

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081524\
 Data File : BM047210.D
 Acq On : 15 Aug 2024 11:31
 Operator : RC/JU
 Sample : PB162717BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162717BL

Quant Time: Aug 15 19:47:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.363	152	275556	20.000	ng	-0.01
21) Naphthalene-d8	10.122	136	992678	20.000	ng	-0.01
39) Acenaphthene-d10	14.021	164	608749	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1179865	20.000	ng	0.00
76) Chrysene-d12	21.021	240	1082659	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1267377	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	1814978	117.692	ng	0.00
7) Phenol-d6	6.581	99	2283485	107.461	ng	0.00
23) Nitrobenzene-d5	8.510	82	1517389	76.969	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	856179	121.297	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	3261665	82.648	ng	-0.01
79) Terphenyl-d14	19.456	244	4429118	87.660	ng	0.00

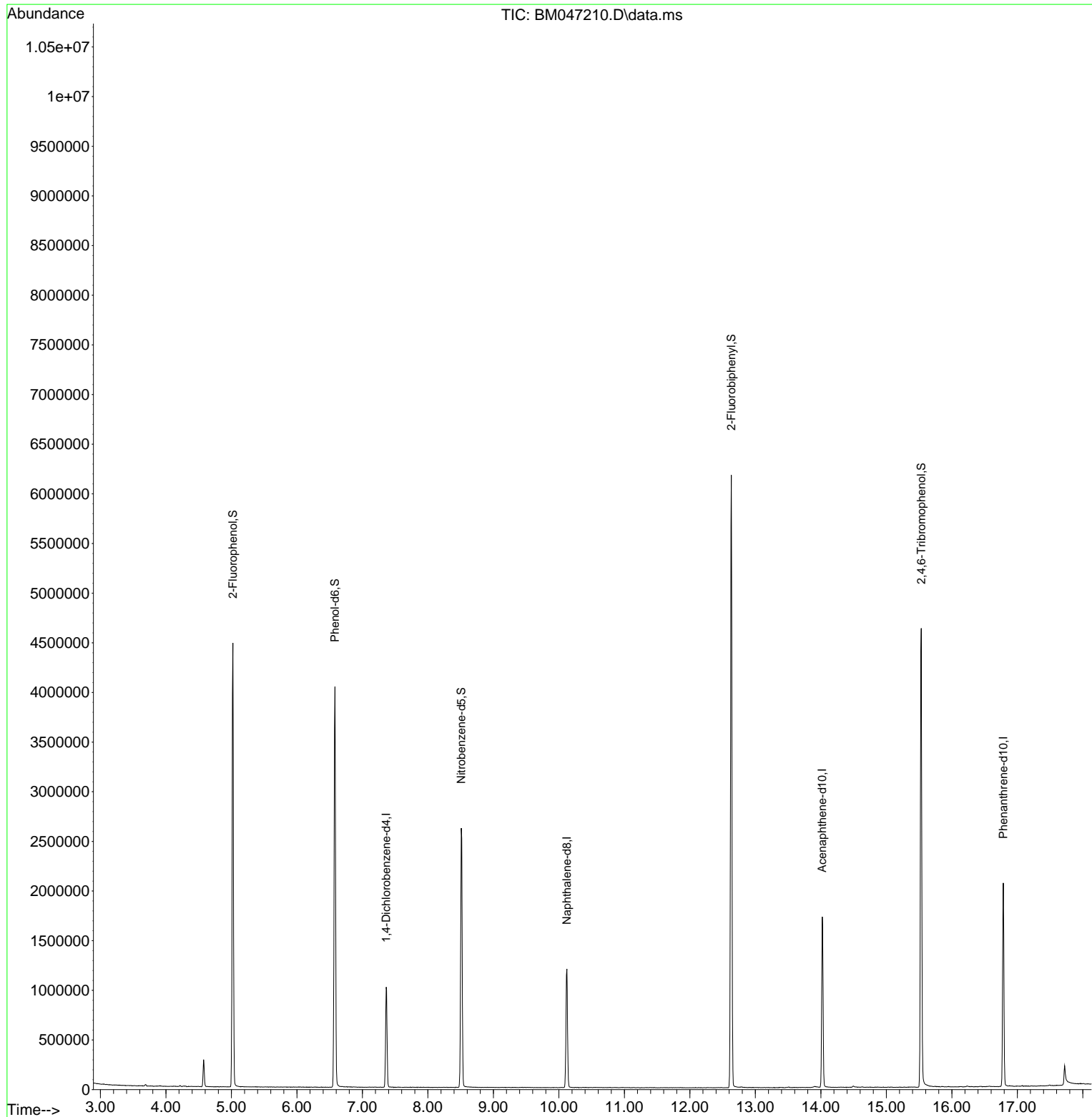
Target Compounds Qvalue

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

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