

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012827.D
 Acq On : 02 Dec 2020 19:32
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.419

Quant Time: Dec 03 00:19:33 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 12:19:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	823	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	3342	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1908	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.83	188	3888	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	3640	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	3805	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	2038	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	3922	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.23	128	4187	0.416	ng/ul	97
5) 2-Methylnaphthalene	11.86	142	2926	0.456	ng/ul	99
7) Acenaphthylene	13.79	152	4148	0.446	ng/ul	99
8) Acenaphthene	14.14	153	3206	0.409	ng/ul	99
9) Fluorene	15.13	166	3550	0.412	ng/ul	99
11) Pentachlorophenol	16.49	266	573	0.464	ng/ul	96
12) Phenanthrene	16.87	178	5542	0.410	ng/ul	98
13) Anthracene	16.97	178	5028	0.428	ng/ul	98
15) Fluoranthene	18.90	202	6907	0.442	ng/ul	99
17) Pyrene	19.27	202	7104	0.436	ng/ul	99
18) Benzo(a)anthracene	21.03	228	7078	0.543	ng/ul	98
19) Chrysene	21.09	228	8097	0.497	ng/ul	100
21) Benzo(b)fluoranthene	22.54	252	8505	0.486	ng/ul	94
22) Benzo(k)fluoranthene	22.57	252	8969	0.444	ng/ul	94
23) Benzo(a)pyrene	23.07	252	7942	0.482	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.22	276	10528	0.489	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.24	278	8407	0.486	ng/ul	92
26) Benzo(g,h,i)perylene	25.86	276	8962	0.462	ng/ul	96

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

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