

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040720\
 Data File : BM025954.D
 Acq On : 07 Apr 2020 12:23
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.443

Quant Time: Apr 07 12:56:43 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 02 10:49:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1972	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	8647	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5298	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	13137	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	12911	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	12381	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	5861	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	16112	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	9254	0.364	ng/ul	100
5) 2-Methylnaphthalene	12.00	142	6825	0.384	ng/ul	99
7) Acenaphthylene	13.92	152	9720	0.429	ng/ul	98
8) Acenaphthene	14.26	153	7590	0.404	ng/ul	99
9) Fluorene	15.26	166	9193	0.394	ng/ul	99
11) Pentachlorophenol	16.61	266	1025	0.318	ng/ul	91
12) Phenanthrene	16.99	178	15742	0.380	ng/ul	99
13) Anthracene	17.08	178	13736	0.376	ng/ul	100
15) Fluoranthene	19.01	202	18781	0.367	ng/ul	99
17) Pyrene	19.37	202	18950	0.410	ng/ul	97
18) Benzo(a)anthracene	21.12	228	17454	0.378	ng/ul	99
19) Chrysene	21.18	228	19305	0.383	ng/ul	100
21) Benzo(b)fluoranthene	22.66	252	18248	0.400	ng/ul	98
22) Benzo(k)fluoranthene	22.70	252	19860	0.406	ng/ul	98
23) Benzo(a)pyrene	23.20	252	16696	0.410	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.41	276	18551	0.340	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.42	278	15264	0.335	ng/ul	97
26) Benzo(g,h,i)perylene	26.05	276	15489	0.336	ng/ul	98

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

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