

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006181.D
 Acq On : 07 Jun 2019 23:22
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.401

Quant Time: Jun 08 02:04:44 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 : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	6058	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	22884	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10515	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	20376	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	16860	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	18639	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	12443	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	21905	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	24628	0.390	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	15716	0.357	ng/ul	99
7) Acenaphthylene	14.01	152	22571	0.386	ng/ul	99
8) Acenaphthene	14.36	153	16003	0.395	ng/ul	100
9) Fluorene	15.35	166	16538	0.349	ng/ul	96
11) Pentachlorophenol	16.71	266	1871	0.361	ng/ul	99
12) Phenanthrene	17.09	178	24541	0.375	ng/ul	99
13) Anthracene	17.18	178	22407	0.359	ng/ul	99
15) Fluoranthene	19.12	202	29029	0.402	ng/ul	96
17) Pyrene	19.49	202	30307	0.341	ng/ul	100
18) Benzo(a)anthracene	21.24	228	26493	0.355	ng/ul	99
19) Chrysene	21.30	228	29904	0.427	ng/ul	99
21) Benzo(b)fluoranthene	22.82	252	27703	0.356	ng/ul	97
22) Benzo(k)fluoranthene	22.86	252	29498	0.390	ng/ul	95
23) Benzo(a)pyrene	23.39	252	28310	0.391	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.72	276	32521	0.415	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.73	278	25915	0.414	ng/ul	94
26) Benzo(g,h,i)perylene	26.40	276	27568	0.420	ng/ul	95

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

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