

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080219\
 Data File : BM021768.D
 Acq On : 02 Aug 2019 21:35
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.420

Quant Time: Aug 03 05:52:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 03 05:33:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	1002	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4366	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	1903	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	3568	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	2430	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	2314	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	2467	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	3201	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	4778	0.391	ng/ul	100
5) 2-Methylnaphthalene	12.14	142	2971	0.423	ng/ul	100
7) Acenaphthylene	14.03	152	3647	0.382	ng/ul	100
8) Acenaphthene	14.38	153	3213	0.396	ng/ul	100
9) Fluorene	15.38	166	3253	0.384	ng/ul	100
11) Pentachlorophenol	16.76	266	162	0.324	ng/ul	100
12) Phenanthrene	17.12	178	4650	0.389	ng/ul	100
13) Anthracene	17.22	178	3655	0.402	ng/ul	100
15) Fluoranthene	19.13	202	4348	0.366	ng/ul	100
17) Pyrene	19.50	202	4636	0.431	ng/ul	100
18) Benzo(a)anthracene	21.25	228	2880	0.396	ng/ul	100
19) Chrysene	21.30	228	4522	0.370	ng/ul	100
21) Benzo(b)fluoranthene	22.82	252	3630	0.352	ng/ul	100
22) Benzo(k)fluoranthene	22.86	252	4157	0.382	ng/ul	100
23) Benzo(a)pyrene	23.40	252	3100	0.382	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.73	276	3601	0.384	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.75	278	2882	0.395	ng/ul	100
26) Benzo(g,h,i)perylene	26.41	276	3426	0.380	ng/ul	100

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

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