

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004283.D
 Acq On : 29 Dec 2018 03:16
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.436

Quant Time: Dec 29 04:07:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 22:16:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2424	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	11061	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6109	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	15311	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	15944	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	15087	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.21	152	6312	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	16317	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.67	128	12128	0.389	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	8258	0.382	ng/ul	99
7) Acenaphthylene	14.19	152	12292	0.402	ng/ul	100
8) Acenaphthene	14.53	153	9293	0.379	ng/ul	99
9) Fluorene	15.52	166	10611	0.382	ng/ul	99
11) Pentachlorophenol	16.89	266	896	0.404	ng/ul	96
12) Phenanthrene	17.26	178	18909	0.380	ng/ul	100
13) Anthracene	17.35	178	17332	0.383	ng/ul	100
15) Fluoranthene	19.27	202	22485	0.375	ng/ul	99
17) Pyrene	19.64	202	23486	0.431	ng/ul	99
18) Benzo(a)anthracene	21.38	228	21655	0.400	ng/ul	100
19) Chrysene	21.43	228	22483	0.387	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	23471	0.410	ng/ul	100
22) Benzo(k)fluoranthene	23.06	252	23667	0.390	ng/ul	100
23) Benzo(a)pyrene	23.62	252	21322	0.392	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.09	276	21833	0.387	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.10	278	17790	0.384	ng/ul	99
26) Benzo(g,h,i)perylene	26.82	276	17880	0.377	ng/ul	100

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

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