

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061719\
 Data File : BN006341.D
 Acq On : 17 Jun 2019 10:09
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.420

Quant Time: Jun 17 10:41:31 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 : Mon Jun 17 00:47:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	3793	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	14779	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7070	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	13186	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	9723	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	11164	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	8595	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14063	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	17006	0.417	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	10725	0.377	ng/ul	98
7) Acenaphthylene	14.01	152	14801	0.377	ng/ul	99
8) Acenaphthene	14.36	153	11300	0.415	ng/ul	100
9) Fluorene	15.35	166	11122	0.349	ng/ul	96
11) Pentachlorophenol	16.72	266	1515	0.452	ng/ul	98
12) Phenanthrene	17.09	178	16541	0.391	ng/ul	99
13) Anthracene	17.19	178	13900	0.344	ng/ul	99
15) Fluoranthene	19.12	202	18546	0.397	ng/ul	95
17) Pyrene	19.49	202	19625	0.383	ng/ul	99
18) Benzo(a)anthracene	21.25	228	15375	0.358	ng/ul	99
19) Chrysene	21.31	228	18574	0.460	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	17527	0.377	ng/ul	94
22) Benzo(k)fluoranthene	22.88	252	18565	0.409	ng/ul	96
23) Benzo(a)pyrene	23.41	252	17368	0.401	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.75	276	19379	0.413	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.75	278	14694	0.392	ng/ul	94
26) Benzo(g,h,i)perylene	26.43	276	17314	0.441	ng/ul	96

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

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