Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030416\
Data File : BF085202.D
Acq On : 4 Mar 2016 15:01
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 04 23:48:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 04 00:54:58 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	59	-0.01
2		1,4-Dioxane	0.610	0.612	-0.3	62	0.00
3		Pyridine	1.526	1.589	-4.1	64	0.00
4		n-Nitrosodimethylamine	0.653	0.602	7.8	54	-0.01
5	S	2-Fluorophenol	1.192	1.313	-10.2	72	0.00
6		Aniline	2.189	2.170	0.9	58	-0.01
7	S	Phenol-d6	1.562	1.508	3.5	61	-0.01
8		2-Chlorophenol	1.435	1.448	-0.9	62	-0.01
9		Benzaldehyde	0.804	0.775	3.6	65	-0.01
10	C	Phenol	1.673	1.650	1.4	62	-0.01
11		bis(2-Chloroethyl)ether	1.390	1.480	-6.5	60	-0.01
12		1,3-Dichlorobenzene	1.541	1.612	-4.6	63	-0.01
13	C	1,4-Dichlorobenzene	1.545	1.648	-6.7	69	0.00
14		1,2-Dichlorobenzene	1.401	1.394	0.5	58	-0.01
15		Benzyl Alcohol	1.147	1.125	1.9	60	-0.01
16		2,2'-oxybis(1-Chloropropane	2.003	1.941	3.1	61	-0.01
17		2-Methylphenol	1.171	1.220	-4.2	60	-0.01
18		Hexachloroethane	0.570	0.572	-0.4	60	-0.01
19	P	n-Nitroso-di-n-propylamine	1.018	0.936	8.1	55	-0.02
20		3+4-Methylphenols	1.387	1.381	0.4	65	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	63	0.00
22		Acetophenone	0.488	0.476	2.5	61	-0.01
23	S	Nitrobenzene-d5	0.374	0.361	3.5	59	-0.01
24		Nitrobenzene	0.380	0.369	2.9	60	-0.01
25		Isophorone	0.693	0.647	6.6	58	-0.01
26	C	2-Nitrophenol	0.185	0.203	-9.7	65	-0.01
27		2,4-Dimethylphenol	0.338	0.351	-3.8	62	-0.01
28		bis(2-Chloroethoxy)methane	0.386	0.353	8.5	59	-0.01
29	C	2,4-Dichlorophenol	0.272	0.281	-3.3	64	-0.01
30		1,2,4-Trichlorobenzene	0.294	0.294	0.0	63	-0.01
31		Naphthalene	0.983	1.002	-1.9	69	0.00
32		Benzoic acid	0.224	0.168	25.0	43#	-0.05
33		4-Chloroaniline	0.398	0.381	4.3	64	-0.01
34	C	Hexachlorobutadiene	0.153	0.164	-7.2	66	-0.01
35		Caprolactam	0.089	0.080	10.1	56	-0.02
36	C	4-Chloro-3-methylphenol	0.309	0.289	6.5	60	-0.01
37		2-Methylnaphthalene	0.631	0.606	4.0	59	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	57	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.572	0.647	-13.1	70	0.00
40	P	Hexachlorocyclopentadiene	0.308	0.311	-1.0	57	0.00
41	S	2,4,6-Tribromophenol	0.171	0.174	-1.8	56	-0.01
42	C	2,4,6-Trichlorophenol	0.426	0.451	-5.9	61	-0.01
43		2,4,5-Trichlorophenol	0.404	0.415	-2.7	58	-0.01
44	S	2-Fluorobiphenyl	1.315	1.374	-4.5	64	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.709	1.888	-10.5	71	0.00
46		2-Chloronaphthalene	1.302	1.368	-5.1	62	-0.01
47		2-Nitroaniline	0.404	0.449	-11.1	62	0.00
48		Acenaphthylene	2.027	2.070	-2.1	61	-0.01
49		Dimethylphthalate	1.535	1.656	-7.9	61	-0.01
50		2,6-Dinitrotoluene	0.328	0.326	0.6	55	-0.01
51	C	Acenaphthene	1.231	1.259	-2.3	62	-0.01
52		3-Nitroaniline	0.393	0.438	-11.5	58	-0.01
53	P	2,4-Dinitrophenol	0.137	0.131	4.4	53	0.00
54		Dibenzofuran	1.613	1.592	1.3	58	-0.01
55	P	4-Nitrophenol	0.309	0.326	-5.5	55	-0.01
56		2,4-Dinitrotoluene	0.420	0.407	3.1	55	-0.01
57		Fluorene	1.267	1.293	-2.1	58	-0.01
58		2,3,4,6-Tetrachlorophenol	0.284	0.281	1.1	54	-0.01
59		Diethylphthalate	1.490	1.571	-5.4	61	-0.01
60		4-Chlorophenyl-phenylether	0.616	0.641	-4.1	63	0.00
61		4-Nitroaniline	0.367	0.347	5.4	52	-0.01
62		Azobenzene	1.468	1.486	-1.2	57	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	57	-0.01
64		4,6-Dinitro-2-methylphenol	0.120	0.114	5.0	48#	-0.01
65	c	n-Nitrosodiphenylamine	0.622	0.683	-9.8	66	0.00
66		4-Bromophenyl-phenylether	0.194	0.200	-3.1	57	-0.01
67		Hexachlorobenzene	0.207	0.209	-1.0	56	-0.01
68		Atrazine	0.173	0.212	-22.5	81	0.00
69	C	Pentachlorophenol	0.115	0.124	-7.8	61	0.00
70		Phenanthrene	1.048	1.102	-5.2	66	-0.01
71		Anthracene	1.113	1.133	-1.8	62	-0.01
72		Carbazole	0.976	0.904	7.4	54	-0.01
73		Di-n-butylphthalate	1.233	1.303	-5.7	64	0.00
74	C	Fluoranthene	1.052	1.025	2.6	58	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	68	-0.01
76		Benzidine	0.595	0.489	17.8	48#	-0.01
77		Pyrene	1.505	1.496	0.6	60	-0.01
78	S	Terphenyl-d14	0.862	0.931	-8.0	72	0.00
79		Butylbenzylphthalate	0.683	0.706	-3.4	65	0.00
80		Benzo(a)anthracene	1.197	1.126	5.9	63	0.00
81		3,3'-Dichlorobenzidine	0.378	0.417	-10.3	69	0.00
82		Chrysene	1.106	1.037	6.2	57	-0.01
83		Bis(2-ethylhexyl)phthalate	0.899	0.874	2.8	63	0.00
84	c	Di-n-octyl phthalate	1.369	1.228	10.3	55	-0.01
85		Indeno(1,2,3-cd)pyrene	0.863	0.962	-11.5	63	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	60	-0.01
87		Benzo(b)fluoranthene	1.333	1.228	7.9	52	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.075	1.194	-11.1	76	0.00
89 C	Benzo(a)pyrene	1.105	1.101	0.4	59	-0.01
90	Dibenzo(a,h)anthracene	0.943	1.118	-18.6	64	0.00
91	Benzo(g,h,i)perylene	0.948	1.129	-19.1	63	0.00

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