

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121518\
 Data File : BF111483.D
 Acq On : 15 Dec 2018 22:01
 Operator : JU/SJ
 Sample : J6373-10 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-WATER

Quant Time: Dec 16 02:01:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	131276	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	462850	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	183411	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	322712	20.00	ng	0.00
76) Chrysene-d12	13.95	240	317093	20.00	ng	0.00
87) Perylene-d12	15.39	264	219996	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	42892	5.44	ng	0.00
7) Phenol-d6	6.43	99	32683	3.11	ng	-0.01
23) Nitrobenzene-d5	7.35	82	61978	7.97	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	18133	11.04	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	117654	9.91	ng	0.00
79) Terphenyl-d14	12.90	244	115509	8.55	ng	0.00
Target Compounds						
10) Phenol	6.45	94	233232	19.952	ng	90
22) Acetophenone	7.20	105	32162	3.113	ng	# 82

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

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