

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090624\
 Data File : BM047463.D
 Acq On : 06 Sep 2024 19:41
 Operator : RC/JU
 Sample : PB163179BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB163179BL

Quant Time: Sep 07 01:32:13 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM090624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 07 01:23:53 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.340	152	226889	20.000	ng	0.00
21) Naphthalene-d8	10.098	136	830958	20.000	ng	0.00
39) Acenaphthene-d10	14.004	164	491342	20.000	ng	0.00
64) Phenanthrene-d10	16.786	188	902658	20.000	ng	0.01
76) Chrysene-d12	21.027	240	722678	20.000	ng	0.00
86) Perylene-d12	23.739	264	824687	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.004	112	1804166	136.999	ng	0.00
7) Phenol-d6	6.569	99	2274881	136.483	ng	0.00
23) Nitrobenzene-d5	8.492	82	1524739	95.327	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	707902	132.283	ng	0.00
45) 2-Fluorobiphenyl	12.604	172	2933621	92.406	ng	0.00
79) Terphenyl-d14	19.439	244	3402129	93.409	ng	0.00

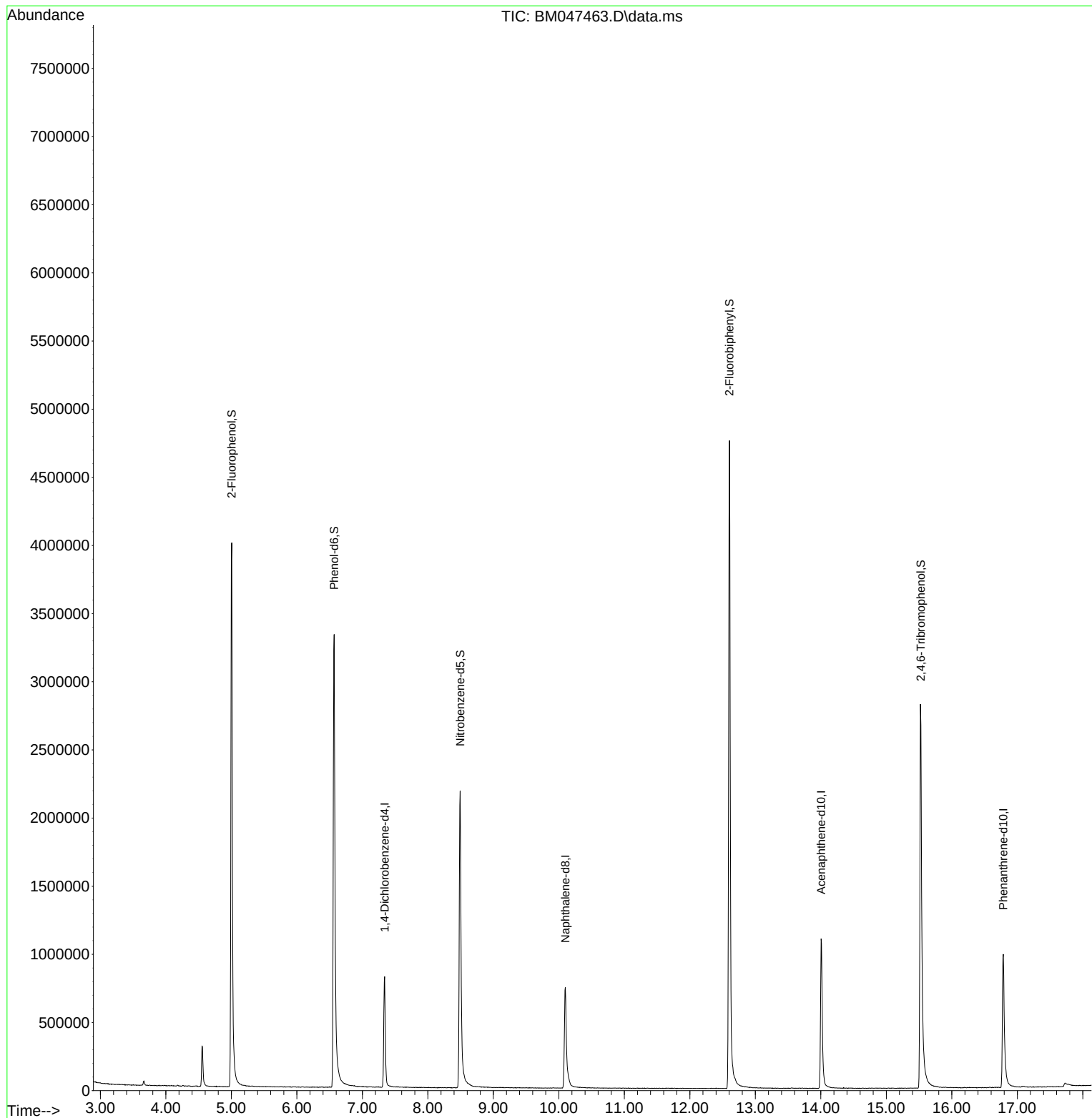
Target Compounds	Qvalue

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

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