

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051115\
 Data File : BM001239.D
 Acq On : 11 May 2015 10:29
 Operator : TP/IZ
 Sample : SSTD0.157
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.157

Quant Time: May 12 02:56:54 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM051115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 12 02:23:42 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	1177	0.40	ng/ul	0.02
2) Naphthalene-d8	10.25	136	4065	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.13	164	2233	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	4690	0.40	ng/ul	0.02
16) Chrysene-d12	21.09	240	3842	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	3429	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.85	152	522	0.10	ng/ul	0.01
14) Fluoranthene-d10	18.92	212	1053	0.10	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.29	128	1059	0.10	ng/ul#	86
5) 2-Methylnaphthalene	11.92	142	686	0.10	ng/ul	98
7) Acenaphthylene	13.84	152	978	0.10	ng/ul#	90
8) Acenaphthene	14.18	153	785	0.10	ng/ul	91
9) Fluorene	15.19	166	886	0.10	ng/ul	88
12) Phenanthrene	16.92	178	1369	0.10	ng/ul#	94
13) Anthracene	17.01	178	1229	0.10	ng/ul#	92
15) Fluoranthene	18.95	202	1496	0.10	ng/ul	90
17) Pyrene	19.31	202	1578	0.11	ng/ul#	93
18) Benzo(a)anthracene	21.08	228	1185	0.10	ng/ul	95
19) Chrysene	21.13	228	1390m	0.10	ng/ul	
21) Benzo(b)fluoranthene	22.58	252	1181	0.10	ng/ul	78
22) Benzo(k)fluoranthene	22.62	252	1307	0.10	ng/ul#	81
23) Benzo(a)pyrene	23.12	252	1216	0.10	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.31	276	1243	0.09	ng/ul	99
25) Dibenzo(a,h)anthracene	25.32	278	979	0.09	ng/ul#	79
26) Benzo(g,h,i)perylene	25.94	276	1117	0.09	ng/ul#	88

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

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