

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025432.D
 Acq On : 10 Mar 2020 02:05
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Mar 10 02:37:09 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	7.66	152	2904	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	12024	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	8264	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	17551	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	11539	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	10937	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	7437	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14784	0.31	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.48	128	12116	0.340	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	8647	0.340	ng/ul	100
7) Acenaphthylene	14.01	152	11144	0.330	ng/ul	98
8) Acenaphthene	14.35	153	9766	0.330	ng/ul	99
9) Fluorene	15.34	166	11640	0.323	ng/ul	99
11) Pentachlorophenol	16.69	266	1223	0.297	ng/ul	99
12) Phenanthrene	17.08	178	18166	0.327	ng/ul	99
13) Anthracene	17.17	178	15571	0.322	ng/ul	100
15) Fluoranthene	19.09	202	18311	0.304	ng/ul	98
17) Pyrene	19.45	202	17759	0.343	ng/ul	99
18) Benzo(a)anthracene	21.21	228	13594	0.330	ng/ul	99
19) Chrysene	21.26	228	15299	0.330	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	13753	0.318	ng/ul	98
22) Benzo(k)fluoranthene	22.80	252	14612	0.322	ng/ul	99
23) Benzo(a)pyrene	23.31	252	11858	0.330	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.58	276	15949	0.327	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.59	278	13189	0.321	ng/ul	98
26) Benzo(g,h,i)perylene	26.24	276	13435	0.317	ng/ul	97

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

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