

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120318\
 Data File : BN003681.D
 Acq On : 03 Dec 2018 16:41
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.481

Quant Time: Dec 04 01:20:06 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 01:17:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	6329	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	24486	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	13016	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	30645	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	29947	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	28664	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	14446	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	32927	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	25458	0.370	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	17214	0.358	ng/ul	100
7) Acenaphthylene	14.20	152	24540	0.419	ng/ul	100
8) Acenaphthene	14.54	153	19769	0.385	ng/ul	99
9) Fluorene	15.52	166	21444	0.358	ng/ul	99
11) Pentachlorophenol	16.86	266	2984	0.541	ng/ul	99
12) Phenanthrene	17.26	178	37442	0.372	ng/ul	99
13) Anthracene	17.35	178	29560	0.358	ng/ul	100
15) Fluoranthene	19.27	202	42205	0.355	ng/ul	97
17) Pyrene	19.64	202	43527	0.405	ng/ul	99
18) Benzo(a)anthracene	21.38	228	35084	0.371	ng/ul	99
19) Chrysene	21.43	228	40814	0.377	ng/ul	99
21) Benzo(b)fluoranthene	23.01	252	42121	0.355	ng/ul	99
22) Benzo(k)fluoranthene	23.06	252	41416	0.359	ng/ul	99
23) Benzo(a)pyrene	23.61	252	36541	0.356	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	41204	0.350	ng/ul	98
25) Dibenzo(a,h)anthracene	26.08	278	32021	0.338	ng/ul	99
26) Benzo(g,h,i)perylene	26.78	276	35092	0.343	ng/ul	98

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

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