

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025514.D
 Acq On : 12 Mar 2020 09:33
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Mar 13 02:50:03 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	2643	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	10495	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	7112	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	14905	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	10062	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	9812	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7379	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	14607	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	12008	0.387	ng/ul	98
5) 2-Methylnaphthalene	12.09	142	8633	0.389	ng/ul	100
7) Acenaphthylene	14.00	152	10269	0.353	ng/ul#	86
8) Acenaphthene	14.35	153	9861	0.387	ng/ul	99
9) Fluorene	15.34	166	11898	0.383	ng/ul	100
11) Pentachlorophenol	16.69	266	1220	0.349	ng/ul	99
12) Phenanthrene	17.07	178	18257	0.387	ng/ul	99
13) Anthracene	17.16	178	15344	0.373	ng/ul	100
15) Fluoranthene	19.09	202	18017	0.352	ng/ul	99
17) Pyrene	19.45	202	17355	0.384	ng/ul	99
18) Benzo(a)anthracene	21.20	228	13747	0.382	ng/ul	99
19) Chrysene	21.25	228	15894	0.393	ng/ul	99
21) Benzo(b)fluoranthene	22.75	252	14982	0.386	ng/ul	99
22) Benzo(k)fluoranthene	22.79	252	15862	0.390	ng/ul	99
23) Benzo(a)pyrene	23.30	252	12795	0.397	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.57	276	17249	0.394	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	14356	0.390	ng/ul	100
26) Benzo(g,h,i)perylene	26.23	276	15357	0.404	ng/ul	99

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

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