

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101718.D
 Acq On : 26 Dec 2017 16:27
 Operator : SJ/JU
 Sample : I6846-03DL 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-34972326-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:45 PM

Quant Time: Dec 26 16:58:10 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	137639	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	508529	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	206286	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	323946	20.00	ng	0.00
75) Chrysene-d12	13.97	240	267210	20.00	ng	0.00
86) Perylene-d12	15.43	264	291726	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	147131	15.92	ng	0.02
7) Phenol-d6	6.43	99	163938	14.26	ng	0.00
23) Nitrobenzene-d5	7.36	82	125736	13.76	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	32158	16.18	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	248630	18.70	ng	0.00
78) Terphenyl-d14	12.89	244	178825	16.13	ng	0.00
Target Compounds						
10) Phenol	6.45	94	1286262	98.136	ng	98
17) 2-Methylphenol	7.05	107	331847	37.115	ng	99
20) 3+4-Methylphenols	7.21	107	934847	98.668	ng	86
27) 2,4-Dimethylphenol	7.73	122	194493	26.356	ng	97
31) Naphthalene	8.09	128	1932591	75.488	ng	99
37) 2-Methylnaphthalene	8.79	142	106142	6.364	ng	98
48) Acenaphthylene	9.69	152	84315	3.813	ng	100
54) Dibenzofuran	10.04	168	43642	2.585	ng	98
57) Fluorene	10.38	166	38039	2.664	ng	99
70) Phenanthrene	11.34	178	42211	2.343	ng	99
72) Carbazole	11.56	167	84828	4.924	ng	98

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

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