Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012716\
Data File : BF084622.D
Acq On : 27 Jan 2016 15:50
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 27 18:12:06 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 25 16:24:07 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	82	0.00
2		1,4-Dioxane	0.586	0.527	10.1	71	-0.01
3		Pyridine	1.633	1.494	8.5	69	-0.01
4		n-Nitrosodimethylamine	0.838	0.779	7.0	72	0.00
5	S	2-Fluorophenol	1.236	1.242	-0.5	88	0.00
6		Aniline	2.272	1.966	13.5	69	0.00
7	S	Phenol-d6	1.553	1.427	8.1	75	0.00
8		2-Chlorophenol	1.403	1.400	0.2	83	0.01
9		Benzaldehyde	1.011	0.836	17.3	68	0.00
10	C	Phenol	1.766	1.582	10.4	68	0.00
11		bis(2-Chloroethyl)ether	1.368	1.385	-1.2	87	0.00
12		1,3-Dichlorobenzene	1.447	1.474	-1.9	86	0.00
13	C	1,4-Dichlorobenzene	1.451	1.454	-0.2	79	0.00
14		1,2-Dichlorobenzene	1.390	1.228	11.7	78	0.00
15		Benzyl Alcohol	1.306	1.208	7.5	72	0.00
16		2,2'-oxybis(1-Chloropropane	2.491	2.168	13.0	73	0.00
17		2-Methylphenol	1.176	1.072	8.8	70	0.00
18		Hexachloroethane	0.580	0.592	-2.1	81	0.00
19	P	n-Nitroso-di-n-propylamine	1.051	0.919	12.6	74	0.00
20		3+4-Methylphenols	1.382	1.307	5.4	83	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
22		Acetophenone	0.481	0.453	5.8	75	0.00
23	S	Nitrobenzene-d5	0.359	0.350	2.5	84	0.00
24		Nitrobenzene	0.364	0.328	9.9	69	0.00
25		Isophorone	0.687	0.667	2.9	82	0.01
26	C	2-Nitrophenol	0.166	0.184	-10.8	96	0.00
27		2,4-Dimethylphenol	0.336	0.294	12.5	70	0.00
28		bis(2-Chloroethoxy)methane	0.420	0.392	6.7	85	0.00
29	C	2,4-Dichlorophenol	0.280	0.252	10.0	78	0.00
30		1,2,4-Trichlorobenzene	0.289	0.290	-0.3	82	0.00
31		Naphthalene	0.907	0.923	-1.8	87	0.00
32		Benzoic acid	0.198	0.226	-14.1	92	0.01
33		4-Chloroaniline	0.416	0.438	-5.3	92	0.00
34	C	Hexachlorobutadiene	0.166	0.151	9.0	77	0.00
35		Caprolactam	0.094	0.097	-3.2	78	0.01
36	C	4-Chloro-3-methylphenol	0.313	0.287	8.3	81	0.00
37		2-Methylnaphthalene	0.651	0.579	11.1	78	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	67	0.00
39		1,2,4,5-Tetrachlorobenzene	0.575	0.738	-28.3#	86	0.00
40	P	Hexachlorocyclopentadiene	0.347	0.398	-14.7	75	0.00
41	S	2,4,6-Tribromophenol	0.202	0.227	-12.4	71	0.01
42	C	2,4,6-Trichlorophenol	0.430	0.478	-11.2	77	0.00
43		2,4,5-Trichlorophenol	0.412	0.510	-23.8	80	0.00
44	S	2-Fluorobiphenyl	1.317	1.526	-15.9	88	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	1.590	1.743	-9.6	79	0.00
46		2-Chloronaphthalene	1.249	1.367	-9.4	82	0.00
47		2-Nitroaniline	0.412	0.530	-28.6#	90	0.00
48		Acenaphthylene	1.845	1.688	8.5	59	0.00
49		Dimethylphthalate	1.430	1.924	-34.5#	87	0.01
50		2,6-Dinitrotoluene	0.314	0.452	-43.9#	87	0.01
51	C	Acenaphthene	1.237	1.261	-1.9	64	0.00
52		3-Nitroaniline	0.389	0.398	-2.3	64	0.00
53	P	2,4-Dinitrophenol	0.093	0.187	-101.1#	95	0.00
54		Dibenzofuran	1.812	1.506	16.9	58	0.00
55	P	4-Nitrophenol	0.282	0.335	-18.8	67	0.01
56		2,4-Dinitrotoluene	0.397	0.477	-20.2	69	0.01
57		Fluorene	1.400	1.168	16.6	60	0.00
58		2,3,4,6-Tetrachlorophenol	0.352	0.379	-7.7	69	0.00
59		Diethylphthalate	1.410	1.374	2.6	64	0.00
60		4-Chlorophenyl-phenylether	0.564	0.678	-20.2	81	0.00
61		4-Nitroaniline	0.346	0.394	-13.9	78	0.01
62		Azobenzene	1.546	1.599	-3.4	68	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	69	0.00
64		4,6-Dinitro-2-methylphenol	0.109	0.145	-33.0#	88	0.01
65	c	n-Nitrosodiphenylamine	0.639	0.665	-4.1	71	0.01
66		4-Bromophenyl-phenylether	0.215	0.227	-5.6	68	0.00
67		Hexachlorobenzene	0.242	0.268	-10.7	81	0.00
68		Atrazine	0.209	0.188	10.0	55	0.00
69	C	Pentachlorophenol	0.126	0.139	-10.3	67	0.00
70		Phenanthrene	1.046	1.027	1.8	62	0.00
71		Anthracene	1.026	1.129	-10.0	79	0.01
72		Carbazole	1.003	0.932	7.1	61	0.00
73		Di-n-butylphthalate	1.111	1.245	-12.1	75	0.00
74	C	Fluoranthene	1.093	1.163	-6.4	76	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	70	0.01
76		Benzidine	0.717	0.675	5.9	65	0.01
77		Pyrene	1.569	1.509	3.8	76	0.01
78	S	Terphenyl-d14	0.896	0.919	-2.6	83	0.00
79		Butylbenzylphthalate	0.679	0.693	-2.1	69	0.00
80		Benzo(a)anthracene	1.174	1.242	-5.8	78	0.00
81		3,3'-Dichlorobenzidine	0.410	0.475	-15.9	79	0.00
82		Chrysene	1.128	1.232	-9.2	77	0.00
83		Bis(2-ethylhexyl)phthalate	0.858	0.937	-9.2	79	0.01
84	c	Di-n-octyl phthalate	1.449	1.459	-0.7	66	0.00
85		Indeno(1,2,3-cd)pyrene	0.895	0.963	-7.6	76	0.01
86	I	Perylene-d12	1.000	1.000	0.0	81	0.01
87		Benzo(b)fluoranthene	1.308	1.228	6.1	73	0.00

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88	Benzo(k)fluoranthene	1.070	1.152	-7.7	88	0.01
89 C	Benzo(a)pyrene	1.110	1.077	3.0	75	0.00
90	Dibenzo(a,h)anthracene	0.955	0.852	10.8	75	0.01
91	Benzo(g,h,i)perylene	0.989	1.124	-13.7	98	0.01

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