

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081624\
 Data File : BM047230.D
 Acq On : 16 Aug 2024 11:47
 Operator : RC/JU
 Sample : PB162729BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB162729BL

Quant Time: Aug 16 12:44:56 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	308565	20.000	ng	-0.01
21) Naphthalene-d8	10.116	136	1161346	20.000	ng	-0.02
39) Acenaphthene-d10	14.022	164	805377	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1638549	20.000	ng	0.00
76) Chrysene-d12	21.021	240	1162658	20.000	ng	-0.01
86) Perylene-d12	23.715	264	1176341	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	2045484	118.450	ng	0.00
7) Phenol-d6	6.581	99	2668101	112.129	ng	0.00
23) Nitrobenzene-d5	8.510	82	1755996	76.136	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	1148413	122.976	ng	0.00
45) 2-Fluorobiphenyl	12.628	172	4151407	79.511	ng	-0.02
79) Terphenyl-d14	19.451	244	5736054	105.715	ng	-0.01

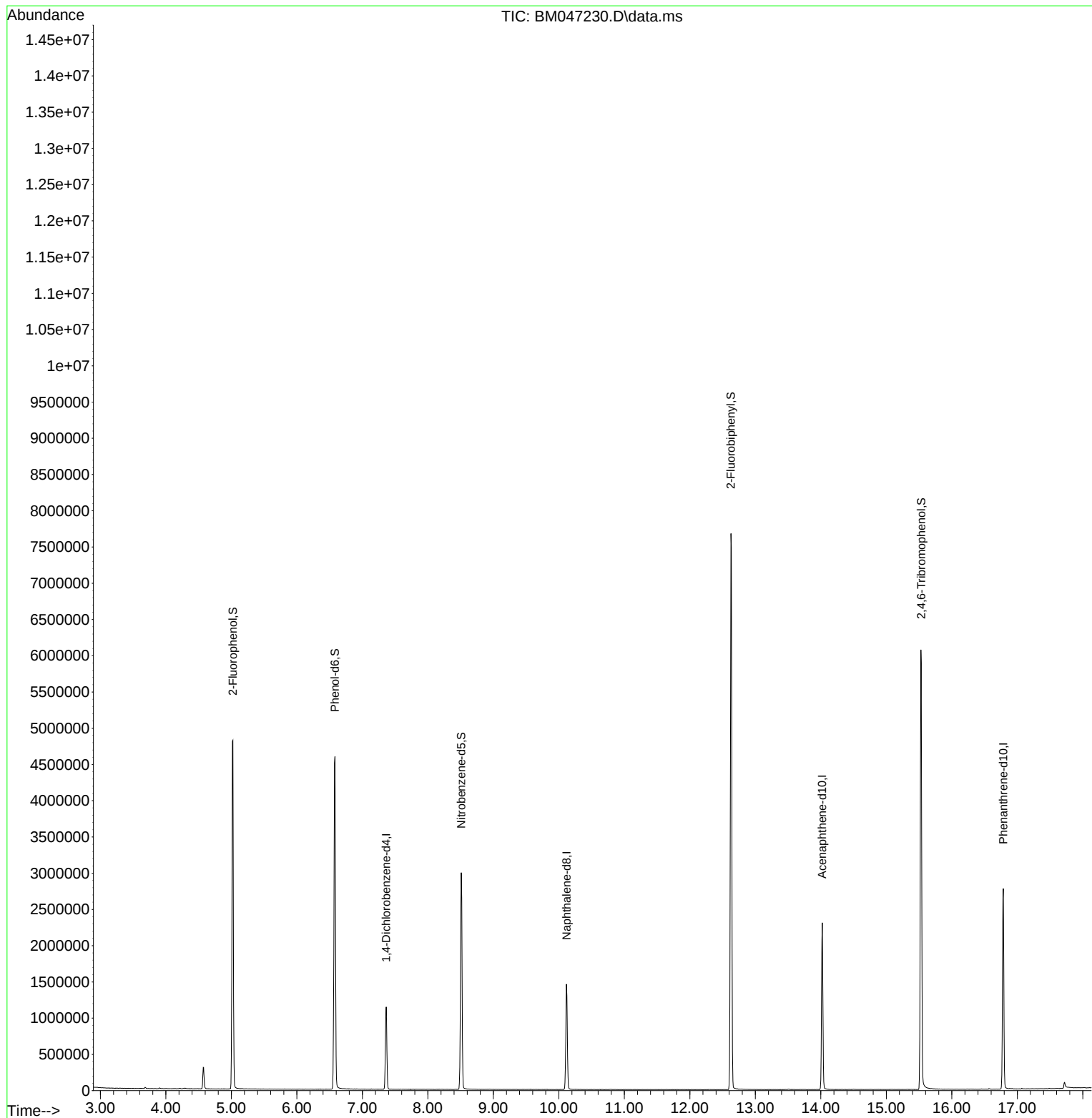
Target Compounds	Qvalue

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

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