Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF102816\
Data File : BF090920.D
Acq On : 28 Oct 2016 20:44
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 28 23:17:02 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 28 15:18:12 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	90	0.01
2		1,4-Dioxane	0.585	0.646	-10.4	96	0.02
3		Pyridine	1.586	1.748	-10.2	93	0.03
4		n-Nitrosodimethylamine	0.447	0.597	-33.6#	123	0.03
5	S	2-Fluorophenol	1.144	1.324	-15.7	105	0.02
6		Aniline	1.763	1.992	-13.0	96	0.02
7	S	Phenol-d6	1.543	1.638	-6.2	91	0.01
8		2-Chlorophenol	1.411	1.408	0.2	83	0.01
9		Benzaldehyde	0.797	0.913	-14.6	103	0.01
10	C	Phenol	1.701	1.669	1.9	90	0.01
11		bis(2-Chloroethyl)ether	1.407	1.449	-3.0	84	0.01
12		1,3-Dichlorobenzene	1.443	1.526	-5.8	90	0.01
13	C	1,4-Dichlorobenzene	1.518	1.521	-0.2	84	0.02
14		1,2-Dichlorobenzene	1.455	1.438	1.2	89	0.01
15		Benzyl Alcohol	1.014	1.135	-11.9	92	0.01
16		2,2'-oxybis(1-Chloropropane	1.400	1.925	-37.5#	108	0.01
17		2-Methylphenol	1.073	1.229	-14.5	95	0.02
18		Hexachloroethane	0.552	0.590	-6.9	97	0.01
19	P	n-Nitroso-di-n-propylamine	0.892	1.022	-14.6	92	0.01
20		3+4-Methylphenols	1.243	1.409	-13.4	90	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.01
22		Acetophenone	0.411	0.458	-11.4	98	0.02
23	S	Nitrobenzene-d5	0.306	0.345	-12.7	96	0.01
24		Nitrobenzene	0.323	0.428	-32.5#	125	0.01
25		Isophorone	0.620	0.670	-8.1	101	0.01
26	C	2-Nitrophenol	0.173	0.172	0.6	75	0.01
27		2,4-Dimethylphenol	0.280	0.296	-5.7	84	0.02
28		bis(2-Chloroethoxy)methane	0.353	0.402	-13.9	94	0.02
29	C	2,4-Dichlorophenol	0.260	0.261	-0.4	87	0.01
30		1,2,4-Trichlorobenzene	0.300	0.294	2.0	87	0.01
31		Naphthalene	0.935	1.008	-7.8	106	0.01
32		Benzoic acid	0.203	0.214	-5.4	91	0.02
33		4-Chloroaniline	0.378	0.418	-10.6	89	0.02
34	C	Hexachlorobutadiene	0.176	0.151	14.2	71	0.01
35		Caprolactam	0.081	0.079	2.5	79	0.02
36	C	4-Chloro-3-methylphenol	0.282	0.295	-4.6	91	0.02
37		2-Methylnaphthalene	0.604	0.581	3.8	80	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.01
39		1,2,4,5-Tetrachlorobenzene	0.549	0.584	-6.4	89	0.01
40	P	Hexachlorocyclopentadiene	0.252	0.135	46.4#	43#	0.01
41	S	2,4,6-Tribromophenol	0.206	0.213	-3.4	84	0.02
42	C	2,4,6-Trichlorophenol	0.354	0.355	-0.3	82	0.01
43		2,4,5-Trichlorophenol	0.365	0.398	-9.0	80	0.02
44	S	2-Fluorobiphenyl	1.139	1.194	-4.8	85	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.441	1.480	-2.7	81	0.01
46		2-Chloronaphthalene	1.096	1.141	-4.1	80	0.01
47		2-Nitroaniline	0.309	0.466	-50.8#	121	0.02
48		Acenaphthylene	1.659	1.699	-2.4	77	0.01
49		Dimethylphthalate	1.336	1.356	-1.5	82	0.02
50		2,6-Dinitrotoluene	0.295	0.282	4.4	69	0.01
51	C	Acenaphthene	1.143	1.154	-1.0	80	0.01
52		3-Nitroaniline	0.314	0.341	-8.6	82	0.02
53	P	2,4-Dinitrophenol	0.155	0.104	32.9#	52	0.01
54		Dibenzofuran	1.475	1.495	-1.4	79	0.01
55	P	4-Nitrophenol	0.192	0.225	-17.2	86	0.02
56		2,4-Dinitrotoluene	0.367	0.408	-11.2	81	0.02
57		Fluorene	1.208	1.231	-1.9	78	0.01
58		2,3,4,6-Tetrachlorophenol	0.326	0.318	2.5	82	0.01
59		Diethylphthalate	1.327	1.423	-7.2	90	0.02
60		4-Chlorophenyl-phenylether	0.590	0.645	-9.3	94	0.02
61		4-Nitroaniline	0.311	0.372	-19.6	92	0.02
62		Azobenzene	1.305	1.598	-22.5	112	0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	80	0.02
64		4,6-Dinitro-2-methylphenol	0.133	0.103	22.6	55	0.01
65	c	n-Nitrosodiphenylamine	0.613	0.639	-4.2	83	0.01
66		4-Bromophenyl-phenylether	0.219	0.219	0.0	81	0.02
67		Hexachlorobenzene	0.259	0.231	10.8	66	0.01
68		Atrazine	0.190	0.193	-1.6	96	0.02
69	C	Pentachlorophenol	0.139	0.108	22.3#	59	0.01
70		Phenanthrene	1.018	1.008	1.0	76	0.01
71		Anthracene	1.072	1.192	-11.2	94	0.02
72		Carbazole	0.947	1.071	-13.1	90	0.02
73		Di-n-butylphthalate	1.221	1.304	-6.8	89	0.02
74	C	Fluoranthene	1.115	1.061	4.8	75	0.02
75	I	Chrysene-d12	1.000	1.000	0.0	72	0.02
76		Benzidine	0.437	0.315	27.9#	51	0.02
77		Pyrene	1.266	1.409	-11.3	78	0.02
78	S	Terphenyl-d14	0.794	0.849	-6.9	74	0.02
79		Butylbenzylphthalate	0.574	0.734	-27.9#	96	0.02
80		Benzo(a)anthracene	1.061	1.133	-6.8	79	0.02
81		3,3'-Dichlorobenzidine	0.412	0.396	3.9	64	0.02
82		Chrysene	1.074	0.983	8.5	68	0.02
83		Bis(2-ethylhexyl)phthalate	0.748	0.901	-20.5	85	0.02
84	c	Di-n-octyl phthalate	1.256	1.447	-15.2	85	0.02
85		Indeno(1,2,3-cd)pyrene	0.951	1.083	-13.9	90	0.06
86	I	Perylene-d12	1.000	1.000	0.0	71	0.03
87		Benzo(b)fluoranthene	1.345	1.194	11.2	55	0.02

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88	Benzo(k)fluoranthene	1.063	1.333	-25.4#	93	0.03
89 C	Benzo(a)pyrene	1.156	1.166	-0.9	68	0.03
90	Dibenzo(a,h)anthracene	0.978	1.240	-26.8#	90	0.06
91	Benzo(g,h,i)perylene	0.921	1.154	-25.3#	91	0.07

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