Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF030416\
Data File : BF085227.D
Acq On : 5 Mar 2016 2:39
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 05 03:51:11 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	67	-0.01
2		1,4-Dioxane	0.610	0.585	4.1	67	0.01
3		Pyridine	1.526	1.522	0.3	69	0.00
4		n-Nitrosodimethylamine	0.653	0.597	8.6	60	-0.01
5	S	2-Fluorophenol	1.192	1.104	7.4	68	-0.01
6		Aniline	2.189	2.095	4.3	64	-0.01
7	S	Phenol-d6	1.562	1.524	2.4	69	-0.01
8		2-Chlorophenol	1.435	1.464	-2.0	70	-0.01
9		Benzaldehyde	0.804	0.760	5.5	72	-0.01
10	C	Phenol	1.673	1.803	-7.8	77	-0.01
11		bis(2-Chloroethyl)ether	1.390	1.330	4.3	61	-0.01
12		1,3-Dichlorobenzene	1.541	1.557	-1.0	69	-0.01
13	C	1,4-Dichlorobenzene	1.545	1.455	5.8	69	-0.01
14		1,2-Dichlorobenzene	1.401	1.457	-4.0	68	-0.01
15		Benzyl Alcohol	1.147	1.169	-1.9	70	-0.01
16		2,2'-oxybis(1-Chloropropane	2.003	1.793	10.5	63	-0.01
17		2-Methylphenol	1.171	1.160	0.9	65	-0.01
18		Hexachloroethane	0.570	0.563	1.2	66	-0.01
19	P	n-Nitroso-di-n-propylamine	1.018	0.876	13.9	58	-0.02
20		3+4-Methylphenols	1.387	1.324	4.5	70	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	66	-0.01
22		Acetophenone	0.488	0.520	-6.6	70	-0.01
23	S	Nitrobenzene-d5	0.374	0.357	4.5	61	-0.01
24		Nitrobenzene	0.380	0.422	-11.1	72	-0.01
25		Isophorone	0.693	0.711	-2.6	67	-0.01
26	C	2-Nitrophenol	0.185	0.193	-4.3	65	-0.01
27		2,4-Dimethylphenol	0.338	0.340	-0.6	63	-0.02
28		bis(2-Chloroethoxy)methane	0.386	0.394	-2.1	69	-0.01
29	C	2,4-Dichlorophenol	0.272	0.278	-2.2	66	-0.02
30		1,2,4-Trichlorobenzene	0.294	0.308	-4.8	70	-0.01
31		Naphthalene	0.983	0.978	0.5	70	-0.01
32		Benzoic acid	0.224	0.123	45.1#	33#	-0.03
33		4-Chloroaniline	0.398	0.397	0.3	70	-0.01
34	C	Hexachlorobutadiene	0.153	0.170	-11.1	72	-0.01
35		Caprolactam	0.089	0.084	5.6	62	-0.02
36	C	4-Chloro-3-methylphenol	0.309	0.330	-6.8	72	-0.01
37		2-Methylnaphthalene	0.631	0.671	-6.3	69	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	67	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.572	0.565	1.2	71	-0.01
40	P	Hexachlorocyclopentadiene	0.308	0.293	4.9	62	-0.01
41	S	2,4,6-Tribromophenol	0.171	0.171	0.0	65	-0.01
42	C	2,4,6-Trichlorophenol	0.426	0.437	-2.6	68	-0.01
43		2,4,5-Trichlorophenol	0.404	0.369	8.7	60	-0.01
44	S	2-Fluorobiphenyl	1.315	1.342	-2.1	73	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.709	1.620	5.2	71	-0.01
46		2-Chloronaphthalene	1.302	1.305	-0.2	68	-0.01
47		2-Nitroaniline	0.404	0.388	4.0	62	0.00
48		Acenaphthylene	2.027	2.023	0.2	70	-0.01
49		Dimethylphthalate	1.535	1.549	-0.9	67	-0.01
50		2,6-Dinitrotoluene	0.328	0.349	-6.4	68	-0.01
51	C	Acenaphthene	1.231	1.159	5.8	66	-0.01
52		3-Nitroaniline	0.393	0.406	-3.3	63	-0.01
53	P	2,4-Dinitrophenol	0.137	0.066	51.8#	31#	-0.01
54		Dibenzofuran	1.613	1.631	-1.1	69	-0.01
55	P	4-Nitrophenol	0.309	0.279	9.7	54	-0.01
56		2,4-Dinitrotoluene	0.420	0.429	-2.1	68	-0.01
57		Fluorene	1.267	1.283	-1.3	68	-0.01
58		2,3,4,6-Tetrachlorophenol	0.284	0.289	-1.8	65	-0.01
59		Diethylphthalate	1.490	1.500	-0.7	68	-0.01
60		4-Chlorophenyl-phenylether	0.616	0.597	3.1	68	-0.01
61		4-Nitroaniline	0.367	0.317	13.6	55	-0.01
62		Azobenzene	1.468	1.296	11.7	58	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	66	-0.01
64		4,6-Dinitro-2-methylphenol	0.120	0.086	28.3#	42#	-0.01
65	c	n-Nitrosodiphenylamine	0.622	0.602	3.2	68	-0.01
66		4-Bromophenyl-phenylether	0.194	0.196	-1.0	65	-0.01
67		Hexachlorobenzene	0.207	0.208	-0.5	65	-0.01
68		Atrazine	0.173	0.172	0.6	76	-0.01
69	C	Pentachlorophenol	0.115	0.108	6.1	61	-0.01
70		Phenanthrene	1.048	0.975	7.0	68	-0.01
71		Anthracene	1.113	1.115	-0.2	71	-0.01
72		Carbazole	0.976	0.937	4.0	65	-0.01
73		Di-n-butylphthalate	1.233	1.240	-0.6	71	0.00
74	C	Fluoranthene	1.052	1.019	3.1	67	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	72	-0.01
76		Benzidine	0.595	0.626	-5.2	65	-0.01
77		Pyrene	1.505	1.551	-3.1	66	-0.01
78	S	Terphenyl-d14	0.862	0.910	-5.6	74	-0.01
79		Butylbenzylphthalate	0.683	0.759	-11.1	74	0.00
80		Benzo(a)anthracene	1.197	1.161	3.0	68	0.00
81		3,3'-Dichlorobenzidine	0.378	0.387	-2.4	68	-0.01
82		Chrysene	1.106	1.155	-4.4	67	-0.01
83		Bis(2-ethylhexyl)phthalate	0.899	0.916	-1.9	70	-0.01
84	c	Di-n-octyl phthalate	1.369	1.382	-0.9	65	-0.01
85		Indeno(1,2,3-cd)pyrene	0.863	0.952	-10.3	66	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	65	-0.01
87		Benzo(b)fluoranthene	1.333	1.427	-7.1	66	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.075	0.974	9.4	68	-0.01
89 C	Benzo(a)pyrene	1.105	1.134	-2.6	66	-0.01
90	Dibenzo(a,h)anthracene	0.943	1.059	-12.3	66	-0.02
91	Benzo(g,h,i)perylene	0.948	1.071	-13.0	66	-0.02

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