

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100815\
 Data File : BM002878.D
 Acq On : 09 Oct 2015 00:00
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 09 07:12:29 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM100815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 09 06:47:52 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	65	0.00
2	1,4-Dioxane	0.499	0.452	9.4	64	0.00
3	n-Nitrosodimethylamine	0.389	0.380	2.3	62	0.00
4 S	2-Fluorophenol	1.069	1.030	3.6	64	0.02
5 S	Phenol-d6	1.321	1.202	9.0	62	0.02
6	bis(2-Chloroethyl)ether	1.303	1.275	2.1	66	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	67	0.00
8 S	Nitrobenzene-d5	0.251	0.233	7.2	63	0.01
9	Nitrobenzene	0.266	0.242	9.0	63	0.00
10	Naphthalene	1.110	1.090	1.8	68	0.00
11	Hexachlorobutadiene	0.176	0.173	1.7	68	0.00
12	2-Methylnaphthalene	0.738	0.741	-0.4	69	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	71	0.00
14 S	2,4,6-Tribromophenol	0.166	0.163	1.8	75	0.00
15 S	2-Fluorobiphenyl	1.514	1.500	0.9	70	0.00
16	Acenaphthylene	2.122	1.991	6.2	67	0.00
17	Acenaphthene	1.370	1.297	5.3	70	0.00
18	Fluorene	1.678	1.654	1.4	71	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
20	4-Bromophenyl-phenylether	0.237	0.227	4.2	74	0.00
21	Hexachlorobenzene	0.292	0.264	9.6	73	0.00
22	Pentachlorophenol	0.027	0.020	25.9#	69	0.00
23	Phenanthrene	1.463	1.369	6.4	76	0.00
24	Anthracene	1.237	1.125	9.1	75	0.00
25	Fluoranthene	1.621	1.610	0.7	84	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	90	0.00
27	Benzydine	0.688	0.590	14.2	88	0.00
28	Pvrene	1.349	1.273	5.6	83	0.00
29 S	Terphenyl-d14	0.667	0.643	3.6	89	-0.02
30	Benzo(a)anthracene	1.435	1.357	5.4	91	0.00
31	3,3'-Dichlorobenzidine	0.551	0.522	5.3	89	0.00
32	Chrysene	1.344	1.390	-3.4	94	0.00
33	Bis(2-ethylhexyl)phthalate	0.874	0.708	19.0	62	-0.02
34	Indeno(1,2,3-cd)pyrene	1.611	1.767	-9.7	102	0.00
35 I	Perylene-d12	1.000	1.000	0.0	98	0.00
36	Benzo(b)fluoranthene	1.431	1.456	-1.7	98	0.00
37	Benzo(k)fluoranthene	1.382	1.282	7.2	97	0.00
38 C	Benzo(a)pyrene	1.316	1.269	3.6	97	-0.01
39	Dibenzo(a,h)anthracene	1.282	1.290	-0.6	102	0.00
40	Benzo(a,h,i)perylene	1.324	1.340	-1.2	102	-0.01

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(#) = Out of Range

SPCC's out = 0 CCC's out = 0