

Data Path : Z:\HPCHEM1\BNA M\DATA\BM041816\
 Data File : BM004968.D
 Acq On : 18 Apr 2016 11:15
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.446

Quant Time: Apr 19 11:01:10 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM041816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 19 08:25:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1355	0.40	ng/ul	0.00
4) Naphthalene-d8	10.61	136	5728	0.40	ng/ul	0.00
8) Acenaphthene-d10	14.45	164	3113	0.40	ng/ul	0.00
12) Phenanthrene-d10	17.19	188	6797	0.40	ng/ul	0.00
18) Chrysene-d12	21.39	240	4371	0.40	ng/ul	0.00
22) Perylene-d12	23.67	264	3457	0.40	ng/ul	0.00
System Monitoring Compounds						
2) 1,4-Dioxane-d8	3.31	96	1273	1.95	ng/uL	0.00
6) 2-Methylnaphthalene-d10	12.19	152	2816	0.41	ng/ul	0.00
16) Fluoranthene-d10	19.23	212	5438	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) 1,4-Dioxane	3.34	88	1539	1.91	ng/ul#	23
5) Naphthalene	10.66	128	5556	0.41	ng/ul	95
7) 2-Methylnaphthalene	12.27	142	3745	0.42	ng/ul	98
9) Acenaphthylene	14.16	152	5184	0.41	ng/ul	96
10) Acenaphthene	14.50	153	4295	0.41	ng/ul	87
11) Fluorene	15.50	166	4774	0.42	ng/ul#	83
13) Pentachlorophenol	16.87	266	230	0.29	ng/ul	97
14) Phenanthrene	17.33	178	5210	0.28	ng/ul	97
15) Anthracene	17.33	178	6283	0.39	ng/ul	96
17) Fluoranthene	19.25	202	7564	0.39	ng/ul	95
19) Pyrene	19.62	202	7558	0.42	ng/ul	98
20) Benzo(a)anthracene	21.37	228	4909	0.42	ng/ul#	94
21) Chrysene	21.42	228	5560	0.40	ng/ul	94
23) Benzo(b)fluoranthene	22.98	252	4839	0.39	ng/ul	91
24) Benzo(k)fluoranthene	23.02	252	4615	0.39	ng/ul#	92
25) Benzo(a)pyrene	23.57	252	4234	0.40	ng/ul#	88
26) Indeno(1,2,3-cd)pyrene	25.99	276	4421	0.40	ng/ul#	79
27) Dibenzo(a,h)anthracene	26.01	278	3342	0.40	ng/ul#	84
28) Benzo(g,h,i)perylene	26.69	276	4085	0.41	ng/ul#	86

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

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