

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122120\
 Data File : BN013232.D
 Acq On : 21 Dec 2020 09:35
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.446

Quant Time: Dec 21 10:25:50 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 21 10:25:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	659	0.40	ng/ul	0.00
2) Naphthalene-d8	10.10	136	2224	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.00	164	1090	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	2144	0.40	ng/ul	0.00
16) Chrysene-d12	21.00	240	1687	0.40	ng/ul	0.00
20) Perylene-d12	23.08	264	1986	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.70	152	1338	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	2351	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.15	128	2878	0.447	ng/ul	98
5) 2-Methylnaphthalene	11.78	142	1786	0.422	ng/ul	99
7) Acenaphthylene	13.72	152	2375	0.460	ng/ul	99
8) Acenaphthene	14.06	153	1873	0.443	ng/ul	99
9) Fluorene	15.07	166	1972	0.406	ng/ul	98
11) Pentachlorophenol	16.43	266	269	0.433	ng/ul	91
12) Phenanthrene	16.80	178	3075	0.428	ng/ul	99
13) Anthracene	16.90	178	2649	0.433	ng/ul	100
15) Fluoranthene	18.84	202	3565	0.417	ng/ul	99
17) Pyrene	19.21	202	3707	0.465	ng/ul	96
18) Benzo(a)anthracene	20.98	228	3028	0.434	ng/ul	99
19) Chrysene	21.03	228	3708	0.481	ng/ul	100
21) Benzo(b)fluoranthene	22.47	252	3758	0.409	ng/ul	96
22) Benzo(k)fluoranthene	22.51	252	4326	0.426	ng/ul	99
23) Benzo(a)pyrene	22.99	252	3775	0.443	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.12	276	4878	0.436	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.13	278	3752	0.424	ng/ul	97
26) Benzo(g,h,i)perylene	25.73	276	4365	0.431	ng/ul	98

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

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