

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101717\
 Data File : BG029332.D
 Acq On : 18 Oct 2017 13:49
 Operator : SJ/JU
 Sample : I5906-13DL 10X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WH-TWS-03DL

Manual Integrations
 APPROVED

Sohil
 10/18/2017 7:29:04 PM

Quant Time: Oct 18 15:55:35 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	63874	20.00	ng	0.00
21) Naphthalene-d8	10.98	136	310620	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	198809	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	447145	20.00	ng	0.00
75) Chrysene-d12	21.79	240	471202	20.00	ng	0.00
86) Perylene-d12	25.10	264	463207	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	50503	13.21	ng	0.00
7) Phenol-d6	7.32	99	67244	12.53	ng	0.00
23) Nitrobenzene-d5	9.33	82	46303	9.47	ng	0.00
41) 2,4,6-Tribromophenol	16.26	330	29035	11.31	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	138063	10.19	ng	0.00
78) Terphenyl-d14	20.11	244	211186	10.20	ng	-0.01
Target Compounds						
10) Phenol	7.35	94	100722	17.408	ng	93
17) 2-Methylphenol	8.60	107	36970	9.619	ng	92
20) 3+4-Methylphenols	8.93	107	142206	26.258	ng	93
27) 2,4-Dimethylphenol	10.14	122	57397	12.077	ng	93
31) Naphthalene	11.04	128	735622	47.519	ng	99
37) 2-Methylnaphthalene	12.63	142	47516m	4.211	ng	
70) Phenanthrene	17.56	178	50490	2.186	ng	97
72) Carbazole	17.92	167	103469	4.644	ng	98

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

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