

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031220\
 Data File : BM025540.D
 Acq On : 13 Mar 2020 01:50
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.440

Quant Time: Mar 13 02:52:06 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	3635	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15023	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	10317	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	21532	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	14066	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	14368	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	10651	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	21134	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	16984	0.382	ng/ul	99
5) 2-Methylnaphthalene	12.09	142	12380	0.389	ng/ul	100
7) Acenaphthylene	14.00	152	15054	0.357	ng/ul#	86
8) Acenaphthene	14.35	153	14331	0.388	ng/ul	99
9) Fluorene	15.33	166	17141	0.380	ng/ul	100
11) Pentachlorophenol	16.69	266	1992	0.394	ng/ul	97
12) Phenanthrene	17.07	178	26248	0.385	ng/ul	100
13) Anthracene	17.16	178	22757	0.383	ng/ul	100
15) Fluoranthene	19.08	202	26149	0.354	ng/ul	98
17) Pyrene	19.45	202	25282	0.400	ng/ul	98
18) Benzo(a)anthracene	21.20	228	20374	0.405	ng/ul	99
19) Chrysene	21.25	228	22494	0.398	ng/ul	98
21) Benzo(b)fluoranthene	22.75	252	21457	0.377	ng/ul	99
22) Benzo(k)fluoranthene	22.79	252	22701	0.381	ng/ul	99
23) Benzo(a)pyrene	23.30	252	18843	0.399	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.57	276	25354	0.396	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.58	278	20915	0.388	ng/ul	99
26) Benzo(g,h,i)perylene	26.22	276	21654	0.389	ng/ul	98

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

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