

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003593.D
 Acq On : 19 Nov 2018 14:35
 Operator : MJ/SJ
 Sample : SSTD1.665
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD1.665

Quant Time: Nov 19 15:22:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 19 13:56:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	4904	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	21454	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	11880	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	27692	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	32316	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	27578	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	54790	1.60	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	135960	1.75	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.69	128	97666	1.623	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	69987	1.617	ng/ul	100
7) Acenaphthylene	14.20	152	90157	1.747	ng/ul	99
8) Acenaphthene	14.54	153	76227	1.653	ng/ul	99
9) Fluorene	15.53	166	89594	1.650	ng/ul	100
11) Pentachlorophenol	16.87	266	8804	1.587	ng/ul	99
12) Phenanthrene	17.26	178	150946	1.662	ng/ul	99
13) Anthracene	17.35	178	131464	1.776	ng/ul	100
15) Fluoranthene	19.28	202	183845	1.801	ng/ul	99
17) Pyrene	19.64	202	176399	1.501	ng/ul	100
18) Benzo(a)anthracene	21.38	228	170750	1.685	ng/ul	99
19) Chrysene	21.43	228	184229	1.600	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	189911	1.598	ng/ul	98
22) Benzo(k)fluoranthene	23.05	252	189086	1.643	ng/ul	98
23) Benzo(a)pyrene	23.61	252	169273	1.706	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.06	276	195763	1.843	ng/ul	100
25) Dibenzo(a,h)anthracene	26.07	278	158955	1.836	ng/ul	98
26) Benzo(g,h,i)perylene	26.78	276	164865	1.810	ng/ul	99

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

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