

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111619.D
 Acq On : 20 Dec 2018 4:36
 Operator : JU/SJ
 Sample : J6386-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS006-1218-01

Quant Time: Dec 20 08:01:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	184661	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	643575	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	303664	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	557230	20.00	ng	0.00
76) Chrysene-d12	14.05	240	499996	20.00	ng	0.00
87) Perylene-d12	15.53	264	341663	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	964640	89.04	ng	0.00
7) Phenol-d6	6.52	99	1211075	87.84	ng	0.00
23) Nitrobenzene-d5	7.45	82	719429	65.07	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	307828	85.79	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1246620	59.84	ng	0.00
79) Terphenyl-d14	13.00	244	1288732	48.88	ng	0.00
Target Compounds						
10) Phenol	6.53	94	73665	4.735	ng	95
50) Dimethylphthalate	9.64	163	109091	4.913	ng	98
71) Phenanthrene	11.44	178	62990	2.131	ng	95
78) Pyrene	12.85	202	85311	2.416	ng	96

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

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