

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025408.D
 Acq On : 09 Mar 2020 09:22
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.421

Quant Time: Mar 09 13:35:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 13:32:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	3095	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	13375	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9527	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	22232	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	15834	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	14271	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	8368	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	20236	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	13539	0.342	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	9842	0.348	ng/ul	100
7) Acenaphthylene	14.01	152	12855	0.330	ng/ul	98
8) Acenaphthene	14.35	153	11350	0.332	ng/ul	100
9) Fluorene	15.34	166	14001	0.337	ng/ul	99
11) Pentachlorophenol	16.69	266	1509	0.289	ng/ul	98
12) Phenanthrene	17.07	178	23051	0.327	ng/ul	100
13) Anthracene	17.17	178	20057	0.327	ng/ul	100
15) Fluoranthene	19.09	202	25545	0.334	ng/ul	98
17) Pyrene	19.46	202	24639	0.347	ng/ul	99
18) Benzo(a)anthracene	21.20	228	19253	0.340	ng/ul	99
19) Chrysene	21.26	228	21174	0.333	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	18317	0.324	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	18926	0.320	ng/ul	99
23) Benzo(a)pyrene	23.31	252	15273	0.325	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	19355	0.304	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.59	278	16083	0.300	ng/ul	99
26) Benzo(g,h,i)perylene	26.24	276	16566	0.300	ng/ul	98

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

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