

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020615\
 Data File : BM000437.D
 Acq On : 06 Feb 2015 17:08
 Operator : TP/IZ
 Sample : SSTD0.282
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD0.282

Quant Time: Feb 07 00:44:09 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	1892	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	6203	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4112	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7620	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	6504	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	5819	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	1794	0.08	ng/ul	0.01
14) Fluoranthene-d10	19.15	212	3705	0.24	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.56	128	3614	0.25	ng/ul	97
5) 2-Methylnaphthalene	12.18	142	2471	0.27	ng/ul#	100
7) Acenaphthylene	14.10	152	3362	0.06	ng/ul	98
8) Acenaphthene	14.43	153	2653	0.07	ng/ul	99
9) Fluorene	15.43	166	3142	0.24	ng/ul	99
11) Pentachlorophenol	16.79	266	59	0.17	ng/ul#	100
12) Phenanthrene	17.16	178	4991	0.08	ng/ul	97
13) Anthracene	17.26	178	4441	0.07	ng/ul	96
15) Fluoranthene	19.19	202	5743	0.08	ng/ul	98
17) Pyrene	19.55	202	5890	0.11	ng/ul	98
18) Benzo(a)anthracene	21.30	228	4409	0.09	ng/ul#	91
19) Chrysene	21.35	228	5301	0.10	ng/ul	92
21) Benzo(b)fluoranthene	22.88	252	5054	0.04	ng/ul	90
22) Benzo(k)fluoranthene	22.93	252	4652	0.04	ng/ul	91
23) Benzo(a)pyrene	23.46	252	4513	0.04	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.82	276	5180	0.12	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.84	278	4029	0.30	ng/ul	93
26) Benzo(g,h,i)perylene	26.51	276	4628	0.10	ng/ul	95

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

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