

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020615\
 Data File : BM000438.D
 Acq On : 06 Feb 2015 17:43
 Operator : TP/IZ
 Sample : SSTD0.483
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.483

Quant Time: Feb 07 00:44:26 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM020615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 07 00:33:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2221	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7159	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4576	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	8596	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	6863	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	5949	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.11	152	4120	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	7923	0.46	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.56	128	8255	0.49	ng/ul	100
5) 2-Methylnaphthalene	12.18	142	5632	0.53	ng/ul#	100
7) Acenaphthylene	14.10	152	7896	0.13	ng/ul	100
8) Acenaphthene	14.43	153	6062	0.15	ng/ul	100
9) Fluorene	15.43	166	7221	0.49	ng/ul	100
11) Pentachlorophenol	16.78	266	167	0.42	ng/ul#	100
12) Phenanthrene	17.16	178	11255	0.15	ng/ul	100
13) Anthracene	17.26	178	9986	0.14	ng/ul	100
15) Fluoranthene	19.19	202	12406	0.16	ng/ul	100
17) Pyrene	19.55	202	12702	0.22	ng/ul	100
18) Benzo(a)anthracene	21.29	228	9168	0.17	ng/ul	100
19) Chrysene	21.34	228	10833	0.19	ng/ul	100
21) Benzo(b)fluoranthene	22.88	252	9336	0.08	ng/ul	100
22) Benzo(k)fluoranthene	22.93	252	10313	0.08	ng/ul	100
23) Benzo(a)pyrene	23.46	252	9043	0.08	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.82	276	10451	0.23	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.84	278	8325	0.60	ng/ul	100
26) Benzo(g,h,i)perylene	26.51	276	9126	0.19	ng/ul	100

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

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