

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111124\
 Data File : BE101567.D
 Acq On : 11 Nov 2024 12:09
 Operator : RC/JU
 Sample : P4756-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-B4

Quant Time: Nov 11 12:07:51 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	7.559	152	48103	20.000	ng	0.00
21) Naphthalene-d8	10.321	136	196681	20.000	ng	0.00
39) Acenaphthene-d10	14.163	164	126914	20.000	ng	0.00
64) Phenanthrene-d10	16.901	188	296210	20.000	ng	0.00
76) Chrysene-d12	21.067	240	364666	20.000	ng	0.00
86) Perylene-d12	23.347	264	446133	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.315	112	458141	168.404	ng	0.00
7) Phenol-d6	6.942	99	565733	151.671	ng	0.00
23) Nitrobenzene-d5	8.770	82	341649	109.722	ng	0.00
42) 2,4,6-Tribromophenol	15.673	330	397603	166.491	ng	0.00
45) 2-Fluorobiphenyl	12.788	172	819672	105.454	ng	0.00
79) Terphenyl-d14	19.498	244	1865492	113.237	ng	0.00

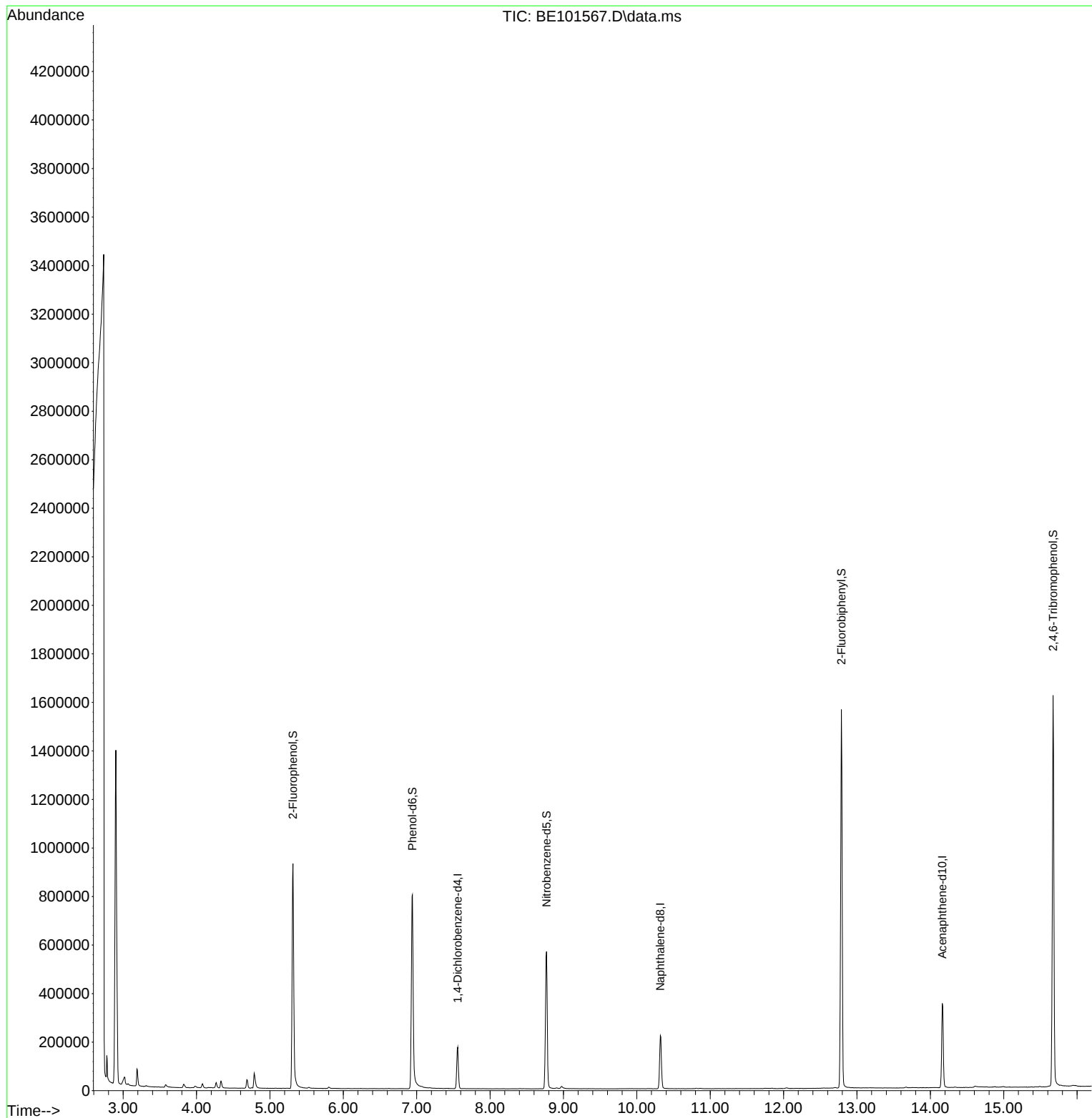
Target Compounds	Qvalue

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

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