

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN081020\
 Data File : BN011403.D
 Acq On : 10 Aug 2020 17:49
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.422

Quant Time: Aug 10 18:18:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 10 13:35:01 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	818	0.40	ng/ul	0.00
2) Naphthalene-d8	10.20	136	2742	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.09	164	1262	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	2721	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	2553	0.40	ng/ul	0.00
20) Perylene-d12	23.18	264	2521	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.80	152	2088	0.30	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	4693	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.25	128	2859	0.351	ng/ul	97
5) 2-Methylnaphthalene	11.88	142	1737	0.318	ng/ul	100
7) Acenaphthylene	13.80	152	2345	0.389	ng/ul	97
8) Acenaphthene	14.15	153	1817	0.379	ng/ul	97
9) Fluorene	15.15	166	2039	0.362	ng/ul	98
11) Pentachlorophenol	16.51	266	256	0.384	ng/ul	96
12) Phenanthrene	16.89	178	3465	0.384	ng/ul	97
13) Anthracene	16.98	178	2921	0.371	ng/ul	95
15) Fluoranthene	18.92	202	4013	0.361	ng/ul	99
17) Pyrene	19.28	202	3804	0.328	ng/ul	99
18) Benzo(a)anthracene	21.05	228	3176	0.324	ng/ul	98
19) Chrysene	21.10	228	4001	0.356	ng/ul	97
21) Benzo(b)fluoranthene	22.56	252	3891	0.370	ng/ul	91
22) Benzo(k)fluoranthene	22.60	252	4536	0.385	ng/ul	93
23) Benzo(a)pyrene	23.09	252	3539	0.387	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.24	276	5048	0.525	ng/ul	94
25) Dibenzo(a,h)anthracene	25.26	278	4006	0.523	ng/ul	94
26) Benzo(g,h,i)perylene	25.86	276	4059	0.467	ng/ul	98

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

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