

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110724\
 Data File : BE101530.D
 Acq On : 7 Nov 2024 14:31
 Operator : RC/JU
 Sample : P4711-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 CF-613-COMP-17

Quant Time: Nov 07 14:13:47 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	7.559	152	52015	20.000	ng	0.00
21) Naphthalene-d8	10.326	136	217168	20.000	ng	0.00
39) Acenaphthene-d10	14.169	164	126302	20.000	ng	0.00
64) Phenanthrene-d10	16.907	188	320318	20.000	ng	0.00
76) Chrysene-d12	21.067	240	355224	20.000	ng	-0.01
86) Perylene-d12	23.352	264	115918	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.315	112	439838	149.517	ng	0.00
7) Phenol-d6	6.942	99	486198	120.545	ng	0.00
23) Nitrobenzene-d5	8.769	82	359492	104.561	ng	0.00
42) 2,4,6-Tribromophenol	15.679	330	402839	169.501	ng	0.00
45) 2-Fluorobiphenyl	12.794	172	896499	115.897	ng	0.00
79) Terphenyl-d14	19.504	244	1991901	124.124	ng	0.00

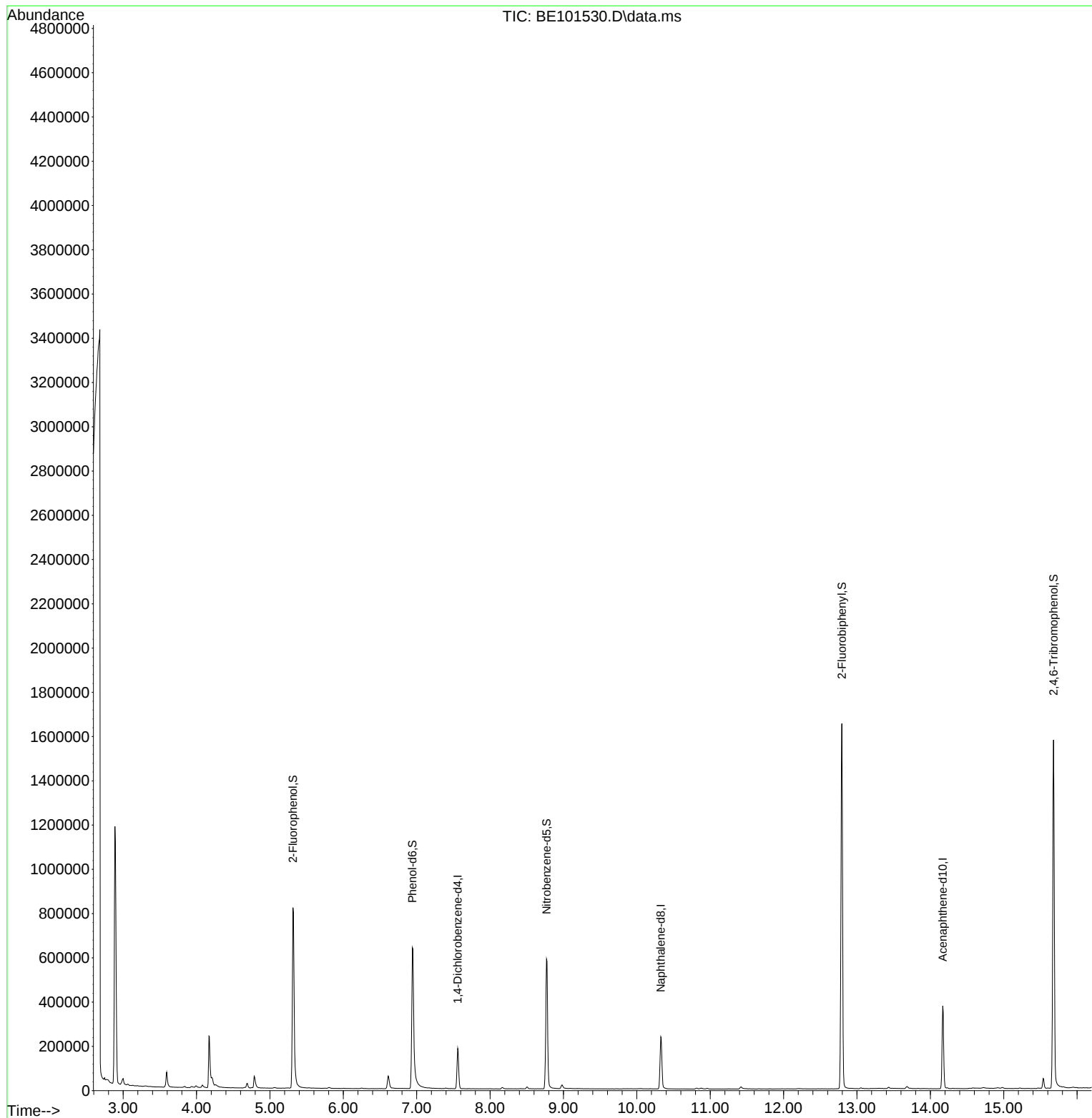
Target Compounds	Qvalue

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

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