

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061318\
 Data File : BF106402.D
 Acq On : 13 Jun 2018 20:15
 Operator : JU/SJ
 Sample : J3451-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VB-44949-SLUDGEBOX

Quant Time: Jun 14 03:46:36 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 13 10:51:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	145722	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	571592	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	271107	20.00	ng	-0.01
63) Phenanthrene-d10	11.50	188	488399	20.00	ng	0.00
75) Chrysene-d12	14.15	240	368506	20.00	ng	-0.01
86) Perylene-d12	15.68	264	368580	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	289933	33.67	ng	0.00
7) Phenol-d6	6.62	99	227948	20.96	ng	0.00
23) Nitrobenzene-d5	7.53	82	555697	57.19	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	222396	85.26	ng	-0.01
44) 2-Fluorobiphenyl	9.32	172	990509	58.19	ng	0.00
78) Terphenyl-d14	13.08	244	862672	58.15	ng	0.00
Target Compounds						
10) Phenol	6.64	94	55174	4.646	ng	99
20) 3+4-Methylphenols	7.39	107	52240	4.945	ng	85
49) Dimethylphthalate	9.71	163	150267	7.758	ng	99

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

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