

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091924\
 Data File : BM047575.D
 Acq On : 19 Sep 2024 17:53
 Operator : RC/JU
 Sample : PB163525BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB163525BL

Quant Time: Sep 20 02:31:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Sep 15 08:56:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	321878	20.000	ng	-0.01
21) Naphthalene-d8	10.634	136	1220647	20.000	ng	-0.01
39) Acenaphthene-d10	14.469	164	605405	20.000	ng	-0.01
64) Phenanthrene-d10	17.204	188	1062846	20.000	ng	-0.01
76) Chrysene-d12	21.427	240	741692	20.000	ng	-0.02
86) Perylene-d12	24.450	264	785495	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.416	112	3220419	159.303	ng	0.00
7) Phenol-d6	7.004	99	3991702	140.210	ng	0.00
23) Nitrobenzene-d5	8.992	82	2404920	98.601	ng	-0.01
42) 2,4,6-Tribromophenol	15.951	330	835950	142.854	ng	-0.01
45) 2-Fluorobiphenyl	13.098	172	4320386	98.305	ng	-0.01
79) Terphenyl-d14	19.821	244	5412069	130.756	ng	-0.01

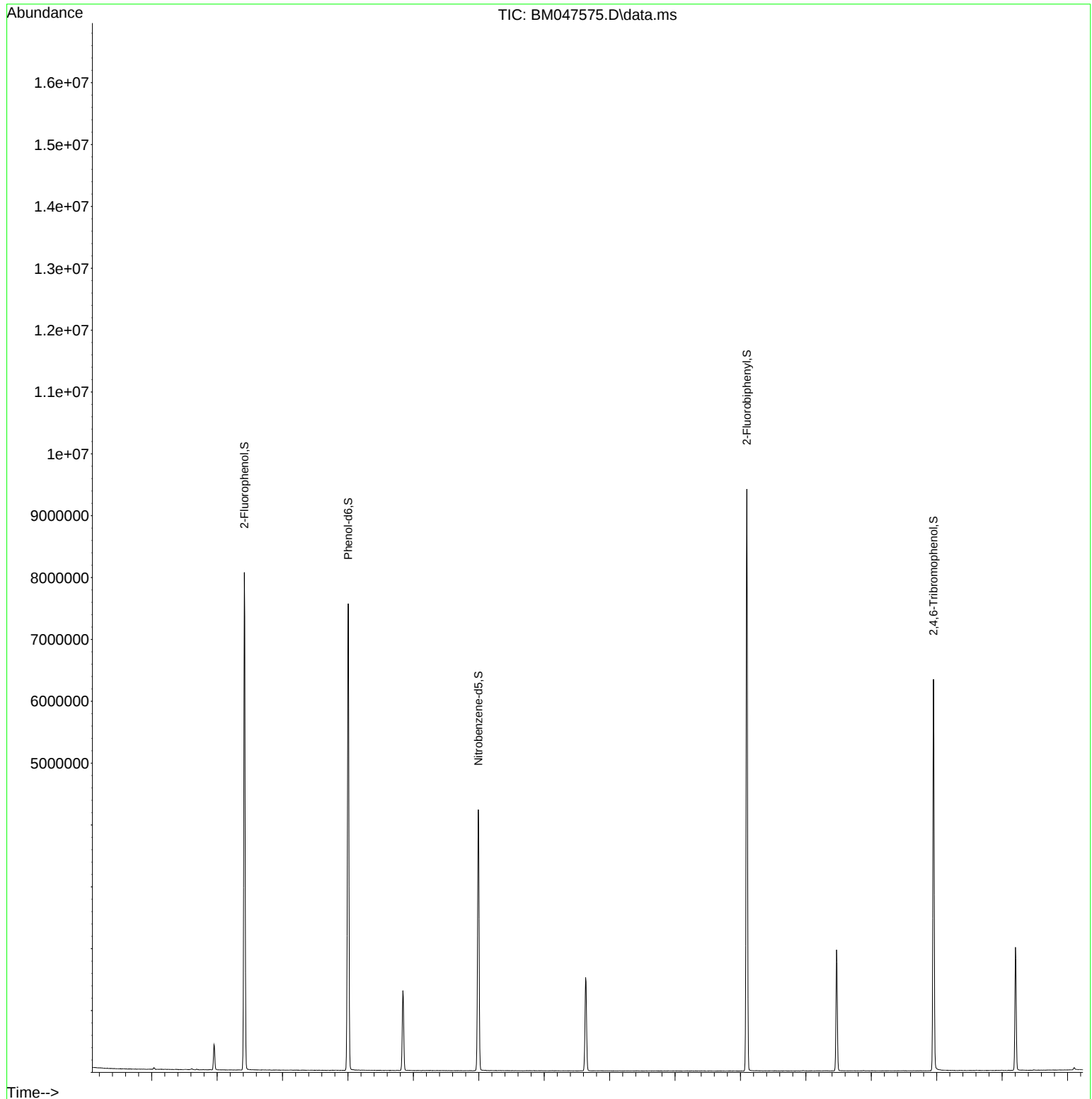
Target Compounds Qvalue

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

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