

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030617\
 Data File : BM009082.D
 Acq On : 06 Mar 2017 15:38
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.442

Quant Time: Mar 07 03:30:30 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 03 23:57:02 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	168	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	480	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.52	164	271	0.40	ng/ul	-0.02
10) Phenanthrene-d10	17.28	188	654	0.40	ng/ul	0.00
16) Chrysene-d12	21.45	240	712	0.40	ng/ul	0.00
20) Perylene-d12	23.78	264	655	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	321	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	796	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	560	0.40	ng/ul#	80
5) 2-Methylnaphthalene	12.35	142	393	0.39	ng/ul	98
7) Acenaphthylene	14.25	152	578	0.41	ng/ul#	82
8) Acenaphthene	14.59	153	430	0.40	ng/ul	97
9) Fluorene	15.58	166	495	0.39	ng/ul#	95
11) Pentachlorophenol	16.92	266	121	0.38	ng/ul	94
12) Phenanthrene	17.31	178	820	0.40	ng/ul#	92
13) Anthracene	17.40	178	773	0.38	ng/ul#	92
15) Fluoranthene	19.33	202	1014	0.38	ng/ul	96
17) Pyrene	19.69	202	1026	0.42	ng/ul	94
18) Benzo(a)anthracene	21.43	228	943	0.40	ng/ul	94
19) Chrysene	21.48	228	908	0.40	ng/ul	93
21) Benzo(b)fluoranthene	23.07	252	1063	0.41	ng/ul	90
22) Benzo(k)fluoranthene	23.12	252	905	0.40	ng/ul#	88
23) Benzo(a)pyrene	23.68	252	930	0.40	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.16	276	1096	0.37	ng/ul#	86
25) Dibenzo(a,h)anthracene	26.18	278	904	0.37	ng/ul#	92
26) Benzo(a,h,i)perylene	26.90	276	920	0.37	ng/ul#	88

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

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