

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101701.D
 Acq On : 23 Dec 2017 13:59
 Operator : SJ/JU
 Sample : I6848-03
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-9871-01

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:38:13 PM

Quant Time: Dec 25 03:13:34 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 16:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	146639	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	579255	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	257658	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	426751	20.00	ng	0.00
75) Chrysene-d12	13.96	240	291671	20.00	ng	0.00
86) Perylene-d12	15.43	264	273320	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1296622	131.71	ng	0.12
7) Phenol-d6	6.48	99	1154261	94.21	ng	0.05
23) Nitrobenzene-d5	7.37	82	986340	94.79	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	323204	130.18	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1522865	91.68	ng	0.00
78) Terphenyl-d14	12.90	244	1308910	108.19	ng	0.00
Target Compounds						
3) Pyridine	3.49	79	7920665	563.584	ng	98
10) Phenol	6.52	94	4476079	320.544	ng	96
17) 2-Methylphenol	7.06	107	876486	92.013	ng	96
20) 3+4-Methylphenols	7.23	107	2713278	268.797	ng	89
27) 2,4-Dimethylphenol	7.74	122	635374	75.589	ng	99
31) Naphthalene	8.09	128	975140	33.439	ng	100
37) 2-Methylnaphthalene	8.78	142	72697	3.826	ng	98
48) Acenaphthylene	9.69	152	78619	2.847	ng	98
54) Dibenzofuran	10.03	168	53602	2.542	ng	98
57) Fluorene	10.37	166	83521	4.683	ng	100
70) Phenanthrene	11.34	178	461337	19.439	ng	98
71) Anthracene	11.39	178	86057m	3.550	ng	
72) Carbazole	11.56	167	1085675	47.834	ng	100
74) Fluoranthene	12.53	202	90327	3.626	ng	97
77) Pyrene	12.76	202	53624	2.450	ng	98

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

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