

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022221\
 Data File : BM028865.D
 Acq On : 22 Feb 2021 15:11
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.443

Quant Time: Feb 22 16:06:44 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 22 11:00:59 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	330	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	1400	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	783	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	1546	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	1229	0.40	ng/ul	0.00
20) Perylene-d12	23.61	264	1154	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	892	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	1642	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1534	0.440	ng/ul	97
5) 2-Methylnaphthalene	12.30	142	1034	0.443	ng/ul	98
7) Acenaphthylene	14.21	152	1335	0.418	ng/ul	98
8) Acenaphthene	14.55	153	1100	0.434	ng/ul	99
9) Fluorene	15.54	166	1241	0.441	ng/ul	97
11) Pentachlorophenol	16.89	266	144	0.356	ng/ul	95
12) Phenanthrene	17.27	178	1914	0.443	ng/ul	99
13) Anthracene	17.36	178	1800	0.437	ng/ul	99
15) Fluoranthene	19.28	202	2090	0.425	ng/ul	97
17) Pyrene	19.64	202	2109	0.472	ng/ul	99
18) Benzo(a)anthracene	21.38	228	1812	0.432	ng/ul	99
19) Chrysene	21.43	228	1840	0.441	ng/ul	99
21) Benzo(b)fluoranthene	22.95	252	1801	0.456	ng/ul	89
22) Benzo(k)fluoranthene	22.99	252	1766	0.448	ng/ul#	89
23) Benzo(a)pyrene	23.52	252	1609	0.457	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.86	276	1956	0.427	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.87	278	1630	0.424	ng/ul	98
26) Benzo(g,h,i)perylene	26.55	276	1654	0.434	ng/ul	99

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

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