

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003788.D
 Acq On : 10 Dec 2018 21:14
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.495

Quant Time: Dec 11 03:46:02 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 11 03:42:06 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	4283	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	16953	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	8566	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	19087	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	16226	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	13863	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	9487	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	19090	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	18307	0.384	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	12587	0.379	ng/ul	99
7) Acenaphthylene	14.19	152	14821	0.384	ng/ul	100
8) Acenaphthene	14.53	153	13075	0.387	ng/ul	99
9) Fluorene	15.52	166	14592	0.370	ng/ul	100
11) Pentachlorophenol	16.86	266	1258	0.366	ng/ul	97
12) Phenanthrene	17.26	178	24056	0.384	ng/ul	99
13) Anthracene	17.35	178	19623	0.381	ng/ul	100
15) Fluoranthene	19.27	202	26280	0.355	ng/ul	96
17) Pyrene	19.63	202	25637	0.441	ng/ul	99
18) Benzo(a)anthracene	21.38	228	20149	0.393	ng/ul	99
19) Chrysene	21.43	228	22676	0.387	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	22378	0.390	ng/ul	98
22) Benzo(k)fluoranthene	23.05	252	21394	0.383	ng/ul	97
23) Benzo(a)pyrene	23.60	252	19117	0.385	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.05	276	20849	0.366	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.07	278	16411	0.358	ng/ul	97
26) Benzo(g,h,i)perylene	26.78	276	18198	0.368	ng/ul	98

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

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