

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120618\
 Data File : BN003756.D
 Acq On : 07 Dec 2018 07:04
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.491

Quant Time: Dec 07 07:50:47 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 : Fri Dec 07 04:08:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5242	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	21295	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	10829	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	23427	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	20432	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	17306	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	12966	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	26023	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	24807	0.415	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	17128	0.410	ng/ul	100
7) Acenaphthylene	14.19	152	21029	0.431	ng/ul	100
8) Acenaphthene	14.53	153	17902	0.419	ng/ul	100
9) Fluorene	15.52	166	19863	0.399	ng/ul	99
11) Pentachlorophenol	16.86	266	1836	0.435	ng/ul	98
12) Phenanthrene	17.26	178	32197	0.419	ng/ul	100
13) Anthracene	17.34	178	27374	0.433	ng/ul	100
15) Fluoranthene	19.27	202	35736	0.394	ng/ul	97
17) Pyrene	19.63	202	34901	0.476	ng/ul	98
18) Benzo(a)anthracene	21.38	228	29140	0.452	ng/ul	99
19) Chrysene	21.43	228	30617	0.415	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	30326	0.424	ng/ul	99
22) Benzo(k)fluoranthene	23.05	252	28788	0.413	ng/ul	99
23) Benzo(a)pyrene	23.60	252	26095	0.421	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.05	276	28155	0.396	ng/ul	98
25) Dibenzo(a,h)anthracene	26.07	278	22477	0.393	ng/ul	99
26) Benzo(g,h,i)perylene	26.77	276	23658	0.383	ng/ul	98

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

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