

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111224\
 Data File : BE101576.D
 Acq On : 12 Nov 2024 11:11
 Operator : RC/JU
 Sample : PB164886BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164886BL

Quant Time: Nov 12 12:01:17 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	43722	20.000	ng	0.00
21) Naphthalene-d8	10.322	136	176416	20.000	ng	0.00
39) Acenaphthene-d10	14.165	164	107077	20.000	ng	0.00
64) Phenanthrene-d10	16.903	188	234410	20.000	ng	0.00
76) Chrysene-d12	21.069	240	303518	20.000	ng	0.00
86) Perylene-d12	23.348	264	454322	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.305	112	377504	152.668	ng	0.00
7) Phenol-d6	6.932	99	456197	134.560	ng	-0.01
23) Nitrobenzene-d5	8.765	82	273748	98.014	ng	0.00
42) 2,4,6-Tribromophenol	15.675	330	251995	125.068	ng	0.00
45) 2-Fluorobiphenyl	12.790	172	676726	103.193	ng	0.00
79) Terphenyl-d14	19.500	244	1360432	99.217	ng	0.00

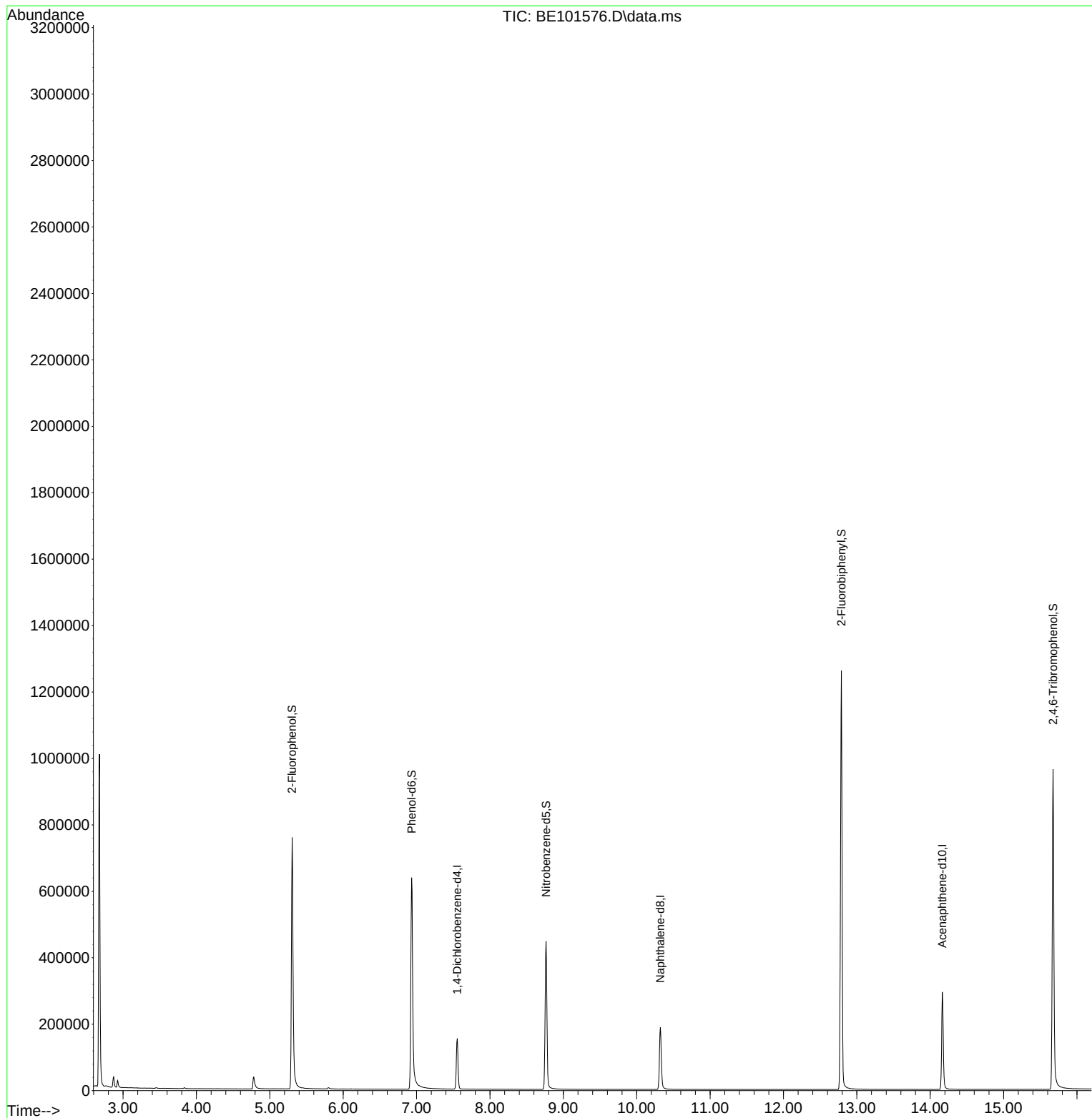
Target Compounds	Qvalue

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

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