Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF010816\
Data File : BF084327.D
Acq On : 8 Jan 2016 14:50
Operator : SJ/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 08 23:08:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 12:22:40 2016
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	99	-0.05
2		1,4-Dioxane	0.586	0.573	2.2	92	-0.09
3		Pyridine	1.633	1.657	-1.5	92	-0.10
4		n-Nitrosodimethylamine	0.838	0.852	-1.7	95	-0.09
5	S	2-Fluorophenol	1.236	1.144	7.4	97	-0.03
6		Aniline	2.272	2.102	7.5	89	-0.03
7	S	Phenol-d6	1.553	1.494	3.8	94	-0.02
8		2-Chlorophenol	1.403	1.346	4.1	96	-0.03
9		Benzaldehyde	1.011	0.968	4.3	95	-0.04
10	C	Phenol	1.766	1.711	3.1	89	-0.02
11		bis(2-Chloroethyl)ether	1.368	1.346	1.6	101	-0.03
12		1,3-Dichlorobenzene	1.447	1.378	4.8	97	-0.05
13	C	1,4-Dichlorobenzene	1.451	1.380	4.9	90	-0.05
14		1,2-Dichlorobenzene	1.390	1.399	-0.6	106	-0.05
15		Benzyl Alcohol	1.306	1.296	0.8	93	-0.03
16		2,2'-oxybis(1-Chloropropane	2.491	2.324	6.7	95	-0.05
17		2-Methylphenol	1.176	1.109	5.7	87	-0.03
18		Hexachloroethane	0.580	0.587	-1.2	96	-0.05
19	P	n-Nitroso-di-n-propylamine	1.051	0.906	13.8	87	-0.03
20		3+4-Methylphenols	1.382	1.407	-1.8	107	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	95	-0.05
22		Acetophenone	0.481	0.449	6.7	83	-0.05
23	S	Nitrobenzene-d5	0.359	0.375	-4.5	102	-0.03
24		Nitrobenzene	0.364	0.338	7.1	80	-0.05
25		Isophorone	0.687	0.698	-1.6	96	-0.03
26	C	2-Nitrophenol	0.166	0.204	-22.9#	120	-0.03
27		2,4-Dimethylphenol	0.336	0.348	-3.6	93	-0.03
28		bis(2-Chloroethoxy)methane	0.420	0.425	-1.2	104	-0.03
29	C	2,4-Dichlorophenol	0.280	0.270	3.6	94	-0.03
30		1,2,4-Trichlorobenzene	0.289	0.289	0.0	92	-0.05
31		Naphthalene	0.907	0.852	6.1	90	-0.05
32		Benzoic acid	0.198	0.140	29.3#	64	-0.02
33		4-Chloroaniline	0.416	0.447	-7.5	106	-0.03
34	C	Hexachlorobutadiene	0.166	0.174	-4.8	100	-0.05
35		Caprolactam	0.094	0.089	5.3	79	-0.03
36	C	4-Chloro-3-methylphenol	0.313	0.345	-10.2	110	-0.02
37		2-Methylnaphthalene	0.651	0.682	-4.8	104	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	-0.05
39		1,2,4,5-Tetrachlorobenzene	0.575	0.527	8.3	88	-0.05
40	P	Hexachlorocyclopentadiene	0.347	0.317	8.6	86	-0.05
41	S	2,4,6-Tribromophenol	0.202	0.176	12.9	79	-0.05
42	C	2,4,6-Trichlorophenol	0.430	0.390	9.3	90	-0.05
43		2,4,5-Trichlorophenol	0.412	0.415	-0.7	94	-0.03
44	S	2-Fluorobiphenyl	1.317	1.156	12.2	96	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.590	1.609	-1.2	105	-0.03
46		2-Chloronaphthalene	1.249	1.237	1.0	106	-0.03
47		2-Nitroaniline	0.412	0.402	2.4	98	-0.03
48		Acenaphthylene	1.845	1.753	5.0	88	-0.05
49		Dimethylphthalate	1.430	1.427	0.2	93	-0.03
50		2,6-Dinitrotoluene	0.314	0.358	-14.0	100	-0.03
51	C	Acenaphthene	1.237	1.196	3.3	87	-0.05
52		3-Nitroaniline	0.389	0.435	-11.8	100	-0.03
53	P	2,4-Dinitrophenol	0.093	0.116	-24.7	84	-0.03
54		Dibenzofuran	1.812	1.544	14.8	86	-0.05
55	P	4-Nitrophenol	0.282	0.256	9.2	73	-0.02
56		2,4-Dinitrotoluene	0.397	0.466	-17.4	97	-0.03
57		Fluorene	1.400	1.270	9.3	93	-0.04
58		2,3,4,6-Tetrachlorophenol	0.352	0.330	6.2	86	-0.05
59		Diethylphthalate	1.410	1.326	6.0	89	-0.05
60		4-Chlorophenyl-phenylether	0.564	0.518	8.2	89	-0.05
61		4-Nitroaniline	0.346	0.397	-14.7	113	-0.02
62		Azobenzene	1.546	1.381	10.7	85	-0.05
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.03
64		4,6-Dinitro-2-methylphenol	0.109	0.102	6.4	91	-0.03
65	c	n-Nitrosodiphenylamine	0.639	0.577	9.7	90	-0.03
66		4-Bromophenyl-phenylether	0.215	0.192	10.7	85	-0.05
67		Hexachlorobenzene	0.242	0.241	0.4	106	-0.03
68		Atrazine	0.209	0.196	6.2	84	-0.05
69	C	Pentachlorophenol	0.126	0.120	4.8	85	-0.03
70		Phenanthrene	1.046	0.959	8.3	85	-0.05
71		Anthracene	1.026	1.017	0.9	104	-0.04
72		Carbazole	1.003	0.831	17.1	80	-0.05
73		Di-n-butylphthalate	1.111	0.998	10.2	89	-0.05
74	C	Fluoranthene	1.093	1.102	-0.8	106	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	93	-0.03
76		Benzidine	0.717	0.623	13.1	80	-0.05
77		Pyrene	1.569	1.547	1.4	104	-0.03
78	S	Terphenyl-d14	0.896	0.871	2.8	105	-0.03
79		Butylbenzylphthalate	0.679	0.616	9.3	83	-0.05
80		Benzo(a)anthracene	1.174	1.095	6.7	92	-0.03
81		3,3'-Dichlorobenzidine	0.410	0.455	-11.0	102	-0.03
82		Chrysene	1.128	1.026	9.0	86	-0.05
83		Bis(2-ethylhexyl)phthalate	0.858	0.778	9.3	88	-0.05
84	c	Di-n-octyl phthalate	1.449	1.379	4.8	84	-0.05
85		Indeno(1,2,3-cd)pyrene	0.895	0.914	-2.1	97	-0.07
86	I	Perylene-d12	1.000	1.000	0.0	91	-0.06
87		Benzo(b)fluoranthene	1.308	1.421	-8.6	94	-0.05

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.070	0.963	10.0	82	-0.05
89 C	Benzo(a)pyrene	1.110	1.125	-1.4	88	-0.06
90	Dibenzo(a,h)anthracene	0.955	0.966	-1.2	96	-0.07
91	Benzo(g,h,i)perylene	0.989	1.013	-2.4	100	-0.08

(#) = Out of Range SPCC's out = 0 CCC's out = 1