

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008446.D
 Acq On : 05 Nov 2019 14:19
 Operator : JU
 Sample : SSTD1.606
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.606

Quant Time: Nov 07 10:51:00 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 10:46:19 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1659	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	7288	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	4062	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	9020	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	6955	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	5586	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	19050	1.55	ng/ul	0.00
14) Fluoranthene-d10	18.99	212	44026	1.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	32863	1.584	ng/ul	97
5) 2-Methylnaphthalene	12.00	142	22781	1.601	ng/ul	98
7) Acenaphthylene	13.91	152	32388	1.674	ng/ul	97
8) Acenaphthene	14.26	153	25279	1.680	ng/ul	99
9) Fluorene	15.25	166	28674	1.678	ng/ul	98
11) Pentachlorophenol	16.62	266	3408	1.562	ng/ul	100
12) Phenanthrene	16.99	178	44499	1.685	ng/ul	99
13) Anthracene	17.08	178	41633	1.791	ng/ul	97
15) Fluoranthene	19.01	202	51971	1.490	ng/ul	98
17) Pyrene	19.38	202	52013	1.878	ng/ul	98
18) Benzo(a)anthracene	21.14	228	41290	1.633	ng/ul	97
19) Chrysene	21.19	228	46832	1.509	ng/ul	98
21) Benzo(b)fluoranthene	22.68	252	35310	1.942	ng/ul	78
22) Benzo(k)fluoranthene	22.72	252	39306	1.908	ng/ul#	88
23) Benzo(a)pyrene	23.23	252	33822	1.767	ng/ul#	65
24) Indeno(1,2,3-cd)pyrene	25.47	276	38100	1.691	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.48	278	30658	1.721	ng/ul#	88
26) Benzo(g,h,i)perylene	26.12	276	33406	1.616	ng/ul	95

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

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