

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110524\
 Data File : BE101473.D
 Acq On : 5 Nov 2024 11:33
 Operator : RC/JU
 Sample : PB164639BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PB164639BL

Quant Time: Nov 05 11:18:06 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	69685	20.000	ng	0.00
21) Naphthalene-d8	10.337	136	281108	20.000	ng	0.00
39) Acenaphthene-d10	14.179	164	180533	20.000	ng	0.00
64) Phenanthrene-d10	16.917	188	412460	20.000	ng	0.00
76) Chrysene-d12	21.077	240	536197	20.000	ng	0.00
86) Perylene-d12	23.369	264	732979	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.319	112	561372	141.192	ng	0.00
7) Phenol-d6	6.947	99	693279	125.752	ng	0.00
23) Nitrobenzene-d5	8.780	82	408173	91.200	ng	0.00
42) 2,4,6-Tribromophenol	15.689	330	450681	142.433	ng	0.00
45) 2-Fluorobiphenyl	12.805	172	1041592	97.709	ng	0.00
79) Terphenyl-d14	19.514	244	2161061	92.761	ng	0.00

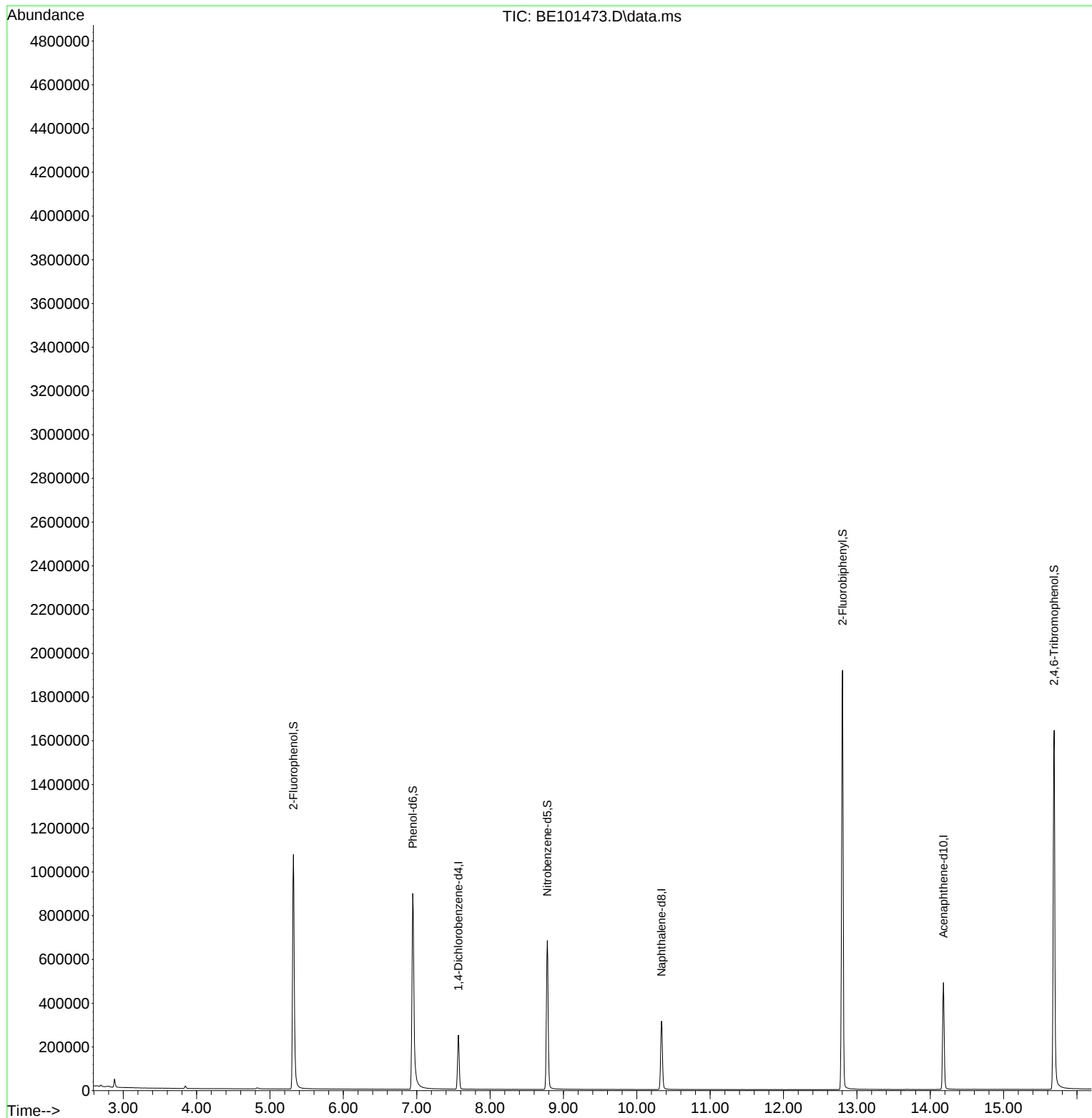
Target Compounds	Qvalue

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

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