

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111519\
 Data File : BN008590.D
 Acq On : 15 Nov 2019 17:30
 Operator : JU
 Sample : SSTD0.804
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.8041

Quant Time: Nov 15 18:01:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 15 17:05:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2779	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	9361	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	5011	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	12186	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	12843	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	12053	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	11694	0.79	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	28234	0.78	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.43	128	21343	0.801	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	14980	0.844	ng/ul	100
7) Acenaphthylene	13.96	152	21412	0.917	ng/ul	98
8) Acenaphthene	14.31	153	15832	0.838	ng/ul	99
9) Fluorene	15.30	166	18596	0.899	ng/ul	100
11) Pentachlorophenol	16.65	266	3516	1.325	ng/ul	99
12) Phenanthrene	17.03	178	31325	0.846	ng/ul	98
13) Anthracene	17.12	178	29529	0.910	ng/ul	98
15) Fluoranthene	19.05	202	38165	0.899	ng/ul	99
17) Pyrene	19.41	202	38478	0.614	ng/ul	97
18) Benzo(a)anthracene	21.17	228	37874	0.847	ng/ul	100
19) Chrysene	21.22	228	37240	0.654	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	37678	0.839	ng/ul	95
22) Benzo(k)fluoranthene	22.76	252	36603	0.717	ng/ul#	93
23) Benzo(a)pyrene	23.26	252	35387	0.782	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.50	276	41723	0.812	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.51	278	33770	0.829	ng/ul	95
26) Benzo(g,h,i)perylene	26.15	276	35086	0.761	ng/ul	96

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

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