

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
 Data File : BM017861.D
 Acq On : 01 Dec 2018 17:46
 Operator : JU/SJ
 Sample : J6060-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 20180945-COMPOSITE-4

Quant Time: Dec 03 00:50:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 12 18:37:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	225387	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	920371	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	527874	20.00	ng	-0.04
64) Phenanthrene-d10	16.99	188	1139175	20.00	ng	-0.03
76) Chrysene-d12	21.19	240	927411	20.00	ng	-0.03
87) Perylene-d12	23.37	264	872186	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	1676481	125.53	ng	-0.03
7) Phenol-d6	6.77	99	2268961	121.64	ng	-0.03
23) Nitrobenzene-d5	8.74	82	1416447	79.45	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	863956	113.91	ng	-0.03
45) 2-Fluorobiphenyl	12.84	172	2964795	72.86	ng	-0.04
79) Terphenyl-d14	19.63	244	3355844	82.01	ng	-0.02
Target Compounds						
10) Phenol	6.80	94	120070	6.133	ng	95
50) Dimethylphthalate	13.70	163	390149	9.091	ng	98
71) Phenanthrene	17.03	178	236612	3.828	ng	99
75) Fluoranthene	19.06	202	269088	3.847	ng	97
78) Pyrene	19.42	202	293450	5.128	ng	98
81) Benzo(a)anthracene	21.17	228	107302	2.008	ng	96
83) Chrysene	21.23	228	115469	2.225	ng	99

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

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