

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091619\
 Data File : BM022684.D
 Acq On : 16 Sep 2019 10:10
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.466

Quant Time: Sep 16 11:16:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM091019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 10 15:06:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	8351	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	32129	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	17674	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	37391	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	32339	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	26667	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	22687	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.23	212	47346	0.46	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	39132	0.427	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	26843	0.428	ng/ul	100
7) Acenaphthylene	14.19	152	32791	0.401	ng/ul	100
8) Acenaphthene	14.54	153	27145	0.396	ng/ul	99
9) Fluorene	15.52	166	30917	0.394	ng/ul	99
11) Pentachlorophenol	16.86	266	2893	0.381	ng/ul	98
12) Phenanthrene	17.25	178	52064	0.437	ng/ul	99
13) Anthracene	17.34	178	43620	0.427	ng/ul	100
15) Fluoranthene	19.26	202	59478	0.463	ng/ul	99
17) Pyrene	19.62	202	59714	0.407	ng/ul	99
18) Benzo(a)anthracene	21.37	228	50748	0.410	ng/ul	99
19) Chrysene	21.42	228	56545	0.434	ng/ul	100
21) Benzo(b)fluoranthene	22.98	252	51583	0.445	ng/ul	99
22) Benzo(k)fluoranthene	23.02	252	50037	0.423	ng/ul	99
23) Benzo(a)pyrene	23.56	252	42688	0.421	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.98	276	48199	0.390	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.99	278	39015	0.390	ng/ul	98
26) Benzo(g,h,i)perylene	26.68	276	41765	0.424	ng/ul	99

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

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