

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122416\
 Data File : BM008643.D
 Acq On : 24 Dec 2016 21:11
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.466

Quant Time: Dec 26 00:02:21 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 25 23:48:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	406	0.40	ng/ul	0.05
2) Naphthalene-d8	10.48	136	1664	0.40	ng/ul	0.02
6) Acenaphthene-d10	14.31	164	995	0.40	ng/ul	0.02
10) Phenanthrene-d10	17.07	188	1926	0.40	ng/ul	0.02
16) Chrysene-d12	21.26	240	1911	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	1861	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.08	152	973	0.38	ng/ul	0.03
14) Fluoranthene-d10	19.10	212	2176	0.41	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.52	128	1838	0.39	ng/ul#	87
5) 2-Methylnaphthalene	12.15	142	1198	0.38	ng/ul	97
7) Acenaphthylene	14.04	152	1748	0.39	ng/ul#	93
8) Acenaphthene	14.37	153	1365	0.40	ng/ul	95
9) Fluorene	15.39	166	1497	0.39	ng/ul#	95
11) Pentachlorophenol	16.76	266	239	0.40	ng/ul	96
12) Phenanthrene	17.11	178	2417	0.41	ng/ul	98
13) Anthracene	17.21	178	2376	0.42	ng/ul	97
15) Fluoranthene	19.13	202	3140	0.44	ng/ul	97
17) Pyrene	19.50	202	3265	0.44	ng/ul	98
18) Benzo(a)anthracene	21.25	228	2581	0.41	ng/ul	96
19) Chrysene	21.30	228	2909	0.42	ng/ul	94
21) Benzo(b)fluoranthene	22.83	252	2736	0.37	ng/ul	97
22) Benzo(k)fluoranthene	22.87	252	2995	0.42	ng/ul	96
23) Benzo(a)pyrene	23.40	252	2817	0.41	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.74	276	3262	0.40	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.74	278	2563	0.40	ng/ul#	89
26) Benzo(g,h,i)perylene	26.43	276	2806	0.40	ng/ul	95

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

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