

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012318\
 Data File : BG032230.D
 Acq On : 23 Jan 2018 14:06
 Operator : SJ/JU
 Sample : J1239-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-14-01

Quant Time: Jan 24 04:22:14 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG011218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 18 10:13:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.73	152	45590	20.00	ng	-0.01
21) Naphthalene-d8	11.61	136	196064	20.00	ng	-0.02
38) Acenaphthene-d10	15.36	164	137974	20.00	ng	-0.02
63) Phenanthrene-d10	18.10	188	341518	20.00	ng	-0.02
75) Chrysene-d12	22.43	240	444333	20.00	ng	-0.02
86) Perylene-d12	26.12	264	468601	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.16	112	310597	124.60	ng	0.00
7) Phenol-d6	7.84	99	370130	117.90	ng	0.00
23) Nitrobenzene-d5	9.92	82	270676	93.40	ng	-0.02
41) 2,4,6-Tribromophenol	16.84	330	317494	139.06	ng	-0.02
44) 2-Fluorobiphenyl	13.99	172	941774	84.64	ng	-0.02
78) Terphenyl-d14	20.63	244	1764745	87.70	ng	-0.02
Target Compounds						
31) Naphthalene	11.66	128	966397	99.500	ng	99
37) 2-Methylnaphthalene	13.23	142	83115	10.779	ng	93
51) Acenaphthene	15.43	154	58887	6.356	ng	99
57) Fluorene	16.40	166	39039	3.444	ng	95
70) Phenanthrene	18.14	178	94123	4.950	ng	98
72) Carbazole	18.49	167	40659	2.471	ng	97

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

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