

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101019\
 Data File : BP000617.D
 Acq On : 10 Oct 2019 15:39
 Operator : JU
 Sample : K5292-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20190855-AMENDED-COMP-B

Quant Time: Oct 11 04:43:02 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	266502	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	966242	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	523580	20.00	ng	-0.02
64) Phenanthrene-d10	11.30	188	939803	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	780849	20.00	ng	-0.01
87) Perylene-d12	15.45	264	851414	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	237076	14.30	ng	-0.02
7) Phenol-d6	6.39	99	676174	33.84	ng	-0.02
23) Nitrobenzene-d5	7.31	82	861881	54.48	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	31766	5.69	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1665518	51.47	ng	-0.02
79) Terphenyl-d14	12.90	244	1971309	51.12	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	74856	3.659	ng	97
50) Dimethylphthalate	9.51	163	234670	6.450	ng	# 97
75) Fluoranthene	12.52	202	119336	2.251	ng	97
78) Pyrene	12.75	202	163759	3.041	ng	97
81) Benzo(a)anthracene	13.96	228	112718	2.366	ng	# 85
83) Chrysene	13.99	228	108163	2.313	ng	# 87

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

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