

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103024\
 Data File : BE101417.D
 Acq On : 30 Oct 2024 10:35
 Operator : RC/JU
 Sample : PB164401BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164401BL

Quant Time: Nov 04 07:00:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE102824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 29 01:26:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	120181	20.000	ng	0.00
21) Naphthalene-d8	10.337	136	516854	20.000	ng	0.00
39) Acenaphthene-d10	14.179	164	338564	20.000	ng	0.00
64) Phenanthrene-d10	16.917	188	823184	20.000	ng	0.00
76) Chrysene-d12	21.071	240	1040671	20.000	ng	0.00
86) Perylene-d12	23.362	264	1417194	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	1063066	155.033	ng	0.00
7) Phenol-d6	6.946	99	1349771	141.962	ng	0.00
23) Nitrobenzene-d5	8.780	82	762023	92.603	ng	0.00
42) 2,4,6-Tribromophenol	15.689	330	820123	138.209	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	1948901	97.486	ng	0.00
79) Terphenyl-d14	19.514	244	3846837	85.077	ng	0.00

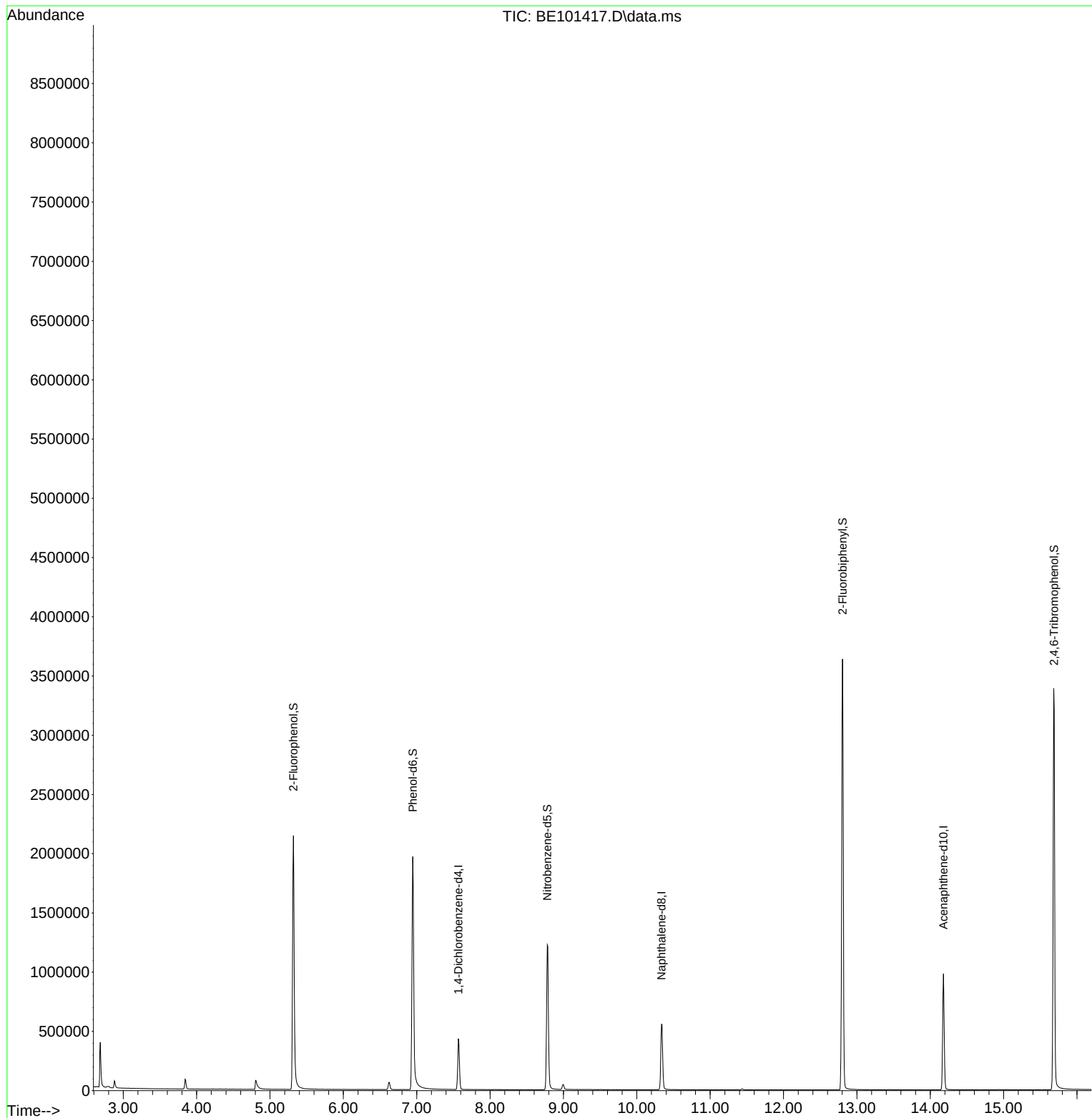
Target Compounds	Qvalue

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

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