

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004101.D
 Acq On : 19 Dec 2018 05:34
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.430

Quant Time: Dec 19 06:24:50 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 : Wed Dec 19 04:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1477	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	7512	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4601	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	11960	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	13599	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	12770	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	4810	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	15343	0.44	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.68	128	8576	0.406	ng/ul	98
5) 2-Methylnaphthalene	12.29	142	6413	0.435	ng/ul	99
7) Acenaphthylene	14.19	152	9920	0.479	ng/ul	99
8) Acenaphthene	14.53	153	7517	0.414	ng/ul	100
9) Fluorene	15.52	166	8911	0.421	ng/ul	100
11) Pentachlorophenol	16.88	266	822	0.382	ng/ul	99
12) Phenanthrene	17.26	178	15871	0.404	ng/ul	99
13) Anthracene	17.35	178	15328	0.475	ng/ul	99
15) Fluoranthene	19.27	202	20473	0.442	ng/ul	93
17) Pyrene	19.64	202	21043	0.431	ng/ul	97
18) Benzo(a)anthracene	21.38	228	20150	0.469	ng/ul	98
19) Chrysene	21.44	228	19541	0.398	ng/ul	99
21) Benzo(b)fluoranthene	23.02	252	19515	0.369	ng/ul	97
22) Benzo(k)fluoranthene	23.06	252	19735	0.383	ng/ul#	97
23) Benzo(a)pyrene	23.62	252	18533	0.405	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.09	276	21407	0.408	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.09	278	17344	0.411	ng/ul	97
26) Benzo(g,h,i)perylene	26.82	276	17958	0.394	ng/ul	96

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

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