

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006164.D
 Acq On : 07 Jun 2019 13:32
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.451

Quant Time: Jun 07 15:04:48 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	5728	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	21913	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	10767	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	22063	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	18141	0.40	ng/ul	0.02
20) Perylene-d12	23.54	264	19990	0.40	ng/ul	0.05
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	12289	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	24136	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	23597	0.390	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	15207	0.361	ng/ul	97
7) Acenaphthylene	14.02	152	23103	0.386	ng/ul	99
8) Acenaphthene	14.36	153	16357	0.394	ng/ul	99
9) Fluorene	15.36	166	17361	0.357	ng/ul	98
11) Pentachlorophenol	16.72	266	1947	0.347	ng/ul	98
12) Phenanthrene	17.09	178	26701	0.377	ng/ul	99
13) Anthracene	17.19	178	24424	0.361	ng/ul	100
15) Fluoranthene	19.13	202	32383	0.414	ng/ul	98
17) Pyrene	19.49	202	33428	0.350	ng/ul	97
18) Benzo(a)anthracene	21.27	228	28061	0.350	ng/ul	99
19) Chrysene	21.33	228	32683	0.434	ng/ul	99
21) Benzo(b)fluoranthene	22.87	252	30050	0.361	ng/ul	95
22) Benzo(k)fluoranthene	22.92	252	31585	0.389	ng/ul	95
23) Benzo(a)pyrene	23.45	252	30370	0.391	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.79	276	34881	0.415	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.80	278	27894	0.415	ng/ul	93
26) Benzo(g,h,i)perylene	26.47	276	29620	0.421	ng/ul	96

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

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