

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006350.D
 Acq On : 17 Jun 2019 16:25
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.421

Quant Time: Jun 18 01:51:36 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4037	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17084	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9269	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	17668	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	10711	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	11563	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10480	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	17149	0.36	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	19663	0.417	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	13041	0.397	ng/ul	98
7) Acenaphthylene	14.01	152	19255	0.374	ng/ul	99
8) Acenaphthene	14.36	153	14821	0.415	ng/ul	100
9) Fluorene	15.35	166	15285	0.366	ng/ul	98
11) Pentachlorophenol	16.72	266	2114	0.471	ng/ul	99
12) Phenanthrene	17.09	178	22038	0.389	ng/ul	99
13) Anthracene	17.19	178	18388	0.339	ng/ul	99
15) Fluoranthene	19.12	202	22477	0.359	ng/ul	98
17) Pyrene	19.48	202	23305	0.413	ng/ul	99
18) Benzo(a)anthracene	21.25	228	17165	0.362	ng/ul	98
19) Chrysene	21.31	228	20374	0.458	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	18154	0.377	ng/ul	95
22) Benzo(k)fluoranthene	22.89	252	19736	0.420	ng/ul	96
23) Benzo(a)pyrene	23.41	252	18441	0.411	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.76	276	20729	0.426	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.76	278	16266	0.419	ng/ul	93
26) Benzo(g,h,i)perylene	26.43	276	17938	0.441	ng/ul	96

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

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