

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031215\
 Data File : BM000697.D
 Acq On : 12 Mar 2015 13:55
 Operator : TP/IZ
 Sample : SSTD0.171
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.171

Quant Time: Mar 13 09:04:44 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 13 08:51:18 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4411	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15578	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	2415	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19976	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	17734	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	15696	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	2027	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	4455	0.10	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.48	128	4069	0.10	ng/ul	95
5) 2-Methylnaphthalene	12.11	142	2740	0.10	ng/ul	100
7) Acenaphthylene	14.02	152	3756	0.23	ng/ul	97
8) Acenaphthene	14.36	153	2842	0.22	ng/ul	99
9) Fluorene	15.35	166	3514	0.23	ng/ul	97
12) Phenanthrene	17.08	178	5805	0.10	ng/ul	97
13) Anthracene	17.17	178	5170	0.09	ng/ul	96
15) Fluoranthene	19.11	202	6560	0.10	ng/ul	97
17) Pyrene	19.48	202	6800	0.09	ng/ul	99
18) Benzo(a)anthracene	21.23	228	5206	0.09	ng/ul	98
19) Chrysene	21.28	228	6610	0.11	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	5578	0.09	ng/ul	85
22) Benzo(k)fluoranthene	22.82	252	5742	0.10	ng/ul#	81
23) Benzo(a)pyrene	23.34	252	5390	0.10	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.65	276	5360	0.10	ng/ul	99
25) Dibenzo(a,h)anthracene	25.66	278	4246	0.10	ng/ul#	86
26) Benzo(g,h,i)perylene	26.31	276	4778	0.11	ng/ul	94

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

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