

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101716.D
 Acq On : 26 Dec 2017 15:33
 Operator : SJ/JU
 Sample : I6844-03DL 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-9974103-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:40 PM

Quant Time: Dec 26 16:05:29 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	142457	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	543164	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	227849	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	366995	20.00	ng	0.00
75) Chrysene-d12	13.97	240	290742	20.00	ng	0.00
86) Perylene-d12	15.43	264	286524	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	150676m	15.76	ng	0.02
7) Phenol-d6	6.43	99	158527	13.32	ng	0.00
23) Nitrobenzene-d5	7.36	82	113729	11.66	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	29870	13.61	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	223518	15.22	ng	0.00
78) Terphenyl-d14	12.89	244	173108	14.35	ng	0.00
Target Compounds						
10) Phenol	6.45	94	1092872	80.561	ng	99
17) 2-Methylphenol	7.05	107	240234	25.960	ng	100
20) 3+4-Methylphenols	7.21	107	768965	78.415	ng	87
27) 2,4-Dimethylphenol	7.73	122	140750	17.857	ng	98
31) Naphthalene	8.09	128	1579582	57.765	ng	99
37) 2-Methylnaphthalene	8.79	142	61213	3.436	ng	99
48) Acenaphthylene	9.69	152	72060	2.951	ng	99
72) Carbazole	11.56	167	79500	4.073	ng	99

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

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