

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100120\
 Data File : BP003439.D
 Acq On : 01 Oct 2020 15:52
 Operator : CG/JU
 Sample : PB132022TB
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB132022TB

Quant Time: Oct 01 17:49:54 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	78407	20.00	ng	0.00
21) Naphthalene-d8	10.310	136	314297	20.00	ng	# 0.00
39) Acenaphthene-d10	14.193	164	193472	20.00	ng	0.00
64) Phenanthrene-d10	16.951	188	442791	20.00	ng	# 0.00
76) Chrysene-d12	21.051	240	507629	20.00	ng	-0.01
86) Perylene-d12	23.239	264	522588	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.146	112	650562	144.88	ng	0.00
7) Phenol-d6	6.734	99	815556	136.26	ng	0.00
23) Nitrobenzene-d5	8.681	82	544625	101.80	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	308756	104.63	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	1159647	87.66	ng	0.00
79) Terphenyl-d14	19.528	244	1919293	78.78	ng	0.00

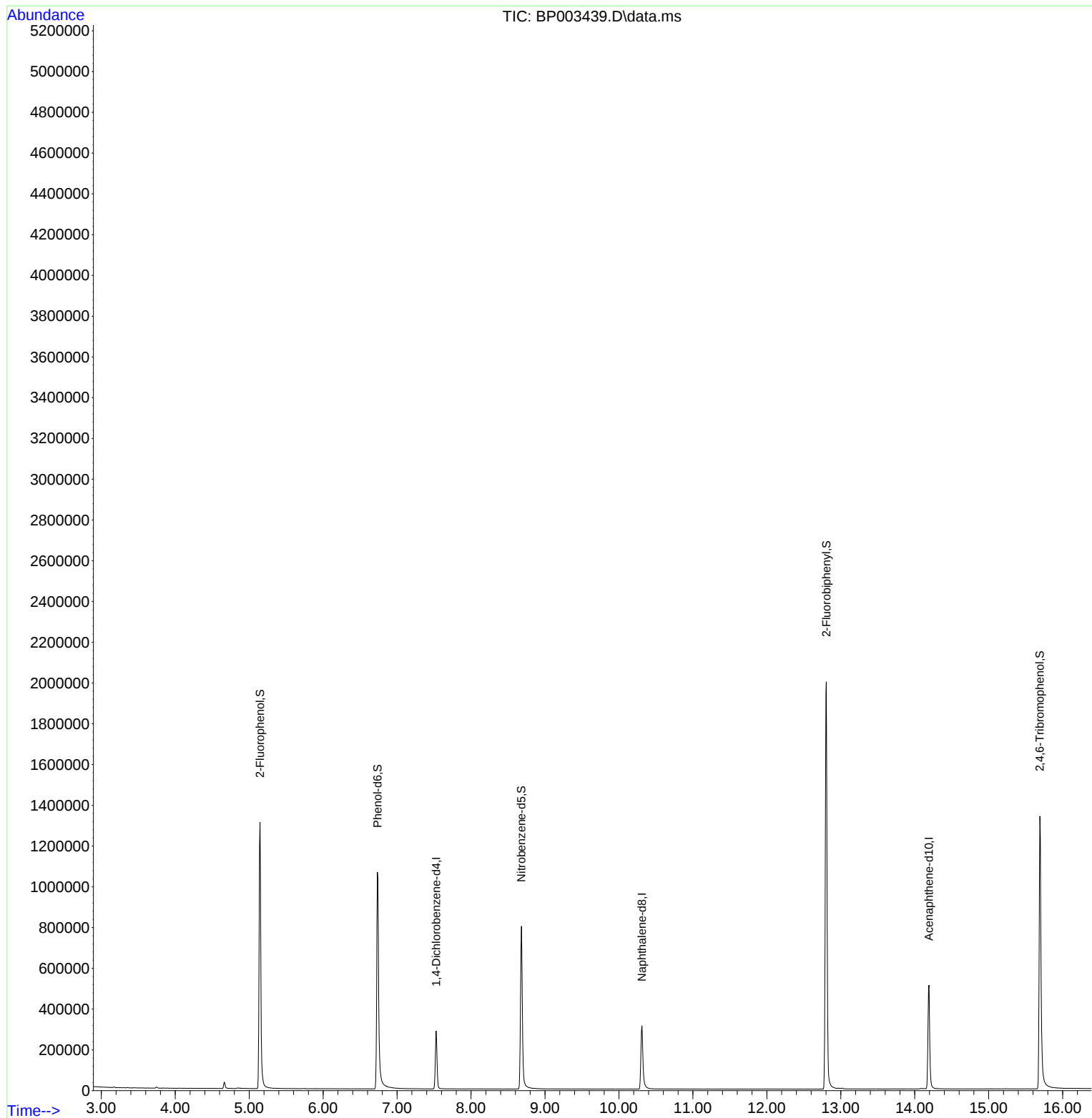
Target Compounds	Qvalue

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

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