

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120318\
 Data File : BN003690.D
 Acq On : 03 Dec 2018 22:20
 Operator : JU/SJ
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.482

Quant Time: Dec 04 01:27:24 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 : Tue Dec 04 01:23:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	6957	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	27776	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	14897	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	34107	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	32599	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	29653	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	16519	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	37066	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	28944	0.371	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	19582	0.359	ng/ul	100
7) Acenaphthylene	14.19	152	28977	0.432	ng/ul	100
8) Acenaphthene	14.54	153	22629	0.385	ng/ul	99
9) Fluorene	15.52	166	24351	0.355	ng/ul	100
11) Pentachlorophenol	16.86	266	3354	0.546	ng/ul	99
12) Phenanthrene	17.26	178	41324	0.369	ng/ul	99
13) Anthracene	17.35	178	33948	0.369	ng/ul	99
15) Fluoranthene	19.27	202	46887	0.355	ng/ul	99
17) Pyrene	19.63	202	48327	0.413	ng/ul	98
18) Benzo(a)anthracene	21.38	228	39298	0.382	ng/ul	99
19) Chrysene	21.43	228	44147	0.375	ng/ul	99
21) Benzo(b)fluoranthene	23.00	252	43604	0.355	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	43631	0.365	ng/ul	100
23) Benzo(a)pyrene	23.61	252	38093	0.359	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	41066	0.337	ng/ul	98
25) Dibenzo(a,h)anthracene	26.07	278	32623	0.333	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	34589	0.327	ng/ul	99

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

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