

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112818\
 Data File : BN003651.D
 Acq On : 28 Nov 2018 12:17
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.476

Quant Time: Nov 28 13:13:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 28 01:42:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4264	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	16396	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	8680	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	20653	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	19765	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	17631	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	9883	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	23056	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	18852	0.409	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	13130	0.408	ng/ul	99
7) Acenaphthylene	14.20	152	16063	0.411	ng/ul	100
8) Acenaphthene	14.54	153	14215	0.415	ng/ul	100
9) Fluorene	15.52	166	16300	0.408	ng/ul	99
11) Pentachlorophenol	16.87	266	1436	0.386	ng/ul	99
12) Phenanthrene	17.26	178	28354	0.418	ng/ul	100
13) Anthracene	17.35	178	23046	0.414	ng/ul	100
15) Fluoranthene	19.28	202	32072	0.401	ng/ul	98
17) Pyrene	19.64	202	31534	0.445	ng/ul	98
18) Benzo(a)anthracene	21.39	228	25455	0.408	ng/ul	98
19) Chrysene	21.44	228	29740	0.416	ng/ul	100
21) Benzo(b)fluoranthene	23.02	252	30802	0.422	ng/ul	98
22) Benzo(k)fluoranthene	23.06	252	29063	0.409	ng/ul	99
23) Benzo(a)pyrene	23.62	252	26306	0.417	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.07	276	28389	0.392	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.09	278	21881	0.375	ng/ul	99
26) Benzo(g,h,i)perylene	26.80	276	24786	0.394	ng/ul	98

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

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