

Data Path : Z:\HPCHEM1\BNA M\DATA\BM053017\
 Data File : BM010307.D
 Acq On : 30 May 2017 12:25
 Operator : SJ/MA
 Sample : SSTD3.212
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: May 30 13:06:26 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	471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.73	136	1310	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	688	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	1465	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	1428	0.40	ng/ul	0.00
20) Perylene-d12	23.82	264	1527	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	6097	3.07	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	13005	3.00	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.78	128	10981	3.15	ng/ul#	94
5) 2-Methylnaphthalene	12.38	142	8010	3.14	ng/ul	99
7) Acenaphthylene	14.28	152	11317	3.38	ng/ul	96
8) Acenaphthene	14.62	153	8301	3.11	ng/ul	90
9) Fluorene	15.62	166	9901	3.12	ng/ul	93
11) Pentachlorophenol	16.95	266	1884	3.32	ng/ul	97
12) Phenanthrene	17.35	178	15342	3.51	ng/ul#	93
13) Anthracene	17.43	178	14965	3.45	ng/ul	93
15) Fluoranthene	19.35	202	19378	3.36	ng/ul	91
17) Pyrene	19.71	202	19765	4.16	ng/ul	96
18) Benzo(a)anthracene	21.45	228	18793	3.94	ng/ul	97
19) Chrysene	21.50	228	18252	4.07	ng/ul	97
21) Benzo(b)fluoranthene	23.09	252	22453	4.26	ng/ul	83
22) Benzo(k)fluoranthene	23.14	252	19580	3.87	ng/ul#	85
23) Benzo(a)pyrene	23.70	252	20632	4.07	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.20	276	27159	4.66	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	21874	4.56	ng/ul#	84
26) Benzo(g,h,i)perylene	26.95	276	23086	4.63	ng/ul#	86

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

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