

Data Path : Z:\HPCHEM1\BNA M\Data\BM060217\
 Data File : BM010385.D
 Acq On : 02 Jun 2017 13:50
 Operator : SJ/MA
 Sample : I3164-03MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHSGMSD

Quant Time: Jun 02 18:21:28 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	7.91	152	782	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.72	136	2397	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.55	164	1280	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2820	0.40	ng/ul	0.00
16) Chrysene-d12	21.46	240	2980	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	146916	0.40	ng/ul	-0.15
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	682	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	1746	0.23	ng/ul	0.00
Target Compounds						
					Ovalue	
3) Naphthalene	10.77	128	168	0.03	ng/ul#	18
8) Acenaphthene	14.61	153	104477	21.62	ng/ul	95
9) Fluorene	15.55	166	249	0.04	ng/ul#	34
11) Pentachlorophenol	16.94	266	29578	27.43	ng/ul	97
12) Phenanthrene	17.33	178	268	0.03	ng/ul#	41
15) Fluoranthene	19.35	202	355	0.03	ng/ul#	27
17) Pyrene	19.71	202	274472	21.62	ng/ul	95
18) Benzo(a)anthracene	21.43	228	416	0.03	ng/ul#	30
19) Chrysene	21.49	228	478	0.04	ng/ul#	27

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

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