

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011760.D
 Acq On : 09 Sep 2020 05:45
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.499

Quant Time: Sep 09 07:13:47 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 09 07:05:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2436	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	9186	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	5044	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	10611	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	8828	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	8364	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	6233	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	12032	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	11691	0.430	ng/ul	98
5) 2-Methylnaphthalene	12.18	142	7827	0.432	ng/ul	100
7) Acenaphthylene	14.09	152	10363	0.426	ng/ul	99
8) Acenaphthene	14.44	153	8374	0.429	ng/ul	99
9) Fluorene	15.43	166	9450	0.419	ng/ul	97
11) Pentachlorophenol	16.77	266	1219	0.381	ng/ul	99
12) Phenanthrene	17.16	178	15357	0.430	ng/ul	99
13) Anthracene	17.25	178	13795	0.434	ng/ul	99
15) Fluoranthene	19.18	202	16924	0.427	ng/ul	99
17) Pyrene	19.54	202	17026	0.434	ng/ul	99
18) Benzo(a)anthracene	21.29	228	14869	0.433	ng/ul	99
19) Chrysene	21.34	228	16302	0.425	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	16378	0.418	ng/ul	98
22) Benzo(k)fluoranthene	22.92	252	17106	0.425	ng/ul	98
23) Benzo(a)pyrene	23.46	252	14859	0.439	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.81	276	19309	0.434	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.83	278	15708	0.433	ng/ul	97
26) Benzo(g,h,i)perylene	26.50	276	16136	0.415	ng/ul	98

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

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