

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025417.D
 Acq On : 09 Mar 2020 17:00
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.482

Quant Time: Mar 09 17:34:14 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	3172	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	13019	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9196	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	20725	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	14850	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	13571	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	8092	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	18307	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	13081	0.339	ng/ul	100
5) 2-Methylnaphthalene	12.10	142	9360	0.340	ng/ul	100
7) Acenaphthylene	14.01	152	12105	0.322	ng/ul	98
8) Acenaphthene	14.35	153	10877	0.330	ng/ul	100
9) Fluorene	15.34	166	13196	0.329	ng/ul	100
11) Pentachlorophenol	16.70	266	1441	0.296	ng/ul	98
12) Phenanthrene	17.07	178	21434	0.326	ng/ul	100
13) Anthracene	17.17	178	18437	0.323	ng/ul	99
15) Fluoranthene	19.09	202	22681	0.319	ng/ul	98
17) Pyrene	19.46	202	22033	0.331	ng/ul	98
18) Benzo(a)anthracene	21.21	228	17608	0.332	ng/ul	99
19) Chrysene	21.26	228	19605	0.329	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	17403	0.324	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	18445	0.328	ng/ul	99
23) Benzo(a)pyrene	23.31	252	14588	0.327	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.59	276	18769	0.310	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.60	278	15709	0.308	ng/ul	99
26) Benzo(g,h,i)perylene	26.24	276	15987	0.304	ng/ul	97

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

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