

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000134.D
 Acq On : 07 Jan 2015 18:49
 Operator : TP/IZ
 Sample : SSTD0.420
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.420

Quant Time: Jan 08 11:59:27 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-SIM-BM010715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 10:41:47 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3004	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	9855	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5925	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	10692	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	9452	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8559	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	5097	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	9443	0.41	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	10289	0.39	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	6860	0.39	ng/ul	99
7) Acenaphthylene	14.20	152	11178	0.38	ng/ul	98
8) Acenaphthene	14.52	153	7059	0.38	ng/ul#	80
9) Fluorene	15.52	166	8261	0.39	ng/ul	97
11) Pentachlorophenol	16.87	266	544	0.33	ng/ul	90
12) Phenanthrene	17.26	178	13017	0.39	ng/ul	98
13) Anthracene	17.35	178	12308	0.40	ng/ul	98
15) Fluoranthene	19.27	202	14487	0.40	ng/ul	97
17) Pyrene	19.63	202	14828	0.38	ng/ul	98
18) Benzo(a)anthracene	21.39	228	11649	0.37	ng/ul	98
19) Chrysene	21.44	228	13320	0.39	ng/ul	98
21) Benzo(b)fluoranthene	23.00	252	12773	0.38	ng/ul	99
22) Benzo(k)fluoranthene	23.04	252	12640	0.39	ng/ul	96
23) Benzo(a)pyrene	23.60	252	11504	0.38	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	26.02	276	13461	0.39	ng/ul	99
25) Dibenzo(a,h)anthracene	26.04	278	10788	0.39	ng/ul	98
26) Benzo(g,h,i)perylene	26.73	276	11884	0.38	ng/ul	98

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

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