

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101692.D
 Acq On : 23 Dec 2017 9:56
 Operator : SJ/JU
 Sample : I6839-03
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-41446854-01

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:37:56 PM

Quant Time: Dec 26 10:59:53 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	162494	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	523817	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	279869	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	486346	20.00	ng	0.00
75) Chrysene-d12	13.96	240	327905	20.00	ng	0.00
86) Perylene-d12	15.43	264	297492	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.43	112	1260344	115.54	ng	0.02
7) Phenol-d6	6.44	99	1305100	96.13	ng	0.00
23) Nitrobenzene-d5	7.36	82	997946	106.05	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	348106	129.08	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1616297	89.59	ng	0.00
78) Terphenyl-d14	12.90	244	1383048	101.68	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Pyridine	3.58	79	68486m	4.398	ng	
10) Phenol	6.46	94	1871769	120.964	ng	97
17) 2-Methylphenol	7.05	107	423270	40.099	ng	99
20) 3+4-Methylphenols	7.22	107	1389939	124.262	ng	85
27) 2,4-Dimethylphenol	7.73	122	332943	43.801	ng	99
31) Naphthalene	8.11	128	6107369	231.595	ng	98
37) 2-Methylnaphthalene	8.78	142	335102	19.504	ng	98
45) 1,1'-Biphenyl	9.25	154	78965	3.181	ng	97
48) Acenaphthylene	9.69	152	262889	8.764	ng	99
51) Acenaphthene	9.86	154	56562	3.139	ng	98
54) Dibenzofuran	10.03	168	214951	9.385	ng	99
57) Fluorene	10.37	166	174077	8.985	ng	98
70) Phenanthrene	11.34	178	241731	8.938	ng	99
72) Carbazole	11.55	167	695304	26.881	ng	99

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

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