

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN020421\
 Data File : BN013570.D
 Acq On : 04 Feb 2021 14:19
 Operator : CG/JU
 Sample : SSTD3.201
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD3.201

Quant Time: Feb 04 14:48:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN020421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 04 13:01:49 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	4673	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	17401	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	9318	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.94	188	18861	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18079	0.40	ng/ul	0.00
20) Perylene-d12	23.31	264	13776	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.91	152	77040	2.95	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	152275	3.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	137476	2.731	ng/ul	99
5) 2-Methylnaphthalene	11.98	142	93199	2.916	ng/ul	100
7) Acenaphthylene	13.90	152	114156	2.973	ng/ul	100
8) Acenaphthene	14.25	153	94598	2.768	ng/ul	99
9) Fluorene	15.24	166	108751	2.878	ng/ul	99
11) Pentachlorophenol	16.59	266	9379	3.372	ng/ul	100
12) Phenanthrene	16.98	178	172740	2.797	ng/ul	99
13) Anthracene	17.07	178	155973	3.079	ng/ul	99
15) Fluoranthene	19.00	202	195556	2.898	ng/ul	99
17) Pyrene	19.37	202	191297	2.585	ng/ul	98
18) Benzo(a)anthracene	21.12	228	183850	3.081	ng/ul	100
19) Chrysene	21.18	228	188087	2.528	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	180575	3.122	ng/ul	97
22) Benzo(k)fluoranthene	22.71	252	182242	2.959	ng/ul	95
23) Benzo(a)pyrene	23.21	252	147510	3.006	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.46	276	198805	3.050	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.47	278	163753	3.110	ng/ul	97
26) Benzo(g,h,i)perylene	26.11	276	168779	2.925	ng/ul	99

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

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