

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE091824\
 Data File : BE101367.D
 Acq On : 18 Sep 2024 17:16
 Operator : RC/JU
 Sample : PB163110BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB163110BL

Quant Time: Sep 18 18:05:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE091724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 17 16:03:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.575	152	146861	20.000	ng	0.00
21) Naphthalene-d8	10.342	136	633980	20.000	ng	0.00
39) Acenaphthene-d10	14.185	164	431261	20.000	ng	0.00
64) Phenanthrene-d10	16.923	188	1053064	20.000	ng	0.00
76) Chrysene-d12	21.083	240	1310107	20.000	ng	0.00
86) Perylene-d12	23.368	264	1774064	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.325	112	1083013	128.938	ng	0.00
7) Phenol-d6	6.952	99	1385894	116.390	ng	0.00
23) Nitrobenzene-d5	8.785	82	780202	78.020	ng	0.00
42) 2,4,6-Tribromophenol	15.689	330	845253	116.528	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	2034924	81.540	ng	0.00
79) Terphenyl-d14	19.514	244	4122600	69.774	ng	0.00

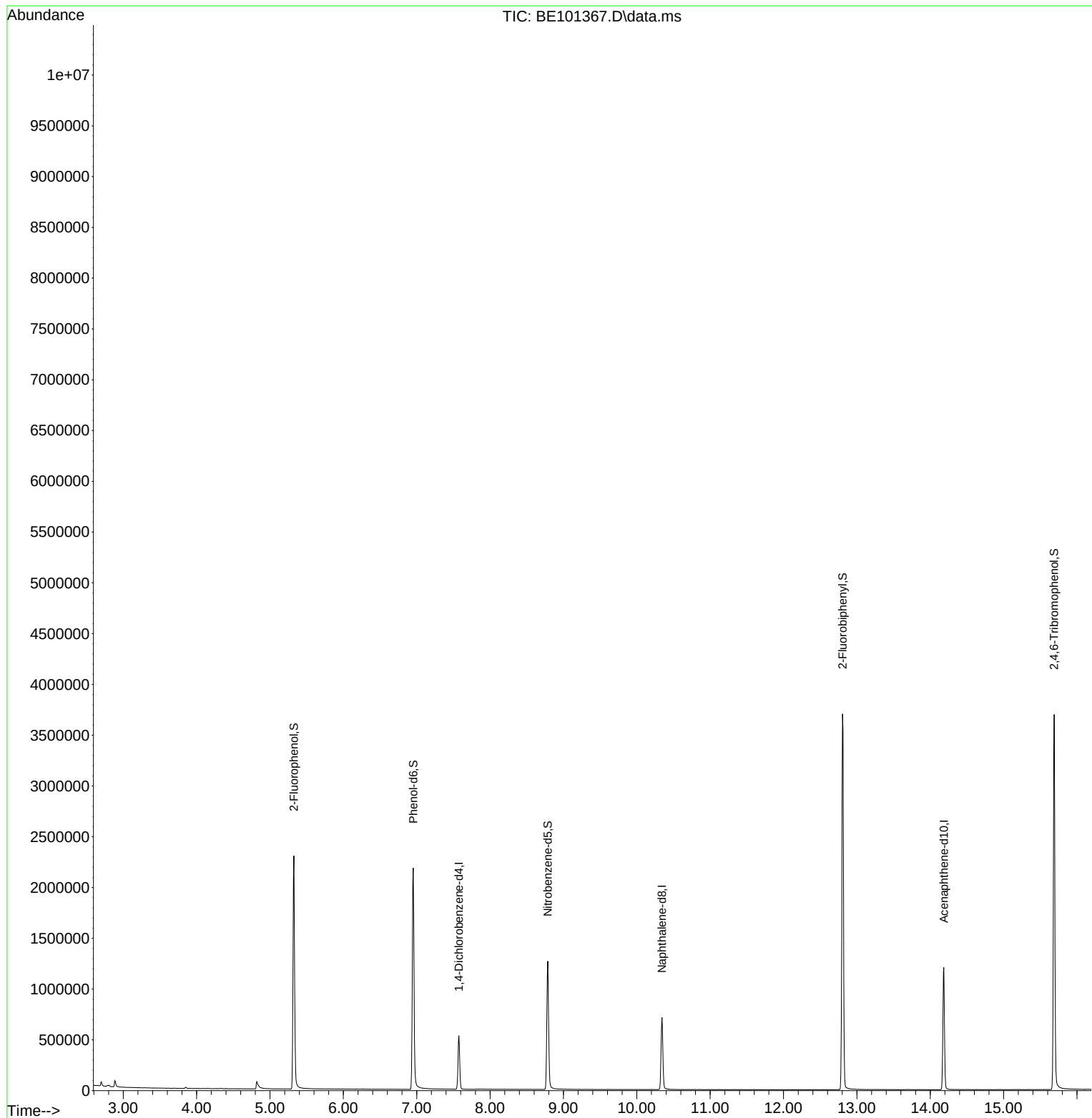
Target Compounds	Qvalue

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

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