

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031318\
 Data File : BG033132.D
 Acq On : 14 Mar 2018 2:13
 Operator : SJ/JU
 Sample : J1874-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MHC-WCIS

Quant Time: Mar 14 04:21:24 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 15:10:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.92	152	50090	20.00	ng	0.00
21) Naphthalene-d8	11.80	136	218715	20.00	ng	0.00
38) Acenaphthene-d10	15.53	164	119754	20.00	ng	0.00
63) Phenanthrene-d10	18.26	188	255428	20.00	ng	0.00
75) Chrysene-d12	22.61	240	242227	20.00	ng	0.00
86) Perylene-d12	26.42	264	215152	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.34	112	437976	139.89	ng	0.01
7) Phenol-d6	8.02	99	530968	121.67	ng	0.02
23) Nitrobenzene-d5	10.12	82	409948	124.81	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	197800	157.09	ng	0.00
44) 2-Fluorobiphenyl	14.16	172	845343	101.81	ng	0.00
78) Terphenyl-d14	20.77	244	1157190	107.33	ng	0.00
Target Compounds						
32) Benzoic acid	11.01	122	35403	23.711	ng	90
58) 2,3,4,6-Tetrachlorophenol	16.13	232	5116	2.540	ng	# 86
69) Pentachlorophenol	17.90	266	17032	9.264	ng	89
76) Benzidine	20.41	184	35533	4.304	ng	97

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

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