Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100417\
Data File : BF099066.D
Acq On : 4 Oct 2017 14:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 04 18:44:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 04 18:42:14 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	76	-0.05
2		1,4-Dioxane	0.600	0.576	4.0	72	-0.13
3		Pyridine	1.624	1.595	1.8	77	-0.12
4		n-Nitrosodimethylamine	0.688	0.622	9.6	69	-0.13
5	S	2-Fluorophenol	1.256	1.264	-0.6	77	-0.06
6		Aniline	2.092	2.082	0.5	77	-0.05
7	S	Phenol-d6	1.550	1.569	-1.2	77	-0.05
8		2-Chlorophenol	1.423	1.439	-1.1	78	-0.06
9		Benzaldehyde	0.899	0.815	9.3	71	-0.06
10	C	Phenol	1.733	1.864	-7.6	83	-0.05
11		bis(2-Chloroethyl)ether	1.348	1.347	0.1	77	-0.05
12		1,3-Dichlorobenzene	1.490	1.511	-1.4	79	-0.05
13	C	1,4-Dichlorobenzene	1.516	1.511	0.3	78	-0.05
14		1,2-Dichlorobenzene	1.401	1.427	-1.9	78	-0.05
15		Benzyl Alcohol	1.137	1.121	1.4	75	-0.06
16		2,2'-oxybis(1-Chloropropane	2.099	2.036	3.0	76	-0.05
17		2-Methylphenol	1.164	1.158	0.5	77	-0.05
18		Hexachloroethane	0.545	0.542	0.6	78	-0.06
19	P	n-Nitroso-di-n-propylamine	0.990	0.944	4.6	76	-0.05
20		3+4-Methylphenols	1.473	1.436	2.5	75	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	78	-0.05
22		Acetophenone	0.485	0.452	6.8	75	-0.05
23	S	Nitrobenzene-d5	0.312	0.313	-0.3	78	-0.05
24		Nitrobenzene	0.354	0.342	3.4	76	-0.05
25		Isophorone	0.645	0.618	4.2	77	-0.05
26	C	2-Nitrophenol	0.170	0.186	-9.4	82	-0.06
27		2,4-Dimethylphenol	0.321	0.311	3.1	77	-0.05
28		bis(2-Chloroethoxy)methane	0.402	0.395	1.7	79	-0.05
29	C	2,4-Dichlorophenol	0.276	0.277	-0.4	79	-0.05
30		1,2,4-Trichlorobenzene	0.276	0.276	0.0	79	-0.05
31		Naphthalene	0.939	0.927	1.3	79	-0.05
32		Benzoic acid	0.264	0.211	20.1	61	-0.05
33		4-Chloroaniline	0.392	0.384	2.0	77	-0.05
34	C	Hexachlorobutadiene	0.142	0.141	0.7	80	-0.05
35		Caprolactam	0.088	0.087	1.1	79	-0.05
36	C	4-Chloro-3-methylphenol	0.310	0.300	3.2	76	-0.05
37		2-Methylnaphthalene	0.611	0.606	0.8	79	-0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	-0.06
39		1,2,4,5-Tetrachlorobenzene	0.596	0.601	-0.8	80	-0.05
40	P	Hexachlorocyclopentadiene	0.273	0.297	-8.8	83	-0.06
41	S	2,4,6-Tribromophenol	0.164	0.169	-3.0	79	-0.05
42	C	2,4,6-Trichlorophenol	0.423	0.417	1.4	78	-0.05
43		2,4,5-Trichlorophenol	0.422	0.416	1.4	77	-0.05
44	S	2-Fluorobiphenyl	1.198	1.213	-1.3	79	-0.05

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.701	1.0	79	-0.05
46		2-Chloronaphthalene	1.291	1.291	0.0	80	-0.06
47		2-Nitroaniline	0.395	0.389	1.5	75	-0.05
48		Acenaphthylene	1.976	1.951	1.3	79	-0.06
49		Dimethylphthalate	1.509	1.500	0.6	80	-0.05
50		2,6-Dinitrotoluene	0.329	0.339	-3.0	79	-0.05
51	C	Acenaphthene	1.251	1.172	6.3	75	-0.06
52		3-Nitroaniline	0.383	0.406	-6.0	80	-0.05
53	P	2,4-Dinitrophenol	0.148	0.154	-4.1	78	-0.05
54		Dibenzofuran	1.666	1.657	0.5	79	-0.06
55	P	4-Nitrophenol	0.324	0.308	4.9	73	-0.04
56		2,4-Dinitrotoluene	0.426	0.452	-6.1	81	-0.05
57		Fluorene	1.339	1.343	-0.3	81	-0.05
58		2,3,4,6-Tetrachlorophenol	0.291	0.282	3.1	75	-0.05
59		Diethylphthalate	1.475	1.450	1.7	79	-0.06
60		4-Chlorophenyl-phenylether	0.602	0.598	0.7	80	-0.05
61		4-Nitroaniline	0.370	0.396	-7.0	82	-0.05
62		Azobenzene	1.356	1.288	5.0	76	-0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	76	-0.05
64		4,6-Dinitro-2-methylphenol	0.122	0.139	-13.9	86	-0.05
65	c	n-Nitrosodiphenylamine	0.693	0.702	-1.3	80	-0.05
66		4-Bromophenyl-phenylether	0.196	0.200	-2.0	79	-0.06
67		Hexachlorobenzene	0.209	0.212	-1.4	79	-0.05
68		Atrazine	0.208	0.211	-1.4	78	-0.05
69	C	Pentachlorophenol	0.130	0.117	10.0	66	-0.05
70		Phenanthrene	1.071	1.070	0.1	79	-0.06
71		Anthracene	1.095	1.102	-0.6	79	-0.06
72		Carbazole	1.073	1.083	-0.9	79	-0.05
73		Di-n-butylphthalate	1.291	1.303	-0.9	78	-0.06
74	C	Fluoranthene	1.084	1.098	-1.3	80	-0.06
75	I	Chrysene-d12	1.000	1.000	0.0	77	-0.05
76		Benzidine	0.740	0.764	-3.2	81	-0.06
77		Pyrene	1.525	1.534	-0.6	79	-0.06
78	S	Terphenyl-d14	0.879	0.907	-3.2	81	-0.06
79		Butylbenzylphthalate	0.717	0.733	-2.2	78	-0.06
80		Benzo(a)anthracene	1.211	1.216	-0.4	80	-0.05
81		3,3'-Dichlorobenzidine	0.444	0.451	-1.6	79	-0.06
82		Chrysene	1.153	1.160	-0.6	80	-0.06
83		Bis(2-ethylhexyl)phthalate	0.987	1.014	-2.7	79	-0.06
84	c	Di-n-octyl phthalate	1.589	1.595	-0.4	80	-0.06
85		Indeno(1,2,3-cd)pyrene	0.922	0.894	3.0	82	-0.11
86	I	Perylene-d12	1.000	1.000	0.0	76	-0.08
87		Benzo(b)fluoranthene	1.230	1.314	-6.8	80	-0.07

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88	Benzo(k)fluoranthene	1.215	1.176	3.2	75	-0.06
89 C	Benzo(a)pyrene	1.129	1.146	-1.5	78	-0.08
90	Dibenzo(a,h)anthracene	0.903	0.942	-4.3	82	-0.11
91	Benzo(g,h,i)perylene	0.893	0.948	-6.2	83	-0.11

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