

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110817\
 Data File : BF100300.D
 Acq On : 8 Nov 2017 19:10
 Operator : SJ/JU
 Sample : I6358-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-WATER-FROM-BERM

Quant Time: Nov 09 00:24:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 08 14:33:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	202838	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	798659	20.00	ng	0.00
38) Acenaphthene-d10	9.94	164	360952	20.00	ng	0.00
63) Phenanthrene-d10	11.43	188	590128	20.00	ng	0.00
75) Chrysene-d12	14.06	240	394256	20.00	ng	0.00
86) Perylene-d12	15.54	264	410129	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	702598	55.52	ng	0.00
7) Phenol-d6	6.52	99	553054	35.44	ng	0.00
23) Nitrobenzene-d5	7.46	82	1073194	88.84	ng	0.00
41) 2,4,6-Tribromophenol	10.73	330	393717	107.04	ng	0.00
44) 2-Fluorobiphenyl	9.26	172	1853260	78.18	ng	0.00
78) Terphenyl-d14	13.02	244	1415008	82.47	ng	0.00
Target Compounds						
10) Phenol	6.53	94	38066	2.324	ng	98
49) Dimethylphthalate	9.65	163	188254	6.831	ng	99
83) Bis(2-ethylhexyl)phthalate	14.04	149	168902	11.205	ng	99

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

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