

Data Path : Z:\HPCHEM1\BNA M\DATA\BM053017\
 Data File : BM010308.D
 Acq On : 30 May 2017 13:01
 Operator : SJ/MA
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV13

Quant Time: May 30 13:31:48 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 13:18:35 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	787	0.40	ng/ul	0.00
2) Naphthalene-d8	10.73	136	2191	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	1191	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	2676	0.40	ng/ul	0.00
16) Chrysene-d12	21.46	240	2639	0.40	ng/ul	0.00
20) Perylene-d12	23.82	264	2830	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.31	152	1287	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	2914	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.78	128	2333	0.40	ng/ul	97
5) 2-Methylnaphthalene	12.38	142	1651	0.41	ng/ul	99
7) Acenaphthylene	14.28	152	2402	0.40	ng/ul	99
8) Acenaphthene	14.62	153	1810	0.40	ng/ul	93
9) Fluorene	15.62	166	2141	0.40	ng/ul	95
11) Pentachlorophenol	16.95	266	403	0.39	ng/ul	93
12) Phenanthrene	17.35	178	3450	0.40	ng/ul	93
13) Anthracene	17.43	178	3160	0.39	ng/ul	98
15) Fluoranthene	19.36	202	4273	0.40	ng/ul	95
17) Pyrene	19.72	202	4457	0.40	ng/ul	100
18) Benzo(a)anthracene	21.45	228	4285	0.40	ng/ul	96
19) Chrysene	21.51	228	4248	0.41	ng/ul	99
21) Benzo(b)fluoranthene	23.09	252	4877	0.39	ng/ul	89
22) Benzo(k)fluoranthene	23.14	252	5050	0.43	ng/ul#	86
23) Benzo(a)pyrene	23.71	252	4881	0.42	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	26.21	276	6362	0.41	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.23	278	5096	0.42	ng/ul#	92
26) Benzo(g,h,i)perylene	26.96	276	5517	0.42	ng/ul#	92

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

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