

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102218\
 Data File : BF110059.D
 Acq On : 22 Oct 2018 19:53
 Operator : JU/SJ
 Sample : J5604-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Oct 23 04:21:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	192103	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	711151	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	278733	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	436396	20.00	ng	-0.01
76) Chrysene-d12	14.14	240	350754	20.00	ng	-0.02
87) Perylene-d12	15.67	264	245953	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1167558	96.34	ng	0.00
7) Phenol-d6	6.59	99	1451902	96.48	ng	-0.01
23) Nitrobenzene-d5	7.53	82	826484	78.29	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	235773	97.70	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	1160881	66.46	ng	-0.01
79) Terphenyl-d14	13.09	244	864280	51.74	ng	-0.02
Target Compounds						
10) Phenol	6.60	94	127101	7.823	ng	# 63
20) 3+4-Methylphenols	7.36	107	40411	2.909	ng	# 76
50) Dimethylphthalate	9.72	163	169160	8.571	ng	100
78) Pyrene	12.95	202	53997	2.096	ng	97

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

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