

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122118\
 Data File : BM018355.D
 Acq On : 21 Dec 2018 11:53
 Operator : JU/SJ
 Sample : J6389-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS014-1218-01

Quant Time: Dec 21 13:04:52 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	116161	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	456488	20.00	ng	-0.02
39) Acenaphthene-d10	14.15	164	264413	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	604418	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	653886	20.00	ng	0.00
87) Perylene-d12	23.30	264	561648	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	913233	131.33	ng	0.00
7) Phenol-d6	6.70	99	1258645	130.23	ng	0.00
23) Nitrobenzene-d5	8.65	82	810209	91.04	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	488862	125.97	ng	0.00
45) 2-Fluorobiphenyl	12.75	172	1543053	75.59	ng	-0.02
79) Terphenyl-d14	19.57	244	2428609	84.00	ng	0.00
Target Compounds						
6) Aniline	6.84	93	45536	3.756	ng	91
10) Phenol	6.73	94	104109	10.171	ng	96
50) Dimethylphthalate	13.62	163	77581	3.529	ng	100
71) Phenanthrene	16.95	178	160941	4.902	ng	97
75) Fluoranthene	18.99	202	95113	2.492	ng	96
78) Pyrene	19.36	202	162991	4.183	ng	99

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

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