

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE103124\
 Data File : BE101429.D
 Acq On : 31 Oct 2024 12:51
 Operator : RC/JU
 Sample : P4616-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 BP-F9-MOVED

Quant Time: Oct 31 14:12:53 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	96511	20.000	ng	0.00
21) Naphthalene-d8	10.338	136	394946	20.000	ng	0.00
39) Acenaphthene-d10	14.174	164	265035	20.000	ng	0.00
64) Phenanthrene-d10	16.912	188	678201	20.000	ng	0.00
76) Chrysene-d12	21.072	240	948781	20.000	ng	0.00
86) Perylene-d12	23.358	264	1233314	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.320	112	876360	159.149	ng	0.00
7) Phenol-d6	6.948	99	977708	128.050	ng	0.00
23) Nitrobenzene-d5	8.781	82	706075	112.289	ng	0.00
42) 2,4,6-Tribromophenol	15.684	330	842539	181.378	ng	0.00
45) 2-Fluorobiphenyl	12.800	172	1693906	108.238	ng	0.00
79) Terphenyl-d14	19.509	244	4061108	98.514	ng	0.00

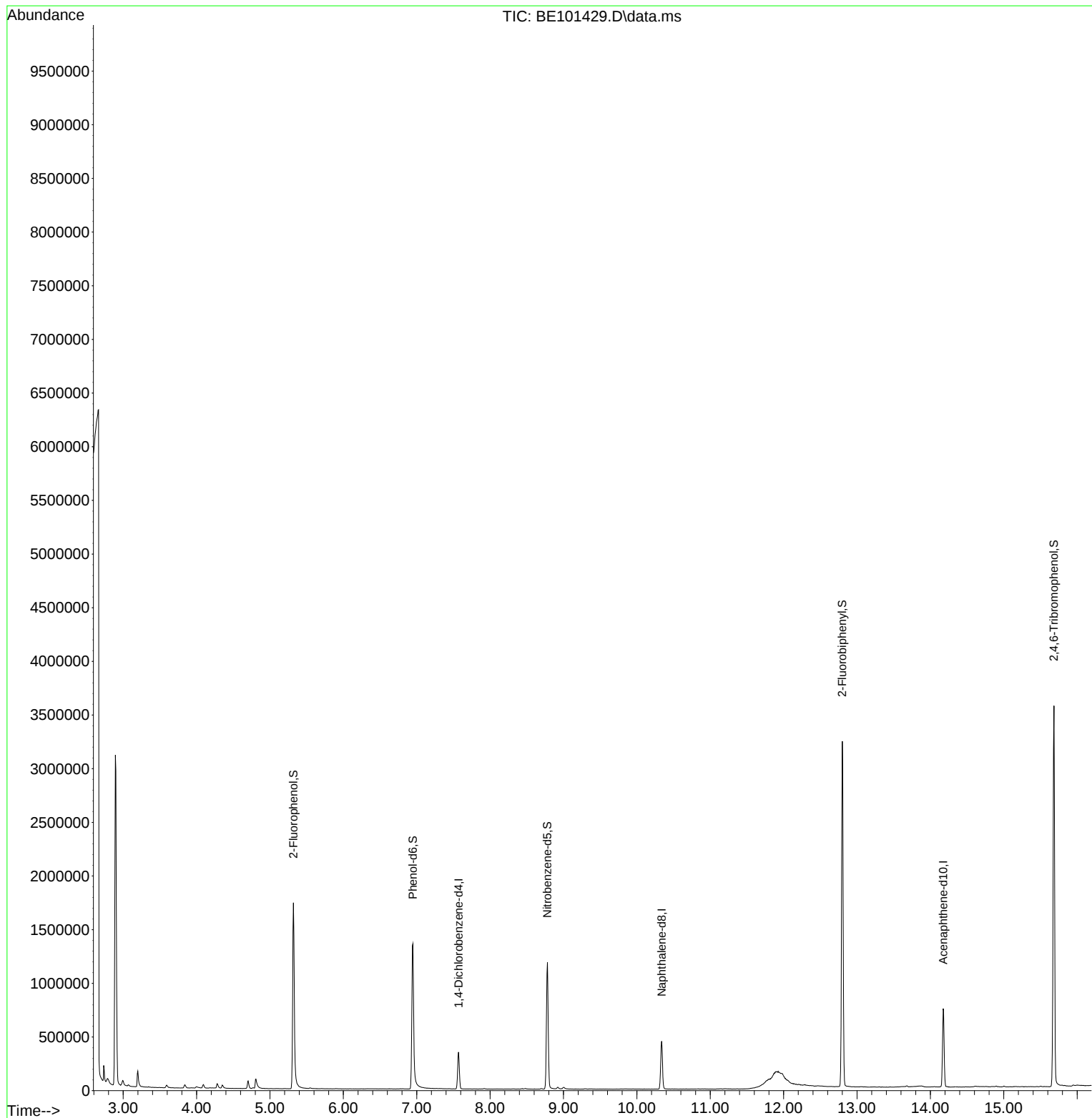
Target Compounds	Qvalue

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

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