

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102116\
 Data File : BF090714.D
 Acq On : 21 Oct 2016 15:46
 Operator : UM/SJ
 Sample : PB94239BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB94239BL

Quant Time: Oct 21 20:15:38 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 21 17:36:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	272783	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	1165353	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	577949	20.00	ng	-0.01
63) Phenanthrene-d10	11.38	188	960361	20.00	ng	-0.01
75) Chrysene-d12	14.02	240	815079	20.00	ng	0.00
86) Perylene-d12	15.46	264	600953	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	1898389	112.24	ng	0.02
7) Phenol-d6	6.50	99	2450717	106.97	ng	0.00
23) Nitrobenzene-d5	7.42	82	1667151	78.46	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	656780	143.58	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	2475059	66.48	ng	-0.01
78) Terphenyl-d14	12.98	244	2791725	78.50	ng	0.01
Target Compounds						
19) n-Nitroso-di-n-propylamine	7.42	70	222312	13.03	ng	# 50
50) 2,6-Dinitrotoluene	9.89	165	74913	9.59	ng	# 39
56) 2,4-Dinitrotoluene	9.89	165	75820	7.94	ng	# 33
64) 4,6-Dinitro-2-methylphenol	10.68	198	937	4.32	ng	# 1
66) 4-Bromophenyl-phenylether	10.68	248	40123	4.31	ng	# 1

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

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