

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110524\
 Data File : BE101483.D
 Acq On : 5 Nov 2024 17:28
 Operator : RC/JU
 Sample : P4643-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-F9-ADDITIONAL

Quant Time: Nov 06 00:02:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE102824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 29 01:26:30 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.570	152	61606	20.000	ng	0.00
21) Naphthalene-d8	10.338	136	260314	20.000	ng	0.00
39) Acenaphthene-d10	14.180	164	172720	20.000	ng	0.00
64) Phenanthrene-d10	16.918	188	405128	20.000	ng	0.00
76) Chrysene-d12	21.078	240	501969	20.000	ng	0.00
86) Perylene-d12	23.364	264	663771	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.320	112	480529	136.708	ng	0.00
7) Phenol-d6	6.947	99	571379	117.233	ng	0.00
23) Nitrobenzene-d5	8.781	82	388714	93.790	ng	0.00
42) 2,4,6-Tribromophenol	15.690	330	445365	147.120	ng	0.00
45) 2-Fluorobiphenyl	12.799	172	953335	93.475	ng	0.00
79) Terphenyl-d14	19.509	244	2163568	99.201	ng	0.00

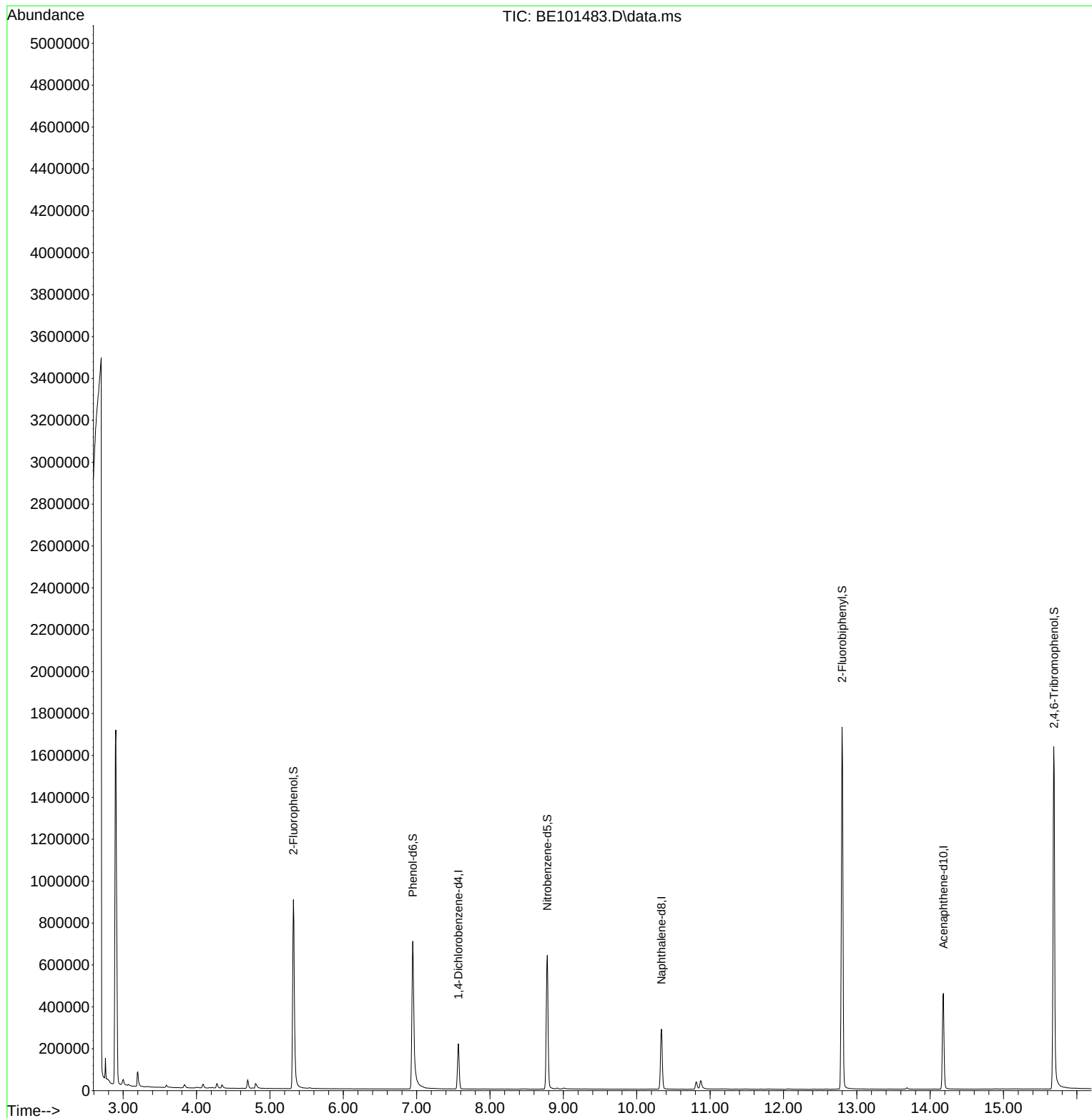
Target Compounds	Qvalue

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

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