Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA_F\Data\BF062317\
Data File : BF096157.D
Acq On : 23 Jun 2017 13:51
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 28 11:58:41 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 20 16:05:47 2017
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	70	-0.04
2		1,4-Dioxane	0.597	0.667	-11.7	75	-0.02
3		Pyridine	1.403	1.253	10.7	60	-0.02
4		n-Nitrosodimethylamine	0.534	0.498	6.7	64	-0.02
5	S	2-Fluorophenol	1.255	1.268	-1.0	65	-0.04
6		Aniline	1.966	2.052	-4.4	69	-0.03
7	S	Phenol-d6	1.503	1.540	-2.5	66	-0.04
8		2-Chlorophenol	1.335	1.409	-5.5	68	-0.03
9		Benzaldehyde	1.014	0.916	9.7	60	-0.02
10	C	Phenol	1.559	1.648	-5.7	67	-0.04
11		bis(2-Chloroethyl)ether	1.235	1.320	-6.9	69	-0.03
12		1,3-Dichlorobenzene	1.511	1.583	-4.8	73	-0.04
13	C	1,4-Dichlorobenzene	1.532	1.605	-4.8	72	-0.04
14		1,2-Dichlorobenzene	1.412	1.494	-5.8	72	-0.03
15		Benzyl Alcohol	1.031	1.107	-7.4	71	-0.03
16		2,2'-oxybis(1-Chloropropane	1.735	1.849	-6.6	72	-0.04
17		2-Methylphenol	1.120	1.185	-5.8	72	-0.03
18		Hexachloroethane	0.506	0.538	-6.3	72	-0.04
19	P	n-Nitroso-di-n-propylamine	0.952	1.023	-7.5	72	-0.02
20		3+4-Methylphenols	1.422	1.616	-13.6	77	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	71	-0.03
22		Acetophenone	0.468	0.482	-3.0	72	-0.02
23	S	Nitrobenzene-d5	0.322	0.329	-2.2	71	-0.03
24		Nitrobenzene	0.344	0.345	-0.3	71	-0.03
25		Isophorone	0.601	0.611	-1.7	72	-0.02
26	C	2-Nitrophenol	0.180	0.188	-4.4	71	-0.02
27		2,4-Dimethylphenol	0.280	0.291	-3.9	73	-0.03
28		bis(2-Chloroethoxy)methane	0.396	0.413	-4.3	72	-0.03
29	C	2,4-Dichlorophenol	0.269	0.270	-0.4	70	-0.03
30		1,2,4-Trichlorobenzene	0.280	0.286	-2.1	72	-0.03
31		Naphthalene	1.004	1.018	-1.4	72	-0.04
32		Benzoic acid	0.161	0.146	9.3	61	0.00
33		4-Chloroaniline	0.392	0.414	-5.6	74	-0.02
34	C	Hexachlorobutadiene	0.145	0.148	-2.1	72	-0.04
35		Caprolactam	0.080	0.079	1.3	66	-0.02
36	C	4-Chloro-3-methylphenol	0.282	0.292	-3.5	72	-0.02
37		2-Methylnaphthalene	0.678	0.710	-4.7	75	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	79	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.599	0.590	1.5	74	-0.03
40	P	Hexachlorocyclopentadiene	0.138	0.152	-10.1	79	-0.04
41	S	2,4,6-Tribromophenol	0.177	0.171	3.4	75	-0.04
42	C	2,4,6-Trichlorophenol	0.385	0.366	4.9	70	-0.02
43		2,4,5-Trichlorophenol	0.409	0.367	10.3	64	-0.02
44	S	2-Fluorobiphenyl	1.437	1.379	4.0	74	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.648	1.617	1.9	74	-0.03
46		2-Chloronaphthalene	1.288	1.259	2.3	74	-0.03
47		2-Nitroaniline	0.377	0.361	4.2	67	-0.02
48		Acenaphthylene	2.202	2.020	8.3	68	-0.04
49		Dimethylphthalate	1.519	1.419	6.6	70	-0.03
50		2,6-Dinitrotoluene	0.331	0.300	9.4	65	-0.03
51	C	Acenaphthene	1.272	1.226	3.6	76	-0.03
52		3-Nitroaniline	0.386	0.372	3.6	75	-0.02
53	P	2,4-Dinitrophenol	0.160	0.132	17.5	63	-0.02
54		Dibenzofuran	1.790	1.734	3.1	77	-0.04
55	P	4-Nitrophenol	0.201	0.166	17.4	60	-0.02
56		2,4-Dinitrotoluene	0.447	0.445	0.4	76	-0.02
57		Fluorene	1.558	1.495	4.0	76	-0.04
58		2,3,4,6-Tetrachlorophenol	0.280	0.269	3.9	73	-0.03
59		Diethylphthalate	1.461	1.418	2.9	76	-0.03
60		4-Chlorophenyl-phenylether	0.677	0.662	2.2	77	-0.03
61		4-Nitroaniline	0.360	0.355	1.4	75	-0.03
62		Azobenzene	1.324	1.278	3.5	76	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	75	-0.03
64		4,6-Dinitro-2-methylphenol	0.131	0.118	9.9	64	-0.02
65	c	n-Nitrosodiphenylamine	0.719	0.700	2.6	74	-0.03
66		4-Bromophenyl-phenylether	0.208	0.207	0.5	74	-0.03
67		Hexachlorobenzene	0.205	0.207	-1.0	75	-0.04
68		Atrazine	0.200	0.190	5.0	73	-0.04
69	C	Pentachlorophenol	0.067	0.055	17.9	61	-0.03
70		Phenanthrene	1.154	1.150	0.3	77	-0.03
71		Anthracene	1.205	1.195	0.8	76	-0.03
72		Carbazole	1.174	1.204	-2.6	76	-0.03
73		Di-n-butylphthalate	1.249	1.294	-3.6	76	-0.03
74	C	Fluoranthene	1.142	1.169	-2.4	77	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	74	-0.02
76		Benzidine	0.768	0.855	-11.3	80	-0.02
77		Pyrene	1.492	1.559	-4.5	77	-0.02
78	S	Terphenyl-d14	0.806	0.842	-4.5	78	-0.02
79		Butylbenzylphthalate	0.652	0.679	-4.1	75	-0.02
80		Benzo(a)anthracene	1.163	1.185	-1.9	75	-0.02
81		3,3'-Dichlorobenzidine	0.413	0.433	-4.8	76	-0.02
82		Chrysene	1.162	1.212	-4.3	76	-0.02
83		Bis(2-ethylhexyl)phthalate	0.919	0.983	-7.0	77	-0.02
84	c	Di-n-octyl phthalate	1.515	1.588	-4.8	76	-0.02
85		Indeno(1,2,3-cd)pyrene	0.994	0.970	2.4	77	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	74	-0.02
87		Benzo(b)fluoranthene	1.252	1.208	3.5	69	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.197	1.285	-7.4	82	-0.02
89 C	Benzo(a)pyrene	1.151	1.157	-0.5	75	-0.02
90	Dibenzo(a,h)anthracene	0.976	0.930	4.7	75	-0.03
91	Benzo(g,h,i)perylene	0.995	0.944	5.1	75	-0.03

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