

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120215\
 Data File : BM003366.D
 Acq On : 02 Dec 2015 21:11
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Dec 03 00:44:55 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM120215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 16:19:00 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	66	0.00
2	1,4-Dioxane	0.448	0.420	6.3	69	0.00
3	n-Nitrosodimethylamine	0.461	0.437	5.2	65	0.00
4 S	2-Fluorophenol	0.997	0.865	13.2	64	0.00
5 S	Phenol-d6	1.227	1.086	11.5	62	0.00
6	bis(2-Chloroethyl)ether	1.193	1.143	4.2	66	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	64	0.00
8 S	Nitrobenzene-d5	0.233	0.208	10.7	62	0.00
9	Nitrobenzene	0.252	0.226	10.3	61	0.00
10	Naphthalene	1.077	1.022	5.1	66	0.00
11	Hexachlorobutadiene	0.164	0.161	1.8	68	0.00
12	2-Methylnaphthalene	0.738	0.697	5.6	63	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	58	0.00
14 S	2,4,6-Tribromophenol	0.104	0.113	-8.7	62	0.00
15 S	2-Fluorobiphenyl	1.505	1.502	0.2	64	0.00
16	Acenaphthylene	1.976	1.815	8.1	54	0.00
17	Acenaphthene	1.321	1.332	-0.8	65	-0.02
18	Fluorene	1.490	1.673	-12.3	69	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	55	0.00
20	4-Bromophenyl-phenylether	0.144	0.169	-17.4	75	-0.02
21	Hexachlorobenzene	0.173	0.215	-24.3	80	0.00
22	Pentachlorophenol	0.055	0.043	21.8	47#	-0.02
23	Phenanthrene	0.964	1.114	-15.6	72	-0.02
24	Anthracene	1.016	1.009	0.7	56	0.00
25	Fluoranthene	0.999	1.082	-8.3	59	-0.01
26 I	Chrysene-d12	1.000	1.000	0.0	67	-0.01
27	Benzydine	0.409	0.263	35.7#	49#	0.00
28	Pvrene	1.812	1.395	23.0	57	0.00
29 S	Terphenyl-d14	0.811	0.675	16.8	60	0.00
30	Benzo(a)anthracene	1.350	1.157	14.3	60	-0.01
31	3,3'-Dichlorobenzidine	0.349	0.212	39.3#	43#	0.00
32	Chrysene	1.563	1.143	26.9#	52	0.01
33	Bis(2-ethylhexyl)phthalate	0.684	0.325	52.5#	32#	-0.01
34	Indeno(1,2,3-cd)pyrene	1.163	0.739	36.5#	55	0.00
35 I	Perylene-d12	1.000	1.000	0.0	54	0.00
36	Benzo(b)fluoranthene	1.522	1.469	3.5	54	-0.01
37	Benzo(k)fluoranthene	1.366	1.245	8.9	51	0.00
38 C	Benzo(a)pyrene	1.249	1.056	15.5	49#	0.00
39	Dibenzo(a,h)anthracene	1.136	0.962	15.3	57	0.00
40	Benzo(a,h,i)perylene	1.199	1.011	15.7	58	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range				
SPCC's out = 0 CCC's out = 0				