Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF060715\
Data File : BF079814.D
Acq On : 7 Jun 2015 14:55
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 09 01:18:28 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 01:08:34 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	75	-0.01
2		1,4-Dioxane	0.528	0.931	-76.3#	135	0.00
3		Pyridine	1.477	2.676	-81.2#	125	-0.01
4		n-Nitrosodimethylamine	0.641	1.187	-85.2#	152#	-0.01
5	S	2-Fluorophenol	1.193	1.133	5.0	72	0.00
6		Aniline	2.237	2.084	6.8	70	0.00
7	S	Phenol-d6	1.541	1.448	6.0	71	-0.01
8		2-Chlorophenol	1.292	1.284	0.6	76	-0.01
9		Benzaldehyde	0.998	0.887	11.1	67	-0.01
10	C	Phenol	1.678	1.501	10.5	67	0.00
11		bis(2-Chloroethyl)ether	1.301	1.272	2.2	74	0.00
12		1,3-Dichlorobenzene	1.617	1.494	7.6	70	-0.01
13	C	1,4-Dichlorobenzene	1.716	1.621	5.5	72	0.00
14		1,2-Dichlorobenzene	1.603	1.414	11.8	68	0.00
15		Benzyl Alcohol	1.274	1.160	8.9	70	0.00
16		2,2'-oxybis(1-Chloropropane	1.968	1.802	8.4	71	0.00
17		2-Methylphenol	1.205	1.015	15.8	65	0.00
18		Hexachloroethane	0.499	0.524	-5.0	82	0.00
19	P	n-Nitroso-di-n-propylamine	1.069	1.046	2.2	77	0.00
20		3+4-Methylphenols	1.535	1.453	5.3	71	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	-0.01
22		Acetophenone	0.477	0.513	-7.5	76	0.00
23	S	Nitrobenzene-d5	0.326	0.332	-1.8	70	0.00
24		Nitrobenzene	0.325	0.333	-2.5	73	-0.01
25		Isophorone	0.665	0.712	-7.1	78	0.00
26	C	2-Nitrophenol	0.137	0.118	13.9	62	0.00
27		2,4-Dimethylphenol	0.308	0.323	-4.9	75	0.00
28		bis(2-Chloroethoxy)methane	0.411	0.361	12.2	64	-0.01
29	C	2,4-Dichlorophenol	0.276	0.266	3.6	69	-0.01
30		1,2,4-Trichlorobenzene	0.335	0.342	-2.1	73	0.00
31		Naphthalene	1.049	1.052	-0.3	74	-0.01
32		Benzoic acid	0.113	0.095	15.9	56	0.00
33		4-Chloroaniline	0.533	0.571	-7.1	75	0.00
34	C	Hexachlorobutadiene	0.204	0.196	3.9	70	-0.01
35		Caprolactam	0.084	0.075	10.7	62	-0.01
36	C	4-Chloro-3-methylphenol	0.309	0.287	7.1	66	0.00
37		2-Methylnaphthalene	0.667	0.719	-7.8	78	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	68	0.00
39		1,2,4,5-Tetrachlorobenzene	0.623	0.709	-13.8	78	0.00
40	P	Hexachlorocyclopentadiene	0.268	0.331	-23.5	84	0.00
41	S	2,4,6-Tribromophenol	0.191	0.203	-6.3	68	-0.01
42	C	2,4,6-Trichlorophenol	0.414	0.412	0.5	68	0.00
43		2,4,5-Trichlorophenol	0.436	0.493	-13.1	77	-0.01
44	S	2-Fluorobiphenyl	1.476	1.487	-0.7	70	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.740	1.654	4.9	66	-0.01
46		2-Chloronaphthalene	1.421	1.401	1.4	70	-0.01
47		2-Nitroaniline	0.298	0.318	-6.7	76	-0.01
48		Acenaphthylene	2.349	2.309	1.7	68	-0.01
49		Dimethylphthalate	1.541	1.740	-12.9	80	-0.01
50		2,6-Dinitrotoluene	0.247	0.301	-21.9	85	0.00
51	C	Acenaphthene	1.330	1.333	-0.2	70	-0.01
52		3-Nitroaniline	0.325	0.348	-7.1	75	-0.01
53	P	2,4-Dinitrophenol	0.062	0.060	3.2	67	-0.01
54		Dibenzofuran	1.884	1.889	-0.3	68	0.00
55	P	4-Nitrophenol	0.243	0.264	-8.6	74	-0.01
56		2,4-Dinitrotoluene	0.293	0.339	-15.7	78	0.00
57		Fluorene	1.630	1.685	-3.4	73	-0.01
58		2,3,4,6-Tetrachlorophenol	0.353	0.341	3.4	62	-0.01
59		Diethylphthalate	1.507	1.728	-14.7	82	0.00
60		4-Chlorophenyl-phenylether	0.739	0.707	4.3	66	-0.01
61		4-Nitroaniline	0.323	0.358	-10.8	79	-0.01
62		Azobenzene	1.522	1.498	1.6	67	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	68	-0.01
64		4,6-Dinitro-2-methylphenol	0.050	0.054	-8.0	82	-0.01
65	c	n-Nitrosodiphenylamine	0.668	0.687	-2.8	72	-0.01
66		4-Bromophenyl-phenylether	0.217	0.240	-10.6	75	0.00
67		Hexachlorobenzene	0.244	0.249	-2.0	74	-0.01
68		Atrazine	0.201	0.206	-2.5	72	-0.01
69	C	Pentachlorophenol	0.104	0.096	7.7	60	-0.01
70		Phenanthrene	1.171	1.189	-1.5	71	0.00
71		Anthracene	1.155	1.192	-3.2	71	-0.01
72		Carbazole	1.097	1.128	-2.8	73	0.00
73		Di-n-butylphthalate	1.244	1.294	-4.0	74	0.00
74	C	Fluoranthene	1.282	1.272	0.8	68	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	65	0.06
76		Benzidine	0.631	0.412	34.7#	45#	0.01
77		Pyrene	1.226	1.258	-2.6	68	0.02
78	S	Terphenyl-d14	0.873	0.915	-4.8	70	0.02
79		Butylbenzylphthalate	0.489	0.478	2.2	64	0.03
80		Benzo(a)anthracene	1.224	1.208	1.3	66	0.06
81		3,3'-Dichlorobenzidine	0.416	0.390	6.2	62	0.05
82		Chrysene	1.132	1.111	1.9	66	0.06
83		Bis(2-ethylhexyl)phthalate	0.772	0.736	4.7	64	0.06
84	c	Di-n-octyl phthalate	1.360	1.201	11.7	60	0.08
85		Indeno(1,2,3-cd)pyrene	1.310	1.189	9.2	61	0.06
86	I	Perylene-d12	1.000	1.000	0.0	61	0.07
87		Benzo(b)fluoranthene	1.242	1.184	4.7	62	0.07

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88	Benzo(k)fluoranthene	1.192	1.293	-8.5	66	0.07
89 C	Benzo(a)pyrene	1.147	1.147	0.0	63	0.06
90	Dibenzo(a,h)anthracene	1.107	1.074	3.0	62	0.06
91	Benzo(g,h,i)perylene	1.144	1.090	4.7	61	0.07

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