

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040420\
 Data File : BM025917.D
 Acq On : 03 Apr 2020 18:08
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.437

Quant Time: Apr 03 18:39:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2804	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	11887	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	7391	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	18785	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	18431	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	18211	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	8213	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	23194	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	12712	0.364	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	9354	0.382	ng/ul	99
7) Acenaphthylene	13.92	152	13216	0.418	ng/ul	98
8) Acenaphthene	14.26	153	10426	0.398	ng/ul	100
9) Fluorene	15.26	166	12718	0.391	ng/ul	100
11) Pentachlorophenol	16.61	266	1385	0.300	ng/ul	96
12) Phenanthrene	16.99	178	22087	0.373	ng/ul	99
13) Anthracene	17.08	178	19292	0.369	ng/ul	100
15) Fluoranthene	19.01	202	26698	0.365	ng/ul	100
17) Pyrene	19.37	202	26661	0.404	ng/ul	97
18) Benzo(a)anthracene	21.13	228	24887	0.378	ng/ul	100
19) Chrysene	21.18	228	27515	0.383	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	26268	0.392	ng/ul	99
22) Benzo(k)fluoranthene	22.70	252	28702	0.399	ng/ul	99
23) Benzo(a)pyrene	23.20	252	24639	0.411	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.41	276	28451	0.355	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.42	278	23722	0.354	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	23089	0.340	ng/ul	98

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

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