

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071415\
 Data File : BM001906.D
 Acq On : 13 Jul 2015 20:57
 Operator : TP/UM
 Sample : SSTD0.243
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.243

Quant Time: Jul 14 04:20:45 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 14 04:19:54 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1187	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4520	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2374	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	5316	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4578	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4095	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	1394	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	2828	0.24	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.49	128	2738	0.22	ng/ul	96
5) 2-Methylnaphthalene	12.12	142	1790	0.23	ng/ul	100
7) Acenaphthylene	14.03	152	2781	0.23	ng/ul#	91
8) Acenaphthene	14.37	153	1956	0.22	ng/ul	99
9) Fluorene	15.37	166	2300	0.23	ng/ul	99
11) Pentachlorophenol	16.72	266	232	0.20	ng/ul	99
12) Phenanthrene	17.09	178	3685	0.22	ng/ul	99
13) Anthracene	17.18	178	3485	0.23	ng/ul	97
15) Fluoranthene	19.12	202	4047	0.23	ng/ul	99
17) Pyrene	19.48	202	4174	0.22	ng/ul	98
18) Benzo(a)anthracene	21.23	228	3565	0.23	ng/ul	99
19) Chrysene	21.28	228	3584	0.22	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	3797	0.23	ng/ul	96
22) Benzo(k)fluoranthene	22.85	252	3056	0.20	ng/ul	94
23) Benzo(a)pyrene	23.37	252	3244	0.22	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.71	276	3489	0.21	ng/ul	100
25) Dibenzo(a,h)anthracene	25.73	278	2801	0.21	ng/ul	95
26) Benzo(g,h,i)perylene	26.40	276	2988	0.21	ng/ul	95

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

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