

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027932.D
 Acq On : 24 Nov 2020 03:04
 Operator : JU/CG
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.441

Quant Time: Nov 24 10:21:34 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 : Mon Nov 23 11:35:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.33	152	1171	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	4516	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2418	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4104	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2350	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2207	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	2835	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3637	0.33	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	5376	0.397	ng/ul	98
5) 2-Methylnaphthalene	12.80	142	3670	0.399	ng/ul	99
7) Acenaphthylene	14.67	152	4770	0.404	ng/ul	98
8) Acenaphthene	15.00	153	3482	0.394	ng/ul	99
9) Fluorene	15.98	166	3814	0.372	ng/ul	100
11) Pentachlorophenol	17.30	266	462	0.331	ng/ul	99
12) Phenanthrene	17.70	178	5153	0.397	ng/ul	99
13) Anthracene	17.79	178	4910	0.393	ng/ul	99
15) Fluoranthene	19.68	202	4955	0.345	ng/ul	96
17) Pyrene	20.03	202	4804	0.403	ng/ul	95
18) Benzo(a)anthracene	21.74	228	3784	0.399	ng/ul	98
19) Chrysene	21.80	228	3723	0.403	ng/ul	99
21) Benzo(b)fluoranthene	23.44	252	3503	0.378	ng/ul	97
22) Benzo(k)fluoranthene	23.49	252	3465	0.381	ng/ul	97
23) Benzo(a)pyrene	24.07	252	3111	0.384	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	26.67	276	3876	0.441	ng/ul#	97
25) Dibenzo(a,h)anthracene	26.68	278	3196	0.437	ng/ul	98
26) Benzo(g,h,i)perylene	27.44	276	3304	0.449	ng/ul	99

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

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