

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101707.D
 Acq On : 26 Dec 2017 11:30
 Operator : SJ/JU
 Sample : I6848-03DL 20X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-9871-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:13 PM

Quant Time: Dec 26 12:00:53 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	124580	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	505880	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	239137	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	484193	20.00	ng	0.00
75) Chrysene-d12	13.96	240	427348	20.00	ng	0.00
86) Perylene-d12	15.43	264	355854	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	68217m	8.16	ng	0.05
7) Phenol-d6	6.43	99	76590	7.36	ng	0.00
23) Nitrobenzene-d5	7.37	82	56216	6.19	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	20109	8.73	ng	0.01
44) 2-Fluorobiphenyl	9.15	172	119207	7.73	ng	0.00
78) Terphenyl-d14	12.89	244	146825	8.28	ng	0.00
Target Compounds						
3) Pyridine	3.25	79	515301	43.158	ng	Qvalue 97
10) Phenol	6.45	94	348730	29.395	ng	98
17) 2-Methylphenol	7.06	107	54938	6.789	ng	98
20) 3+4-Methylphenols	7.22	107	209006	24.372	ng	89
27) 2,4-Dimethylphenol	7.77	122	33574	4.574	ng	# 85
31) Naphthalene	8.09	128	54318	2.133	ng	# 95
72) Carbazole	11.56	167	91998	3.573	ng	96

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

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