

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM102516\
 Data File : BM007726.D
 Acq On : 24 Oct 2016 20:37
 Operator : UM/SJ
 Sample : SSTD0.457
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.457

Quant Time: Oct 25 01:46:41 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 25 01:45:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	910	0.40	ng/ul	0.00
2) Naphthalene-d8	10.55	136	2916	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.40	164	1334	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.14	188	2935	0.40	ng/ul	0.00
16) Chrysene-d12	21.32	240	2804	0.40	ng/ul	0.00
20) Perylene-d12	23.59	264	2741	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.15	152	1428	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.18	212	2810	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.60	128	3091	0.44	ng/ul	100
5) 2-Methylnaphthalene	12.23	142	1960	0.42	ng/ul	100
7) Acenaphthylene	14.11	152	2669	0.45	ng/ul	100
8) Acenaphthene	14.45	153	2051	0.45	ng/ul	100
9) Fluorene	15.46	166	2230	0.41	ng/ul	100
11) Pentachlorophenol	16.83	266	252	0.32	ng/ul	99
12) Phenanthrene	17.18	178	3458	0.43	ng/ul	100
13) Anthracene	17.28	178	3052	0.42	ng/ul	100
15) Fluoranthene	19.21	202	4135	0.41	ng/ul	100
17) Pyrene	19.57	202	4275	0.45	ng/ul	100
18) Benzo(a)anthracene	21.31	228	3487	0.42	ng/ul	100
19) Chrysene	21.37	228	4225	0.46	ng/ul	100
21) Benzo(b)fluoranthene	22.91	252	4359	0.46	ng/ul	100
22) Benzo(k)fluoranthene	22.95	252	4156	0.44	ng/ul	100
23) Benzo(a)pyrene	23.49	252	3839	0.44	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.87	276	4794	0.51	ng/ul	100
25) Dibenzo(a,h)anthracene	25.88	278	3851	0.53	ng/ul	100
26) Benzo(g,h,i)perylene	26.56	276	4368	0.51	ng/ul	100

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

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