

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010715\
 Data File : BM000122.D
 Acq On : 07 Jan 2015 10:57
 Operator : TP/IZ
 Sample : SSTD0.416
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.416

Quant Time: Jan 08 10:12:21 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 10:05:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	3041	0.40	ng/ul	0.00
2) Naphthalene-d8	10.64	136	9993	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5977	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	11128	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	9449	0.40	ng/ul	0.00
20) Perylene-d12	23.70	264	8513	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	5184	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	9388	0.39	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
3) Naphthalene	10.68	128	10397	0.39	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	6928	0.39	ng/ul	100
7) Acenaphthylene	14.20	152	11446	0.86	ng/ul	98
8) Acenaphthene	14.54	153	7147	0.39	ng/ul	100
9) Fluorene	15.52	166	8429	0.40	ng/ul	100
11) Pentachlorophenol	16.88	266	537	0.31	ng/ul	99
12) Phenanthrene	17.26	178	13339	0.39	ng/ul	100
13) Anthracene	17.35	178	12279	0.38	ng/ul	100
15) Fluoranthene	19.27	202	14651	0.39	ng/ul	100
17) Pyrene	19.64	202	15131	0.39	ng/ul	100
18) Benzo(a)anthracene	21.39	228	12314	0.40	ng/ul	100
19) Chrysene	21.44	228	13173	0.39	ng/ul	100
21) Benzo(b)fluoranthene	23.00	252	12248	0.36	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	13078	0.41	ng/ul	100
23) Benzo(a)pyrene	23.60	252	11835	0.39	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	26.03	276	13302	0.38	ng/ul	100
25) Dibenzo(a,h)anthracene	26.06	278	10619	0.38	ng/ul	100
26) Benzo(g,h,i)perylene	26.74	276	11987	0.39	ng/ul	100

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

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