

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM101818\
 Data File : BM017224.D
 Acq On : 19 Oct 2018 10:56
 Operator : MJ/SJ
 Sample : J5204-08
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SU-03-101718-B

Quant Time: Oct 19 12:25:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	300387	20.00	ng	0.00
21) Naphthalene-d8	10.44	136	1146253	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	676347	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	1598026	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1590734	20.00	ng	-0.01
87) Perylene-d12	23.47	264	1430109	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	1949064	109.74	ng	0.00
7) Phenol-d6	6.85	99	2273597	93.99	ng	0.00
23) Nitrobenzene-d5	8.82	82	1818500	92.79	ng	-0.01
42) 2,4,6-Tribromophenol	15.80	330	1274790	139.37	ng	-0.01
45) 2-Fluorobiphenyl	12.93	172	4328472	85.15	ng	0.00
79) Terphenyl-d14	19.70	244	6689308	94.78	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.49	128	698490	12.435	ng	99
37) 2-Methylnaphthalene	12.11	142	138376	3.600	ng	98
38) 1-Methylnaphthalene	12.33	142	258923	6.921	ng	# 99
52) Acenaphthene	14.37	154	131296	3.222	ng	99
58) Fluorene	15.36	166	174357	3.476	ng	99
71) Phenanthrene	17.10	178	398513	4.622	ng	100

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

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