

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020821\
 Data File : BM028676.D
 Acq On : 08 Feb 2021 16:17
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.484

Quant Time: Feb 08 16:47:57 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 08 13:26:08 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	418	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	1723	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	931	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	1880	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	1756	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	1774	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1036	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	2009	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	1796	0.419	ng/ul	99
5) 2-Methylnaphthalene	12.38	142	1211	0.422	ng/ul	98
7) Acenaphthylene	14.28	152	1520	0.400	ng/ul	100
8) Acenaphthene	14.61	153	1227	0.407	ng/ul	100
9) Fluorene	15.60	166	1391	0.416	ng/ul	99
11) Pentachlorophenol	16.95	266	190	0.387	ng/ul	97
12) Phenanthrene	17.33	178	2172	0.413	ng/ul	99
13) Anthracene	17.42	178	2032	0.406	ng/ul	99
15) Fluoranthene	19.33	202	2440	0.408	ng/ul	98
17) Pyrene	19.69	202	2503	0.392	ng/ul	99
18) Benzo(a)anthracene	21.42	228	2429	0.405	ng/ul	98
19) Chrysene	21.47	228	2451	0.411	ng/ul	97
21) Benzo(b)fluoranthene	23.00	252	2542	0.419	ng/ul	92
22) Benzo(k)fluoranthene	23.04	252	2557	0.422	ng/ul#	92
23) Benzo(a)pyrene	23.57	252	2278	0.421	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.93	276	2925	0.415	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.95	278	2430	0.411	ng/ul	93
26) Benzo(g,h,i)perylene	26.63	276	2478	0.423	ng/ul	96

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

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