Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031017\
Data File : BF093545.D
Acq On : 11 Mar 2017 2:10
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 03:46:21 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 07 15:55:22 2017
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	89	0.02
2		1,4-Dioxane	0.662	0.698	-5.4	91	0.03
3		Pyridine	1.688	1.799	-6.6	87	0.05
4		n-Nitrosodimethylamine	0.806	0.846	-5.0	99	0.05
5	S	2-Fluorophenol	1.202	1.351	-12.4	100	0.02
6		Aniline	2.080	1.986	4.5	91	0.02
7	S	Phenol-d6	1.515	1.548	-2.2	95	0.03
8		2-Chlorophenol	1.346	1.338	0.6	90	0.03
9		Benzaldehyde	1.017	0.893	12.2	87	0.02
10	C	Phenol	1.857	1.933	-4.1	102	0.02
11		bis(2-Chloroethyl)ether	1.501	1.497	0.3	93	0.02
12		1,3-Dichlorobenzene	1.541	1.551	-0.6	94	0.02
13	C	1,4-Dichlorobenzene	1.578	1.578	0.0	93	0.02
14		1,2-Dichlorobenzene	1.397	1.384	0.9	90	0.02
15		Benzyl Alcohol	1.080	1.059	1.9	94	0.02
16		2,2'-oxybis(1-Chloropropane	2.493	2.633	-5.6	95	0.01
17		2-Methylphenol	1.139	1.115	2.1	91	0.02
18		Hexachloroethane	0.532	0.563	-5.8	97	0.02
19	P	n-Nitroso-di-n-propylamine	1.042	1.072	-2.9	102	0.02
20		3+4-Methylphenols	1.477	1.610	-9.0	98	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	86	0.01
22		Acetophenone	0.481	0.501	-4.2	93	0.02
23	S	Nitrobenzene-d5	0.360	0.363	-0.8	89	0.01
24		Nitrobenzene	0.405	0.453	-11.9	94	0.02
25		Isophorone	0.727	0.752	-3.4	92	0.02
26	C	2-Nitrophenol	0.185	0.184	0.5	82	0.02
27		2,4-Dimethylphenol	0.301	0.304	-1.0	81	0.02
28		bis(2-Chloroethoxy)methane	0.459	0.450	2.0	89	0.01
29	C	2,4-Dichlorophenol	0.283	0.303	-7.1	93	0.02
30		1,2,4-Trichlorobenzene	0.301	0.303	-0.7	87	0.02
31		Naphthalene	1.002	0.973	2.9	88	0.01
32		Benzoic acid	0.156	0.120	23.1	70	0.00
33		4-Chloroaniline	0.407	0.406	0.2	86	0.02
34	C	Hexachlorobutadiene	0.155	0.162	-4.5	94	0.02
35		Caprolactam	0.097	0.101	-4.1	87	0.02
36	C	4-Chloro-3-methylphenol	0.302	0.296	2.0	84	0.02
37		2-Methylnaphthalene	0.638	0.646	-1.3	88	0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.01
39		1,2,4,5-Tetrachlorobenzene	0.546	0.570	-4.4	83	0.01
40	P	Hexachlorocyclopentadiene	0.098	0.138	-40.8#	111	0.01
41	S	2,4,6-Tribromophenol	0.130	0.139	-6.9	86	0.02
42	C	2,4,6-Trichlorophenol	0.341	0.376	-10.3	88	0.01
43		2,4,5-Trichlorophenol	0.366	0.381	-4.1	79	0.02
44	S	2-Fluorobiphenyl	1.075	1.077	-0.2	80	0.01

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Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.457	1.527	-4.8	87	0.01
46		2-Chloronaphthalene	1.115	1.207	-8.3	82	0.01
47		2-Nitroaniline	0.394	0.446	-13.2	89	0.01
48		Acenaphthylene	1.839	1.881	-2.3	81	0.01
49		Dimethylphthalate	1.326	1.412	-6.5	90	0.01
50		2,6-Dinitrotoluene	0.291	0.282	3.1	86	0.01
51	C	Acenaphthene	1.088	1.077	1.0	78	0.01
52		3-Nitroaniline	0.350	0.382	-9.1	98	0.02
53	P	2,4-Dinitrophenol	0.132	0.016#	87.9#	8#	0.05
54		Dibenzofuran	1.361	1.552	-14.0	87	0.01
55	P	4-Nitrophenol	0.170	0.065	61.8#	28#	0.03
56		2,4-Dinitrotoluene	0.357	0.400	-12.0	98	0.01
57		Fluorene	1.154	1.241	-7.5	80	0.01
58		2,3,4,6-Tetrachlorophenol	0.258	0.287	-11.2	81	0.02
59		Diethylphthalate	1.267	1.408	-11.1	90	0.02
60		4-Chlorophenyl-phenylether	0.517	0.609	-17.8	87	0.02
61		4-Nitroaniline	0.357	0.378	-5.9	93	0.02
62		Azobenzene	1.440	1.690	-17.4	97	0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.01
64		4,6-Dinitro-2-methylphenol	0.130	0.081	37.7#	56	0.02
65	c	n-Nitrosodiphenylamine	0.668	0.649	2.8	85	0.02
66		4-Bromophenyl-phenylether	0.211	0.199	5.7	87	0.01
67		Hexachlorobenzene	0.202	0.194	4.0	84	0.02
68		Atrazine	0.195	0.196	-0.5	90	0.02
69	C	Pentachlorophenol	0.070	0.095	-35.7#	120	0.02
70		Phenanthrene	1.142	1.093	4.3	88	0.02
71		Anthracene	1.155	1.119	3.1	96	0.02
72		Carbazole	1.085	1.090	-0.5	96	0.01
73		Di-n-butylphthalate	1.255	1.191	5.1	90	0.01
74	C	Fluoranthene	1.103	1.098	0.5	92	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	87	0.02
76		Benzidine	0.658	0.410	37.7#	55	0.02
77		Pyrene	1.455	1.457	-0.1	90	0.02
78	S	Terphenyl-d14	0.769	0.872	-13.4	90	0.02
79		Butylbenzylphthalate	0.612	0.673	-10.0	90	0.02
80		Benzo(a)anthracene	1.181	1.201	-1.7	88	0.03
81		3,3'-Dichlorobenzidine	0.414	0.401	3.1	81	0.02
82		Chrysene	1.093	1.235	-13.0	87	0.02
83		Bis(2-ethylhexyl)phthalate	0.853	0.963	-12.9	92	0.02
84	c	Di-n-octyl phthalate	1.422	1.431	-0.6	84	0.02
85		Indeno(1,2,3-cd)pyrene	0.915	0.897	2.0	85	0.06
86	I	Perylene-d12	1.000	1.000	0.0	79	0.02
87		Benzo(b)fluoranthene	1.218	1.286	-5.6	75	0.02

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88	Benzo(k)fluoranthene	1.175	1.148	2.3	92	0.02
89 C	Benzo(a)pyrene	1.113	1.133	-1.8	80	0.03
90	Dibenzo(a,h)anthracene	0.921	0.938	-1.8	84	0.05
91	Benzo(g,h,i)perylene	0.945	0.983	-4.0	86	0.06

(#) = Out of Range SPCC's out = 1 CCC's out = 1