

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM071415\
 Data File : BM001905.D
 Acq On : 13 Jul 2015 20:21
 Operator : TP/UM
 Sample : SSTD0.142
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.142

Quant Time: Jul 14 04:22:35 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	1161	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	4378	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	2283	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	5115	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	4339	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	4077	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	686	0.12	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	1346	0.12	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	1350	0.11	ng/ul#	90
5) 2-Methylnaphthalene	12.12	142	891	0.12	ng/ul	100
7) Acenaphthylene	14.03	152	1344	0.12	ng/ul#	92
8) Acenaphthene	14.37	153	970	0.11	ng/ul	98
9) Fluorene	15.37	166	1089	0.11	ng/ul#	98
12) Phenanthrene	17.09	178	1819	0.11	ng/ul	94
13) Anthracene	17.18	178	1709	0.12	ng/ul#	92
15) Fluoranthene	19.12	202	1992	0.12	ng/ul	94
17) Pyrene	19.48	202	2071	0.12	ng/ul	95
18) Benzo(a)anthracene	21.23	228	1804	0.12	ng/ul	96
19) Chrysene	21.28	228	1840	0.12	ng/ul	97
21) Benzo(b)fluoranthene	22.81	252	1903	0.12	ng/ul	86
22) Benzo(k)fluoranthene	22.85	252	1685	0.11	ng/ul#	86
23) Benzo(a)pyrene	23.38	252	1722	0.12	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.71	276	1996	0.12	ng/ul	98
25) Dibenzo(a,h)anthracene	25.73	278	1584	0.12	ng/ul#	86
26) Benzo(g,h,i)perylene	26.40	276	1650	0.11	ng/ul#	89

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

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