

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121619\
 Data File : BF118217.D
 Acq On : 17 Dec 2019 00:54
 Operator : JU
 Sample : K6267-02
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WO-4

Quant Time: Dec 17 04:44:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 17:34:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	99178	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	383793	20.00	ng	0.00
39) Acenaphthene-d10	9.91	164	191172	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	312552	20.00	ng	0.00
76) Chrysene-d12	14.06	240	215252	20.00	ng	-0.01
87) Perylene-d12	15.55	264	264789	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	575274	101.59	ng	0.00
7) Phenol-d6	6.49	99	726336	106.05	ng	0.00
23) Nitrobenzene-d5	7.43	82	425809	62.42	ng	-0.01
42) 2,4,6-Tribromophenol	10.70	330	206177	88.78	ng	-0.01
45) 2-Fluorobiphenyl	9.23	172	881585	70.17	ng	-0.01
79) Terphenyl-d14	12.99	244	760712	60.77	ng	-0.01
Target Compounds						
10) Phenol	6.50	94	24732	3.166	ng	85
50) Dimethylphthalate	9.62	163	37012	2.621	ng	98
78) Pyrene	12.85	202	37166	2.191	ng	97
84) Bis(2-ethylhexyl)phthalate	14.03	149	1474173	145.529	ng	99

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

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