

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN061419\
 Data File : BN006326.D
 Acq On : 14 Jun 2019 09:20
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Jun 14 16:08:37 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	4182	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	16350	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7962	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	15563	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	11747	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	13053	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	9393	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	16563	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.47	128	18441	0.409	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	11684	0.372	ng/ul	97
7) Acenaphthylene	14.01	152	16264	0.367	ng/ul	99
8) Acenaphthene	14.36	153	12602	0.410	ng/ul	99
9) Fluorene	15.35	166	12870	0.358	ng/ul	98
11) Pentachlorophenol	16.72	266	1047	0.265	ng/ul	97
12) Phenanthrene	17.09	178	19445	0.389	ng/ul	100
13) Anthracene	17.18	178	16227	0.340	ng/ul	99
15) Fluoranthene	19.12	202	21768	0.395	ng/ul	96
17) Pyrene	19.49	202	22769	0.368	ng/ul	99
18) Benzo(a)anthracene	21.25	228	18383	0.354	ng/ul	98
19) Chrysene	21.30	228	21597	0.443	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	20646	0.379	ng/ul	95
22) Benzo(k)fluoranthene	22.87	252	21578	0.407	ng/ul	95
23) Benzo(a)pyrene	23.40	252	20679	0.408	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.74	276	23349	0.426	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.74	278	18395	0.419	ng/ul	93
26) Benzo(g,h,i)perylene	26.42	276	20222	0.440	ng/ul	95

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

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