

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100518\
 Data File : BF109617.D
 Acq On : 5 Oct 2018 19:11
 Operator : JU/SJ
 Sample : J5244-04
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-WATER

Quant Time: Oct 06 06:32:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	97963	20.00	ng	-0.01
21) Naphthalene-d8	7.98	136	371756	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	179368	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	333339	20.00	ng	0.00
75) Chrysene-d12	13.83	240	208746	20.00	ng	-0.01
86) Perylene-d12	15.23	264	220124	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	259994	43.71	ng	0.00
7) Phenol-d6	6.35	99	207374	27.32	ng	0.00
23) Nitrobenzene-d5	7.26	82	485485	77.09	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	191632	108.38	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	879700	73.97	ng	-0.02
78) Terphenyl-d14	12.79	244	746340	87.74	ng	-0.01
Target Compounds						
10) Phenol	6.36	94	22571	2.626	ng	85
49) Dimethylphthalate	9.45	163	167422	12.833	ng	99

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

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