

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE091724\
 Data File : BE101353.D
 Acq On : 17 Sep 2024 16:12
 Operator : RC/JU
 Sample : PB163307TB
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SLEB307

Quant Time: Sep 17 17:18:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE091724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 17 16:03:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.573	152	162016	20.000	ng	0.00
21) Naphthalene-d8	10.346	136	685347	20.000	ng	0.00
39) Acenaphthene-d10	14.183	164	464519	20.000	ng	0.00
64) Phenanthrene-d10	16.921	188	1111912	20.000	ng	0.00
76) Chrysene-d12	21.081	240	1415522	20.000	ng	0.00
86) Perylene-d12	23.372	264	1928976	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.323	112	1195209	128.985	ng	0.00
7) Phenol-d6	6.950	99	1515408	115.362	ng	0.00
23) Nitrobenzene-d5	8.784	82	839505	77.659	ng	0.00
42) 2,4,6-Tribromophenol	15.693	330	910316	116.513	ng	0.00
45) 2-Fluorobiphenyl	12.808	172	2194500	81.638	ng	0.00
79) Terphenyl-d14	19.518	244	4421634	69.262	ng	0.00

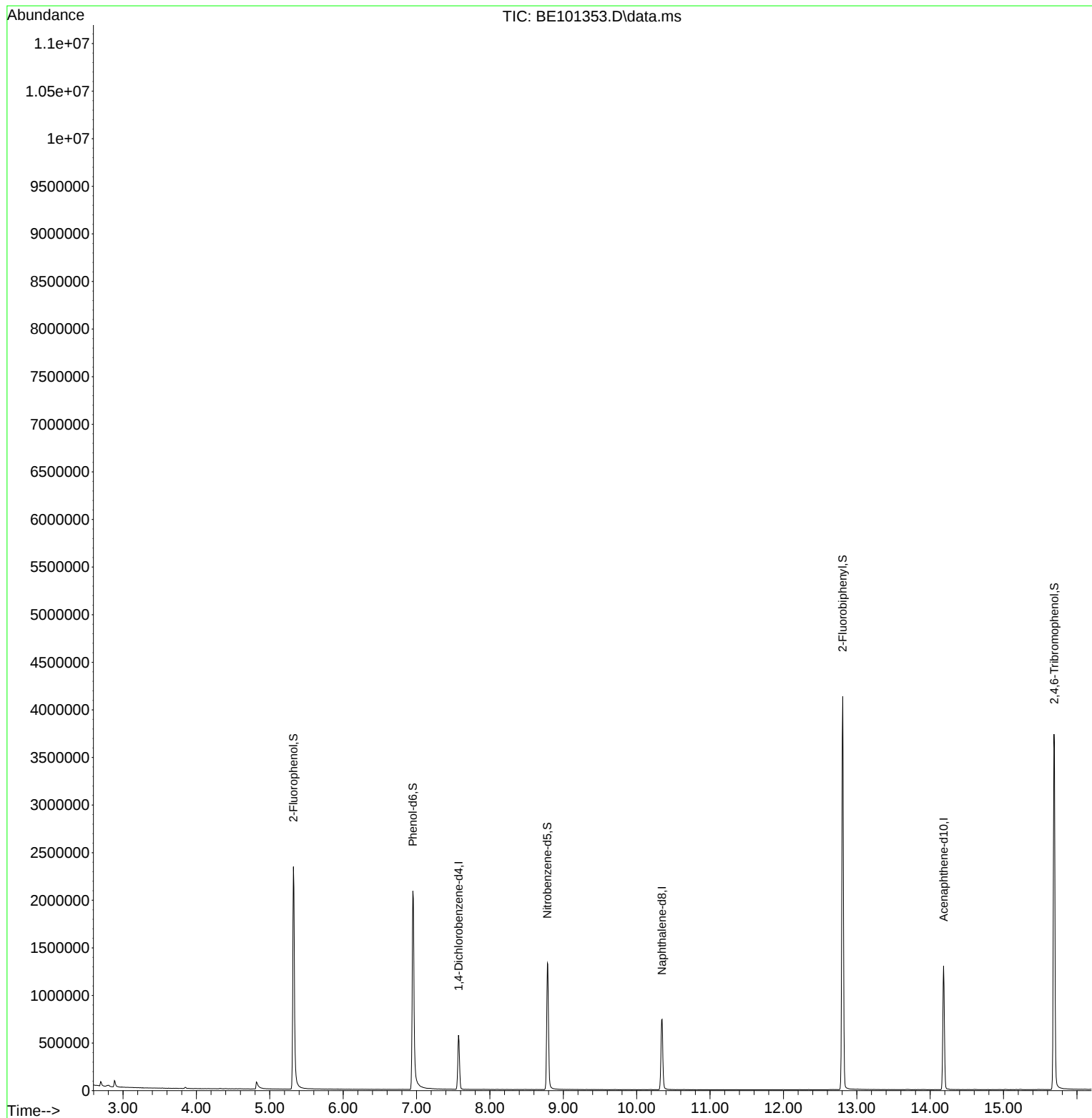
Target Compounds	Qvalue

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

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