

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010719\
 Data File : BF111901.D
 Acq On : 8 Jan 2019 2:31
 Operator : JU/SJ
 Sample : K1064-05 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-1419-W

Quant Time: Jan 08 06:51:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	104765	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	305465	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	153490	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	273206	20.00	ng	0.00
76) Chrysene-d12	14.06	240	313336	20.00	ng	0.00
87) Perylene-d12	15.54	264	207908	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	28312	4.62	ng	0.00
7) Phenol-d6	6.53	99	21251	2.69	ng	0.00
23) Nitrobenzene-d5	7.45	82	85977	15.04	ng	-0.01
42) 2,4,6-Tribromophenol	10.73	330	16455	8.24	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	72887	6.92	ng	0.00
79) Terphenyl-d14	13.00	244	73544	4.46	ng	0.00
Target Compounds						
10) Phenol	6.55	94	20022	2.334	ng	88
20) 3+4-Methylphenols	7.30	107	42403	6.234	ng	# 64
22) Acetophenone	7.29	105	407437	52.685	ng	# 50
31) Naphthalene	8.20	128	376708	25.979	ng	# 92
37) 2-Methylnaphthalene	8.89	142	405262	45.103	ng	97
38) 1-Methylnaphthalene	8.99	142	208693	24.216	ng	# 98
46) 1,1'-Biphenyl	9.35	154	69018	5.376	ng	92

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

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