

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031887.D
 Acq On : 3 Jan 2018 12:51
 Operator : SJ/JU
 Sample : I7106-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EPE-106-2(25-27)

Quant Time: Jan 03 14:27:53 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 15:51:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.79	152	43737	20.00	ng	0.00
21) Naphthalene-d8	11.67	136	187555	20.00	ng	0.00
38) Acenaphthene-d10	15.41	164	137800	20.00	ng	-0.01
63) Phenanthrene-d10	18.15	188	387495	20.00	ng	-0.01
75) Chrysene-d12	22.49	240	448740	20.00	ng	-0.02
86) Perylene-d12	26.22	264	464101	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	277718	115.65	ng	0.00
7) Phenol-d6	7.89	99	363303	116.53	ng	0.00
23) Nitrobenzene-d5	9.98	82	192815	77.63	ng	0.00
41) 2,4,6-Tribromophenol	16.89	330	269563	128.20	ng	0.00
44) 2-Fluorobiphenyl	14.05	172	700261	63.78	ng	0.00
78) Terphenyl-d14	20.68	244	1384188	70.47	ng	-0.01
Target Compounds						
10) Phenol	7.92	94	16629	5.283	ng	90
49) Dimethylphthalate	14.84	163	54628	4.500	ng	# 97
83) Bis(2-ethylhexyl)phthalate	22.31	149	33849	2.428	ng	91

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

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