

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062719\
 Data File : BF115246.D
 Acq On : 27 Jun 2019 20:55
 Operator : HP/JU
 Sample : K3529-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Jun 28 06:57:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 28 05:26:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	221786	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	768055	20.00	ng	0.00
39) Acenaphthene-d10	9.62	164	377126	20.00	ng	0.00
64) Phenanthrene-d10	11.09	188	751807	20.00	ng	0.00
76) Chrysene-d12	13.71	240	658398	20.00	ng	-0.01
87) Perylene-d12	15.07	264	705665	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	1443383	100.43	ng	0.00
7) Phenol-d6	6.24	99	1743231	99.93	ng	0.00
23) Nitrobenzene-d5	7.16	82	1200948	96.19	ng	0.00
42) 2,4,6-Tribromophenol	10.40	330	447197	112.45	ng	0.00
45) 2-Fluorobiphenyl	8.95	172	2064581	92.99	ng	0.00
79) Terphenyl-d14	12.68	244	2404459	77.65	ng	0.00
Target Compounds						
10) Phenol	6.25	94	425534	20.825	ng	86
20) 3+4-Methylphenols	7.01	107	374904	24.340	ng	# 59
50) Dimethylphthalate	9.35	163	64637	2.330	ng	96

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

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