

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101710.D
 Acq On : 26 Dec 2017 12:51
 Operator : SJ/JU
 Sample : I6836-03DL 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-70-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:22 PM

Quant Time: Dec 26 13:21:52 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	149912	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	594900	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	290574	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	578983	20.00	ng	0.00
75) Chrysene-d12	13.96	240	530716	20.00	ng	0.00
86) Perylene-d12	15.43	264	467321	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	127314m	12.65	ng	0.03
7) Phenol-d6	6.43	99	141703	11.31	ng	0.00
23) Nitrobenzene-d5	7.36	82	105820	9.90	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	40760	14.56	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	223058	11.91	ng	0.00
78) Terphenyl-d14	12.89	244	271728	12.34	ng	0.00
Target Compounds						
10) Phenol	6.45	94	644803	45.168	ng	99
17) 2-Methylphenol	7.06	107	104172	10.697	ng	98
20) 3+4-Methylphenols	7.21	107	348726	33.793	ng	88
27) 2,4-Dimethylphenol	7.75	122	49082	5.686	ng	97
31) Naphthalene	8.09	128	1154212	38.539	ng	100
48) Acenaphthylene	9.69	152	71775	2.305	ng	99
72) Carbazole	11.56	167	115003	3.735	ng	96

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

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