Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF071117\
Data File : BF096635.D
Acq On : 11 Jul 2017 10:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 17:03:56 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 07 13:02:33 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	80	-0.01
2		1,4-Dioxane	0.643	0.577	10.3	69	-0.03
3		Pyridine	1.763	1.406	20.2	64	-0.03
4		n-Nitrosodimethylamine	0.870	0.763	12.3	69	-0.04
5	S	2-Fluorophenol	1.355	1.250	7.7	75	-0.01
6		Aniline	2.153	1.981	8.0	74	-0.01
7	S	Phenol-d6	1.666	1.588	4.7	77	-0.01
8		2-Chlorophenol	1.445	1.456	-0.8	81	-0.01
9		Benzaldehyde	1.043	0.916	12.2	70	-0.01
10	C	Phenol	1.898	1.751	7.7	74	-0.01
11		bis(2-Chloroethyl)ether	1.467	1.405	4.2	76	-0.01
12		1,3-Dichlorobenzene	1.515	1.522	-0.5	82	0.00
13	C	1,4-Dichlorobenzene	1.514	1.499	1.0	78	-0.01
14		1,2-Dichlorobenzene	1.390	1.358	2.3	80	-0.01
15		Benzyl Alcohol	1.114	1.074	3.6	77	-0.01
16		2,2'-oxybis(1-Chloropropane	2.982	2.900	2.7	77	0.00
17		2-Methylphenol	1.152	1.085	5.8	75	0.00
18		Hexachloroethane	0.549	0.561	-2.2	81	-0.01
19	P	n-Nitroso-di-n-propylamine	0.992	0.953	3.9	77	-0.01
20		3+4-Methylphenols	1.284	1.294	-0.8	80	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	77	-0.01
22		Acetophenone	0.437	0.451	-3.2	82	-0.01
23	S	Nitrobenzene-d5	0.333	0.346	-3.9	79	-0.01
24		Nitrobenzene	0.349	0.359	-2.9	78	-0.01
25		Isophorone	0.645	0.640	0.8	76	-0.01
26	C	2-Nitrophenol	0.185	0.196	-5.9	79	0.00
27		2,4-Dimethylphenol	0.315	0.300	4.8	73	-0.01
28		bis(2-Chloroethoxy)methane	0.426	0.438	-2.8	78	-0.01
29	C	2,4-Dichlorophenol	0.271	0.287	-5.9	81	-0.01
30		1,2,4-Trichlorobenzene	0.256	0.272	-6.3	81	-0.01
31		Naphthalene	0.944	0.935	1.0	77	0.00
32		Benzoic acid	0.264	0.239	9.5	67	0.00
33		4-Chloroaniline	0.392	0.408	-4.1	80	0.00
34	C	Hexachlorobutadiene	0.131	0.142	-8.4	84	-0.01
35		Caprolactam	0.094	0.085	9.6	72	0.00
36	C	4-Chloro-3-methylphenol	0.300	0.289	3.7	75	-0.01
37		2-Methylnaphthalene	0.601	0.607	-1.0	79	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	83	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.519	0.520	-0.2	84	-0.01
40	P	Hexachlorocyclopentadiene	0.253	0.223	11.9	69	-0.01
41	S	2,4,6-Tribromophenol	0.156	0.166	-6.4	89	-0.01
42	C	2,4,6-Trichlorophenol	0.411	0.380	7.5	77	0.00
43		2,4,5-Trichlorophenol	0.406	0.385	5.2	79	0.00
44	S	2-Fluorobiphenyl	1.170	1.135	3.0	80	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.714	1.582	7.7	78	-0.01
46		2-Chloronaphthalene	1.277	1.166	8.7	77	-0.01
47		2-Nitroaniline	0.465	0.414	11.0	73	0.00
48		Acenaphthylene	2.024	1.870	7.6	78	-0.01
49		Dimethylphthalate	1.493	1.424	4.6	81	-0.01
50		2,6-Dinitrotoluene	0.330	0.320	3.0	79	0.00
51	C	Acenaphthene	1.247	1.122	10.0	76	-0.01
52		3-Nitroaniline	0.412	0.381	7.5	77	0.00
53	P	2,4-Dinitrophenol	0.175	0.167	4.6	76	-0.01
54		Dibenzofuran	1.647	1.618	1.8	82	-0.01
55	P	4-Nitrophenol	0.341	0.304	10.9	72	0.00
56		2,4-Dinitrotoluene	0.418	0.423	-1.2	81	-0.01
57		Fluorene	1.163	1.198	-3.0	85	0.00
58		2,3,4,6-Tetrachlorophenol	0.281	0.290	-3.2	85	-0.01
59		Diethylphthalate	1.411	1.443	-2.3	86	-0.01
60		4-Chlorophenyl-phenylether	0.506	0.505	0.2	84	-0.01
61		4-Nitroaniline	0.374	0.388	-3.7	86	0.00
62		Azobenzene	1.365	1.329	2.6	80	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.138	0.0	80	-0.01
65	c	n-Nitrosodiphenylamine	0.702	0.681	3.0	83	-0.01
66		4-Bromophenyl-phenylether	0.195	0.203	-4.1	89	-0.01
67		Hexachlorobenzene	0.213	0.219	-2.8	87	-0.01
68		Atrazine	0.200	0.200	0.0	85	0.00
69	C	Pentachlorophenol	0.126	0.127	-0.8	80	0.00
70		Phenanthrene	1.081	1.049	3.0	85	-0.01
71		Anthracene	1.102	1.090	1.1	85	-0.01
72		Carbazole	1.108	1.082	2.3	84	-0.01
73		Di-n-butylphthalate	1.342	1.278	4.8	82	0.00
74	C	Fluoranthene	1.060	1.040	1.9	84	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	86	-0.01
76		Benzidine	0.960	0.827	13.9	74	0.00
77		Pyrene	1.605	1.549	3.5	84	-0.01
78	S	Terphenyl-d14	0.936	0.921	1.6	89	-0.01
79		Butylbenzylphthalate	0.813	0.807	0.7	85	-0.01
80		Benzo(a)anthracene	1.158	1.087	6.1	83	-0.01
81		3,3'-Dichlorobenzidine	0.476	0.480	-0.8	86	-0.01
82		Chrysene	1.213	1.196	1.4	86	-0.01
83		Bis(2-ethylhexyl)phthalate	0.907	0.934	-3.0	91	-0.01
84	c	Di-n-octyl phthalate	1.786	1.880	-5.3	91	-0.01
85		Indeno(1,2,3-cd)pyrene	0.957	0.918	4.1	87	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	90	-0.02
87		Benzo(b)fluoranthene	1.253	1.198	4.4	84	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.121	1.106	1.3	92	-0.01
89 C	Benzo(a)pyrene	1.119	1.092	2.4	89	-0.02
90	Dibenzo(a,h)anthracene	0.880	0.814	7.5	86	-0.02
91	Benzo(g,h,i)perylene	0.912	0.895	1.9	91	-0.02

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