

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006271.D
 Acq On : 12 Jun 2019 06:09
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Jun 12 06:52:31 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 : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5225	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	21079	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10382	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18887	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	12691	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	14255	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	12770	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	19550	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	24223	0.416	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	15951	0.394	ng/ul	99
7) Acenaphthylene	14.01	152	22749	0.394	ng/ul	99
8) Acenaphthene	14.36	153	16724	0.418	ng/ul	100
9) Fluorene	15.35	166	17175	0.367	ng/ul	98
11) Pentachlorophenol	16.72	266	522	0.109	ng/ul	93
12) Phenanthrene	17.09	178	23857	0.394	ng/ul	99
13) Anthracene	17.18	178	21129	0.365	ng/ul	99
15) Fluoranthene	19.12	202	25574	0.382	ng/ul	96
17) Pyrene	19.49	202	26855	0.402	ng/ul	99
18) Benzo(a)anthracene	21.25	228	21010	0.374	ng/ul	99
19) Chrysene	21.30	228	24892	0.472	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	22090	0.372	ng/ul	96
22) Benzo(k)fluoranthene	22.87	252	24709	0.427	ng/ul	98
23) Benzo(a)pyrene	23.40	252	22762	0.411	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.73	276	26514	0.442	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.74	278	20865	0.436	ng/ul	96
26) Benzo(g,h,i)perylene	26.41	276	22764	0.454	ng/ul	97

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

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