

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091219\
 Data File : BG042798.D
 Acq On : 13 Sep 2019 1:19
 Operator : HP/JU
 Sample : K4712-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-4-14-16

Quant Time: Sep 13 08:24:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:06:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	87543	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	341980	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	234796	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	543289	20.00	ng	0.00
76) Chrysene-d12	21.75	240	558951	20.00	ng	-0.01
87) Perylene-d12	25.02	264	611783	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	335619	66.65	ng	0.00
7) Phenol-d6	7.26	99	317907	43.99	ng	0.00
23) Nitrobenzene-d5	9.28	82	615679	93.73	ng	0.00
42) 2,4,6-Tribromophenol	16.22	330	432840	138.56	ng	0.00
45) 2-Fluorobiphenyl	13.37	172	1336954	90.13	ng	0.00
79) Terphenyl-d14	20.09	244	2273484	90.15	ng	0.00
Target Compounds						
10) Phenol	7.29	94	20109	2.713	ng	93
50) Dimethylphthalate	14.18	163	152960	8.696	ng	# 98
52) Acenaphthene	14.80	154	29348	2.162	ng	99

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

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