Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
Data File : BF088588.D
Acq On : 1 Jul 2016 22:11
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 02 03:32:27 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 30 18:01:55 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	98	0.00
2		1,4-Dioxane	0.662	0.595	10.1	91	-0.02
3		Pyridine	1.920	1.731	9.8	92	-0.02
4		n-Nitrosodimethylamine	0.733	0.647	11.7	89	-0.02
5	S	2-Fluorophenol	1.334	1.166	12.6	85	-0.01
6		Aniline	2.513	2.163	13.9	84	-0.01
7	S	Phenol-d6	1.724	1.574	8.7	93	0.00
8		2-Chlorophenol	1.434	1.366	4.7	100	0.00
9		Benzaldehyde	1.168	1.071	8.3	96	0.00
10	C	Phenol	1.968	1.838	6.6	91	0.00
11		bis(2-Chloroethyl)ether	1.468	1.332	9.3	95	-0.01
12		1,3-Dichlorobenzene	1.578	1.500	4.9	102	0.00
13	C	1,4-Dichlorobenzene	1.567	1.456	7.1	97	0.00
14		1,2-Dichlorobenzene	1.466	1.382	5.7	103	-0.01
15		Benzyl Alcohol	1.269	1.058	16.6	80	-0.01
16		2,2'-oxybis(1-Chloropropane	2.125	1.827	14.0	85	0.00
17		2-Methylphenol	1.170	1.079	7.8	91	0.00
18		Hexachloroethane	0.606	0.483	20.3	80	-0.01
19	P	n-Nitroso-di-n-propylamine	1.158	1.080	6.7	106	0.00
20		3+4-Methylphenols	1.404	1.180	16.0	88	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	89	-0.01
22		Acetophenone	0.485	0.445	8.2	83	-0.01
23	S	Nitrobenzene-d5	0.382	0.380	0.5	94	-0.01
24		Nitrobenzene	0.385	0.370	3.9	90	-0.01
25		Isophorone	0.707	0.657	7.1	87	-0.01
26	C	2-Nitrophenol	0.158	0.160	-1.3	98	0.00
27		2,4-Dimethylphenol	0.329	0.310	5.8	83	-0.01
28		bis(2-Chloroethoxy)methane	0.460	0.460	0.0	97	-0.01
29	C	2,4-Dichlorophenol	0.266	0.276	-3.8	99	0.00
30		1,2,4-Trichlorobenzene	0.291	0.294	-1.0	90	-0.01
31		Naphthalene	0.986	0.936	5.1	84	-0.01
32		Benzoic acid	0.239	0.168	29.7#	63	-0.01
33		4-Chloroaniline	0.439	0.450	-2.5	93	0.00
34	C	Hexachlorobutadiene	0.156	0.155	0.6	89	-0.01
35		Caprolactam	0.090	0.086	4.4	82	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.309	1.6	89	0.00
37		2-Methylnaphthalene	0.635	0.640	-0.8	102	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
39		1,2,4,5-Tetrachlorobenzene	0.665	0.694	-4.4	84	-0.01
40	P	Hexachlorocyclopentadiene	0.327	0.075	77.1#	19#	-0.01
41	S	2,4,6-Tribromophenol	0.186	0.158	15.1	68	-0.01
42	C	2,4,6-Trichlorophenol	0.439	0.469	-6.8	97	0.00
43		2,4,5-Trichlorophenol	0.377	0.380	-0.8	81	0.00
44	S	2-Fluorobiphenyl	1.266	1.441	-13.8	105	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.656	1.691	-2.1	83	-0.01
46		2-Chloronaphthalene	1.300	1.294	0.5	81	-0.01
47		2-Nitroaniline	0.461	0.458	0.7	75	-0.01
48		Acenaphthylene	2.069	2.052	0.8	82	-0.01
49		Dimethylphthalate	1.425	1.663	-16.7	106	0.00
50		2,6-Dinitrotoluene	0.316	0.348	-10.1	98	0.00
51	C	Acenaphthene	1.268	1.210	4.6	83	0.00
52		3-Nitroaniline	0.401	0.372	7.2	68	-0.01
53	P	2,4-Dinitrophenol	0.139	0.020#	85.6#	12#	-0.01
54		Dibenzofuran	1.660	1.605	3.3	81	-0.01
55	P	4-Nitrophenol	0.305	0.285	6.6	71	0.00
56		2,4-Dinitrotoluene	0.413	0.419	-1.5	89	0.00
57		Fluorene	1.279	1.113	13.0	70	-0.01
58		2,3,4,6-Tetrachlorophenol	0.292	0.282	3.4	78	0.00
59		Diethylphthalate	1.430	1.506	-5.3	88	-0.01
60		4-Chlorophenyl-phenylether	0.594	0.599	-0.8	92	0.00
61		4-Nitroaniline	0.386	0.369	4.4	86	-0.01
62		Azobenzene	1.631	1.618	0.8	87	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
64		4,6-Dinitro-2-methylphenol	0.107	0.023	78.5#	15#	-0.01
65	c	n-Nitrosodiphenylamine	0.677	0.687	-1.5	75	-0.01
66		4-Bromophenyl-phenylether	0.202	0.211	-4.5	82	0.00
67		Hexachlorobenzene	0.223	0.221	0.9	76	-0.01
68		Atrazine	0.211	0.178	15.6	62	-0.01
69	C	Pentachlorophenol	0.132	0.113	14.4	62	0.00
70		Phenanthrene	1.093	1.064	2.7	78	0.00
71		Anthracene	1.087	0.973	10.5	68	-0.01
72		Carbazole	1.023	0.989	3.3	82	0.00
73		Di-n-butylphthalate	1.190	1.275	-7.1	86	0.00
74	C	Fluoranthene	1.108	0.914	17.5	60	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	75	0.00
76		Benzidine	1.011	0.894	11.6	65	-0.01
77		Pyrene	1.696	1.434	15.4	60	-0.01
78	S	Terphenyl-d14	0.898	0.763	15.0	65	-0.01
79		Butylbenzylphthalate	0.756	0.690	8.7	69	-0.01
80		Benzo(a)anthracene	1.288	1.194	7.3	69	0.00
81		3,3'-Dichlorobenzidine	0.484	0.527	-8.9	85	0.00
82		Chrysene	1.274	1.227	3.7	73	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	0.851	1.5	70	-0.01
84	c	Di-n-octyl phthalate	1.467	1.580	-7.7	77	-0.01
85		Indeno(1,2,3-cd)pyrene	0.980	0.925	5.6	74	0.00
86	I	Perylene-d12	1.000	1.000	0.0	76	0.00
87		Benzo(b)fluoranthene	1.255	1.438	-14.6	91	-0.01

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88	Benzo(k)fluoranthene	1.092	0.855	21.7	59	0.00
89 C	Benzo(a)pyrene	1.062	1.048	1.3	74	-0.01
90	Dibenzo(a,h)anthracene	0.887	0.867	2.3	75	-0.01
91	Benzo(g,h,i)perylene	0.918	0.840	8.5	71	-0.01

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