

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101512.D
 Acq On : 19 Dec 2017 6:58
 Operator : SJ/JU
 Sample : I6910-04
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121317-TIDO-E-U-M

Quant Time: Dec 19 07:47:41 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:38:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	134834	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	491852	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	179121	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	276431	20.00	ng	0.00
75) Chrysene-d12	13.99	240	307848	20.00	ng	0.00
86) Perylene-d12	15.46	264	283737	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	388615	42.93	ng	0.01
7) Phenol-d6	6.45	99	276461	24.54	ng	0.00
23) Nitrobenzene-d5	7.38	82	795154	89.99	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	148660	86.13	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1250654	108.31	ng	0.00
78) Terphenyl-d14	12.92	244	966555	75.69	ng	0.00
Target Compounds						
51) Acenaphthene	9.88	154	23077	2.001	ng	95
70) Phenanthrene	11.36	178	60716	3.950	ng	97

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

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