

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091219\
 Data File : BM022674.D
 Acq On : 12 Sep 2019 20:14
 Operator : HP/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.465

Quant Time: Sep 13 01:15:50 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	7.85	152	7780	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	30528	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	17121	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	36197	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	32081	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	29026	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	21576	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	45505	0.46	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.69	128	37306	0.429	ng/ul	100
5) 2-Methylnaphthalene	12.30	142	25658	0.430	ng/ul	100
7) Acenaphthylene	14.20	152	31760	0.401	ng/ul	99
8) Acenaphthene	14.54	153	26269	0.396	ng/ul	99
9) Fluorene	15.52	166	30433	0.400	ng/ul	99
11) Pentachlorophenol	16.86	266	2708	0.369	ng/ul	99
12) Phenanthrene	17.25	178	50850	0.441	ng/ul	100
13) Anthracene	17.34	178	42625	0.431	ng/ul	100
15) Fluoranthene	19.27	202	57307	0.460	ng/ul	100
17) Pyrene	19.62	202	56822	0.391	ng/ul	98
18) Benzo(a)anthracene	21.37	228	50702	0.413	ng/ul	99
19) Chrysene	21.42	228	55710	0.431	ng/ul	100
21) Benzo(b)fluoranthene	22.98	252	55003	0.436	ng/ul	99
22) Benzo(k)fluoranthene	23.02	252	53916	0.419	ng/ul	100
23) Benzo(a)pyrene	23.57	252	46913	0.425	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.98	276	55527	0.412	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.99	278	45455	0.418	ng/ul	98
26) Benzo(g,h,i)perylene	26.68	276	45841	0.427	ng/ul	99

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

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