

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM021215\
 Data File : BM000501.D
 Acq On : 12 Feb 2015 14:08
 Operator : TP/IZ
 Sample : SSTD0.298
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.298

Quant Time: Feb 12 16:34:43 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:25:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	3244	0.40	ng/ul	0.00
2) Naphthalene-d8	10.49	136	10934	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.35	164	6454	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.09	188	12876	0.40	ng/ul	0.00
16) Chrysene-d12	21.29	240	10869	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	9577	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.09	152	2757	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	5404	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.54	128	5060	0.19	ng/ul	98
5) 2-Methylnaphthalene	12.16	142	3402	0.18	ng/ul	99
7) Acenaphthylene	14.07	152	4715	0.18	ng/ul	98
8) Acenaphthene	14.41	153	3561	0.18	ng/ul	99
9) Fluorene	15.40	166	4411	0.18	ng/ul	100
11) Pentachlorophenol	16.76	266	314	0.24	ng/ul	96
12) Phenanthrene	17.13	178	6987	0.19	ng/ul	99
13) Anthracene	17.23	178	6218	0.19	ng/ul	98
15) Fluoranthene	19.16	202	7653	0.18	ng/ul	98
17) Pyrene	19.53	202	7846	0.20	ng/ul	100
18) Benzo(a)anthracene	21.27	228	6013	0.19	ng/ul	99
19) Chrysene	21.32	228	7175	0.19	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	6612	0.20	ng/ul	93
22) Benzo(k)fluoranthene	22.89	252	6446	0.19	ng/ul#	91
23) Benzo(a)pyrene	23.41	252	6049	0.20	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.75	276	6663	0.20	ng/ul	98
25) Dibenzo(a,h)anthracene	25.77	278	5319	0.20	ng/ul	95
26) Benzo(g,h,i)perylene	26.44	276	5949	0.21	ng/ul	98

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

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