

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027877.D
 Acq On : 21 Nov 2020 12:32
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.435

Quant Time: Nov 21 13:03:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 17:20:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	748	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	2941	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1779	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3647	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2447	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	1928	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	1920	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3704	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	3488	0.396	ng/ul	100
5) 2-Methylnaphthalene	12.80	142	2477	0.413	ng/ul	98
7) Acenaphthylene	14.67	152	3438	0.396	ng/ul	100
8) Acenaphthene	15.01	153	2549	0.392	ng/ul	99
9) Fluorene	15.98	166	2958	0.392	ng/ul	99
11) Pentachlorophenol	17.31	266	463	0.373	ng/ul	98
12) Phenanthrene	17.70	178	4518	0.392	ng/ul	99
13) Anthracene	17.79	178	4293	0.386	ng/ul	99
15) Fluoranthene	19.68	202	4929	0.386	ng/ul	100
17) Pyrene	20.04	202	4913	0.395	ng/ul	100
18) Benzo(a)anthracene	21.75	228	3903	0.395	ng/ul	100
19) Chrysene	21.80	228	3817	0.397	ng/ul	99
21) Benzo(b)fluoranthene	23.44	252	3243	0.400	ng/ul	99
22) Benzo(k)fluoranthene	23.49	252	3101	0.391	ng/ul	99
23) Benzo(a)pyrene	24.08	252	2771	0.391	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.68	276	3237	0.422	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.69	278	2672	0.418	ng/ul	99
26) Benzo(g,h,i)perylene	27.45	276	2743	0.426	ng/ul	100

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

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