

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120220\
 Data File : BN012814.D
 Acq On : 02 Dec 2020 11:43
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.418

Quant Time: Dec 02 12:18:54 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 : Tue Dec 01 10:54:12 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	628	0.40	ng/ul	0.00
2) Naphthalene-d8	10.18	136	2553	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1478	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	3096	0.40	ng/ul	0.00
16) Chrysene-d12	21.05	240	2955	0.40	ng/ul	0.00
20) Perylene-d12	23.16	264	3003	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	1563	0.42	ng/ul	0.00
14) Fluoranthene-d10	18.87	212	3169	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.23	128	3229	0.420	ng/ul	99
5) 2-Methylnaphthalene	11.86	142	2230	0.454	ng/ul	100
7) Acenaphthylene	13.79	152	3203	0.444	ng/ul	99
8) Acenaphthene	14.14	153	2479	0.408	ng/ul	99
9) Fluorene	15.14	166	2767	0.414	ng/ul	97
11) Pentachlorophenol	16.49	266	437	0.444	ng/ul	95
12) Phenanthrene	16.87	178	4433	0.411	ng/ul	99
13) Anthracene	16.97	178	3930	0.420	ng/ul	99
15) Fluoranthene	18.91	202	5510	0.443	ng/ul	96
17) Pyrene	19.27	202	5620	0.425	ng/ul	99
18) Benzo(a)anthracene	21.04	228	5526	0.522	ng/ul	98
19) Chrysene	21.09	228	6590	0.498	ng/ul	100
21) Benzo(b)fluoranthene	22.54	252	6692	0.484	ng/ul	94
22) Benzo(k)fluoranthene	22.58	252	7348	0.461	ng/ul	94
23) Benzo(a)pyrene	23.07	252	6335	0.487	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	25.23	276	8474	0.498	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.24	278	6745	0.494	ng/ul	95
26) Benzo(g,h,i)perylene	25.86	276	7289	0.476	ng/ul	97

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

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