

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021721\
 Data File : BM028764.D
 Acq On : 17 Feb 2021 10:09
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.434

Quant Time: Feb 17 11:34:56 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 : Wed Feb 17 11:34:13 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	396	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1585	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	806	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	1569	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1507	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	1514	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	910	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1645	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	1679	0.425	ng/ul	98
5) 2-Methylnaphthalene	12.36	142	1091	0.413	ng/ul	99
7) Acenaphthylene	14.26	152	1271	0.386	ng/ul	97
8) Acenaphthene	14.60	153	1081	0.415	ng/ul	98
9) Fluorene	15.58	166	1192	0.412	ng/ul	99
11) Pentachlorophenol	16.94	266	121	0.295	ng/ul	97
12) Phenanthrene	17.31	178	1831	0.418	ng/ul	100
13) Anthracene	17.41	178	1628	0.389	ng/ul	98
15) Fluoranthene	19.32	202	2018	0.404	ng/ul	98
17) Pyrene	19.68	202	2087	0.381	ng/ul	99
18) Benzo(a)anthracene	21.40	228	1944	0.378	ng/ul	98
19) Chrysene	21.45	228	2108	0.412	ng/ul	98
21) Benzo(b)fluoranthene	22.98	252	2201	0.425	ng/ul	89
22) Benzo(k)fluoranthene	23.02	252	2202	0.426	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	1876	0.406	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.90	276	2567	0.427	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	2143	0.425	ng/ul#	89
26) Benzo(g,h,i)perylene	26.59	276	2230	0.446	ng/ul	94

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

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