

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101717.D
 Acq On : 26 Dec 2017 16:00
 Operator : SJ/JU
 Sample : I6845-03DL 10X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-10257-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:42 PM

Quant Time: Dec 26 16:38:24 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	141981	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	544200	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	230635	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	366653	20.00	ng	0.00
75) Chrysene-d12	13.97	240	294457	20.00	ng	0.00
86) Perylene-d12	15.43	264	282166	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	125181	13.13	ng	0.02
7) Phenol-d6	6.43	99	142789	12.04	ng	0.00
23) Nitrobenzene-d5	7.36	82	112072	11.46	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	28987	13.04	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	219398	14.76	ng	0.00
78) Terphenyl-d14	12.89	244	171450	14.04	ng	0.00
Target Compounds						
10) Phenol	6.45	94	531309	39.297	ng	98
17) 2-Methylphenol	7.05	107	264027	28.627	ng	99
20) 3+4-Methylphenols	7.21	107	688314	70.426	ng	85
27) 2,4-Dimethylphenol	7.73	122	232282	29.414	ng	99
31) Naphthalene	8.10	128	2270526	82.875	ng	99
37) 2-Methylnaphthalene	8.78	142	110070	6.166	ng	99
48) Acenaphthylene	9.69	152	154019	6.231	ng	99
54) Dibenzofuran	10.03	168	42406	2.247	ng	98
57) Fluorene	10.37	166	38083	2.385	ng	99
70) Phenanthrene	11.34	178	40865	2.004	ng	98
72) Carbazole	11.56	167	111135	5.699	ng	98

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

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