

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121016\
 Data File : BM008256.D
 Acq On : 11 Dec 2016 05:28
 Operator : UM/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Manual Integrations
 APPROVED

UMANGI
 12/12/2016 6:50:45 PM

Quant Time: Dec 12 07:06:42 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 12 03:51:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	565	0.40	ng/ul	0.00
2) Naphthalene-d8	10.48	136	2061	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	1191	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	2098m	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	1655	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	1601	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	1188	0.41	ng/ul	-0.01
14) Fluoranthene-d10	19.12	212	2125m	0.36	ng/ul	0.00
Target Compounds						Ovalue
3) Naphthalene	10.52	128	2229	0.39	ng/ul	97
5) 2-Methylnaphthalene	12.14	142	1501	0.40	ng/ul	98
7) Acenaphthylene	14.04	152	2121	0.39	ng/ul	97
8) Acenaphthene	14.38	153	1635	0.40	ng/ul	93
9) Fluorene	15.39	166	1771	0.38	ng/ul	94
11) Pentachlorophenol	16.76	266	269	0.37	ng/ul	93
12) Phenanthrene	17.12	178	2551	0.39	ng/ul	96
13) Anthracene	17.23	178	2519	0.41	ng/ul	95
15) Fluoranthene	19.14	202	3055	0.36	ng/ul	95
17) Pyrene	19.51	202	3001	0.52	ng/ul	97
18) Benzo(a)anthracene	21.26	228	2434	0.48	ng/ul	97
19) Chrysene	21.31	228	2284	0.36	ng/ul	97
21) Benzo(b)fluoranthene	22.84	252	2276	0.36	ng/ul	98
22) Benzo(k)fluoranthene	22.89	252	2535	0.38	ng/ul	93
23) Benzo(a)pyrene	23.42	252	2354	0.39	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.77	276	2855	0.39	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.77	278	2254	0.38	ng/ul	96
26) Benzo(g,h,i)perylene	26.44	276	2424	0.37	ng/ul	95

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

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