

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042715\
 Data File : BM001123.D
 Acq On : 27 Apr 2015 19:09
 Operator : TP/IZ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.499

Quant Time: Apr 28 01:38:26 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM042715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 27 14:06:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	982	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3511	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	1808	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	4026	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	3625	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	3461	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	1796	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	3649	0.40	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.37	128	3671	0.41	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	2382	0.40	ng/ul	100
7) Acenaphthylene	13.91	152	3247	0.39	ng/ul	99
8) Acenaphthene	14.27	153	2551	0.40	ng/ul	97
9) Fluorene	15.26	166	2941	0.40	ng/ul	96
11) Pentachlorophenol	16.63	266	215	0.29	ng/ul	90
12) Phenanthrene	17.00	178	4652	0.40	ng/ul	99
13) Anthracene	17.09	178	4232	0.40	ng/ul	99
15) Fluoranthene	19.02	202	5352	0.40	ng/ul	99
17) Pyrene	19.39	202	5447	0.40	ng/ul	99
18) Benzo(a)anthracene	21.14	228	4498	0.39	ng/ul	99
19) Chrysene	21.20	228	5296	0.42	ng/ul	91
21) Benzo(b)fluoranthene	22.67	252	4888	0.40	ng/ul	100
22) Benzo(k)fluoranthene	22.71	252	5252	0.41	ng/ul	99
23) Benzo(a)pyrene	23.22	252	4865	0.40	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.46	276	5535	0.40	ng/ul	100
25) Dibenzo(a,h)anthracene	25.49	278	4417	0.40	ng/ul	98
26) Benzo(g,h,i)perylene	26.12	276	4873	0.40	ng/ul	99

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

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