

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111124\
 Data File : BE101571.D
 Acq On : 11 Nov 2024 14:33
 Operator : RC/JU
 Sample : P4795-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 LAW-23-00189

Quant Time: Nov 11 14:20:14 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	7.553	152	39706	20.000	ng	0.00
21) Naphthalene-d8	10.320	136	167207	20.000	ng	-0.01
39) Acenaphthene-d10	14.169	164	112501	20.000	ng	0.00
64) Phenanthrene-d10	16.901	188	270240	20.000	ng	-0.01
76) Chrysene-d12	21.067	240	338714	20.000	ng	-0.01
86) Perylene-d12	23.346	264	425347	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.309	112	368633	164.159	ng	0.00
7) Phenol-d6	6.936	99	423859	137.667	ng	-0.01
23) Nitrobenzene-d5	8.763	82	297927	112.546	ng	-0.01
42) 2,4,6-Tribromophenol	15.673	330	360088	170.099	ng	-0.01
45) 2-Fluorobiphenyl	12.788	172	720361	104.551	ng	-0.01
79) Terphenyl-d14	19.498	244	1819130	118.884	ng	0.00

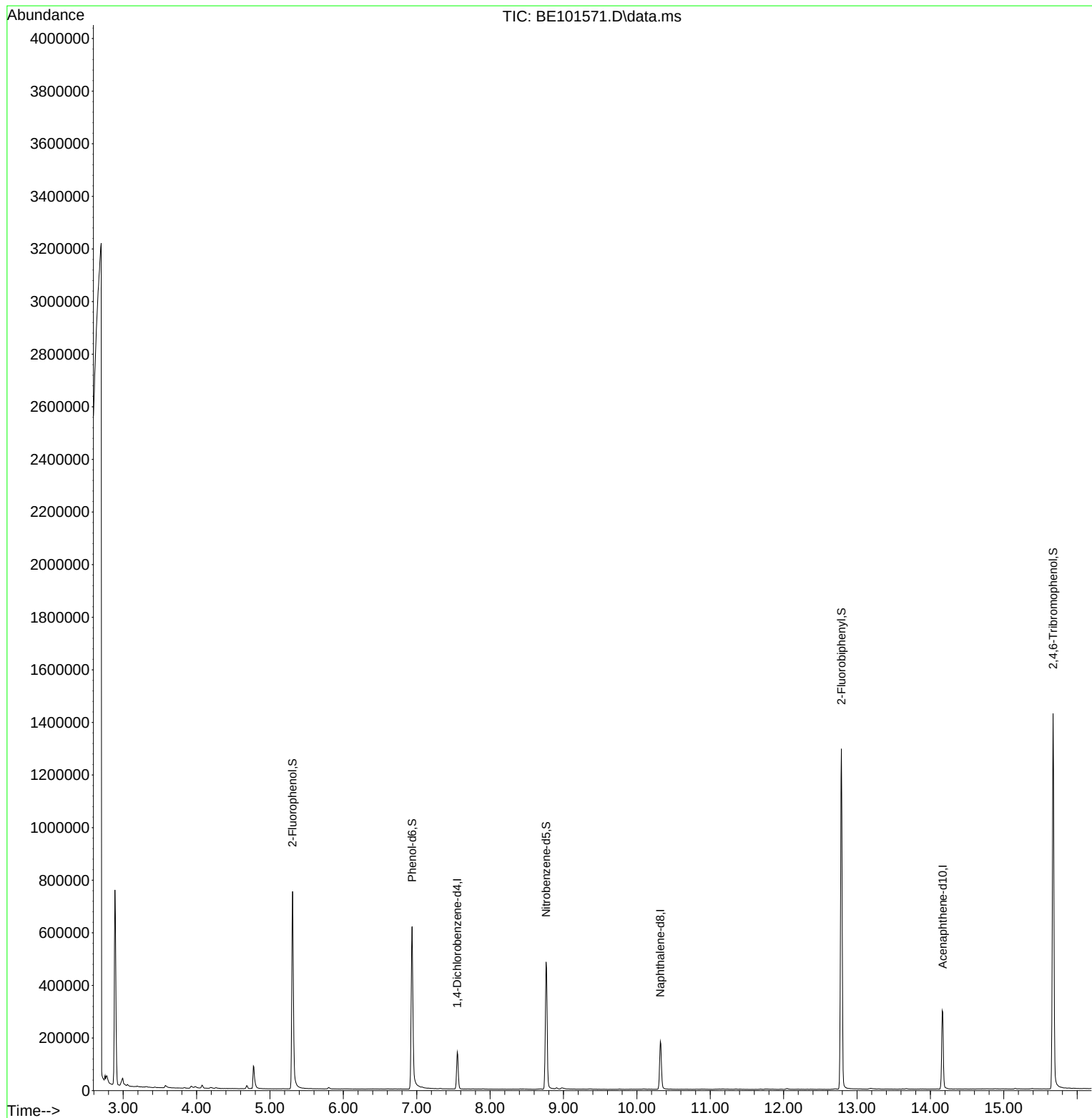
Target Compounds	Qvalue

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

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