

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110624\
 Data File : BE101511.D
 Acq On : 6 Nov 2024 00:39
 Operator : RC/JU
 Sample : P4667-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-F-7

Quant Time: Nov 07 01:06:58 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	7.558	152	50314	20.000	ng	0.00
21) Naphthalene-d8	10.326	136	206452	20.000	ng	0.00
39) Acenaphthene-d10	14.168	164	119151	20.000	ng	0.00
64) Phenanthrene-d10	16.906	188	351377	20.000	ng	0.00
76) Chrysene-d12	21.066	240	415210	20.000	ng	-0.01
86) Perylene-d12	23.346	264	88152	20.000	ng	#-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.314	112	425582	149.562	ng	0.00
7) Phenol-d6	6.941	99	466116	119.473	ng	0.00
23) Nitrobenzene-d5	8.769	82	347114	106.201	ng	0.00
42) 2,4,6-Tribromophenol	15.678	330	418998	186.881	ng	0.00
45) 2-Fluorobiphenyl	12.793	172	857624	117.526	ng	0.00
79) Terphenyl-d14	19.503	244	2263990	120.698	ng	0.00

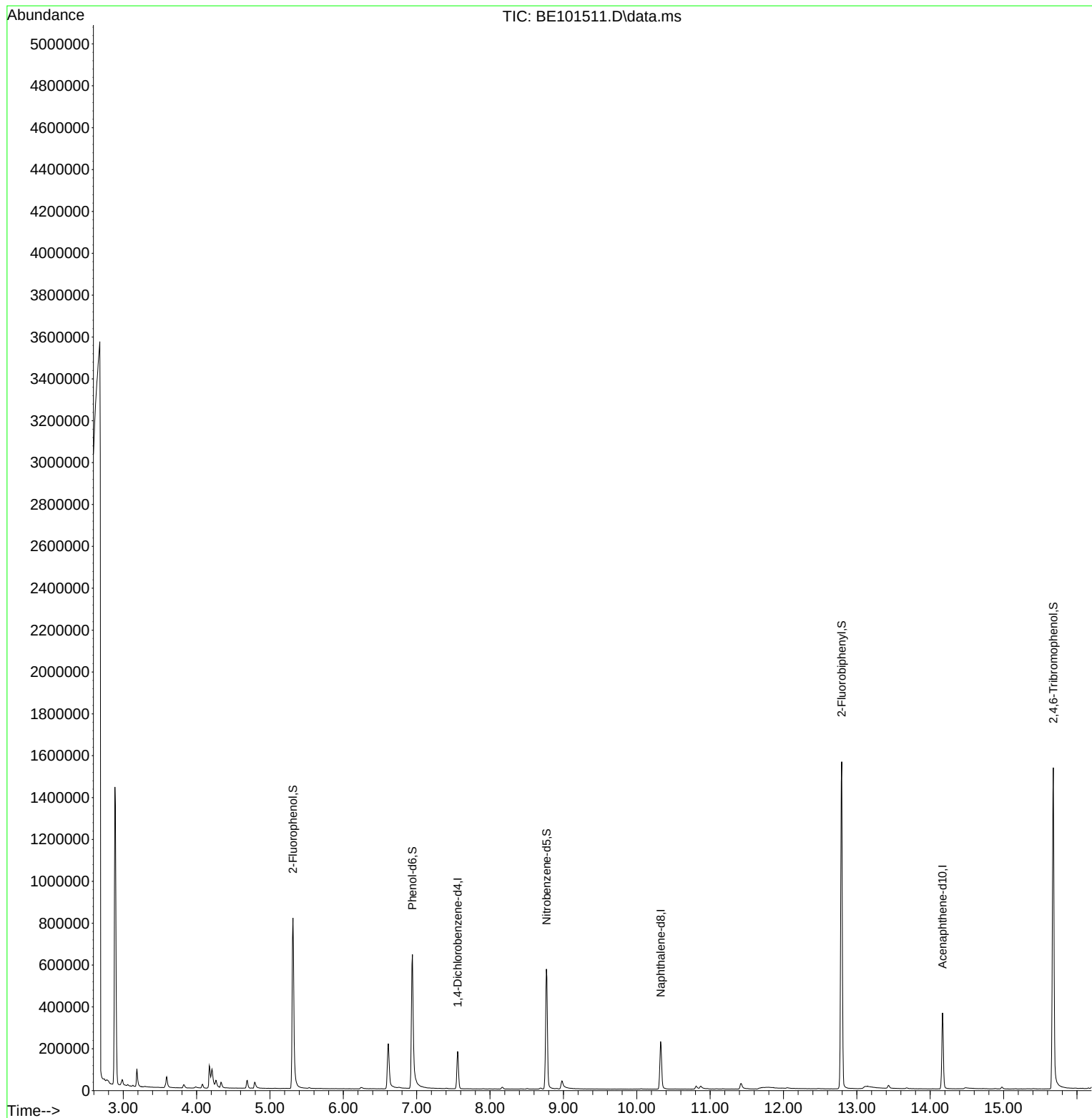
Target Compounds	Qvalue

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

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