

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082718\
 Data File : BF108529.D
 Acq On : 28 Aug 2018 00:34
 Operator : JU/SJ
 Sample : J4659-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Aug 28 03:09:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 22 11:01:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	105683	20.00	ng	0.00
21) Naphthalene-d8	8.28	136	370095	20.00	ng	0.00
38) Acenaphthene-d10	10.04	164	151582	20.00	ng	-0.01
63) Phenanthrene-d10	11.54	188	265389	20.00	ng	-0.01
75) Chrysene-d12	14.19	240	267951	20.00	ng	-0.01
86) Perylene-d12	15.74	264	185018	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	221864	38.01	ng	0.00
7) Phenol-d6	6.61	99	174315	22.47	ng	-0.02
23) Nitrobenzene-d5	7.56	82	687982	103.73	ng	-0.01
41) 2,4,6-Tribromophenol	10.84	330	137480	90.93	ng	-0.01
44) 2-Fluorobiphenyl	9.36	172	1226447	129.90	ng	0.00
78) Terphenyl-d14	13.12	244	1015744	96.38	ng	0.00
Target Compounds						
19) n-Nitroso-di-n-propylamine	7.56	70	92356	17.606	ng	# 83
25) Isophorone	7.56	82	687975	54.422	ng	# 64
49) Dimethylphthalate	9.75	163	35004	3.315	ng	# 99
50) 2,6-Dinitrotoluene	10.04	165	19278	8.726	ng	# 25
56) 2,4-Dinitrotoluene	10.04	165	19278	6.779	ng	# 23
64) 4,6-Dinitro-2-methylphenol	10.84	198	121	5.466	ng	# 1
66) 4-Bromophenyl-phenylether	10.84	248	8919	3.093	ng	# 1

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

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