Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF061715\
Data File : BF080029.D
Acq On : 17 Jun 2015 16:17
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF061715

Quant Time: Jun 17 17:09:01 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 17 16:41:53 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	108	0.00
2		1,4-Dioxane	0.537	0.517	3.7	103	0.00
3		Pyridine	1.428	1.418	0.7	107	0.01
4		n-Nitrosodimethylamine	0.679	0.606	10.8	90	0.00
5	S	2-Fluorophenol	1.130	1.121	0.8	104	0.00
6		Aniline	2.068	2.096	-1.4	104	0.00
7	S	Phenol-d6	1.503	1.573	-4.7	105	0.00
8		2-Chlorophenol	1.306	1.327	-1.6	105	0.00
9		Benzaldehyde	0.868	0.843	2.9	100	0.00
10	C	Phenol	1.641	1.740	-6.0	109	-0.01
11		bis(2-Chloroethyl)ether	1.302	1.343	-3.1	107	0.00
12		1,3-Dichlorobenzene	1.569	1.609	-2.5	106	0.00
13	C	1,4-Dichlorobenzene	1.492	1.504	-0.8	106	0.00
14		1,2-Dichlorobenzene	1.452	1.487	-2.4	107	0.00
15		Benzyl Alcohol	1.229	1.267	-3.1	106	0.00
16		2,2'-oxybis(1-Chloropropane	1.932	1.912	1.0	107	0.00
17		2-Methylphenol	1.115	1.189	-6.6	108	0.00
18		Hexachloroethane	0.497	0.504	-1.4	104	0.00
19	P	n-Nitroso-di-n-propylamine	1.044	1.012	3.1	108	-0.01
20		3+4-Methylphenols	1.440	1.473	-2.3	106	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
22		Acetophenone	0.466	0.500	-7.3	106	0.00
23	S	Nitrobenzene-d5	0.344	0.356	-3.5	105	0.00
24		Nitrobenzene	0.355	0.375	-5.6	108	0.00
25		Isophorone	0.691	0.731	-5.8	107	0.00
26	C	2-Nitrophenol	0.148	0.155	-4.7	107	-0.01
27		2,4-Dimethylphenol	0.309	0.309	0.0	107	0.00
28		bis(2-Chloroethoxy)methane	0.406	0.423	-4.2	106	0.00
29	C	2,4-Dichlorophenol	0.265	0.284	-7.2	107	0.00
30		1,2,4-Trichlorobenzene	0.301	0.319	-6.0	109	0.00
31		Naphthalene	1.046	1.097	-4.9	106	0.00
32		Benzoic acid	0.187	0.205	-9.6	111	-0.01
33		4-Chloroaniline	0.448	0.445	0.7	107	0.00
34	C	Hexachlorobutadiene	0.185	0.200	-8.1	117	0.00
35		Caprolactam	0.086	0.096	-11.6	108	-0.01
36	C	4-Chloro-3-methylphenol	0.299	0.324	-8.4	107	0.00
37		2-Methylnaphthalene	0.676	0.696	-3.0	104	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
39		1,2,4,5-Tetrachlorobenzene	0.582	0.578	0.7	105	-0.01
40	P	Hexachlorocyclopentadiene	0.260	0.267	-2.7	106	0.00
41	S	2,4,6-Tribromophenol	0.196	0.198	-1.0	107	-0.01
42	C	2,4,6-Trichlorophenol	0.396	0.404	-2.0	106	0.00
43		2,4,5-Trichlorophenol	0.456	0.472	-3.5	107	0.00
44	S	2-Fluorobiphenyl	1.402	1.406	-0.3	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.627	1.590	2.3	105	0.00
46		2-Chloronaphthalene	1.320	1.367	-3.6	108	0.00
47		2-Nitroaniline	0.362	0.393	-8.6	108	0.00
48		Acenaphthylene	2.204	2.265	-2.8	105	0.00
49		Dimethylphthalate	1.504	1.521	-1.1	110	0.00
50		2,6-Dinitrotoluene	0.301	0.331	-10.0	108	0.00
51	C	Acenaphthene	1.348	1.394	-3.4	109	0.00
52		3-Nitroaniline	0.348	0.389	-11.8	113	0.00
53	P	2,4-Dinitrophenol	0.106	0.124	-17.0	119	0.00
54		Dibenzofuran	1.913	1.913	0.0	106	0.00
55	P	4-Nitrophenol	0.278	0.284	-2.2	112	-0.01
56		2,4-Dinitrotoluene	0.383	0.395	-3.1	108	0.00
57		Fluorene	1.518	1.513	0.3	106	0.00
58		2,3,4,6-Tetrachlorophenol	0.385	0.417	-8.3	110	0.00
59		Diethylphthalate	1.514	1.598	-5.5	107	0.00
60		4-Chlorophenyl-phenylether	0.729	0.723	0.8	107	0.00
61		4-Nitroaniline	0.359	0.376	-4.7	106	0.00
62		Azobenzene	1.541	1.584	-2.8	105	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
64		4,6-Dinitro-2-methylphenol	0.094	0.102	-8.5	110	0.00
65	c	n-Nitrosodiphenylamine	0.651	0.679	-4.3	108	0.00
66		4-Bromophenyl-phenylether	0.213	0.222	-4.2	110	0.00
67		Hexachlorobenzene	0.236	0.235	0.4	104	0.00
68		Atrazine	0.206	0.209	-1.5	104	0.00
69	C	Pentachlorophenol	0.134	0.139	-3.7	115	0.00
70		Phenanthrene	1.211	1.227	-1.3	110	0.00
71		Anthracene	1.166	1.197	-2.7	112	0.00
72		Carbazole	1.113	1.142	-2.6	109	0.00
73		Di-n-butylphthalate	1.286	1.312	-2.0	114	0.01
74	C	Fluoranthene	1.290	1.295	-0.4	111	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	110	0.06
76		Benzidine	0.559	0.575	-2.9	110	0.02
77		Pyrene	1.231	1.188	3.5	105	0.02
78	S	Terphenyl-d14	0.856	0.829	3.2	107	0.02
79		Butylbenzylphthalate	0.495	0.515	-4.0	109	0.03
80		Benzo(a)anthracene	1.218	1.231	-1.1	112	0.06
81		3,3'-Dichlorobenzidine	0.420	0.433	-3.1	112	0.06
82		Chrysene	1.124	1.149	-2.2	112	0.06
83		Bis(2-ethylhexyl)phthalate	0.782	0.796	-1.8	108	0.06
84	c	Di-n-octyl phthalate	1.339	1.358	-1.4	109	0.07
85		Indeno(1,2,3-cd)pyrene	1.417	1.417	0.0	110	0.08
86	I	Perylene-d12	1.000	1.000	0.0	112	0.08
87		Benzo(b)fluoranthene	1.211	1.253	-3.5	112	0.07

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88	Benzo(k)fluoranthene	1.233	1.195	3.1	108	0.08
89 C	Benzo(a)pyrene	1.150	1.168	-1.6	112	0.08
90	Dibenzo(a,h)anthracene	1.177	1.178	-0.1	111	0.08
91	Benzo(g,h,i)perylene	1.188	1.167	1.8	108	0.08

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