

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020515\
 Data File : BM000428.D
 Acq On : 05 Feb 2015 13:28
 Operator : TP/IZ
 Sample : SSTD3.279
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.279

Quant Time: Feb 05 14:07:05 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 05 13:55:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2304	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7636	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	4716	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	9136	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	8558	0.40	ng/ul	0.00
20) Perylene-d12	23.55	264	7268	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	32232	3.16	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	64901	3.27	ng/ul	0.00

Target Compounds						Qvalue
3) Naphthalene	10.56	128	59814	3.09	ng/ul	97
5) 2-Methylnaphthalene	12.18	142	42343	3.12	ng/ul	99
7) Acenaphthylene	14.10	152	66167	3.37	ng/ul	98
8) Acenaphthene	14.43	153	44291	3.03	ng/ul	100
9) Fluorene	15.43	166	54369	3.13	ng/ul	99
11) Pentachlorophenol	16.78	266	7407	3.57	ng/ul	98
12) Phenanthrene	17.16	178	85253	3.27	ng/ul	97
13) Anthracene	17.16	178	85253	3.55	ng/ul	97
15) Fluoranthene	19.19	202	100535	3.39	ng/ul	99
17) Pyrene	19.55	202	101079	3.08	ng/ul	98
18) Benzo(a)anthracene	21.29	228	86383	3.20	ng/ul	98
19) Chrysene	21.34	228	90571	3.10	ng/ul	99
21) Benzo(b)fluoranthene	22.88	252	82548	3.07	ng/ul	90
22) Benzo(k)fluoranthene	22.92	252	86032	3.32	ng/ul#	88
23) Benzo(a)pyrene	23.46	252	78187	3.17	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.82	276	89813	3.23	ng/ul	100
25) Dibenzo(a,h)anthracene	25.83	278	72448	3.25	ng/ul	94
26) Benzo(g,h,i)perylene	26.50	276	76157	3.14	ng/ul	96

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

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