

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012896.D
 Acq On : 11 Nov 2017 11:42
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02018

Quant Time: Nov 13 00:38:17 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 00:36:35 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.79	152	62729	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	237475	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	138428	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	322371	20.00	ng/ul	0.00
17) Chrysene-d12	21.36	240	407616	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	464206	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	273488	18.99	ng/ul	0.00
10) Fluorene-d10	15.42	176	188020	19.09	ng/ul	0.00
14) Anthracene-d10	17.27	188	296906	19.05	ng/ul	0.00
18) Pyrene-d10	19.56	212	334214	18.33	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.49	264	415155	18.94	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.62	128	229393	19.078	ng/ul	99
4) 2-Methylnaphthalene	12.24	142	169687	19.165	ng/ul	99
5) 1-Methylnaphthalene	12.46	142	166027	19.133	ng/ul#	93
8) Acenaphthylene	14.15	152	268616	19.174	ng/ul	99
9) Acenaphthene	14.49	153	179708	19.153	ng/ul	96
11) Fluorene	15.47	166	217728	19.453	ng/ul	94
13) Phenanthrene	17.22	178	341069	19.155	ng/ul	99
15) Anthracene	17.30	178	355915	19.239	ng/ul	98
16) Fluoranthene	19.23	202	418099	19.088	ng/ul#	90
19) Pyrene	19.59	202	439961	18.803	ng/ul#	94
20) Benzo(a)anthracene	21.34	228	472421	18.836	ng/ul	99
21) Chrysene	21.39	228	425311	18.975	ng/ul	98
23) Benzo(b)fluoranthene	22.94	252	507226	18.375	ng/ul#	96
24) Benzo(k)fluoranthene	22.99	252	514912	19.517	ng/ul#	96
26) Benzo(a)pyrene	23.53	252	503242	18.797	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	25.94	276	641091	18.565	ng/ul	96
28) Dibenzo(a,h)anthracene	25.94	278	540068	18.605	ng/ul#	96
29) Benzo(g,h,i)perylene	26.64	276	538693	18.688	ng/ul#	95

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

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