

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE091824\
 Data File : BE101356.D
 Acq On : 18 Sep 2024 10:18
 Operator : RC/JU
 Sample : PB163109BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB163109BL

Quant Time: Sep 18 10:57:54 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	7.577	152	149016	20.000	ng	0.00
21) Naphthalene-d8	10.344	136	621553	20.000	ng	0.00
39) Acenaphthene-d10	14.187	164	398373	20.000	ng	0.00
64) Phenanthrene-d10	16.924	188	920959	20.000	ng	0.00
76) Chrysene-d12	21.084	240	1163723	20.000	ng	0.00
86) Perylene-d12	23.376	264	1662666	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.320	112	1331803	156.264	ng	0.00
7) Phenol-d6	6.954	99	1657655	137.200	ng	0.00
23) Nitrobenzene-d5	8.787	82	932844	95.150	ng	0.00
42) 2,4,6-Tribromophenol	15.691	330	895188	133.601	ng	0.00
45) 2-Fluorobiphenyl	12.812	172	2347907	101.848	ng	0.00
79) Terphenyl-d14	19.516	244	4270939	81.377	ng	0.00

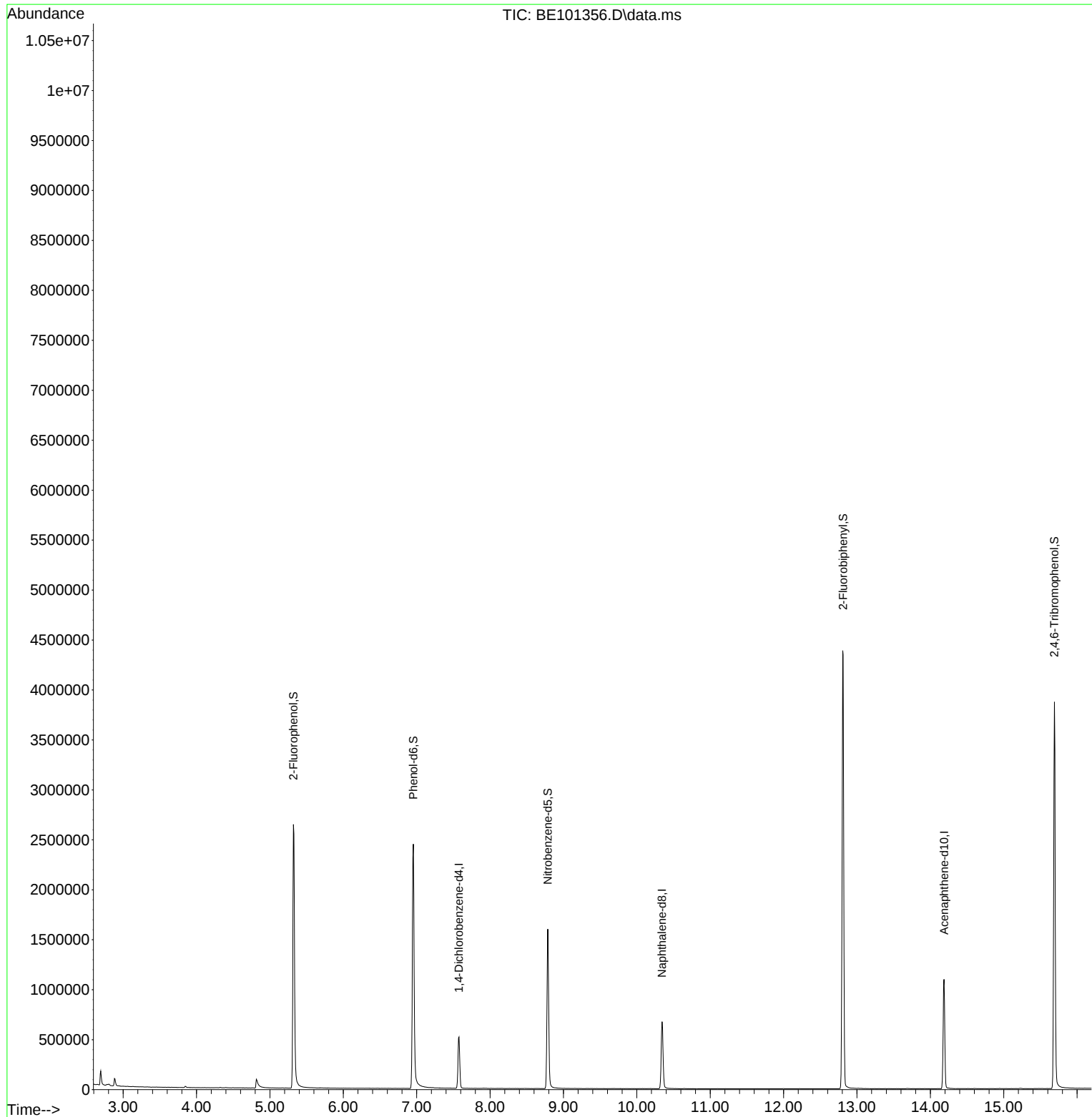
Target Compounds	Qvalue

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

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