

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012886.D
 Acq On : 11 Nov 2017 05:09
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02017

Quant Time: Nov 13 00:29:46 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 22:45:25 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	65200	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	248860	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.43	164	147905	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	345036	20.00	ng/ul	0.00
17) Chrysene-d12	21.36	240	416036	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	451621	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	289151	18.80	ng/ul	0.00
10) Fluorene-d10	15.42	176	200304	19.04	ng/ul	0.00
14) Anthracene-d10	17.27	188	316085	18.95	ng/ul	0.00
18) Pyrene-d10	19.57	212	356610	19.16	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	404859	18.99	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.63	128	239305	18.992	ng/ul	99
4) 2-Methylnaphthalene	12.24	142	178400	19.227	ng/ul	99
5) 1-Methylnaphthalene	12.46	142	172106	18.926	ng/ul#	93
8) Acenaphthylene	14.15	152	286468	19.138	ng/ul	99
9) Acenaphthene	14.49	153	192350	19.187	ng/ul	95
11) Fluorene	15.47	166	231940	19.395	ng/ul	92
13) Phenanthrene	17.22	178	366471	19.229	ng/ul	99
15) Anthracene	17.31	178	377728	19.077	ng/ul	99
16) Fluoranthene	19.23	202	444229	18.949	ng/ul#	91
19) Pyrene	19.60	202	463151	19.393	ng/ul#	93
20) Benzo(a)anthracene	21.34	228	484185	18.914	ng/ul	100
21) Chrysene	21.39	228	436109	19.063	ng/ul	98
23) Benzo(b)fluoranthene	22.94	252	516406	19.228	ng/ul#	97
24) Benzo(k)fluoranthene	22.99	252	501275	19.530	ng/ul#	96
26) Benzo(a)pyrene	23.53	252	493373	18.942	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	25.94	276	628379	18.704	ng/ul	97
28) Dibenzo(a,h)anthracene	25.94	278	530118	18.771	ng/ul#	96
29) Benzo(a,h,i)perylene	26.64	276	522158	18.619	ng/ul#	95

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

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