

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010320\
 Data File : BG043898.D
 Acq On : 3 Jan 2020 21:52
 Operator : JU
 Sample : K6449-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DUPLICATE

Quant Time: Jan 04 05:24:06 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	129817	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	570051	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	378897	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	817239	20.00	ng	0.00
76) Chrysene-d12	21.58	240	687020	20.00	ng	0.00
87) Perylene-d12	24.72	264	797010	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	725634	102.63	ng	0.00
7) Phenol-d6	7.12	99	713844	72.23	ng	0.00
23) Nitrobenzene-d5	9.12	82	962669	90.34	ng	0.00
42) 2,4,6-Tribromophenol	16.08	330	623489	157.24	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	1981454	94.75	ng	0.00
79) Terphenyl-d14	19.95	244	2691664	98.34	ng	0.00
Target Compounds						
10) Phenol	7.15	94	39810	3.910	ng	96
50) Dimethylphthalate	14.05	163	236091	8.776	ng	96
84) Bis(2-ethylhexyl)phthalate	21.50	149	140132	4.757	ng	93

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

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