

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110824\
 Data File : BE101551.D
 Acq On : 8 Nov 2024 14:32
 Operator : RC/JU
 Sample : P4694-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 Z-03A

Quant Time: Nov 08 14:31:47 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)

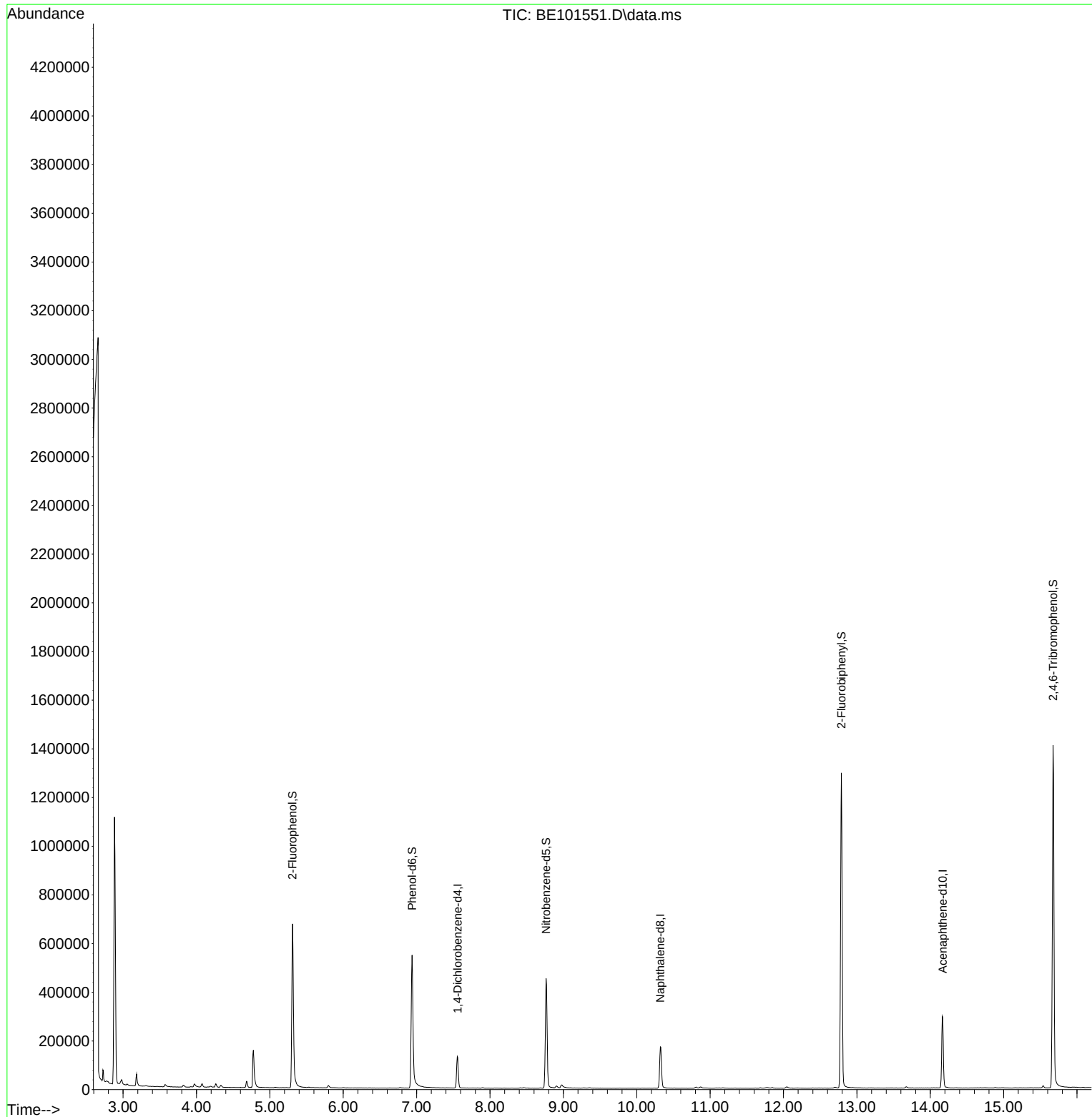
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.560	152	37676	20.000	ng	0.00
21) Naphthalene-d8	10.322	136	159494	20.000	ng	0.00
39) Acenaphthene-d10	14.170	164	109930	20.000	ng	0.00
64) Phenanthrene-d10	16.902	188	283947	20.000	ng	0.00
76) Chrysene-d12	21.068	240	377741	20.000	ng	0.00
86) Perylene-d12	23.348	264	463708	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.310	112	336368	157.861	ng	0.00
7) Phenol-d6	6.938	99	384167	131.498	ng	0.00
23) Nitrobenzene-d5	8.765	82	269661	106.795	ng	0.00
42) 2,4,6-Tribromophenol	15.674	330	360865	174.453	ng	0.00
45) 2-Fluorobiphenyl	12.790	172	701703	104.225	ng	0.00
79) Terphenyl-d14	19.505	244	2011694	117.885	ng	0.00
Target Compounds						
77) Benzidine	19.206	184	56539	6.968	ng	Qvalue 99

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

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