

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM110920\
 Data File : BM027763.D
 Acq On : 09 Nov 2020 19:10
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.483

Quant Time: Nov 10 06:35:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM02.4-EPA-SIM-BM110920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 06:30:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.40	152	645	0.40	ng/ul	0.00
2) Naphthalene-d8	11.23	136	2290	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	1374	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.72	188	2919	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	2355	0.40	ng/ul	0.00
20) Perylene-d12	24.26	264	1919	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2057	0.49	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	4232	0.49	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	11.28	128	3539	0.484	ng/ul#	77
5) 2-Methylnaphthalene	12.87	142	2392	0.474	ng/ul	97
7) Acenaphthylene	14.73	152	3343	0.489	ng/ul#	79
8) Acenaphthene	15.07	153	2577	0.497	ng/ul	92
9) Fluorene	16.04	166	3009	0.492	ng/ul#	94
11) Pentachlorophenol	17.36	266	536	0.452	ng/ul	90
12) Phenanthrene	17.76	178	4709	0.497	ng/ul#	86
13) Anthracene	17.85	178	4512	0.493	ng/ul#	85
15) Fluoranthene	19.73	202	5329	0.494	ng/ul	83
17) Pyrene	20.09	202	5274	0.513	ng/ul#	85
18) Benzo(a)anthracene	21.80	228	4664	0.500	ng/ul#	88
19) Chrysene	21.85	228	4786	0.502	ng/ul#	88
21) Benzo(b)fluoranthene	23.51	252	4490	0.532	ng/ul	74
22) Benzo(k)fluoranthene	23.56	252	4313	0.514	ng/ul#	74
23) Benzo(a)pyrene	24.15	252	3860	0.519	ng/ul#	69
24) Indeno(1,2,3-cd)pyrene	26.79	276	4476	0.508	ng/ul#	89
25) Dibenzo(a,h)anthracene	26.80	278	3715	0.507	ng/ul#	65
26) Benzo(g,h,i)perylene	27.57	276	3785	0.498	ng/ul#	74

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

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