

Data Path : Z:\HPCHEM1\BNA M\DATA\BM053017\
 Data File : BM010304.D
 Acq On : 30 May 2017 10:37
 Operator : SJ/MA
 Sample : SSTD0.409
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.409

Quant Time: May 30 13:04:15 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	728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.73	136	2037	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1085	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2492	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	2507	0.40	ng/ul	0.00
20) Perylene-d12	23.82	264	2751	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	1180	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	2691	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.78	128	2118	0.39	ng/ul	97
5) 2-Methylnaphthalene	12.38	142	1495	0.38	ng/ul	100
7) Acenaphthylene	14.28	152	2178	0.41	ng/ul	98
8) Acenaphthene	14.62	153	1632	0.39	ng/ul	93
9) Fluorene	15.62	166	1960	0.39	ng/ul	98
11) Pentachlorophenol	16.95	266	381	0.39	ng/ul	94
12) Phenanthrene	17.35	178	3168	0.43	ng/ul	93
13) Anthracene	17.43	178	2955	0.40	ng/ul	99
15) Fluoranthene	19.35	202	3960	0.40	ng/ul	96
17) Pyrene	19.71	202	4143	0.50	ng/ul	96
18) Benzo(a)anthracene	21.45	228	3978	0.48	ng/ul	99
19) Chrysene	21.50	228	3906	0.50	ng/ul	96
21) Benzo(b)fluoranthene	23.10	252	4740	0.50	ng/ul	85
22) Benzo(k)fluoranthene	23.14	252	4472	0.49	ng/ul#	86
23) Benzo(a)pyrene	23.71	252	4488	0.49	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	26.20	276	5913	0.56	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.22	278	4705	0.54	ng/ul#	92
26) Benzo(g,h,i)perylene	26.95	276	5073	0.57	ng/ul#	94

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

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