

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110824\
 Data File : BE101549.D
 Acq On : 8 Nov 2024 13:21
 Operator : RC/JU
 Sample : P4700-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 MH-8

Quant Time: Nov 08 13:03:51 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	7.558	152	52937	20.000	ng	0.00
21) Naphthalene-d8	10.326	136	221809	20.000	ng	0.00
39) Acenaphthene-d10	14.168	164	140890	20.000	ng	0.00
64) Phenanthrene-d10	16.906	188	331523	20.000	ng	0.00
76) Chrysene-d12	21.066	240	393207	20.000	ng	-0.01
86) Perylene-d12	23.352	264	509724	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.314	112	406941	135.924	ng	0.00
7) Phenol-d6	6.942	99	496179	120.877	ng	0.00
23) Nitrobenzene-d5	8.769	82	323645	92.165	ng	0.00
42) 2,4,6-Tribromophenol	15.678	330	411747	155.310	ng	0.00
45) 2-Fluorobiphenyl	12.788	172	872227	101.084	ng	-0.01
79) Terphenyl-d14	19.503	244	1982468	111.603	ng	0.00

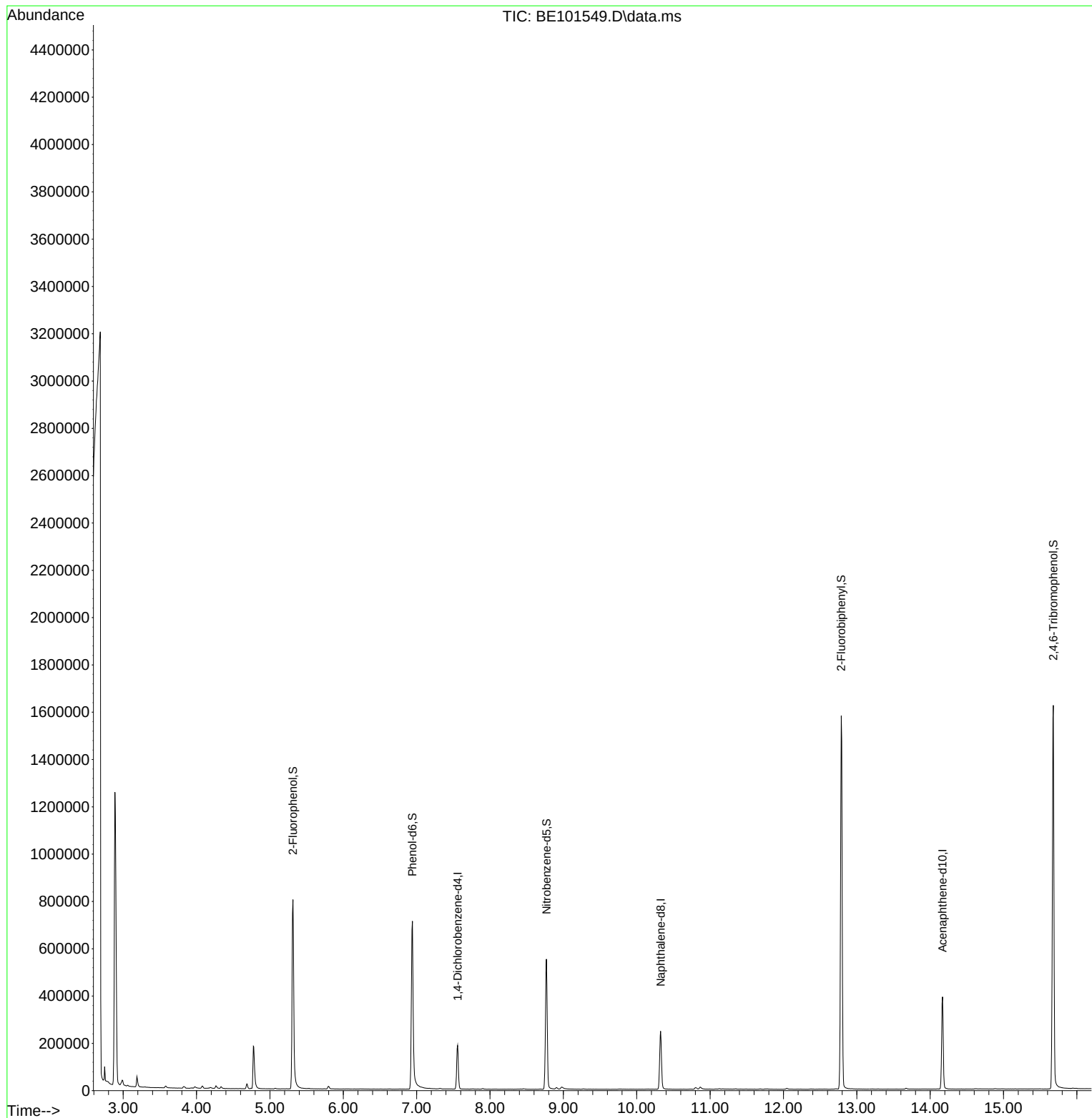
Target Compounds	Qvalue

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

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