

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031320\
 Data File : BM025565.D
 Acq On : 14 Mar 2020 06:14
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.443

Quant Time: Mar 16 06:49:43 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 16 06:44:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	3673	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15830	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	10184	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	20630	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	14505	0.40	ng/ul	0.00
20) Perylene-d12	23.39	264	15201	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	10949	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	21291	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	17779	0.379	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	12786	0.382	ng/ul	100
7) Acenaphthylene	14.00	152	15341	0.369	ng/ul#	87
8) Acenaphthene	14.34	153	14070	0.386	ng/ul	99
9) Fluorene	15.33	166	16543	0.372	ng/ul	100
11) Pentachlorophenol	16.68	266	1978	0.409	ng/ul	98
12) Phenanthrene	17.06	178	25318	0.387	ng/ul	100
13) Anthracene	17.16	178	22451	0.395	ng/ul	99
15) Fluoranthene	19.08	202	26488	0.374	ng/ul	96
17) Pyrene	19.45	202	25697	0.395	ng/ul	96
18) Benzo(a)anthracene	21.20	228	21774	0.420	ng/ul	100
19) Chrysene	21.25	228	22815	0.392	ng/ul	98
21) Benzo(b)fluoranthene	22.74	252	22155	0.368	ng/ul	98
22) Benzo(k)fluoranthene	22.79	252	23025	0.365	ng/ul	98
23) Benzo(a)pyrene	23.30	252	20186	0.404	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.56	276	26008	0.384	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.58	278	21620	0.379	ng/ul	98
26) Benzo(g,h,i)perylene	26.22	276	22257	0.378	ng/ul	97

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

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