

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM041715\
 Data File : BM001025.D
 Acq On : 16 Apr 2015 20:08
 Operator : TP/IZ
 Sample : SSTD0.830
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.830

Quant Time: Apr 17 06:36:26 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM041715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 06:23:20 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1338	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	4636	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	2690	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	4843	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	4185	0.40	ng/ul	-0.01
20) Perylene-d12	23.36	264	3795	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	11.97	152	4635	0.88	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	8305	0.88	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.42	128	9573	0.88	ng/ul	98
5) 2-Methylnaphthalene	12.03	142	6211	0.89	ng/ul	100
7) Acenaphthylene	13.96	152	8053	0.89	ng/ul	98
8) Acenaphthene	14.30	153	6210	0.86	ng/ul	99
9) Fluorene	15.29	166	7179	0.87	ng/ul	99
11) Pentachlorophenol	16.65	266	336	0.61	ng/ul	93
12) Phenanthrene	17.02	178	11144	0.87	ng/ul	98
13) Anthracene	17.12	178	10140	0.90	ng/ul	97
15) Fluoranthene	19.05	202	12425	0.88	ng/ul	97
17) Pyrene	19.42	202	12505	0.84	ng/ul#	92
18) Benzo(a)anthracene	21.18	228	10148	0.88	ng/ul	98
19) Chrysene	21.23	228	11776	0.87	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	10832	0.86	ng/ul	96
22) Benzo(k)fluoranthene	22.75	252	11432	0.87	ng/ul	95
23) Benzo(a)pyrene	23.26	252	10326	0.88	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.53	276	12165	0.89	ng/ul	99
25) Dibenzo(a,h)anthracene	25.54	278	9926	0.90	ng/ul	96
26) Benzo(g,h,i)perylene	26.19	276	10743	0.88	ng/ul	97

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

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