

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000121.D
 Acq On : 07 Jan 2015 10:21
 Operator : TP/IZ
 Sample : SSTD0.215
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.215

Quant Time: Jan 08 10:13:08 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 10:05:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3504	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	11729	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	6985	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11947	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	9082	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8276	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	3190	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	4997	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	6359	0.20	ng/ul	99
5) 2-Methylnaphthalene	12.29	142	4162	0.20	ng/ul	98
7) Acenaphthylene	14.20	152	6829	0.44	ng/ul	99
8) Acenaphthene	14.54	153	4444	0.21	ng/ul	99
9) Fluorene	15.52	166	5009	0.20	ng/ul	99
11) Pentachlorophenol	16.88	266	264	0.14	ng/ul#	88
12) Phenanthrene	17.26	178	7504	0.20	ng/ul	99
13) Anthracene	17.35	178	6831	0.20	ng/ul	99
15) Fluoranthene	19.27	202	7787	0.19	ng/ul	99
17) Pyrene	19.63	202	8057	0.22	ng/ul	98
18) Benzo(a)anthracene	21.39	228	5982	0.20	ng/ul	98
19) Chrysene	21.44	228	6787	0.21	ng/ul	98
21) Benzo(b)fluoranthene	23.00	252	6299	0.19	ng/ul	95
22) Benzo(k)fluoranthene	23.04	252	6472	0.21	ng/ul#	95
23) Benzo(a)pyrene	23.60	252	5879	0.20	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	26.03	276	6644	0.20	ng/ul	99
25) Dibenzo(a,h)anthracene	26.06	278	5252	0.19	ng/ul	93
26) Benzo(g,h,i)perylene	26.74	276	5949	0.20	ng/ul	95

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

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