Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF061215\
Data File : BF079958.D
Acq On : 13 Jun 2015 4:52
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 13 06:05:05 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 22:57:49 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	96	0.00
2		1,4-Dioxane	0.496	0.542	-9.3	103	0.00
3		Pyridine	1.272	1.560	-22.6	126	0.00
4		n-Nitrosodimethylamine	0.627	0.687	-9.6	115	0.00
5	S	2-Fluorophenol	1.091	1.211	-11.0	104	0.00
6		Aniline	2.151	2.229	-3.6	94	0.00
7	S	Phenol-d6	1.486	1.571	-5.7	95	0.00
8		2-Chlorophenol	1.326	1.326	0.0	92	-0.01
9		Benzaldehyde	0.854	0.838	1.9	100	0.00
10	C	Phenol	1.634	1.570	3.9	86	-0.01
11		bis(2-Chloroethyl)ether	1.315	1.229	6.5	83	0.00
12		1,3-Dichlorobenzene	1.546	1.615	-4.5	96	0.00
13	C	1,4-Dichlorobenzene	1.606	1.770	-10.2	103	0.00
14		1,2-Dichlorobenzene	1.427	1.646	-15.3	108	0.00
15		Benzyl Alcohol	1.191	1.230	-3.3	94	0.00
16		2,2'-oxybis(1-Chloropropane	1.948	2.155	-10.6	101	0.00
17		2-Methylphenol	1.158	1.127	2.7	89	0.00
18		Hexachloroethane	0.497	0.505	-1.6	94	0.00
19	P	n-Nitroso-di-n-propylamine	1.062	1.052	0.9	90	0.00
20		3+4-Methylphenols	1.480	1.552	-4.9	94	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
22		Acetophenone	0.472	0.481	-1.9	90	0.00
23	S	Nitrobenzene-d5	0.303	0.361	-19.1	106	0.00
24		Nitrobenzene	0.322	0.369	-14.6	101	0.00
25		Isophorone	0.705	0.658	6.7	82	-0.01
26	C	2-Nitrophenol	0.100	0.139	-39.0#	121	0.00
27		2,4-Dimethylphenol	0.316	0.308	2.5	90	0.00
28		bis(2-Chloroethoxy)methane	0.419	0.420	-0.2	89	0.00
29	C	2,4-Dichlorophenol	0.265	0.313	-18.1	108	0.00
30		1,2,4-Trichlorobenzene	0.337	0.362	-7.4	98	0.00
31		Naphthalene	1.093	1.081	1.1	89	0.00
32		Benzoic acid	0.107	0.098	8.4	78	0.00
33		4-Chloroaniline	0.424	0.552	-30.2#	114	0.00
34	C	Hexachlorobutadiene	0.187	0.225	-20.3#	113	0.00
35		Caprolactam	0.084	0.093	-10.7	88	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.288	4.6	82	0.00
37		2-Methylnaphthalene	0.739	0.740	-0.1	91	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
39		1,2,4,5-Tetrachlorobenzene	0.670	0.667	0.4	98	0.00
40	P	Hexachlorocyclopentadiene	0.211	0.223	-5.7	93	0.00
41	S	2,4,6-Tribromophenol	0.182	0.186	-2.2	88	-0.01
42	C	2,4,6-Trichlorophenol	0.413	0.414	-0.2	93	0.00
43		2,4,5-Trichlorophenol	0.456	0.418	8.3	85	0.00
44	S	2-Fluorobiphenyl	1.428	1.378	3.5	98	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.619	1.679	-3.7	102	0.00
46		2-Chloronaphthalene	1.378	1.265	8.2	91	0.00
47		2-Nitroaniline	0.256	0.279	-9.0	98	0.00
48		Acenaphthylene	2.326	2.185	6.1	91	-0.01
49		Dimethylphthalate	1.602	1.562	2.5	95	0.00
50		2,6-Dinitrotoluene	0.250	0.267	-6.8	89	-0.01
51	C	Acenaphthene	1.322	1.294	2.1	96	0.00
52		3-Nitroaniline	0.319	0.332	-4.1	92	0.00
53	P	2,4-Dinitrophenol	0.037	0.045#	-21.6	113	0.00
54		Dibenzofuran	1.932	1.895	1.9	95	0.00
55	P	4-Nitrophenol	0.215	0.272	-26.5#	108	0.00
56		2,4-Dinitrotoluene	0.314	0.380	-21.0	99	0.00
57		Fluorene	1.684	1.674	0.6	93	0.00
58		2,3,4,6-Tetrachlorophenol	0.345	0.351	-1.7	90	0.00
59		Diethylphthalate	1.602	1.433	10.5	84	0.00
60		4-Chlorophenyl-phenylether	0.752	0.729	3.1	94	0.00
61		4-Nitroaniline	0.326	0.340	-4.3	87	-0.01
62		Azobenzene	1.542	1.436	6.9	89	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	89	-0.03
64		4,6-Dinitro-2-methylphenol	0.038	0.044	-15.8	98	-0.01
65	c	n-Nitrosodiphenylamine	0.646	0.643	0.5	86	0.00
66		4-Bromophenyl-phenylether	0.224	0.236	-5.4	89	-0.01
67		Hexachlorobenzene	0.244	0.256	-4.9	88	-0.01
68		Atrazine	0.199	0.234	-17.6	94	-0.01
69	C	Pentachlorophenol	0.095	0.093	2.1	83	-0.01
70		Phenanthrene	1.192	1.241	-4.1	88	-0.02
71		Anthracene	1.153	1.228	-6.5	91	-0.03
72		Carbazole	1.119	1.191	-6.4	88	-0.05
73		Di-n-butylphthalate	1.219	1.221	-0.2	83	-0.07
74	C	Fluoranthene	1.318	1.377	-4.5	87	-0.13
75	I	Chrysene-d12	1.000	1.000	0.0	89	-0.30
76		Benzidine	0.554	0.567	-2.3	93	-0.14
77		Pyrene	1.256	1.199	4.5	88	-0.15
78	S	Terphenyl-d14	0.867	0.844	2.7	89	-0.16
79		Butylbenzylphthalate	0.424	0.426	-0.5	84	-0.23
80		Benzo(a)anthracene	1.270	1.213	4.5	89	-0.29
81		3,3'-Dichlorobenzidine	0.406	0.394	3.0	84	-0.29
82		Chrysene	1.229	1.130	8.1	88	-0.30
83		Bis(2-ethylhexyl)phthalate	0.679	0.695	-2.4	86	-0.31
84	c	Di-n-octyl phthalate	1.107	1.102	0.5	81	-0.39
85		Indeno(1,2,3-cd)pyrene	1.330	1.389	-4.4	95	-0.42
86	I	Perylene-d12	1.000	1.000	0.0	89	-0.39
87		Benzo(b)fluoranthene	1.225	1.178	3.8	88	-0.38

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.250	1.334	-6.7	91	-0.39
89 C	Benzo(a)pyrene	1.165	1.166	-0.1	89	-0.39
90	Dibenzo(a,h)anthracene	1.146	1.178	-2.8	95	-0.42
91	Benzo(g,h,i)perylene	1.185	1.208	-1.9	94	-0.42

(#) = Out of Range SPCC's out = 1 CCC's out = 2