

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100520\
 Data File : BP003496.D
 Acq On : 05 Oct 2020 18:29
 Operator : CG/JU
 Sample : L4221-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-02-100120

Quant Time: Oct 06 04:18:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP091520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 05 14:33:14 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.499	152	95925	20.00	ng	0.00
21) Naphthalene-d8	10.275	136	383951	20.00	ng	# 0.00
39) Acenaphthene-d10	14.157	164	238714	20.00	ng	0.00
64) Phenanthrene-d10	16.916	188	533890	20.00	ng	# 0.00
76) Chrysene-d12	21.027	240	610821	20.00	ng	0.00
86) Perylene-d12	23.198	264	561405	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.116	112	710532	129.34	ng	0.00
7) Phenol-d6	6.710	99	875939	119.62	ng	0.00
23) Nitrobenzene-d5	8.652	82	733635	112.25	ng	0.00
42) 2,4,6-Tribromophenol	15.663	330	466878	128.23	ng	0.00
45) 2-Fluorobiphenyl	12.775	172	1506035	92.27	ng	0.00
79) Terphenyl-d14	19.504	244	2549775	86.98	ng	0.00

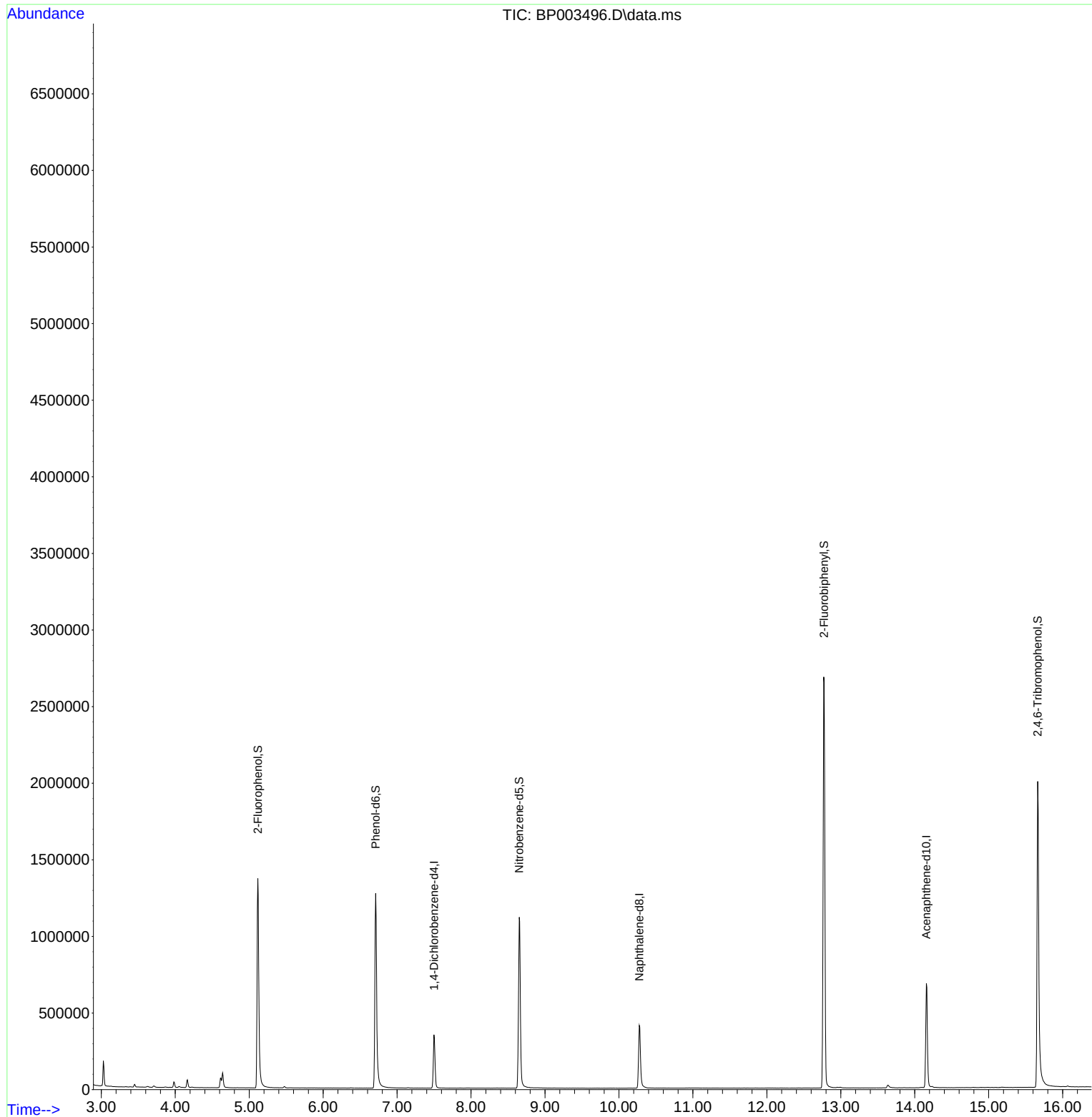
Target Compounds	Qvalue

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

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