Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA B\Data\BB093016\
Data File : BB058572.D
Acq On : 30 Sep 2016 18:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_B
LabSampleId : SSTDCCC040

Quant Time: Sep 30 19:02:01 2016
Quant Method : Z:\HPCHEM1\BNA B\METHOD\8270-BB093016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 30 14:16:00 2016
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	88	0.00
2		1,4-Dioxane	0.520	0.611	-17.5	105	0.00
3		Pyridine	1.308	1.555	-18.9	104	0.00
4		n-Nitrosodimethylamine	0.976	1.165	-19.4	105	0.00
5	S	2-Fluorophenol	1.303	1.545	-18.6	104	0.00
6		Aniline	1.894	2.213	-16.8	103	-0.02
7	S	Phenol-d6	1.533	1.811	-18.1	106	-0.02
8		2-Chlorophenol	1.478	1.738	-17.6	105	0.00
9		Benzaldehyde	0.825	0.795	3.6	84	0.00
10	C	Phenol	1.582	1.874	-18.5	103	-0.02
11		bis(2-Chloroethyl)ether	1.386	1.652	-19.2	105	-0.02
12		1,3-Dichlorobenzene	1.500	1.731	-15.4	101	0.00
13	C	1,4-Dichlorobenzene	1.539	1.774	-15.3	102	0.00
14		1,2-Dichlorobenzene	1.440	1.701	-18.1	103	0.00
15		Benzyl Alcohol	1.110	1.320	-18.9	105	-0.02
16		2,2'-oxybis(1-Chloropropane	2.523	3.130	-24.1	106	0.00
17		2-Methylphenol	1.050	1.228	-17.0	102	-0.02
18		Hexachloroethane	0.678	0.811	-19.6	105	0.00
19	P	n-Nitroso-di-n-propylamine	1.106	1.327	-20.0	108	-0.04
20		3+4-Methylphenols	1.372	1.597	-16.4	104	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.459	0.510	-11.1	101	-0.02
23	S	Nitrobenzene-d5	0.415	0.477	-14.9	105	-0.02
24		Nitrobenzene	0.419	0.479	-14.3	103	-0.02
25		Isophorone	0.803	0.900	-12.1	105	-0.04
26	C	2-Nitrophenol	0.239	0.272	-13.8	102	0.00
27		2,4-Dimethylphenol	0.340	0.387	-13.8	106	-0.02
28		bis(2-Chloroethoxy)methane	0.425	0.502	-18.1	107	0.00
29	C	2,4-Dichlorophenol	0.315	0.354	-12.4	103	0.00
30		1,2,4-Trichlorobenzene	0.319	0.359	-12.5	103	-0.02
31		Naphthalene	0.967	1.075	-11.2	100	-0.02
32		Benzoic acid	0.228	0.280	-22.8	115	-0.09
33		4-Chloroaniline	0.429	0.481	-12.1	103	-0.02
34	C	Hexachlorobutadiene	0.169	0.186	-10.1	100	0.00
35		Caprolactam	0.131	0.150	-14.5	109	-0.11
36	C	4-Chloro-3-methylphenol	0.346	0.391	-13.0	103	-0.02
37		2-Methylnaphthalene	0.633	0.713	-12.6	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	104	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.570	0.588	-3.2	105	0.00
40	P	Hexachlorocyclopentadiene	0.278	0.284	-2.2	100	0.00
41	S	2,4,6-Tribromophenol	0.256	0.265	-3.5	100	-0.02
42	C	2,4,6-Trichlorophenol	0.437	0.438	-0.2	102	0.00
43		2,4,5-Trichlorophenol	0.452	0.440	2.7	102	-0.02
44	S	2-Fluorobiphenyl	1.210	1.220	-0.8	94	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.573	1.606	-2.1	99	-0.02
46		2-Chloronaphthalene	1.248	1.218	2.4	99	-0.02
47		2-Nitroaniline	0.497	0.500	-0.6	100	-0.02
48		Acenaphthylene	1.916	2.007	-4.7	108	0.00
49		Dimethylphthalate	1.463	1.528	-4.4	106	-0.04
50		2,6-Dinitrotoluene	0.419	0.439	-4.8	107	-0.02
51	C	Acenaphthene	1.300	1.282	1.4	97	-0.02
52		3-Nitroaniline	0.441	0.469	-6.3	111	-0.02
53	P	2,4-Dinitrophenol	0.270	0.273	-1.1	104	-0.04
54		Dibenzofuran	1.634	1.664	-1.8	103	0.00
55	P	4-Nitrophenol	0.392	0.393	-0.3	98	-0.04
56		2,4-Dinitrotoluene	0.554	0.551	0.5	104	-0.04
57		Fluorene	1.346	1.435	-6.6	110	0.00
58		2,3,4,6-Tetrachlorophenol	0.335	0.322	3.9	100	0.00
59		Diethylphthalate	1.582	1.643	-3.9	108	-0.04
60		4-Chlorophenyl-phenylether	0.626	0.650	-3.8	102	0.00
61		4-Nitroaniline	0.441	0.469	-6.3	111	-0.02
62		Azobenzene	1.598	1.665	-4.2	111	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
64		4,6-Dinitro-2-methylphenol	0.210	0.221	-5.2	95	-0.04
65	c	n-Nitrosodiphenylamine	0.706	0.822	-16.4	110	-0.02
66		4-Bromophenyl-phenylether	0.222	0.230	-3.6	91	-0.02
67		Hexachlorobenzene	0.272	0.300	-10.3	108	0.00
68		Atrazine	0.229	0.231	-0.9	90	-0.02
69	C	Pentachlorophenol	0.176	0.189	-7.4	95	-0.02
70		Phenanthrene	1.064	1.183	-11.2	108	-0.02
71		Anthracene	1.087	1.174	-8.0	99	-0.02
72		Carbazole	1.079	1.108	-2.7	94	-0.02
73		Di-n-butylphthalate	1.554	1.584	-1.9	100	-0.02
74	C	Fluoranthene	1.065	1.108	-4.0	92	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	93	-0.02
76		Benzidine	0.627	0.608	3.0	92	0.00
77		Pyrene	1.194	1.219	-2.1	101	-0.02
78	S	Terphenyl-d14	0.701	0.682	2.7	91	-0.02
79		Butylbenzylphthalate	0.852	0.860	-0.9	89	-0.02
80		Benzo(a)anthracene	1.069	1.171	-9.5	103	-0.02
81		3,3'-Dichlorobenzidine	0.415	0.435	-4.8	101	-0.02
82		Chrysene	1.021	1.142	-11.9	104	-0.02
83		Bis(2-ethylhexyl)phthalate	1.078	1.134	-5.2	96	0.00
84	c	Di-n-octyl phthalate	1.885	1.995	-5.8	98	0.00
85		Indeno(1,2,3-cd)pyrene	1.015	1.027	-1.2	95	-0.10
86	I	Perylene-d12	1.000	1.000	0.0	91	-0.02
87		Benzo(b)fluoranthene	1.257	1.286	-2.3	93	-0.02

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88	Benzo(k)fluoranthene	1.030	1.251	-21.5	106	-0.04
89 C	Benzo(a)pyrene	1.076	1.211	-12.5	99	-0.02
90	Dibenzo(a,h)anthracene	0.896	0.970	-8.3	95	-0.06
91	Benzo(g,h,i)perylene	0.850	0.904	-6.4	94	-0.06

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