

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101706.D
 Acq On : 26 Dec 2017 11:03
 Operator : SJ/JU
 Sample : I6848-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-9871-01

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:10 PM

Quant Time: Dec 26 11:36:49 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	152232	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	600372	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	290360	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	561893	20.00	ng	0.00
75) Chrysene-d12	13.96	240	529428	20.00	ng	0.00
86) Perylene-d12	15.43	264	485344	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1344881	131.60	ng	0.12
7) Phenol-d6	6.50	99	1182768m	92.99	ng	0.06
23) Nitrobenzene-d5	7.37	82	1032706	95.75	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	411532	147.09	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1685244	90.03	ng	0.00
78) Terphenyl-d14	12.90	244	1842961	83.92	ng	0.00
Target Compounds						
3) Pyridine	3.49	79	8072471	553.283	ng	97
10) Phenol	6.52	94	4593537	316.870	ng	95
17) 2-Methylphenol	7.07	107	894420	90.446	ng	99
20) 3+4-Methylphenols	7.24	107	2784721	265.739	ng	87
27) 2,4-Dimethylphenol	7.74	122	671677	77.097	ng	98
31) Naphthalene	8.09	128	1022357	33.825	ng	99
37) 2-Methylnaphthalene	8.78	142	77835	3.953	ng	98
48) Acenaphthylene	9.69	152	89876	2.888	ng	98
54) Dibenzofuran	10.03	168	64273	2.705	ng	98
57) Fluorene	10.37	166	98372	4.894	ng	98
70) Phenanthrene	11.34	178	596174	19.079	ng	100
71) Anthracene	11.39	178	126139m	3.952	ng	
72) Carbazole	11.56	167	1444759	48.345	ng	99
74) Fluoranthene	12.53	202	149633	4.562	ng	96
77) Pyrene	12.76	202	91477	2.302	ng	100

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

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