

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110824\
 Data File : BE101555.D
 Acq On : 8 Nov 2024 16:55
 Operator : RC/JU
 Sample : P4739-16
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 TP-11

Quant Time: Nov 08 17:25:33 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	7.557	152	38129	20.000	ng	0.00
21) Naphthalene-d8	10.325	136	158412	20.000	ng	0.00
39) Acenaphthene-d10	14.167	164	106535	20.000	ng	0.00
64) Phenanthrene-d10	16.905	188	274316	20.000	ng	0.00
76) Chrysene-d12	21.065	240	372052	20.000	ng	-0.01
86) Perylene-d12	23.351	264	471691	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.307	112	313885	145.560	ng	0.00
7) Phenol-d6	6.935	99	371488	125.647	ng	-0.01
23) Nitrobenzene-d5	8.768	82	251308	100.206	ng	0.00
42) 2,4,6-Tribromophenol	15.677	330	320840	160.047	ng	0.00
45) 2-Fluorobiphenyl	12.787	172	626278	95.986	ng	-0.01
79) Terphenyl-d14	19.502	244	1850655	110.107	ng	0.00

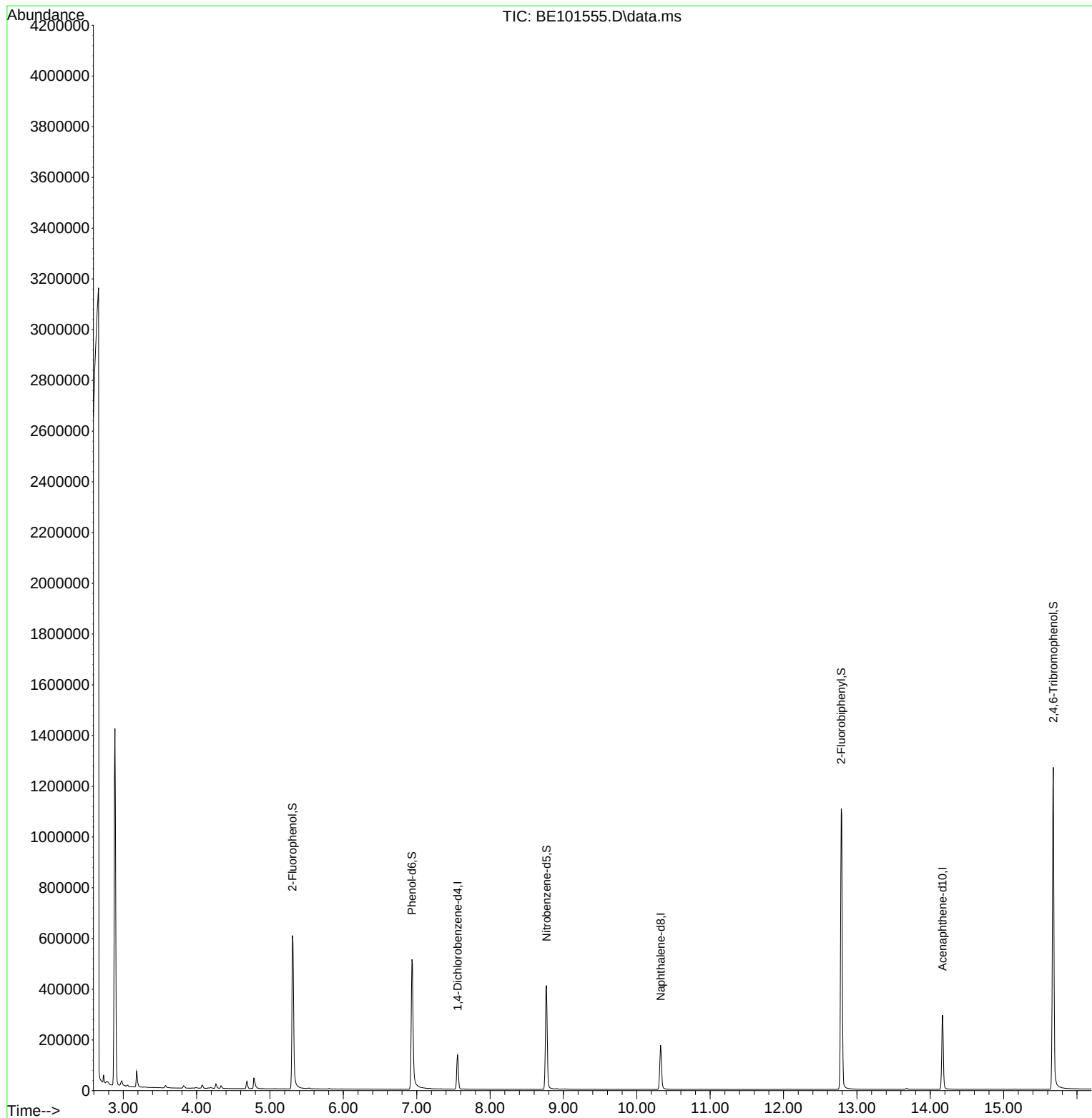
Target Compounds	Qvalue

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

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