

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061119\
 Data File : BN006246.D
 Acq On : 11 Jun 2019 13:25
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Jun 11 14:04:46 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 11 12:32:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4570	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	17688	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8346	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	16026	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	11685	0.40	ng/ul	0.03
20) Perylene-d12	23.55	264	13513	0.40	ng/ul	0.06

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	10196	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	17199	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.48	128	20161	0.413	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	12784	0.376	ng/ul	99
7) Acenaphthylene	14.01	152	17883	0.385	ng/ul	99
8) Acenaphthene	14.36	153	13324	0.414	ng/ul	100
9) Fluorene	15.36	166	13487	0.358	ng/ul	98
11) Pentachlorophenol	16.72	266	1241	0.305	ng/ul	96
12) Phenanthrene	17.09	178	20087	0.390	ng/ul	98
13) Anthracene	17.19	178	17315	0.352	ng/ul	99
15) Fluoranthene	19.13	202	22774	0.401	ng/ul	97
17) Pyrene	19.49	202	23981	0.390	ng/ul	99
18) Benzo(a)anthracene	21.27	228	19152	0.371	ng/ul	99
19) Chrysene	21.33	228	23186	0.478	ng/ul	99
21) Benzo(b)fluoranthene	22.87	252	21258	0.377	ng/ul	98
22) Benzo(k)fluoranthene	22.92	252	23311	0.425	ng/ul	96
23) Benzo(a)pyrene	23.45	252	21805	0.416	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.79	276	24975	0.440	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.80	278	19725	0.434	ng/ul	94
26) Benzo(g,h,i)perylene	26.48	276	21521	0.452	ng/ul	95

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

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