Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF012215\
Data File : BF077035.D
Acq On : 23 Jan 2015 6:51
Operator : TP/IZ
Sample : SSTDCCC040
Misc : S
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jan 23 07:40:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 07:32:16 2015
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	98	0.00
2		1,4-Dioxane	0.521	0.402	22.8	80	0.00
3		Pyridine	1.348	1.213	10.0	74	-0.01
4		n-Nitrosodimethylamine	0.668	0.679	-1.6	99	-0.01
5	S	2-Fluorophenol	1.216	1.194	1.8	92	-0.01
6		Aniline	2.124	2.107	0.8	91	0.00
7	S	Phenol-d6	1.535	1.527	0.5	93	-0.01
8		2-Chlorophenol	1.402	1.369	2.4	90	0.00
9		Benzaldehyde	0.894	0.901	-0.8	113	0.00
10	C	Phenol	1.629	1.666	-2.3	92	-0.01
11		bis(2-Chloroethyl)ether	1.275	1.281	-0.5	96	0.00
12		1,3-Dichlorobenzene	1.595	1.586	0.6	95	0.00
13	C	1,4-Dichlorobenzene	1.692	1.657	2.1	94	-0.01
14		1,2-Dichlorobenzene	1.559	1.538	1.3	92	0.00
15		Benzyl Alcohol	1.242	1.285	-3.5	94	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.697	1.0	94	0.00
17		2-Methylphenol	1.146	1.171	-2.2	94	-0.01
18		Hexachloroethane	0.588	0.587	0.2	93	-0.01
19	P	n-Nitroso-di-n-propylamine	1.074	1.054	1.9	90	0.00
20		3+4-Methylphenols	1.517	1.476	2.7	91	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.490	0.494	-0.8	94	0.00
23	S	Nitrobenzene-d5	0.361	0.363	-0.6	93	-0.01
24		Nitrobenzene	0.368	0.359	2.4	89	0.00
25		Isophorone	0.657	0.646	1.7	91	-0.01
26	C	2-Nitrophenol	0.172	0.186	-8.1	94	-0.01
27		2,4-Dimethylphenol	0.284	0.290	-2.1	92	-0.01
28		bis(2-Chloroethoxy)methane	0.374	0.389	-4.0	94	0.00
29	C	2,4-Dichlorophenol	0.265	0.269	-1.5	92	-0.01
30		1,2,4-Trichlorobenzene	0.294	0.301	-2.4	92	0.00
31		Naphthalene	1.030	1.036	-0.6	93	0.00
32		Benzoic acid	0.158	0.140	11.4	81	-0.01
33		4-Chloroaniline	0.416	0.403	3.1	90	-0.01
34	C	Hexachlorobutadiene	0.175	0.174	0.6	91	0.00
35		Caprolactam	0.078	0.082	-5.1	93	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.302	0.3	92	-0.01
37		2-Methylnaphthalene	0.703	0.709	-0.9	95	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.493	0.2	92	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.186	7.9	71	0.00
41	S	2,4,6-Tribromophenol	0.162	0.170	-4.9	91	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.372	-3.3	90	-0.01
43		2,4,5-Trichlorophenol	0.367	0.393	-7.1	94	0.00
44	S	2-Fluorobiphenyl	1.314	1.305	0.7	90	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.483	1.509	-1.8	90	0.00
46		2-Chloronaphthalene	1.220	1.224	-0.3	92	0.00
47		2-Nitroaniline	0.370	0.380	-2.7	89	0.00
48		Acenaphthylene	2.058	2.052	0.3	90	0.00
49		Dimethylphthalate	1.418	1.391	1.9	90	0.00
50		2,6-Dinitrotoluene	0.283	0.294	-3.9	89	0.00
51	C	Acenaphthene	1.168	1.145	2.0	89	0.00
52		3-Nitroaniline	0.339	0.348	-2.7	87	0.00
53	P	2,4-Dinitrophenol	0.095	0.063	33.7#	56	0.00
54		Dibenzofuran	1.701	1.696	0.3	91	0.00
55	P	4-Nitrophenol	0.222	0.223	-0.5	88	-0.01
56		2,4-Dinitrotoluene	0.384	0.413	-7.6	91	0.00
57		Fluorene	1.441	1.410	2.2	90	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.278	-3.7	91	0.00
59		Diethylphthalate	1.433	1.453	-1.4	92	0.00
60		4-Chlorophenyl-phenylether	0.590	0.588	0.3	91	0.00
61		4-Nitroaniline	0.328	0.344	-4.9	90	0.00
62		Azobenzene	1.494	1.499	-0.3	91	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.082	21.2	59	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.723	-1.4	89	-0.01
66		4-Bromophenyl-phenylether	0.203	0.199	2.0	85	0.00
67		Hexachlorobenzene	0.214	0.214	0.0	91	0.00
68		Atrazine	0.216	0.223	-3.2	86	0.00
69	C	Pentachlorophenol	0.082	0.098	-19.5	103	0.00
70		Phenanthrene	1.247	1.191	4.5	87	0.00
71		Anthracene	1.216	1.217	-0.1	91	0.00
72		Carbazole	1.180	1.199	-1.6	91	0.00
73		Di-n-butylphthalate	1.365	1.421	-4.1	91	0.00
74	C	Fluoranthene	1.278	1.238	3.1	85	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	96	0.00
76		Benzidine	0.640	0.721	-12.7	108	0.00
77		Pyrene	1.371	1.388	-1.2	89	0.00
78	S	Terphenyl-d14	0.869	0.922	-6.1	95	0.00
79		Butylbenzylphthalate	0.621	0.685	-10.3	93	0.00
80		Benzo(a)anthracene	1.255	1.287	-2.5	91	0.00
81		3,3'-Dichlorobenzidine	0.399	0.428	-7.3	95	0.00
82		Chrysene	1.144	1.150	-0.5	89	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.947	-10.2	98	0.00
84	c	Di-n-octyl phthalate	1.491	1.682	-12.8	99	0.01
85		Indeno(1,2,3-cd)pyrene	1.213	1.278	-5.4	92	0.05
86	I	Perylene-d12	1.000	1.000	0.0	95	0.02
87		Benzo(b)fluoranthene	1.347	1.302	3.3	96	0.02

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88	Benzo(k)fluoranthene	1.177	1.279	-8.7	91	0.06
89 C	Benzo(a)pyrene	1.188	1.215	-2.3	92	0.02
90	Dibenzo(a,h)anthracene	1.090	1.131	-3.8	93	0.06
91	Benzo(g,h,i)perylene	1.084	1.149	-6.0	95	0.05

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