

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121119\
 Data File : BG043811.D
 Acq On : 12 Dec 2019 1:19
 Operator : JU
 Sample : K6245-11
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RELAY-FLOOR

Quant Time: Dec 12 03:54:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 09 18:16:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	192661	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	799438	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	513918	20.00	ng	0.00
64) Phenanthrene-d10	17.37	188	972196	20.00	ng	0.00
76) Chrysene-d12	21.63	240	798849	20.00	ng	0.00
87) Perylene-d12	24.79	264	924012	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	958995	94.31	ng	0.00
7) Phenol-d6	7.17	99	1413022	96.36	ng	0.01
23) Nitrobenzene-d5	9.16	82	891727	62.63	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	427799	87.32	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	1824273	62.35	ng	0.00
79) Terphenyl-d14	19.99	244	2132196	64.21	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.08	163	92190	2.513	ng	99
71) Phenanthrene	17.41	178	130288	2.831	ng	98
84) Bis(2-ethylhexyl)phthalate	21.55	149	138679	4.447	ng	98

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

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