

Data Path : Z:\HPCHEM1\BNA M\DATA\BM053017\
 Data File : BM010303.D
 Acq On : 30 May 2017 10:01
 Operator : SJ/MA
 Sample : SSTD0.208
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: May 30 13:02:53 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM053017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 30 13:00:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	842	0.40	ng/ul	0.00
2) Naphthalene-d8	10.74	136	2506	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.55	164	1362	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2996	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	2521	0.40	ng/ul	0.00
20) Perylene-d12	23.82	264	2641	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	708	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	1437	0.16	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.78	128	1311	0.20	ng/ul#	91
5) 2-Methylnaphthalene	12.38	142	907	0.19	ng/ul	99
7) Acenaphthylene	14.28	152	1331	0.20	ng/ul	95
8) Acenaphthene	14.62	153	1016	0.19	ng/ul	93
9) Fluorene	15.62	166	1196	0.19	ng/ul	96
11) Pentachlorophenol	16.95	266	211	0.18	ng/ul	89
12) Phenanthrene	17.35	178	1850	0.21	ng/ul#	93
13) Anthracene	17.43	178	1724	0.19	ng/ul	98
15) Fluoranthene	19.35	202	2154	0.18	ng/ul	98
17) Pyrene	19.73	202	2211	0.26	ng/ul#	94
18) Benzo(a)anthracene	21.45	228	2015	0.24	ng/ul	95
19) Chrysene	21.45	228	2015	0.25	ng/ul	96
21) Benzo(b)fluoranthene	23.10	252	2294	0.25	ng/ul	70
22) Benzo(k)fluoranthene	23.14	252	2185	0.25	ng/ul#	69
23) Benzo(a)pyrene	23.70	252	2188	0.25	ng/ul#	70
24) Indeno(1,2,3-cd)pyrene	26.21	276	2821	0.28	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.22	278	2238	0.27	ng/ul#	83
26) Benzo(a,h,i)perylene	26.95	276	2481	0.29	ng/ul#	93

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

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