

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110419\
 Data File : BG043495.D
 Acq On : 5 Nov 2019 00:15
 Operator : JU
 Sample : PB124488BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB124488BL

Quant Time: Nov 05 17:32:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	39017	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	138018	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	103592	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	249767	20.00	ng	0.00
76) Chrysene-d12	21.66	240	276278	20.00	ng	0.00
87) Perylene-d12	24.85	264	327740	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	217792	102.29	ng	0.00
7) Phenol-d6	7.20	99	192543	62.85	ng	0.00
23) Nitrobenzene-d5	9.18	82	251875	94.08	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	263987	133.91	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	672750	89.74	ng	0.00
79) Terphenyl-d14	20.00	244	1373731	93.28	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propylamine	9.17	70	35546	17.957	ng	# 75
51) 2,6-Dinitrotoluene	14.64	165	13900	7.974	ng	# 26
57) 2,4-Dinitrotoluene	14.64	165	13900	5.580	ng	# 29
67) 4-Bromophenyl-phenylether	16.13	248	16128	4.798	ng	# 1
77) Benzidine	20.00	184	21438	3.856	ng	# 90

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

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