

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110719\
 Data File : BN008479.D
 Acq On : 07 Nov 2019 11:35
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.499

Quant Time: Nov 08 05:42:18 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 14:41:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1167	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5098	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2991	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	5041	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3491	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	3605	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	3067	0.38	ng/ul	0.01
14) Fluoranthene-d10	18.99	212	5834	0.39	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.37	128	5898	0.407	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	3717	0.385	ng/ul	100
7) Acenaphthylene	13.91	152	5082	0.365	ng/ul	99
8) Acenaphthene	14.26	153	4341	0.385	ng/ul	99
9) Fluorene	15.25	166	4491	0.364	ng/ul	98
11) Pentachlorophenol	16.62	266	391	0.356	ng/ul	99
12) Phenanthrene	16.99	178	6355	0.415	ng/ul	98
13) Anthracene	17.08	178	5378	0.401	ng/ul	98
15) Fluoranthene	19.01	202	6893	0.392	ng/ul	99
17) Pyrene	19.38	202	7178	0.421	ng/ul	99
18) Benzo(a)anthracene	21.15	228	4623	0.380	ng/ul	100
19) Chrysene	21.20	228	6739	0.436	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	5215	0.388	ng/ul	94
22) Benzo(k)fluoranthene	22.73	252	6311	0.413	ng/ul	96
23) Benzo(a)pyrene	23.24	252	5433	0.402	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.48	276	6546	0.426	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.49	278	5147	0.422	ng/ul	98
26) Benzo(g,h,i)perylene	26.12	276	6064	0.440	ng/ul	99

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

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