

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121920\
 Data File : BN013224.D
 Acq On : 19 Dec 2020 09:40
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.444

Quant Time: Dec 21 00:43:12 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 14 10:24:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.34	152	328	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.10	136	1032	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.01	164	572	0.40	ng/ul	-0.02
10) Phenanthrene-d10	16.78	188	1164	0.40	ng/ul	-0.02
16) Chrysene-d12	21.00	240	974	0.40	ng/ul	-0.02
20) Perylene-d12	23.09	264	1021	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	582	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.82	212	1359	0.43	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.16	128	1201	0.402	ng/ul#	91
5) 2-Methylnaphthalene	11.80	142	763	0.389	ng/ul	98
7) Acenaphthylene	13.72	152	1212	0.448	ng/ul	96
8) Acenaphthene	14.07	153	983	0.443	ng/ul	95
9) Fluorene	15.08	166	1080	0.424	ng/ul#	96
11) Pentachlorophenol	16.46	266	141	0.418	ng/ul	94
12) Phenanthrene	16.82	178	1694	0.434	ng/ul#	91
13) Anthracene	16.92	178	1381	0.416	ng/ul#	91
15) Fluoranthene	18.85	202	2006	0.433	ng/ul	97
17) Pyrene	19.22	202	2136	0.464	ng/ul	97
18) Benzo(a)anthracene	20.99	228	1542	0.383	ng/ul	96
19) Chrysene	21.04	228	2278	0.512	ng/ul	98
21) Benzo(b)fluoranthene	22.47	252	2108	0.446	ng/ul	98
22) Benzo(k)fluoranthene	22.51	252	2529	0.484	ng/ul	96
23) Benzo(a)pyrene	23.00	252	2041	0.466	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.12	276	2570	0.446	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.14	278	1941	0.426	ng/ul#	90
26) Benzo(g,h,i)perylene	25.75	276	2410	0.463	ng/ul	98

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

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