

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101712.D
 Acq On : 26 Dec 2017 13:45
 Operator : SJ/JU
 Sample : I6838-03DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-7317-01DL

Manual Integrations
 APPROVED

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

Quant Time: Dec 26 14:38:03 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	138169	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	538263	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	236941	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	417385	20.00	ng	0.00
75) Chrysene-d12	13.97	240	326907	20.00	ng	0.00
86) Perylene-d12	15.43	264	293158	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.42	112	300184	32.36	ng	0.01
7) Phenol-d6	6.43	99	337051	29.20	ng	0.00
23) Nitrobenzene-d5	7.36	82	233224	24.12	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	83095	36.40	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	445168	29.14	ng	0.00
78) Terphenyl-d14	12.89	244	403925	29.79	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.45	94	1010202	76.778	ng	99
17) 2-Methylphenol	7.05	107	176056	19.615	ng	99
20) 3+4-Methylphenols	7.21	107	537692	56.533	ng	85
27) 2,4-Dimethylphenol	7.75	122	85965	11.006	ng	99
31) Naphthalene	8.10	128	2116659	78.111	ng	99
37) 2-Methylnaphthalene	8.79	142	69309	3.926	ng	98
48) Acenaphthylene	9.69	152	61835	2.435	ng	98
54) Dibenzofuran	10.03	168	42257	2.179	ng	97
57) Fluorene	10.37	166	33697	2.054	ng	# 97
70) Phenanthrene	11.34	178	49800	2.145	ng	97
72) Carbazole	11.56	167	154882	6.977	ng	98

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

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