

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
 Data File : BF106981.D
 Acq On : 29 Jun 2018 18:54
 Operator : JU/SJ
 Sample : J3629-16
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A0C22-02

Quant Time: Jun 29 23:50:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	69012	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	256830	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	115057	20.00	ng	-0.01
63) Phenanthrene-d10	11.48	188	190830	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	187899	20.00	ng	-0.02
86) Perylene-d12	15.68	264	145582	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	429754	105.45	ng	0.00
7) Phenol-d6	6.60	99	525266	103.20	ng	-0.01
23) Nitrobenzene-d5	7.52	82	320671	69.39	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	138448	103.82	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	599850	88.93	ng	-0.02
78) Terphenyl-d14	13.07	244	570179	73.50	ng	-0.01
Target Compounds						
22) Acetophenone	7.35	105	13702	2.385	ng	# 98
49) Dimethylphthalate	9.70	163	87046	10.087	ng	# 99
83) Bis(2-ethylhexyl)phthalate	14.09	149	81503	12.514	ng	99

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

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