

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112320\
 Data File : BM027953.D
 Acq On : 24 Nov 2020 21:22
 Operator : JU/CG
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.443

Quant Time: Nov 25 00:50:31 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 : Tue Nov 24 15:00:24 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.32	152	1357	0.40	ng/ul	0.00
2) Naphthalene-d8	11.16	136	5171	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	2704	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	4418	0.40	ng/ul	0.00
16) Chrysene-d12	21.76	240	2703	0.40	ng/ul	0.00
20) Perylene-d12	24.18	264	2684	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.72	152	3235	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	3999	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.21	128	6244	0.403	ng/ul	98
5) 2-Methylnaphthalene	12.79	142	4253	0.403	ng/ul	99
7) Acenaphthylene	14.67	152	5611	0.425	ng/ul	98
8) Acenaphthene	15.00	153	3989	0.403	ng/ul	98
9) Fluorene	15.98	166	4348	0.379	ng/ul	100
11) Pentachlorophenol	17.30	266	584	0.388	ng/ul	99
12) Phenanthrene	17.70	178	5729	0.411	ng/ul	99
13) Anthracene	17.79	178	5471	0.406	ng/ul	99
15) Fluoranthene	19.68	202	5623	0.363	ng/ul	96
17) Pyrene	20.03	202	5486	0.400	ng/ul#	94
18) Benzo(a)anthracene	21.74	228	4392	0.403	ng/ul	97
19) Chrysene	21.80	228	4332	0.408	ng/ul	97
21) Benzo(b)fluoranthene	23.44	252	4244	0.376	ng/ul	96
22) Benzo(k)fluoranthene	23.49	252	4203	0.380	ng/ul#	95
23) Benzo(a)pyrene	24.07	252	3808	0.386	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.67	276	4599	0.431	ng/ul#	96
25) Dibenzo(a,h)anthracene	26.69	278	3798	0.427	ng/ul	97
26) Benzo(g,h,i)perylene	27.44	276	3912	0.437	ng/ul	97

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

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