

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061815\
 Data File : BM001726.D
 Acq On : 17 Jun 2015 20:03
 Operator : TP/IZ
 Sample : SSTD0.205
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.205

Quant Time: Jun 18 04:23:44 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM061815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 18 04:21:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	1214	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4291	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1994	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	4047	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3193	0.40	ng/ul	0.00
20) Perylene-d12	23.48	264	2855	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	1037	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.10	212	1680	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.50	128	2235	0.20	ng/ul	97
5) 2-Methylnaphthalene	12.12	142	1412	0.20	ng/ul	99
7) Acenaphthylene	14.03	152	1891	0.21	ng/ul	97
8) Acenaphthene	14.37	153	1437	0.20	ng/ul	99
9) Fluorene	15.37	166	1595	0.20	ng/ul	99
11) Pentachlorophenol	16.72	266	159	0.18	ng/ul	97
12) Phenanthrene	17.11	178	2453	0.21	ng/ul	99
13) Anthracene	17.20	178	2176	0.20	ng/ul#	92
15) Fluoranthene	19.12	202	2487	0.19	ng/ul	88
17) Pyrene	19.48	202	2559	0.20	ng/ul#	91
18) Benzo(a)anthracene	21.23	228	2072	0.21	ng/ul	98
19) Chrysene	21.29	228	2164	0.20	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	2113	0.20	ng/ul	93
22) Benzo(k)fluoranthene	22.84	252	2081	0.20	ng/ul#	94
23) Benzo(a)pyrene	23.37	252	1907	0.19	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.71	276	2222	0.20	ng/ul	99
25) Dibenzo(a,h)anthracene	25.72	278	1798	0.21	ng/ul	93
26) Benzo(g,h,i)perylene	26.39	276	1978	0.20	ng/ul	94

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

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