

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122315\
 Data File : BM003734.D
 Acq On : 23 Dec 2015 06:56
 Operator : UM/SJ
 Sample : G4869-14
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LTCWC-04

Quant Time: Dec 23 10:35:08 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 09 18:52:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	39246	20.00	ng	-0.03
21) Naphthalene-d8	10.48	136	153089	20.00	ng	-0.04
38) Acenaphthene-d10	14.34	164	77757	20.00	ng	-0.03
63) Phenanthrene-d10	17.09	188	167958	20.00	ng	-0.03
75) Chrysene-d12	21.29	240	184117	20.00	ng	-0.03
86) Perylene-d12	23.53	264	187971	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	308274	139.65	ng	-0.02
7) Phenol-d6	6.87	99	366893	119.69	ng	-0.02
23) Nitrobenzene-d5	8.85	82	253107	102.51	ng	-0.03
41) 2,4,6-Tribromophenol	15.84	330	116849	150.47	ng	-0.03
44) 2-Fluorobiphenyl	12.96	172	578484	101.36	ng	-0.04
78) Terphenyl-d14	19.73	244	706454	90.49	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	10.53	128	592468	73.96	ng	99
37) 2-Methylnaphthalene	12.15	142	63212	11.50	ng	97
45) 1,1'-Biphenyl	13.17	154	18608	2.70	ng	99
51) Acenaphthene	14.41	154	71494	14.61	ng	98
54) Dibenzofuran	14.74	168	67755	9.59	ng	95
57) Fluorene	15.39	166	44866	7.66	ng	97
70) Phenanthrene	17.13	178	83026	8.81	ng	100
72) Carbazole	17.50	167	26362	3.23	ng	99

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

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