

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010419\
 Data File : BF111866.D
 Acq On : 5 Jan 2019 1:03
 Operator : JU/SJ
 Sample : K1036-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 13-N

Quant Time: Jan 05 05:29:28 2019
 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	6.89	152	144678	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	533980	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	220756	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	407846	20.00	ng	0.00
76) Chrysene-d12	14.06	240	327362	20.00	ng	0.01
87) Perylene-d12	15.54	264	252445	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	865424	101.96	ng	0.02
7) Phenol-d6	6.54	99	1021061	94.52	ng	0.02
23) Nitrobenzene-d5	7.45	82	636640	69.40	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	267681	102.62	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	1157951	76.46	ng	0.00
79) Terphenyl-d14	13.00	244	918672	53.22	ng	0.00
Target Compounds						
10) Phenol	6.55	94	30176	2.476	ng	99
50) Dimethylphthalate	9.64	163	74192	4.596	ng	98
71) Phenanthrene	11.44	178	44958	2.079	ng	95
75) Fluoranthene	12.63	202	47423	2.165	ng	97

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

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