

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042417\
 Data File : BM009688.D
 Acq On : 24 Apr 2017 12:36
 Operator : SJ/MA
 Sample : SSTD0.215
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.215

Quant Time: Apr 24 14:19:15 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM042417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 24 14:10:59 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	109	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	367	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	222	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.23	188	596	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	748	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	827	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.22	152	112	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	350	0.19	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.68	128	189	0.18	ng/ul#	41
5) 2-Methylnaphthalene	12.29	142	141	0.19	ng/ul	95
7) Acenaphthylene	14.19	152	205	0.18	ng/ul#	58
8) Acenaphthene	14.54	153	172	0.19	ng/ul	93
9) Fluorene	15.53	166	201	0.19	ng/ul#	88
11) Pentachlorophenol	16.88	266	44	0.15	ng/ul#	57
12) Phenanthrene	17.26	178	345	0.18	ng/ul#	74
13) Anthracene	17.35	178	351	0.19	ng/ul#	71
15) Fluoranthene	19.28	202	465	0.19	ng/ul	77
17) Pyrene	19.65	202	473	0.18	ng/ul#	76
18) Benzo(a)anthracene	21.39	228	503	0.20	ng/ul#	84
19) Chrysene	21.44	228	450	0.19	ng/ul#	80
21) Benzo(b)fluoranthene	23.01	252	544	0.17	ng/ul	79
22) Benzo(k)fluoranthene	23.06	252	544	0.19	ng/ul#	73
23) Benzo(a)pyrene	23.62	252	532	0.18	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	26.07	276	608	0.16	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.08	278	505	0.16	ng/ul#	69
26) Benzo(g,h,i)perylene	26.80	276	527	0.17	ng/ul#	73

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

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