

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031519\
 Data File : BM019200.D
 Acq On : 15 Mar 2019 23:34
 Operator : JU/SJ
 Sample : K1911-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RBR200051

Quant Time: Mar 16 01:39:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM031119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 17:30:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	190651	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	816373	20.00	ng	-0.01
39) Acenaphthene-d10	14.30	164	524925	20.00	ng	-0.02
64) Phenanthrene-d10	17.06	188	1185327	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	938160	20.00	ng	-0.01
87) Perylene-d12	23.50	264	826827	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1580306	134.24	ng	0.00
7) Phenol-d6	6.84	99	2091990	121.49	ng	0.00
23) Nitrobenzene-d5	8.82	82	1656053	95.89	ng	-0.01
42) 2,4,6-Tribromophenol	15.81	330	1153623	148.85	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	3840689	95.17	ng	-0.01
79) Terphenyl-d14	19.70	244	6132873	129.78	ng	0.00
Target Compounds						
31) Naphthalene	10.49	128	245761	5.712	ng	98
52) Acenaphthene	14.37	154	107141	3.223	ng	99
58) Fluorene	15.36	166	105944	2.397	ng	99
71) Phenanthrene	17.10	178	267643	3.852	ng	99

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

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