

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103124\
 Data File : BE101442.D
 Acq On : 31 Oct 2024 20:36
 Operator : RC/JU
 Sample : P4595-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PITKIN-COMP

Quant Time: Nov 01 06:16:51 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)

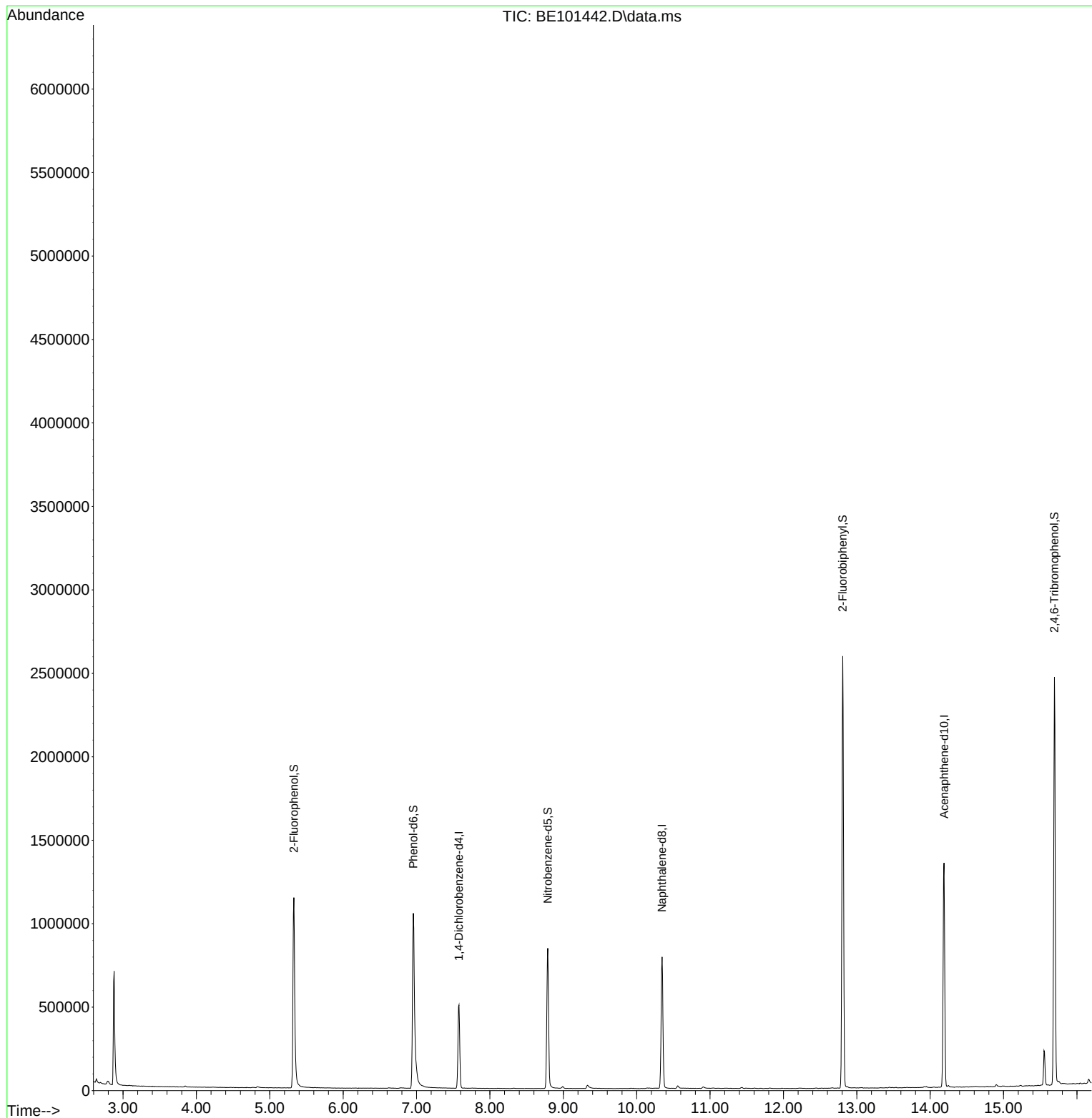
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.579	152	148112	20.000	ng	0.00
21) Naphthalene-d8	10.346	136	698047	20.000	ng	0.00
39) Acenaphthene-d10	14.189	164	479391	20.000	ng	0.00
64) Phenanthrene-d10	16.927	188	1031178	20.000	ng	0.00
76) Chrysene-d12	21.086	240	1039489	20.000	ng	0.00
86) Perylene-d12	23.384	264	1358300	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.328	112	625722	74.044	ng	0.00
7) Phenol-d6	6.956	99	850492	72.582	ng	0.00
23) Nitrobenzene-d5	8.789	82	513706	46.222	ng	0.00
42) 2,4,6-Tribromophenol	15.693	330	593575	70.645	ng	0.00
45) 2-Fluorobiphenyl	12.808	172	1395500	49.299	ng	0.00
79) Terphenyl-d14	19.518	244	2503604	55.433	ng	0.00
Target Compounds						
75) Fluoranthene	18.971	202	160620	2.679	ng	96
78) Pyrene	19.336	202	155894	2.727	ng	99

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

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