

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103124\
 Data File : BE101434.D
 Acq On : 31 Oct 2024 15:50
 Operator : RC/JU
 Sample : P4639-03 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 EO-02-103024

Quant Time: Oct 31 16:51:48 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)

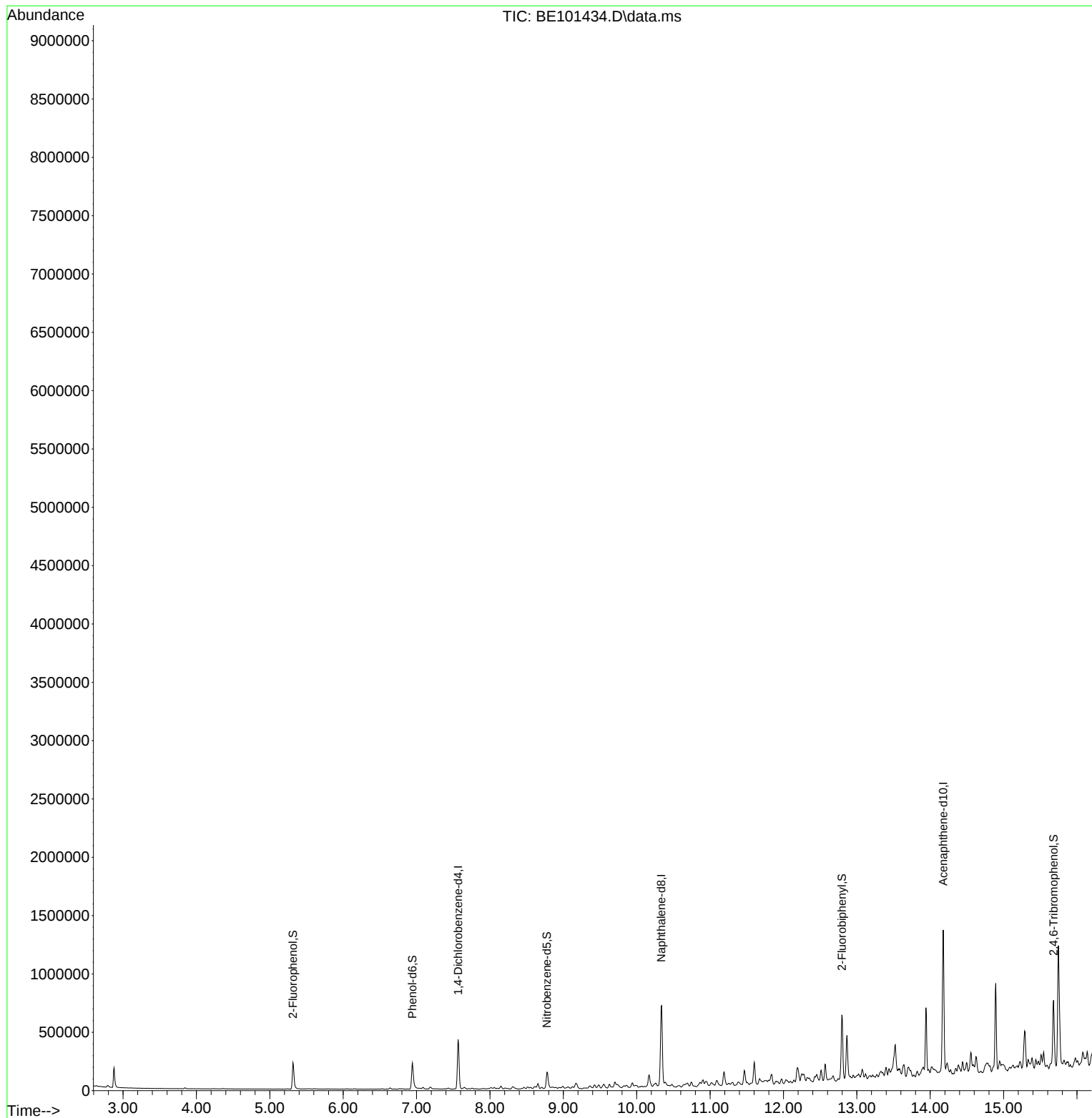
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.567	152	121922	20.000	ng	0.00
21) Naphthalene-d8	10.334	136	577495	20.000	ng	0.00
39) Acenaphthene-d10	14.177	164	429096	20.000	ng	0.00
64) Phenanthrene-d10	16.915	188	1012535	20.000	ng	0.00
76) Chrysene-d12	21.074	240	1092047	20.000	ng	0.00
86) Perylene-d12	23.366	264	1412315	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.316	112	119757	17.215	ng	0.00
7) Phenol-d6	6.944	99	155748	16.147	ng	0.00
23) Nitrobenzene-d5	8.777	82	89935	9.781	ng	0.00
42) 2,4,6-Tribromophenol	15.681	330	131595	17.498	ng	0.00
45) 2-Fluorobiphenyl	12.796	172	294586	11.627	ng	0.00
79) Terphenyl-d14	19.506	244	688531	14.511	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	16.956	178	101855	2.101	ng	98
75) Fluoranthene	18.959	202	135575	2.303	ng	98
78) Pyrene	19.329	202	159975	2.664	ng	98

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

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