

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022720\
 Data File : BM025125.D
 Acq On : 28 Feb 2020 04:14
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.492

Quant Time: Feb 28 04:50:53 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 28 04:39:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	2799	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.46	136	10262	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.32	164	6042	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	13772	0.40	ng/ul	-0.01
16) Chrysene-d12	21.24	240	14140	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	15494	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.05	152	7355	0.38	ng/ul	-0.02
14) Fluoranthene-d10	19.09	212	18073	0.38	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.51	128	12876	0.412	ng/ul#	83
5) 2-Methylnaphthalene	12.13	142	8632	0.383	ng/ul	100
7) Acenaphthylene	14.03	152	12397	0.356	ng/ul	94
8) Acenaphthene	14.38	153	9260	0.396	ng/ul	99
9) Fluorene	15.37	166	10980	0.374	ng/ul	97
11) Pentachlorophenol	16.71	266	1493	0.355	ng/ul	99
12) Phenanthrene	17.10	178	19144	0.395	ng/ul	96
13) Anthracene	17.19	178	16805	0.358	ng/ul	95
15) Fluoranthene	19.11	202	23226	0.330	ng/ul	96
17) Pyrene	19.48	202	23229	0.365	ng/ul	95
18) Benzo(a)anthracene	21.22	228	22554	0.369	ng/ul	98
19) Chrysene	21.27	228	24081	0.396	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	24030	0.396	ng/ul	92
22) Benzo(k)fluoranthene	22.82	252	25934	0.431	ng/ul	98
23) Benzo(a)pyrene	23.34	252	21483	0.387	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.63	276	30196	0.435	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.64	278	25287	0.433	ng/ul	96
26) Benzo(g,h,i)perylene	26.29	276	24673	0.421	ng/ul#	92

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

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