

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110824\
 Data File : BE101554.D
 Acq On : 8 Nov 2024 16:20
 Operator : RC/JU
 Sample : P4739-12
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-B2

Quant Time: Nov 08 16:13:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 00:01:20 2024
 Response via : Initial Calibration

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

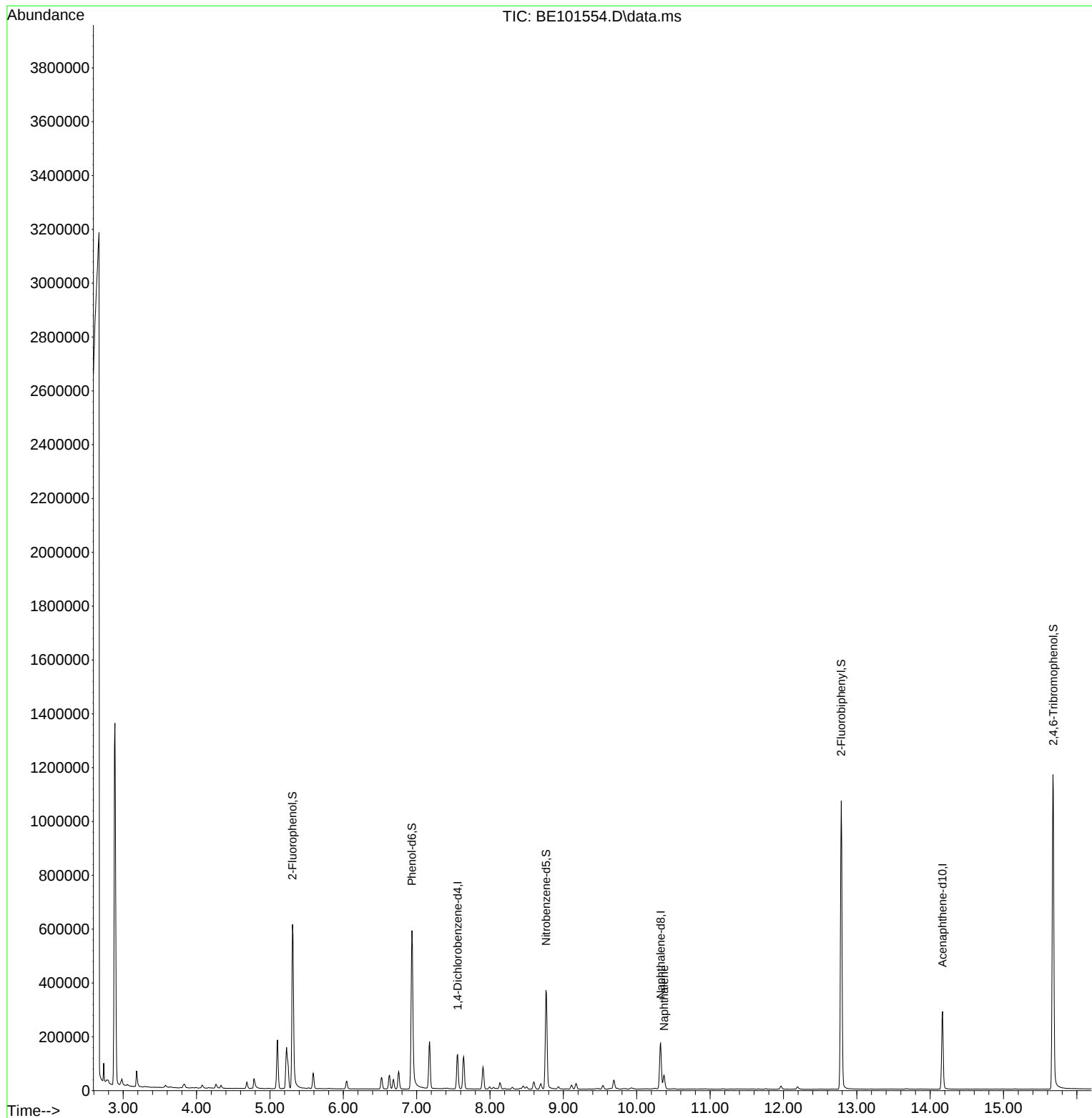
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.559	152	35981	20.000	ng	0.00
21) Naphthalene-d8	10.326	136	152008	20.000	ng	0.00
39) Acenaphthene-d10	14.168	164	104618	20.000	ng	0.00
64) Phenanthrene-d10	16.906	188	262241	20.000	ng	0.00
76) Chrysene-d12	21.066	240	362112	20.000	ng	-0.01
86) Perylene-d12	23.352	264	459208	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.308	112	309909	152.295	ng	0.00
7) Phenol-d6	6.936	99	381111	136.597	ng	-0.01
23) Nitrobenzene-d5	8.763	82	227095	94.366	ng	-0.01
42) 2,4,6-Tribromophenol	15.678	330	316067	160.555	ng	0.00
45) 2-Fluorobiphenyl	12.788	172	574774	89.707	ng	-0.01
79) Terphenyl-d14	19.497	244	1791177	109.493	ng	0.00
Target Compounds						
31) Naphthalene	10.373	128	43586	5.680	ng	Qvalue 99

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

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