

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113527.D
 Acq On : 3 Apr 2019 2:45
 Operator : JU/SJ
 Sample : K2170-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OIL-WATER-SEPERATOR-2

Quant Time: Apr 03 08:20:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	129909	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	446969	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	185301	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	352058	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	330180	20.00	ng	-0.01
87) Perylene-d12	15.24	264	238705	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	789450	99.37	ng	0.00
7) Phenol-d6	6.35	99	907861	90.33	ng	-0.01
23) Nitrobenzene-d5	7.26	82	678502	90.10	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	269462	117.73	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1158557	101.40	ng	-0.02
79) Terphenyl-d14	12.79	244	1087528	72.60	ng	-0.01
Target Compounds						
10) Phenol	6.36	94	47989	4.182	ng	91
15) Benzyl Alcohol	6.86	79	1024084	154.447	ng	97
19) n-Nitroso-di-n-propylamine	7.11	70	34761	5.506	ng	# 77
20) 3+4-Methylphenols	7.12	107	151365	12.882	ng	# 81
32) Benzoic acid	7.92	122	1432975	297.939	ng	100

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

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