

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021924.D
 Acq On : 08 Aug 2019 09:28
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Aug 08 12:35:33 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 08 06:54:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	653	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	2994	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	1671	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	3556	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3238	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	3753	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1905	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4220	0.35	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	3103	0.367	ng/ul	98
5) 2-Methylnaphthalene	12.10	142	2151	0.382	ng/ul	98
7) Acenaphthylene	14.01	152	2810	0.361	ng/ul	99
8) Acenaphthene	14.35	153	2390	0.375	ng/ul	95
9) Fluorene	15.35	166	2522	0.356	ng/ul	99
11) Pentachlorophenol	16.72	266	230	0.351	ng/ul	97
12) Phenanthrene	17.09	178	3841	0.371	ng/ul	98
13) Anthracene	17.19	178	3329	0.377	ng/ul	98
15) Fluoranthene	19.11	202	4652	0.336	ng/ul	98
17) Pyrene	19.48	202	4896	0.368	ng/ul#	94
18) Benzo(a)anthracene	21.23	228	4248	0.367	ng/ul	98
19) Chrysene	21.28	228	5325	0.357	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	4659	0.368	ng/ul	98
22) Benzo(k)fluoranthene	22.84	252	5223	0.365	ng/ul	98
23) Benzo(a)pyrene	23.37	252	4573	0.355	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.69	276	5498	0.358	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.70	278	4369	0.352	ng/ul	98
26) Benzo(g,h,i)perylene	26.37	276	4871	0.352	ng/ul#	94

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

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