

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111918\
 Data File : BN003591.D
 Acq On : 19 Nov 2018 13:25
 Operator : MJ/SJ
 Sample : SSTD0.463
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.463

Quant Time: Nov 19 14:01:54 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	7.84	152	5158	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	17586	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	10570	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	26429	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	24872	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	22923	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	10480	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	30113	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.69	128	19175	0.389	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	13285	0.374	ng/ul	100
7) Acenaphthylene	14.20	152	18560	0.404	ng/ul	100
8) Acenaphthene	14.54	153	16535	0.403	ng/ul	100
9) Fluorene	15.53	166	19316	0.400	ng/ul	100
11) Pentachlorophenol	16.87	266	1610	0.304	ng/ul	100
12) Phenanthrene	17.26	178	34420	0.397	ng/ul	100
13) Anthracene	17.36	178	27778	0.393	ng/ul	100
15) Fluoranthene	19.28	202	39866	0.409	ng/ul	100
17) Pyrene	19.64	202	39163	0.433	ng/ul	100
18) Benzo(a)anthracene	21.38	228	30626	0.393	ng/ul	100
19) Chrysene	21.43	228	35937	0.405	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	37588	0.381	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	36597	0.383	ng/ul	100
23) Benzo(a)pyrene	23.61	252	32627	0.396	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.06	276	37365	0.423	ng/ul	100
25) Dibenzo(a,h)anthracene	26.07	278	29863	0.415	ng/ul	100
26) Benzo(g,h,i)perylene	26.78	276	33253	0.439	ng/ul	100

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

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