

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091019\
 Data File : BM022617.D
 Acq On : 10 Sep 2019 13:25
 Operator : HP/JU
 Sample : SSTD0.855
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.855

Quant Time: Sep 10 14:15:52 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 13:22:48 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	16822	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	65427	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	33063	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	72577	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	57803	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	51069	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	81092	0.75	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	154808	0.63	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.69	128	141482	0.766	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	96471	0.783	ng/ul	100
7) Acenaphthylene	14.20	152	117290	0.761	ng/ul	100
8) Acenaphthene	14.54	153	97714	0.774	ng/ul	99
9) Fluorene	15.52	166	112579	0.803	ng/ul	99
11) Pentachlorophenol	16.86	266	10936	0.817	ng/ul	99
12) Phenanthrene	17.26	178	177440	0.840	ng/ul	100
13) Anthracene	17.35	178	154336	0.857	ng/ul	100
15) Fluoranthene	19.27	202	195322	0.690	ng/ul	100
17) Pyrene	19.63	202	195904	0.825	ng/ul	99
18) Benzo(a)anthracene	21.38	228	171929	0.832	ng/ul	100
19) Chrysene	21.43	228	175799	0.661	ng/ul	100
21) Benzo(b)fluoranthene	22.99	252	172998	1.005	ng/ul	94
22) Benzo(k)fluoranthene	23.03	252	172310	0.884	ng/ul	95
23) Benzo(a)pyrene	23.58	252	149871	0.856	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.99	276	181534	0.868	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.01	278	147197	0.871	ng/ul	97
26) Benzo(g,h,i)perylene	26.70	276	145945	0.775	ng/ul	99

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

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