

Data Path : U:\HPCHEM1\BNA G\DATA\BG122617\
 Data File : BG031789.D
 Acq On : 27 Dec 2017 9:00
 Operator : SJ/JU
 Sample : I7009-07
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121917-TIDO-F-D-B

Quant Time: Dec 27 10:09:29 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 11:08:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.80	152	42002	20.00	ng	-0.04
21) Naphthalene-d8	11.69	136	173322	20.00	ng	-0.04
38) Acenaphthene-d10	15.44	164	129333	20.00	ng	-0.03
63) Phenanthrene-d10	18.17	188	350593	20.00	ng	-0.04
75) Chrysene-d12	22.51	240	420981	20.00	ng	-0.06
86) Perylene-d12	26.26	264	446135	20.00	ng	-0.10
System Monitoring Compounds						
5) 2-Fluorophenol	6.22	112	87029	37.74	ng	-0.03
7) Phenol-d6	7.90	99	71332	23.82	ng	-0.03
23) Nitrobenzene-d5	10.00	82	207305	90.32	ng	-0.04
41) 2,4,6-Tribromophenol	16.90	330	185648	94.07	ng	-0.04
44) 2-Fluorobiphenyl	14.07	172	832974	80.84	ng	-0.03
78) Terphenyl-d14	20.69	244	1542444	83.71	ng	-0.04
Target Compounds						
10) Phenol	7.93	94	7932	2.624	ng	88
49) Dimethylphthalate	14.86	163	25363	2.226	ng	# 94

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

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