

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110724\
 Data File : BE101525.D
 Acq On : 7 Nov 2024 11:26
 Operator : RC/JU
 Sample : PB164711BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164711BL

Quant Time: Nov 07 11:12:24 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	7.566	152	70050	20.000	ng	0.00
21) Naphthalene-d8	10.333	136	282724	20.000	ng	0.00
39) Acenaphthene-d10	14.176	164	172547	20.000	ng	0.00
64) Phenanthrene-d10	16.914	188	381552	20.000	ng	0.00
76) Chrysene-d12	21.074	240	468916	20.000	ng	0.00
86) Perylene-d12	23.359	264	671548	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.322	112	567284	143.192	ng	0.00
7) Phenol-d6	6.949	99	692828	127.550	ng	0.00
23) Nitrobenzene-d5	8.776	82	412982	92.267	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	422295	130.065	ng	0.00
45) 2-Fluorobiphenyl	12.795	172	993875	94.050	ng	0.00
79) Terphenyl-d14	19.505	244	1939141	91.539	ng	0.00

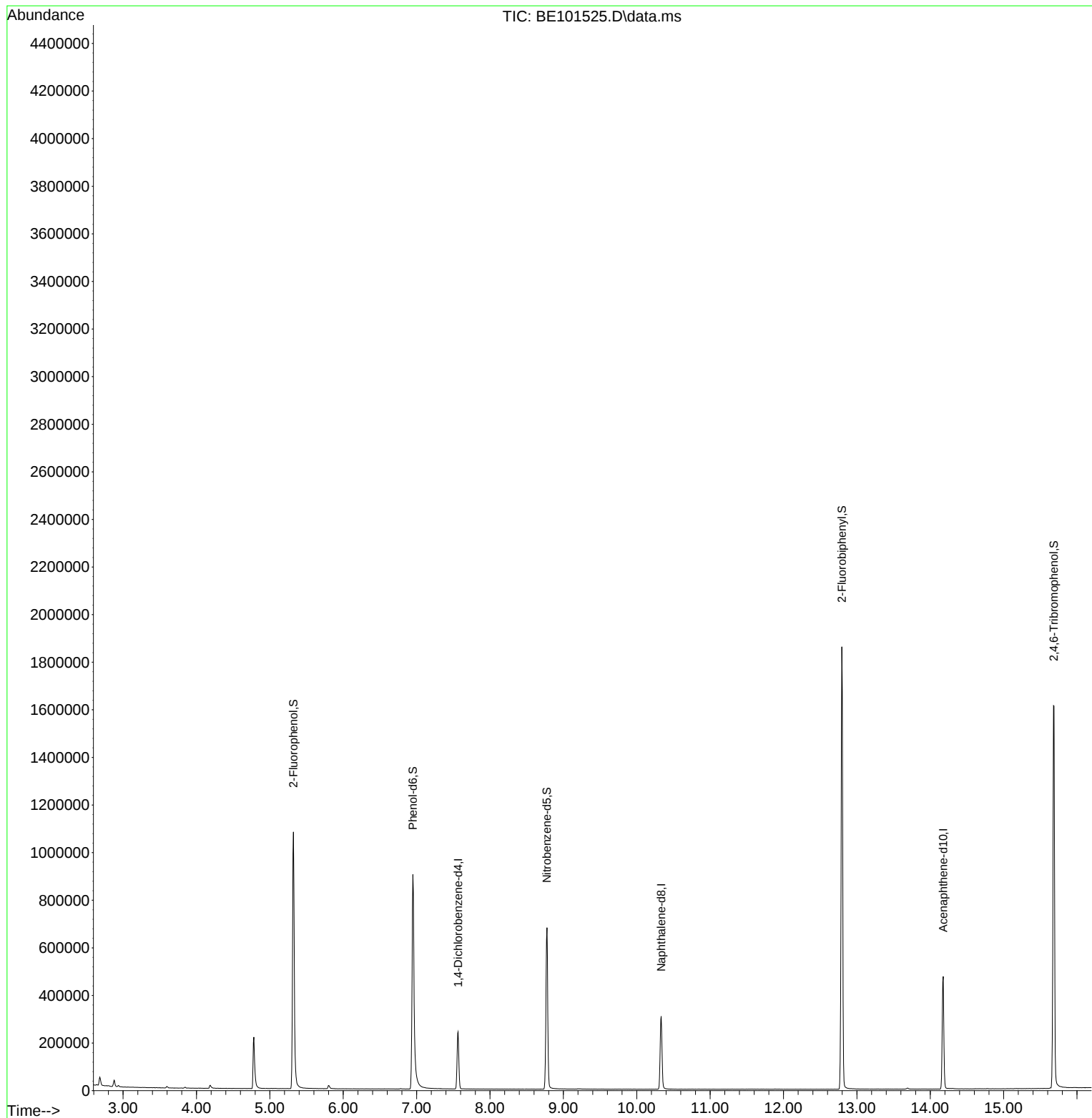
Target Compounds	Qvalue

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

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