

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
 Data File : BF106457.D
 Acq On : 14 Jun 2018 21:58
 Operator : JU/SJ
 Sample : J3469-01 10X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OILY-WATER

Quant Time: Jun 15 03:35:04 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 14 18:39:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	117939	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	454479	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	186070	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	320850	20.00	ng	0.00
75) Chrysene-d12	14.15	240	289303	20.00	ng	0.00
86) Perylene-d12	15.69	264	210018	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	39164	5.62	ng	0.00
7) Phenol-d6	6.62	99	35836	4.07	ng	0.00
23) Nitrobenzene-d5	7.52	82	61539	7.97	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	18982	10.60	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	118695	2.15	ng	0.00
78) Terphenyl-d14	13.08	244	96701	8.30	ng	0.00
Target Compounds						
15) Benzyl Alcohol	7.10	79	71084	9.428	ng	99
20) 3+4-Methylphenols	7.38	107	88614	10.364	ng	86
32) Benzoic acid	8.10	122	113859	27.785	ng	94
83) Bis(2-ethylhexyl)phthalate	14.11	149	47880	4.925	ng	95

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

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061418\
Data File : BF106457.D
Acq On : 14 Jun 2018 21:58
Operator : JU/SJ
Sample : J3469-01 10X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OILY-WATER

Quant Time: Jun 15 03:35:04 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 14 18:39:24 2018
Response via : Initial Calibration

