

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101508.D
 Acq On : 6 Nov 2024 22:51
 Operator : RC/JU
 Sample : P4667-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-F-6

Quant Time: Nov 07 01:06:04 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	7.564	152	50252	20.000	ng	0.00
21) Naphthalene-d8	10.332	136	209129	20.000	ng	0.00
39) Acenaphthene-d10	14.168	164	139188	20.000	ng	0.00
64) Phenanthrene-d10	16.906	188	336796	20.000	ng	0.00
76) Chrysene-d12	21.066	240	442281	20.000	ng	-0.01
86) Perylene-d12	23.352	264	579467	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.314	112	467725	164.575	ng	0.00
7) Phenol-d6	6.942	99	534758	137.236	ng	0.00
23) Nitrobenzene-d5	8.769	82	367897	111.119	ng	0.00
42) 2,4,6-Tribromophenol	15.678	330	449119	171.479	ng	0.00
45) 2-Fluorobiphenyl	12.794	172	933037	109.454	ng	0.00
79) Terphenyl-d14	19.503	244	2271915	113.707	ng	0.00

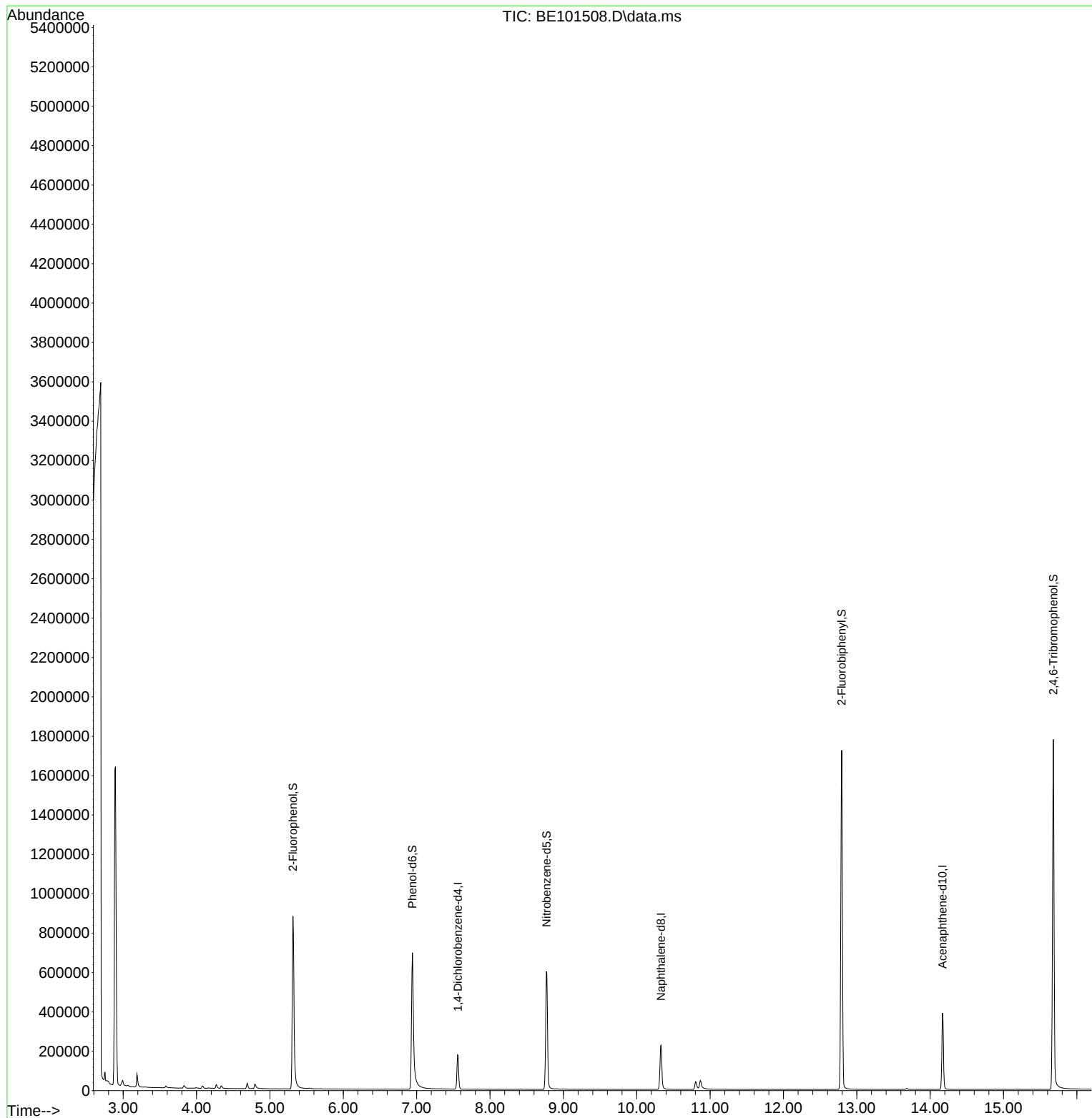
Target Compounds	Qvalue

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

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