

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101715.D
 Acq On : 26 Dec 2017 15:06
 Operator : SJ/JU
 Sample : I6837-03DL 10X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-69-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:37 PM

Quant Time: Dec 26 15:51:09 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	134009	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	512485	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	204871	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	329494	20.00	ng	0.00
75) Chrysene-d12	13.97	240	273409	20.00	ng	0.00
86) Perylene-d12	15.43	264	291102	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	121648m	13.52	ng	0.04
7) Phenol-d6	6.43	99	119129	10.64	ng	0.00
23) Nitrobenzene-d5	7.36	82	102172	11.10	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	26501	13.42	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	200079	15.15	ng	0.00
78) Terphenyl-d14	12.89	244	152415	13.44	ng	0.00
Target Compounds						
3) Pyridine	3.25	79	606204	47.199	ng	Qvalue 96
10) Phenol	6.45	94	373580	29.274	ng	99
17) 2-Methylphenol	7.06	107	51863	5.958	ng	96
20) 3+4-Methylphenols	7.22	107	192893	20.910	ng	84
27) 2,4-Dimethylphenol	7.77	122	20454	2.750	ng	# 91
31) Naphthalene	8.09	128	259842	10.071	ng	98
70) Phenanthrene	11.34	178	45078	2.460	ng	99
72) Carbazole	11.56	167	77681	4.433	ng	99

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

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