

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091019\
 Data File : BM022615.D
 Acq On : 10 Sep 2019 12:12
 Operator : HP/JU
 Sample : SSTD0.253
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.253

Quant Time: Sep 10 13:39:11 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	15949	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	59959	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.48	164	29869	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	66733	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	48598	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	42456	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	19417	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	35345	0.16	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.69	128	34462	0.203	ng/ul	99
5) 2-Methylnaphthalene	12.31	142	23164	0.205	ng/ul	100
7) Acenaphthylene	14.20	152	27189	0.195	ng/ul	99
8) Acenaphthene	14.54	153	23337	0.205	ng/ul	99
9) Fluorene	15.53	166	25989	0.205	ng/ul	99
11) Pentachlorophenol	16.87	266	2359	0.192	ng/ul	98
12) Phenanthrene	17.26	178	42339	0.218	ng/ul	99
13) Anthracene	17.35	178	34923	0.211	ng/ul	99
15) Fluoranthene	19.27	202	44024	0.169	ng/ul	99
17) Pyrene	19.63	202	45032	0.225	ng/ul	98
18) Benzo(a)anthracene	21.38	228	35359	0.203	ng/ul	99
19) Chrysene	21.43	228	38946	0.174	ng/ul	100
21) Benzo(b)fluoranthene	22.99	252	35894	0.251	ng/ul	96
22) Benzo(k)fluoranthene	23.03	252	37923	0.234	ng/ul#	95
23) Benzo(a)pyrene	23.57	252	31440	0.216	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.99	276	38524	0.222	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.01	278	31198	0.222	ng/ul	98
26) Benzo(g,h,i)perylene	26.70	276	30812	0.197	ng/ul	98

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

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