Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF080115\
Data File : BF080780.D
Acq On : 1 Aug 2015 14:52
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 03 02:52:31 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF073015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 31 01:45:22 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	82	0.00
2		1,4-Dioxane	0.575	0.529	8.0	79	0.00
3		Pyridine	1.595	1.369	14.2	68	0.00
4		n-Nitrosodimethylamine	0.912	0.928	-1.8	82	-0.01
5	S	2-Fluorophenol	1.166	1.099	5.7	82	0.00
6		Aniline	2.285	2.082	8.9	75	-0.01
7	S	Phenol-d6	1.587	1.548	2.5	77	0.00
8		2-Chlorophenol	1.307	1.221	6.6	79	0.00
9		Benzaldehyde	1.000	0.859	14.1	72	-0.01
10	C	Phenol	1.837	1.792	2.4	73	0.00
11		bis(2-Chloroethyl)ether	1.312	1.169	10.9	80	0.00
12		1,3-Dichlorobenzene	1.455	1.470	-1.0	84	0.00
13	C	1,4-Dichlorobenzene	1.484	1.444	2.7	78	0.00
14		1,2-Dichlorobenzene	1.357	1.368	-0.8	89	0.00
15		Benzyl Alcohol	1.318	1.021	22.5	60	0.00
16		2,2'-oxybis(1-Chloropropane	2.803	2.664	5.0	81	0.00
17		2-Methylphenol	1.073	0.946	11.8	75	0.00
18		Hexachloroethane	0.554	0.548	1.1	83	0.00
19	P	n-Nitroso-di-n-propylamine	1.069	0.992	7.2	78	0.00
20		3+4-Methylphenols	1.323	1.297	2.0	79	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	71	-0.01
22		Acetophenone	0.483	0.484	-0.2	80	0.00
23	S	Nitrobenzene-d5	0.413	0.414	-0.2	72	-0.01
24		Nitrobenzene	0.414	0.435	-5.1	87	0.00
25		Isophorone	0.734	0.690	6.0	71	0.00
26	C	2-Nitrophenol	0.168	0.181	-7.7	73	0.00
27		2,4-Dimethylphenol	0.316	0.317	-0.3	79	0.00
28		bis(2-Chloroethoxy)methane	0.435	0.447	-2.8	74	0.00
29	C	2,4-Dichlorophenol	0.280	0.283	-1.1	73	0.00
30		1,2,4-Trichlorobenzene	0.306	0.299	2.3	73	0.00
31		Naphthalene	0.998	0.920	7.8	73	0.00
32		Benzoic acid	0.209	0.182	12.9	61	-0.01
33		4-Chloroaniline	0.421	0.374	11.2	67	-0.01
34	C	Hexachlorobutadiene	0.197	0.219	-11.2	84	0.00
35		Caprolactam	0.084	0.078	7.1	67	0.00
36	C	4-Chloro-3-methylphenol	0.302	0.277	8.3	72	0.00
37		2-Methylnaphthalene	0.627	0.600	4.3	69	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	66	0.00
39		1,2,4,5-Tetrachlorobenzene	0.628	0.644	-2.5	78	0.00
40	P	Hexachlorocyclopentadiene	0.386	0.386	0.0	69	-0.01
41	S	2,4,6-Tribromophenol	0.208	0.204	1.9	63	-0.01
42	C	2,4,6-Trichlorophenol	0.437	0.438	-0.2	68	0.00
43		2,4,5-Trichlorophenol	0.433	0.485	-12.0	76	0.00
44	S	2-Fluorobiphenyl	1.341	1.416	-5.6	75	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.534	1.454	5.2	69	-0.01
46		2-Chloronaphthalene	1.253	1.206	3.8	68	-0.01
47		2-Nitroaniline	0.487	0.429	11.9	59	-0.01
48		Acenaphthylene	1.992	1.905	4.4	64	-0.01
49		Dimethylphthalate	1.409	1.412	-0.2	71	0.00
50		2,6-Dinitrotoluene	0.299	0.314	-5.0	72	0.00
51	C	Acenaphthene	1.215	1.199	1.3	66	-0.01
52		3-Nitroaniline	0.345	0.286	17.1	56	-0.01
53	P	2,4-Dinitrophenol	0.159	0.148	6.9	67	0.00
54		Dibenzofuran	1.625	1.496	7.9	62	-0.01
55	P	4-Nitrophenol	0.254	0.210	17.3	53	-0.01
56		2,4-Dinitrotoluene	0.392	0.371	5.4	64	-0.01
57		Fluorene	1.261	1.189	5.7	67	-0.01
58		2,3,4,6-Tetrachlorophenol	0.327	0.324	0.9	64	0.00
59		Diethylphthalate	1.359	1.355	0.3	71	0.00
60		4-Chlorophenyl-phenylether	0.596	0.664	-11.4	84	0.00
61		4-Nitroaniline	0.312	0.293	6.1	65	-0.01
62		Azobenzene	1.472	1.628	-10.6	80	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
64		4,6-Dinitro-2-methylphenol	0.120	0.115	4.2	71	0.00
65	c	n-Nitrosodiphenylamine	0.656	0.607	7.5	70	0.00
66		4-Bromophenyl-phenylether	0.206	0.199	3.4	67	0.00
67		Hexachlorobenzene	0.239	0.233	2.5	76	0.00
68		Atrazine	0.158	0.161	-1.9	67	0.00
69	C	Pentachlorophenol	0.144	0.132	8.3	63	0.00
70		Phenanthrene	1.008	0.954	5.4	72	0.00
71		Anthracene	0.990	0.944	4.6	67	0.00
72		Carbazole	0.859	0.861	-0.2	78	0.00
73		Di-n-butylphthalate	1.066	1.008	5.4	66	0.00
74	C	Fluoranthene	0.974	0.897	7.9	65	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
76		Benzidine	0.728	0.522	28.3#	58	-0.01
77		Pyrene	1.551	1.193	23.1	63	-0.01
78	S	Terphenyl-d14	1.057	0.822	22.2	64	-0.01
79		Butylbenzylphthalate	0.592	0.548	7.4	77	0.00
80		Benzo(a)anthracene	1.203	1.091	9.3	80	0.00
81		3,3'-Dichlorobenzidine	0.411	0.388	5.6	83	0.00
82		Chrysene	1.168	0.907	22.3	65	-0.01
83		Bis(2-ethylhexyl)phthalate	0.772	0.729	5.6	77	-0.01
84	c	Di-n-octyl phthalate	1.295	1.165	10.0	75	-0.01
85		Indeno(1,2,3-cd)pyrene	0.958	1.315	-37.3#	116	0.00
86	I	Perylene-d12	1.000	1.000	0.0	92	0.00
87		Benzo(b)fluoranthene	1.195	1.142	4.4	94	-0.01

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88	Benzo(k)fluoranthene	0.983	0.748	23.9	70	0.00
89 C	Benzo(a)pyrene	0.981	0.932	5.0	90	0.00
90	Dibenzo(a,h)anthracene	0.878	1.080	-23.0	114	-0.01
91	Benzo(g,h,i)perylene	0.864	1.075	-24.4	116	0.00

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