

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111124\
 Data File : BE101570.D
 Acq On : 11 Nov 2024 13:57
 Operator : RC/JU
 Sample : P4793-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 M00-24-00345

Quant Time: Nov 11 13:35:35 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)

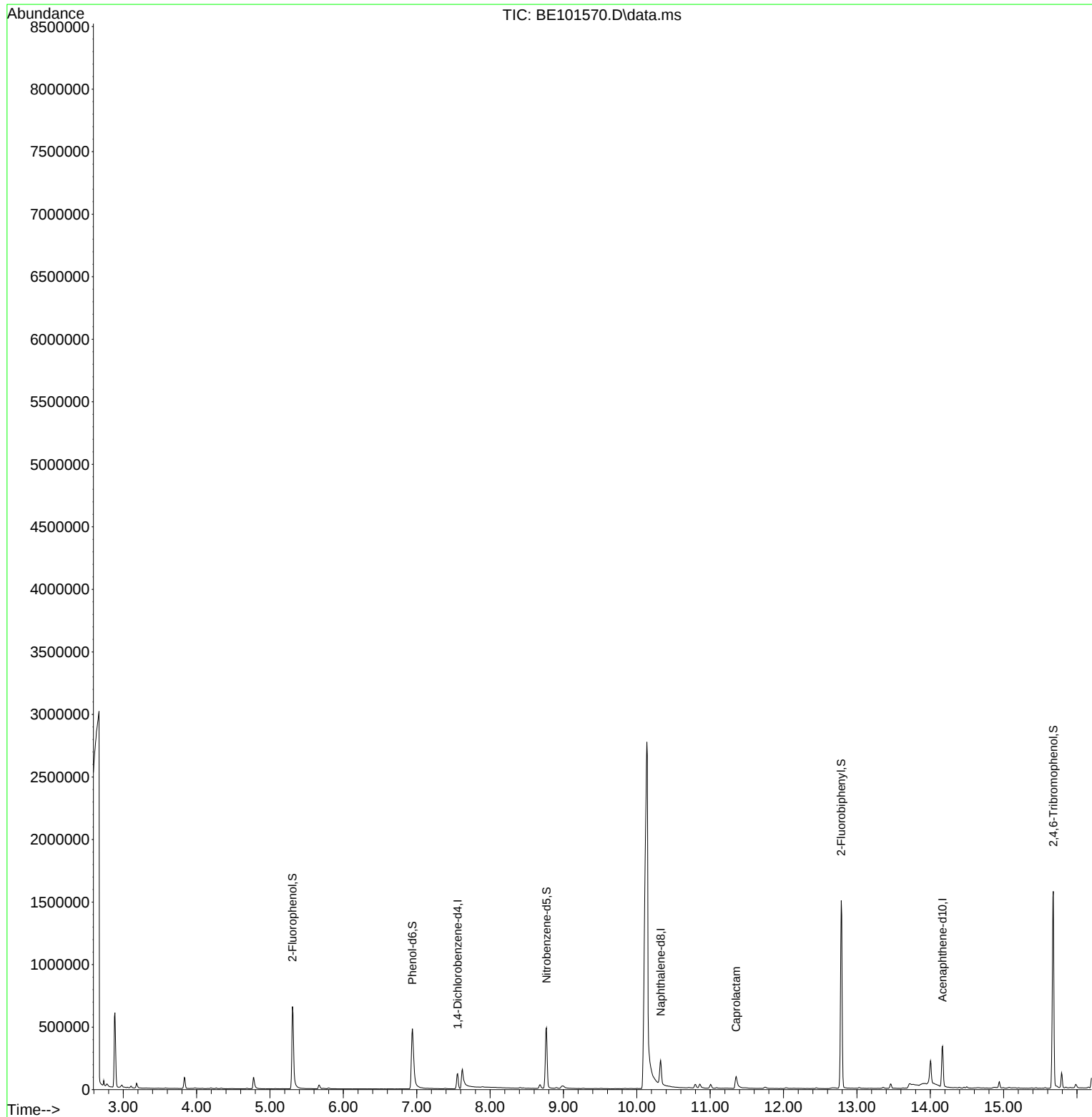
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.557	152	35270	20.000	ng	0.00
21) Naphthalene-d8	10.325	136	166554	20.000	ng	0.00
39) Acenaphthene-d10	14.167	164	117788	20.000	ng	0.00
64) Phenanthrene-d10	16.905	188	293304	20.000	ng	0.00
76) Chrysene-d12	21.065	240	330346	20.000	ng	-0.01
86) Perylene-d12	23.351	264	426782	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.307	112	350177	175.553	ng	0.00
7) Phenol-d6	6.940	99	439022	160.526	ng	0.00
23) Nitrobenzene-d5	8.768	82	295608	112.108	ng	0.00
42) 2,4,6-Tribromophenol	15.677	330	396415	178.854	ng	0.00
45) 2-Fluorobiphenyl	12.787	172	798312	110.664	ng	-0.01
79) Terphenyl-d14	19.502	244	1855088	124.305	ng	0.00
Target Compounds						
35) Caprolactam	11.353	113	24208	27.498	ng	Qvalue 95

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

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