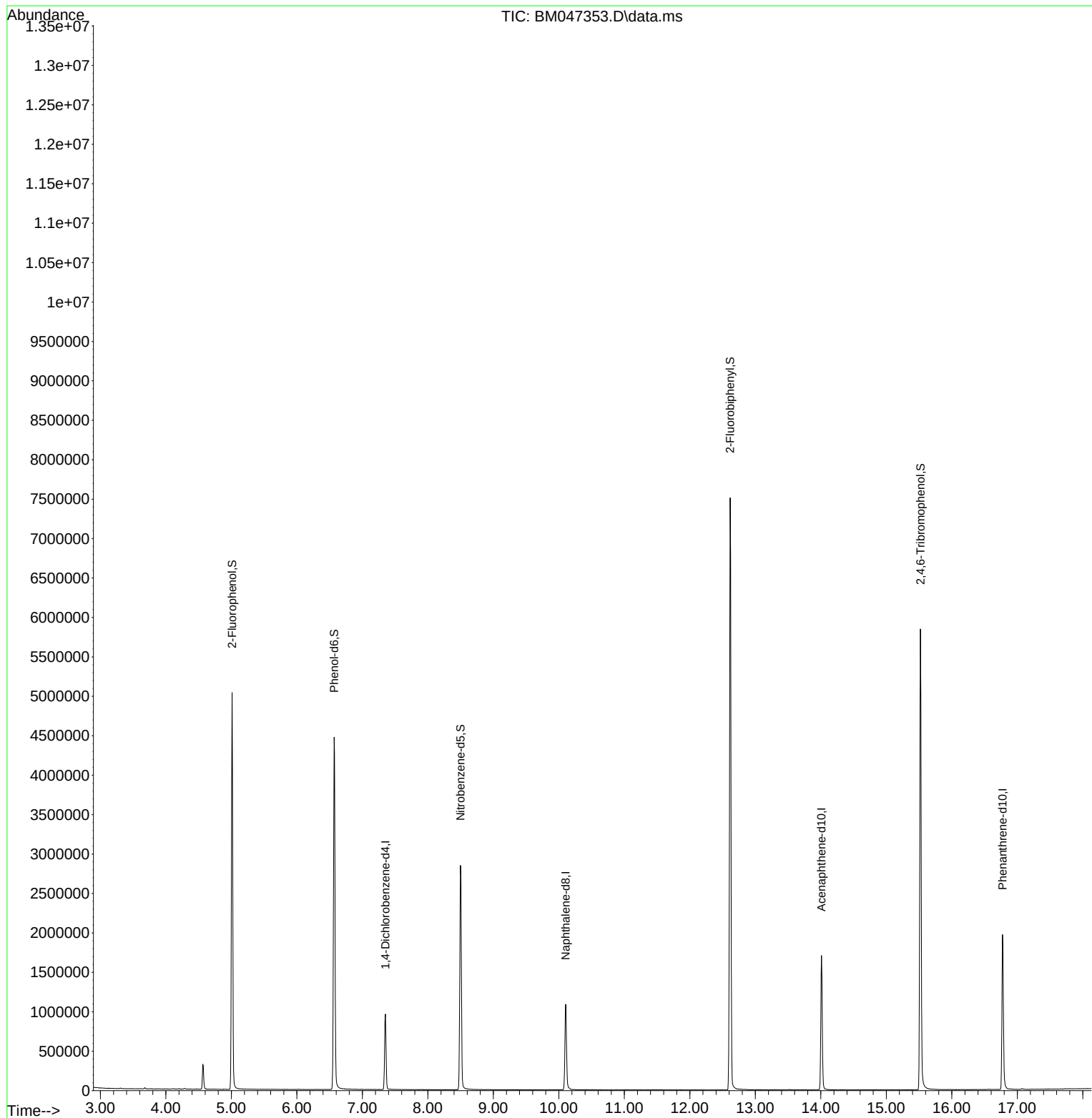
Target Compounds	Qvalue

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

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