

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100220\
 Data File : BP003483.D
 Acq On : 03 Oct 2020 07:24
 Operator : CG/JU
 Sample : L4198-23
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RO-17

Quant Time: Oct 03 08:48:05 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 02 11:17:24 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.528	152	111934	20.00	ng	0.00
21) Naphthalene-d8	10.304	136	445271	20.00	ng	# 0.00
39) Acenaphthene-d10	14.187	164	276977	20.00	ng	0.00
64) Phenanthrene-d10	16.951	188	611432	20.00	ng	# 0.01
76) Chrysene-d12	21.057	240	626748	20.00	ng	0.00
86) Perylene-d12	23.239	264	547769	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.146	112	902657	140.81	ng	0.00
7) Phenol-d6	6.734	99	1048227	122.68	ng	0.00
23) Nitrobenzene-d5	8.681	82	866264	114.29	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	465847	110.27	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	1799592	95.02	ng	0.00
79) Terphenyl-d14	19.527	244	2631313	87.48	ng	0.00

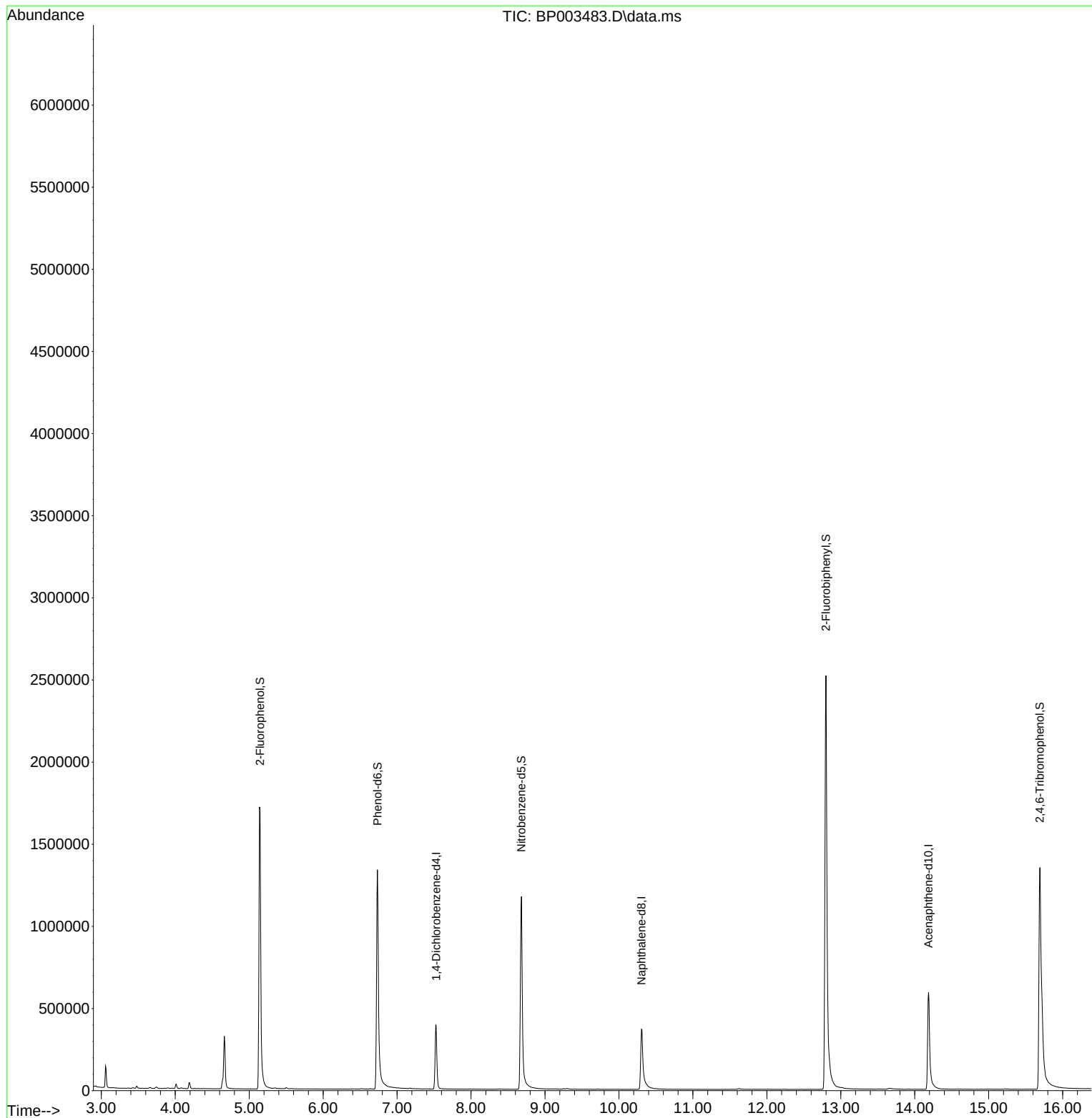
Target Compounds	Qvalue

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

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