

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120418\
 Data File : BM017884.D
 Acq On : 03 Dec 2018 21:17
 Operator : JU/SJ
 Sample : J6149-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ENGINE-WASH

Quant Time: Dec 04 07:23:38 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	190752	20.00	ng	-0.04
21) Naphthalene-d8	10.35	136	818607	20.00	ng	-0.04
39) Acenaphthene-d10	14.23	164	479499	20.00	ng	-0.04
64) Phenanthrene-d10	16.98	188	1062228	20.00	ng	-0.04
76) Chrysene-d12	21.19	240	912684	20.00	ng	-0.03
87) Perylene-d12	23.37	264	820257	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	1171136	103.61	ng	-0.03
7) Phenol-d6	6.77	99	1391910	88.17	ng	-0.03
23) Nitrobenzene-d5	8.73	82	1189466	75.01	ng	-0.04
42) 2,4,6-Tribromophenol	15.73	330	887875	128.87	ng	-0.04
45) 2-Fluorobiphenyl	12.84	172	3045205	82.39	ng	-0.04
79) Terphenyl-d14	19.63	244	3984209	98.94	ng	-0.02
Target Compounds						
10) Phenol	6.80	94	54659	3.299	ng	93
20) 3+4-Methylphenols	8.36	107	67567	4.357	ng	92
32) Benzoic acid	9.71	122	64212	6.637	ng	96

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

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