

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022221\
 Data File : BM028858.D
 Acq On : 22 Feb 2021 09:39
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Feb 22 11:00:48 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 11:00:22 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	346	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	1429	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.49	164	775	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	1480	0.40	ng/ul	0.00
16) Chrysene-d12	21.39	240	1138	0.40	ng/ul	0.00
20) Perylene-d12	23.61	264	1115	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.23	152	895	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	1561	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.68	128	1591	0.447	ng/ul	97
5) 2-Methylnaphthalene	12.30	142	1056	0.443	ng/ul	98
7) Acenaphthylene	14.21	152	1353	0.428	ng/ul	97
8) Acenaphthene	14.55	153	1097	0.437	ng/ul	99
9) Fluorene	15.54	166	1208	0.434	ng/ul	97
11) Pentachlorophenol	16.90	266	138	0.357	ng/ul	98
12) Phenanthrene	17.27	178	1831	0.443	ng/ul	98
13) Anthracene	17.36	178	1761	0.446	ng/ul	99
15) Fluoranthene	19.28	202	1974	0.419	ng/ul	97
17) Pyrene	19.64	202	2019	0.488	ng/ul	98
18) Benzo(a)anthracene	21.37	228	1715	0.441	ng/ul	96
19) Chrysene	21.43	228	1714	0.444	ng/ul	96
21) Benzo(b)fluoranthene	22.94	252	1711	0.448	ng/ul	88
22) Benzo(k)fluoranthene	22.98	252	1716	0.451	ng/ul#	87
23) Benzo(a)pyrene	23.52	252	1565	0.460	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.85	276	1963	0.443	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.86	278	1629	0.439	ng/ul	96
26) Benzo(g,h,i)perylene	26.54	276	1680	0.456	ng/ul	99

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

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