

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000123.D
 Acq On : 07 Jan 2015 11:43
 Operator : TP/IZ
 Sample : SSTD0.817
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.817

Quant Time: Jan 08 10:13:54 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 10:05:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3156	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	10457	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6054	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	10681	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	9058	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8098	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	11316	0.82	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	18499	0.81	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	22755	0.82	ng/ul	99
5) 2-Methylnaphthalene	12.30	142	15193	0.82	ng/ul	100
7) Acenaphthylene	14.20	152	25772	1.91	ng/ul	99
8) Acenaphthene	14.54	153	15497	0.83	ng/ul	99
9) Fluorene	15.52	166	17756	0.82	ng/ul	97
11) Pentachlorophenol	16.87	266	1309	0.78	ng/ul	97
12) Phenanthrene	17.26	178	26754	0.81	ng/ul	99
13) Anthracene	17.35	178	25535	0.82	ng/ul	98
15) Fluoranthene	19.27	202	29261	0.81	ng/ul	98
17) Pyrene	19.64	202	29921	0.80	ng/ul	99
18) Benzo(a)anthracene	21.39	228	24405	0.82	ng/ul	97
19) Chrysene	21.44	228	26551	0.82	ng/ul	97
21) Benzo(b)fluoranthene	23.00	252	26503	0.83	ng/ul	98
22) Benzo(k)fluoranthene	23.04	252	24812	0.82	ng/ul	95
23) Benzo(a)pyrene	23.60	252	24195	0.84	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.03	276	28109	0.85	ng/ul	100
25) Dibenzo(a,h)anthracene	26.05	278	22598	0.85	ng/ul	97
26) Benzo(g,h,i)perylene	26.74	276	24426	0.84	ng/ul	99

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

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