

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061319\
 Data File : BN006301.D
 Acq On : 13 Jun 2019 09:04
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.415

Quant Time: Jun 13 11:27:50 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 : Thu Jun 13 11:20:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4395	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17389	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8329	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15424	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11245	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	12785	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10111	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	16040	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	19459	0.406	ng/ul	100
5) 2-Methylnaphthalene	12.09	142	12614	0.377	ng/ul	98
7) Acenaphthylene	14.01	152	17489	0.378	ng/ul	100
8) Acenaphthene	14.36	153	13211	0.411	ng/ul	100
9) Fluorene	15.35	166	13370	0.356	ng/ul	97
11) Pentachlorophenol	16.72	266	1256	0.320	ng/ul	98
12) Phenanthrene	17.09	178	19153	0.387	ng/ul	99
13) Anthracene	17.18	178	16404	0.347	ng/ul	99
15) Fluoranthene	19.12	202	21168	0.387	ng/ul	98
17) Pyrene	19.49	202	22044	0.372	ng/ul	98
18) Benzo(a)anthracene	21.25	228	18010	0.362	ng/ul	98
19) Chrysene	21.30	228	21003	0.450	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	19935	0.374	ng/ul	95
22) Benzo(k)fluoranthene	22.88	252	20877	0.402	ng/ul	96
23) Benzo(a)pyrene	23.40	252	19926	0.401	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.74	276	22975	0.428	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.75	278	18111	0.422	ng/ul	93
26) Benzo(g,h,i)perylene	26.42	276	19750	0.439	ng/ul	96

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

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