

Data Path : Z:\HPCHEM1\BNA M\Data\BM060217\
 Data File : BM010382.D
 Acq On : 02 Jun 2017 12:02
 Operator : SJ/MA
 Sample : I3163-09MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSF7MSD

Quant Time: Jun 02 18:12:00 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM053017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 02 17:41:09 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	683	0.40	ng/ul	0.00
2) Naphthalene-d8	10.73	136	2110	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1160	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2548	0.40	ng/ul	0.00
16) Chrysene-d12	21.46	240	2597	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	147102	0.40	ng/ul	-0.15
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	686	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	1756	0.26	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.61	153	1053	0.24	ng/ul	96
9) Fluorene	15.55	166	254	0.05	ng/ul#	33
11) Pentachlorophenol	16.95	266	233	0.24	ng/ul#	82
12) Phenanthrene	17.33	178	211	0.03	ng/ul#	40
15) Fluoranthene	19.35	202	384	0.04	ng/ul#	24
17) Pyrene	19.71	202	3500	0.32	ng/ul	96
19) Chrysene	21.50	228	247	0.02	ng/ul#	23

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

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