

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022715\
 Data File : BM000612.D
 Acq On : 27 Feb 2015 18:34
 Operator : TP/IZ
 Sample : SSTD0.233
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.233

Quant Time: Feb 27 22:50:56 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM01.2-EPA-SIM-BM022715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 27 22:41:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	2735	0.40	ng/ul	0.00
4) Naphthalene-d8	10.46	136	9389	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.32	164	5782	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.05	188	9835	0.40	ng/ul	-0.01
18) Chrysene-d12	21.25	240	7135	0.40	ng/ul	0.00
22) Perylene-d12	23.46	264	6140	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.14	96	2113	1.12	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.06	152	2367	0.21	ng/ul	0.01
16) Fluoranthene-d10	19.09	212	3818	0.22	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.17	88	3687	1.16	ng/ul#	83
5) Naphthalene	10.50	128	4914	0.21	ng/ul	97
7) 2-Methylnaphthalene	12.13	142	3211	0.21	ng/ul	98
9) Acenaphthylene	14.04	152	4280	0.18	ng/ul#	96
10) Acenaphthene	14.37	153	3451	0.19	ng/ul#	81
11) Fluorene	15.37	166	3884	0.19	ng/ul	96
13) Pentachlorophenol	16.73	266	150	0.12	ng/ul	95
14) Phenanthrene	17.10	178	5772	0.21	ng/ul	97
15) Anthracene	17.19	178	4956	0.20	ng/ul	96
17) Fluoranthene	19.13	202	5944	0.22	ng/ul	95
19) Pyrene	19.49	202	6125	0.20	ng/ul	98
20) Benzo(a)anthracene	21.23	228	4258	0.21	ng/ul	96
21) Chrysene	21.28	228	5108	0.22	ng/ul	96
23) Benzo(b)fluoranthene	22.80	252	4520	0.21	ng/ul	86
24) Benzo(k)fluoranthene	22.84	252	4458	0.21	ng/ul#	87
25) Benzo(a)pyrene	23.36	252	4144	0.21	ng/ul#	81
26) Indeno(1,2,3-cd)pyrene	25.67	276	4648	0.21	ng/ul	98
27) Dibenzo(a,h)anthracene	25.68	278	3687	0.22	ng/ul#	85
28) Benzo(g,h,i)perylene	26.35	276	4169	0.22	ng/ul	96

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

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