

Data Path : U:\HPCHEM1\BNA F\DATA\BF010518\
 Data File : BF101933.D
 Acq On : 5 Jan 2018 16:48
 Operator : SJ/JU
 Sample : J1032-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OWS-4

Quant Time: Jan 05 17:48:07 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 13:21:45 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	127273	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	467458	20.00	ng	-0.01
38) Acenaphthene-d10	9.81	164	159935	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	256834	20.00	ng	-0.01
75) Chrysene-d12	13.96	240	271310	20.00	ng	-0.02
86) Perylene-d12	15.43	264	217513	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1148146	134.38	ng	0.01
7) Phenol-d6	6.43	99	1213477	114.12	ng	0.00
23) Nitrobenzene-d5	7.35	82	875539	104.26	ng	-0.02
41) 2,4,6-Tribromophenol	10.61	330	178596	115.89	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	1123379	108.96	ng	-0.01
78) Terphenyl-d14	12.88	244	901234	80.08	ng	-0.02
Target Compounds						
20) 3+4-Methylphenols	7.25	107	47035	5.369	ng	89
32) Benzoic acid	7.89	122	119073	29.692	ng	96
49) Dimethylphthalate	9.53	163	66712	5.186	ng	99

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

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