

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM111017\
 Data File : BM012912.D
 Acq On : 11 Nov 2017 21:11
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02019

Quant Time: Nov 13 01:10:57 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM111017MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 01:07:29 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	71554	20.00	ng/ul	0.00
2) Naphthalene-d8	10.57	136	275211	20.00	ng/ul	0.00
6) Acenaphthene-d10	14.42	164	166455	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.17	188	390896	20.00	ng/ul	0.00
17) Chrysene-d12	21.35	240	483184	20.00	ng/ul	0.00
22) Perylene-d12	23.63	264	544189	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.12	160	325362	18.79	ng/ul	0.00
10) Fluorene-d10	15.42	176	226077	19.09	ng/ul	0.00
14) Anthracene-d10	17.27	188	358313	18.96	ng/ul	0.00
18) Pyrene-d10	19.56	212	401717	18.58	ng/ul	0.00
25) Benzo(a)pyrene-d12	23.48	264	490420	19.09	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.62	128	268234	19.249	ng/ul	100
4) 2-Methylnaphthalene	12.23	142	199282	19.421	ng/ul	98
5) 1-Methylnaphthalene	12.46	142	196430	19.533	ng/ul#	94
8) Acenaphthylene	14.15	152	321088	19.061	ng/ul	98
9) Acenaphthene	14.49	153	217531	19.281	ng/ul	97
11) Fluorene	15.47	166	261138	19.403	ng/ul	93
13) Phenanthrene	17.21	178	413780	19.165	ng/ul	99
15) Anthracene	17.30	178	432648	19.287	ng/ul	98
16) Fluoranthene	19.23	202	507795	19.119	ng/ul#	92
19) Pyrene	19.59	202	531159	19.150	ng/ul#	93
20) Benzo(a)anthracene	21.34	228	565980	19.037	ng/ul	99
21) Chrysene	21.39	228	505220	19.015	ng/ul	99
23) Benzo(b)fluoranthene	22.94	252	649627	20.074	ng/ul#	97
24) Benzo(k)fluoranthene	22.99	252	591451	19.123	ng/ul#	97
26) Benzo(a)pyrene	23.53	252	596847	19.017	ng/ul#	97
27) Indeno(1,2,3-cd)pyrene	25.93	276	760758	18.792	ng/ul	96
28) Dibenzo(a,h)anthracene	25.93	278	642436	18.879	ng/ul#	97
29) Benzo(a,h,i)perylene	26.63	276	632519	18.718	ng/ul#	95

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

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