

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102919\
 Data File : BG043378.D
 Acq On : 29 Oct 2019 17:40
 Operator : JU
 Sample : K5564-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CSA-2-1-1

Quant Time: Oct 30 08:21:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 21 16:37:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	45389	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	177004	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	127943	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	284990	20.00	ng	-0.01
76) Chrysene-d12	21.66	240	307214	20.00	ng	-0.01
87) Perylene-d12	24.85	264	371923	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	303079	122.36	ng	0.00
7) Phenol-d6	7.21	99	436657	122.53	ng	0.00
23) Nitrobenzene-d5	9.19	82	268036	78.07	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	265338	108.98	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	728019	78.63	ng	0.00
79) Terphenyl-d14	20.00	244	1300339	79.41	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.10	163	112420	11.344	ng	97
75) Fluoranthene	19.45	202	45747	2.404	ng	98
78) Pyrene	19.81	202	41637	2.190	ng	96
80) Butylbenzylphthalate	20.69	149	60532	8.402	ng	99

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

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