

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061815\
 Data File : BM001728.D
 Acq On : 17 Jun 2015 21:15
 Operator : TP/IZ
 Sample : SSTD0.807
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.807

Quant Time: Jun 18 04:24:06 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	1137	0.40	ng/ul	0.00
2) Naphthalene-d8	10.46	136	4024	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.32	164	1912	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.06	188	3958	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	3393	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	2881	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.04	152	4005	0.79	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6875	0.80	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.50	128	8550	0.83	ng/ul	98
5) 2-Methylnaphthalene	12.12	142	5453	0.81	ng/ul	100
7) Acenaphthylene	14.03	152	7395	0.86	ng/ul	97
8) Acenaphthene	14.37	153	5611	0.83	ng/ul	99
9) Fluorene	15.37	166	6295	0.84	ng/ul	99
11) Pentachlorophenol	16.70	266	656	0.76	ng/ul	98
12) Phenanthrene	17.09	178	9513	0.84	ng/ul#	90
13) Anthracene	17.18	178	8796	0.84	ng/ul	97
15) Fluoranthene	19.12	202	10361	0.83	ng/ul	86
17) Pyrene	19.49	202	10674	0.80	ng/ul	98
18) Benzo(a)anthracene	21.23	228	8819	0.84	ng/ul	99
19) Chrysene	21.28	228	9322	0.79	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	8809	0.84	ng/ul	96
22) Benzo(k)fluoranthene	22.84	252	8774	0.82	ng/ul	94
23) Benzo(a)pyrene	23.37	252	8179	0.82	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.70	276	9293	0.84	ng/ul	99
25) Dibenzo(a,h)anthracene	25.71	278	7557	0.86	ng/ul	96
26) Benzo(g,h,i)perylene	26.39	276	8024	0.82	ng/ul	98

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

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