

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101019\
 Data File : BP000609.D
 Acq On : 10 Oct 2019 12:27
 Operator : JU
 Sample : K5285-07
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 TP-1-D

Quant Time: Oct 11 04:20:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	270791	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	1011832	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	573912	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	1091743	20.00	ng	-0.02
76) Chrysene-d12	13.96	240	909047	20.00	ng	-0.02
87) Perylene-d12	15.44	264	974696	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1446749	85.88	ng	-0.02
7) Phenol-d6	6.39	99	1847427	91.01	ng	-0.02
23) Nitrobenzene-d5	7.31	82	1003525	60.57	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	527048	86.18	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	2062601	58.15	ng	-0.02
79) Terphenyl-d14	12.90	244	2456808	54.72	ng	-0.02
Target Compounds						
10) Phenol	6.40	94	78829	3.793	ng	89
50) Dimethylphthalate	9.51	163	340227	8.532	ng	99

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

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