

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110824\
 Data File : BE101552.D
 Acq On : 8 Nov 2024 15:08
 Operator : RC/JU
 Sample : P4738-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 72-12018

Quant Time: Nov 08 15:04:26 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	7.555	152	40340	20.000	ng	0.00
21) Naphthalene-d8	10.322	136	167454	20.000	ng	0.00
39) Acenaphthene-d10	14.165	164	113482	20.000	ng	0.00
64) Phenanthrene-d10	16.903	188	293109	20.000	ng	0.00
76) Chrysene-d12	21.063	240	380485	20.000	ng	-0.01
86) Perylene-d12	23.348	264	482208	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.311	112	334221	146.495	ng	0.00
7) Phenol-d6	6.938	99	411552	131.569	ng	0.00
23) Nitrobenzene-d5	8.765	82	258705	97.586	ng	0.00
42) 2,4,6-Tribromophenol	15.675	330	357416	167.378	ng	0.00
45) 2-Fluorobiphenyl	12.790	172	658218	94.706	ng	0.00
79) Terphenyl-d14	19.500	244	1757418	102.242	ng	0.00

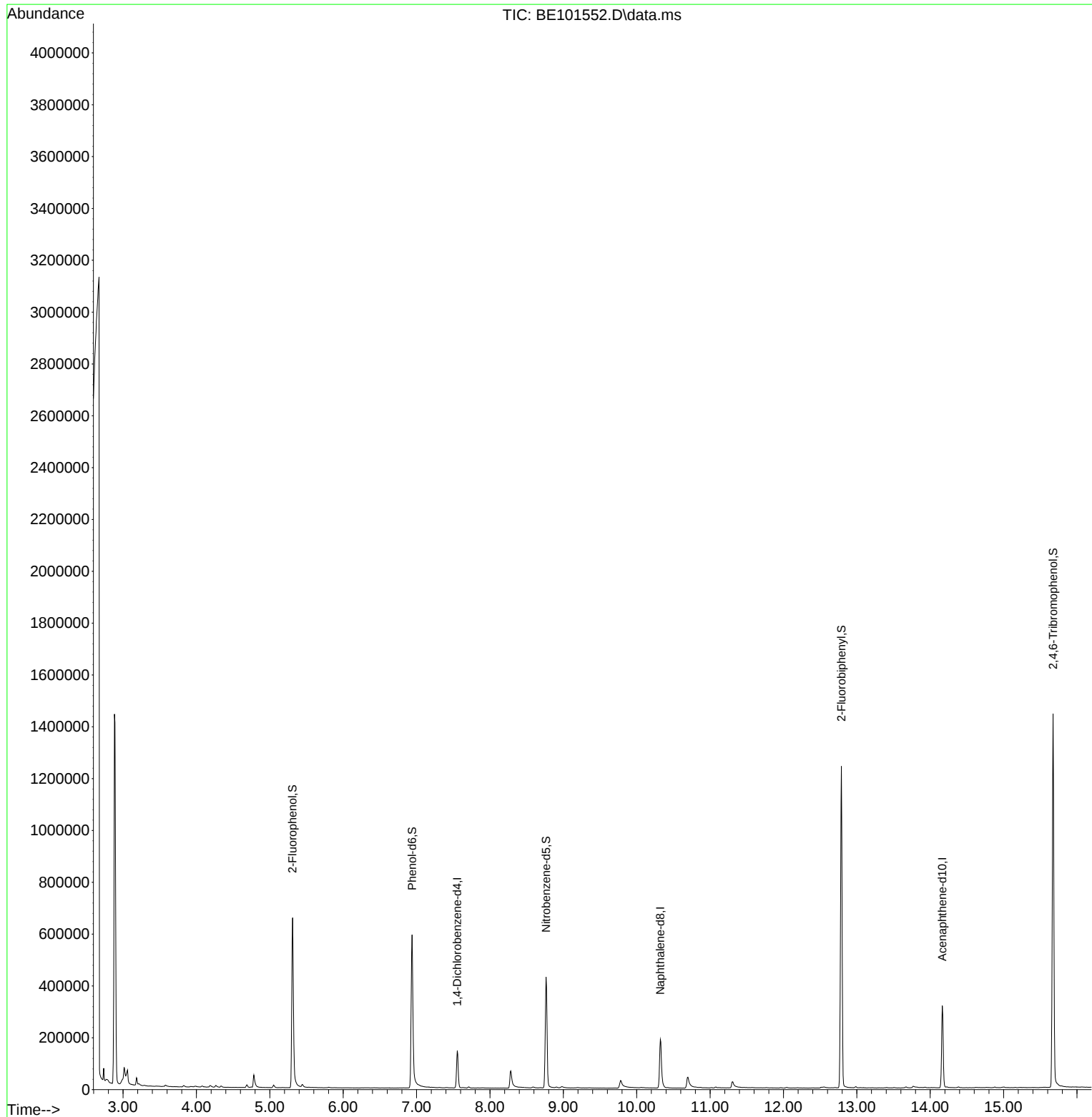
Target Compounds	Qvalue

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

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