

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101714.D
 Acq On : 26 Dec 2017 14:39
 Operator : SJ/JU
 Sample : I6842-03DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-04-01DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:34 PM

Quant Time: Dec 26 15:20:43 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	146025	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	557683	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	234396	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	381957	20.00	ng	0.00
75) Chrysene-d12	13.97	240	307113	20.00	ng	0.00
86) Perylene-d12	15.43	264	286774	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	307962	31.41	ng	0.02
7) Phenol-d6	6.43	99	363341	29.78	ng	0.00
23) Nitrobenzene-d5	7.36	82	234049	23.36	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	67746	29.99	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	448820	29.70	ng	0.00
78) Terphenyl-d14	12.89	244	352311	27.66	ng	0.00
Target Compounds						
3) Pyridine	3.30	79	305172m	21.805	ng	Qvalue
10) Phenol	6.45	94	527853	37.960	ng	98
17) 2-Methylphenol	7.05	107	125787	13.261	ng	97
20) 3+4-Methylphenols	7.21	107	316743	31.511	ng	89
27) 2,4-Dimethylphenol	7.75	122	71096	8.785	ng	90
31) Naphthalene	8.09	128	141393	5.036	ng	98
70) Phenanthrene	11.34	178	46089	2.170	ng	97
72) Carbazole	11.56	167	75998	3.741	ng	99

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

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