

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031419\
 Data File : BM019179.D
 Acq On : 14 Mar 2019 21:07
 Operator : JU/SJ
 Sample : K1899-03DL 25X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-4ADL

Quant Time: Mar 15 06:03:44 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	147927	20.00	ng	-0.01
21) Naphthalene-d8	10.44	136	722996	20.00	ng	-0.01
39) Acenaphthene-d10	14.31	164	469339	20.00	ng	-0.01
64) Phenanthrene-d10	17.06	188	1116744	20.00	ng	-0.01
76) Chrysene-d12	21.26	240	1156182	20.00	ng	-0.01
87) Perylene-d12	23.50	264	1057254	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	23917	2.62	ng	0.00
7) Phenol-d6	6.85	99	19672	1.47	ng	0.00
23) Nitrobenzene-d5	8.83	82	54029	3.53	ng	0.00
42) 2,4,6-Tribromophenol	15.81	330	40764	5.88	ng	-0.01
45) 2-Fluorobiphenyl	12.92	172	145160	4.02	ng	-0.01
79) Terphenyl-d14	19.70	244	241003	4.14	ng	-0.01
Target Compounds						
31) Naphthalene	10.49	128	3723421	97.709	ng	100
37) 2-Methylnaphthalene	12.11	142	960064	34.670	ng	99
38) 1-Methylnaphthalene	12.33	142	1000939	36.672	ng	100
46) 1,1'-Biphenyl	13.13	154	180405	4.401	ng	97
52) Acenaphthene	14.37	154	286209	9.629	ng	98
55) Dibenzofuran	14.71	168	276200	5.762	ng	99
58) Fluorene	15.36	166	165202	4.181	ng	99
71) Phenanthrene	17.10	178	141386	2.160	ng	100

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

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