

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032317\
 Data File : BG026406.D
 Acq On : 23 Mar 2017 17:14
 Operator : SJ/MA
 Sample : I2364-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-COMP-10

Quant Time: Mar 24 01:33:31 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG032017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 20 17:43:27 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	144901	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	588627	20.00	ng	0.00
38) Acenaphthene-d10	14.65	164	348263	20.00	ng	0.00
63) Phenanthrene-d10	17.38	188	832576	20.00	ng	0.00
75) Chrysene-d12	21.63	240	978847	20.00	ng	-0.01
86) Perylene-d12	24.80	264	946958	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	1096684	134.33	ng	0.00
7) Phenol-d6	7.21	99	1371561	125.14	ng	0.00
23) Nitrobenzene-d5	9.20	82	836338	85.43	ng	0.00
41) 2,4,6-Tribromophenol	16.13	330	682218	171.06	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	2239390	90.17	ng	0.00
78) Terphenyl-d14	19.98	244	2907165	72.13	ng	0.00
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	10.02	122	17900	2.17	ng	84
31) Naphthalene	10.89	128	992119	33.34	ng	99
37) 2-Methylnaphthalene	12.49	142	132956	6.10	ng	# 90
48) Acenaphthylene	14.38	152	162959	4.44	ng	99
57) Fluorene	15.69	166	65170	2.30	ng	93
70) Phenanthrene	17.42	178	134114	3.00	ng	# 92
72) Carbazole	17.78	167	94071	2.00	ng	# 94

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

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