

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031620\
 Data File : BN010224.D
 Acq On : 16 Mar 2020 12:39
 Operator : CG/JU
 Sample : SSTD0.205
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.205

Quant Time: Mar 16 14:00:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN031620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 13:58:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1702	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5362	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3167	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	7517	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	6837	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	6971	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	1895	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	4755	0.21	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	3539	0.219	ng/ul	96
5) 2-Methylnaphthalene	12.06	142	2354	0.213	ng/ul	99
7) Acenaphthylene	13.97	152	3358	0.208	ng/ul	96
8) Acenaphthene	14.32	153	2636	0.203	ng/ul	98
9) Fluorene	15.31	166	3105	0.208	ng/ul	98
11) Pentachlorophenol	16.66	266	228	0.111	ng/ul	97
12) Phenanthrene	17.04	178	5455	0.210	ng/ul	99
13) Anthracene	17.13	178	4725	0.205	ng/ul	98
15) Fluoranthene	19.06	202	6524	0.212	ng/ul	98
17) Pyrene	19.42	202	6689	0.204	ng/ul	99
18) Benzo(a)anthracene	21.18	228	5916	0.197	ng/ul	100
19) Chrysene	21.23	228	6989	0.215	ng/ul	100
21) Benzo(b)fluoranthene	22.73	252	6328	0.205	ng/ul	96
22) Benzo(k)fluoranthene	22.77	252	6613	0.188	ng/ul	95
23) Benzo(a)pyrene	23.28	252	5381	0.174	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.52	276	6770	0.196	ng/ul	100
25) Dibenzo(a,h)anthracene	25.54	278	5537	0.203	ng/ul	94
26) Benzo(g,h,i)perylene	26.17	276	6052	0.200	ng/ul	98

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

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