

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100120\
 Data File : BP003441.D
 Acq On : 01 Oct 2020 16:59
 Operator : CG/JU
 Sample : PB132017BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB132017BL

Quant Time: Oct 01 17:50:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 10:05:32 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.528	152	80711	20.00	ng	0.00
21) Naphthalene-d8	10.310	136	322946	20.00	ng	# 0.00
39) Acenaphthene-d10	14.187	164	199371	20.00	ng	0.00
64) Phenanthrene-d10	16.945	188	455911	20.00	ng	# 0.00
76) Chrysene-d12	21.051	240	517579	20.00	ng	-0.01
86) Perylene-d12	23.239	264	538676	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.146	112	647001	139.98	ng	0.00
7) Phenol-d6	6.734	99	814208	132.15	ng	0.00
23) Nitrobenzene-d5	8.681	82	543034	98.79	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	304886	100.26	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	1158101	84.96	ng	0.00
79) Terphenyl-d14	19.528	244	1894878	76.28	ng	0.00

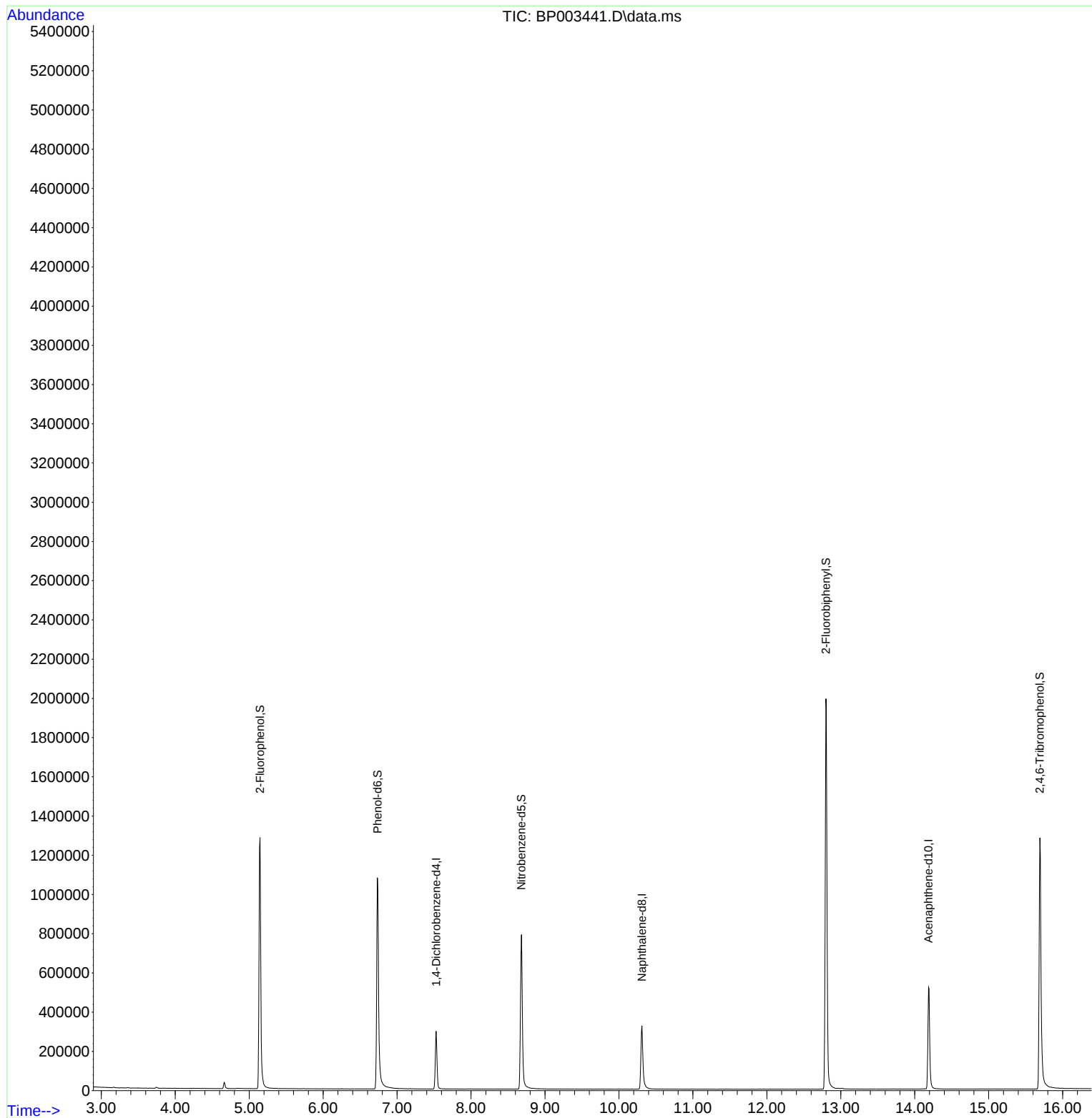
Target Compounds	Qvalue

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

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