

Data Path : Z:\HPCHEM1\BNA N\DATA\BN031518\
 Data File : BN000418.D
 Acq On : 15 Mar 2018 16:33
 Operator : JU/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD02007

Quant Time: Mar 16 05:45:54 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN030918MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 16 05:42:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	195915	20.00	ng/ul	0.00
2) Naphthalene-d8	11.29	136	953294	20.00	ng/ul	0.00
6) Acenaphthene-d10	15.07	164	536151	20.00	ng/ul	0.00
12) Phenanthrene-d10	17.79	188	1182304	20.00	ng/ul	0.00
17) Chrysene-d12	21.91	240	1012358	20.00	ng/ul	0.00
22) Perylene-d12	24.40	264	938616	20.00	ng/ul	0.00
System Monitoring Compounds						
7) Acenaphthylene-d8	14.77	160	1083979	20.37	ng/ul	0.00
10) Fluorene-d10	16.05	176	726022	20.15	ng/ul	0.00
14) Anthracene-d10	17.89	188	1096598	19.85	ng/ul	0.00
18) Pyrene-d10	20.14	212	1140068	23.02	ng/ul	0.00
25) Benzo(a)pyrene-d12	24.24	264	862114	20.11	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.34	128	979441	19.470	ng/ul	99
4) 2-Methylnaphthalene	12.92	142	679338	19.184	ng/ul	97
5) 1-Methylnaphthalene	13.14	142	679806	19.129	ng/ul#	92
8) Acenaphthylene	14.80	152	1070952	20.027	ng/ul	98
9) Acenaphthene	15.13	153	726045	19.508	ng/ul	96
11) Fluorene	16.10	166	813071	19.807	ng/ul	93
13) Phenanthrene	17.83	178	1314138	19.381	ng/ul	99
15) Anthracene	17.92	178	1364024	19.826	ng/ul	99
16) Fluoranthene	19.81	202	1471882	20.091	ng/ul	95
19) Pyrene	20.17	202	1447413	22.646	ng/ul#	89
20) Benzo(a)anthracene	21.90	228	1270368	20.185	ng/ul	99
21) Chrysene	21.95	228	1174415	19.664	ng/ul	100
23) Benzo(b)fluoranthene	23.64	252	1155920	20.652	ng/ul#	96
24) Benzo(k)fluoranthene	23.69	252	1061266	20.257	ng/ul#	96
26) Benzo(a)pyrene	24.29	252	1064749	19.988	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.98	276	1176492	20.484	ng/ul#	79
28) Dibenzo(a,h)anthracene	26.98	278	986518	20.955	ng/ul#	93
29) Benzo(g,h,i)perylene	27.77	276	978288	20.399	ng/ul#	88

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

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