Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF081217\
Data File : BF097654.D
Acq On : 13 Aug 2017 3:30
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 14 12:09:41 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 11 15:55:57 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	87	0.00
2		1,4-Dioxane	0.546	0.544	0.4	92	-0.03
3		Pyridine	1.593	1.633	-2.5	82	-0.02
4		n-Nitrosodimethylamine	0.924	0.942	-1.9	85	-0.02
5	S	2-Fluorophenol	1.258	1.360	-8.1	87	0.02
6		Aniline	2.243	2.216	1.2	89	0.00
7	S	Phenol-d6	1.561	1.531	1.9	87	0.03
8		2-Chlorophenol	1.442	1.472	-2.1	89	0.01
9		Benzaldehyde	1.029	1.059	-2.9	89	0.00
10	C	Phenol	1.709	1.829	-7.0	87	0.04
11		bis(2-Chloroethyl)ether	1.371	1.490	-8.7	96	0.00
12		1,3-Dichlorobenzene	1.615	1.623	-0.5	88	0.00
13	C	1,4-Dichlorobenzene	1.644	1.671	-1.6	90	0.00
14		1,2-Dichlorobenzene	1.500	1.545	-3.0	91	0.00
15		Benzyl Alcohol	1.079	1.148	-6.4	89	0.01
16		2,2'-oxybis(1-Chloropropane	3.078	3.229	-4.9	93	0.00
17		2-Methylphenol	1.118	1.137	-1.7	90	0.04
18		Hexachloroethane	0.563	0.546	3.0	86	0.00
19	P	n-Nitroso-di-n-propylamine	1.086	1.118	-2.9	88	0.00
20		3+4-Methylphenols	1.369	1.473	-7.6	91	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	83	0.00
22		Acetophenone	0.472	0.502	-6.4	88	0.00
23	S	Nitrobenzene-d5	0.319	0.343	-7.5	87	0.00
24		Nitrobenzene	0.348	0.380	-9.2	87	0.00
25		Isophorone	0.588	0.607	-3.2	83	0.00
26	C	2-Nitrophenol	0.176	0.192	-9.1	87	0.00
27		2,4-Dimethylphenol	0.293	0.309	-5.5	86	0.01
28		bis(2-Chloroethoxy)methane	0.425	0.452	-6.4	86	0.00
29	C	2,4-Dichlorophenol	0.259	0.263	-1.5	83	0.02
30		1,2,4-Trichlorobenzene	0.256	0.268	-4.7	86	0.00
31		Naphthalene	1.002	1.006	-0.4	85	0.00
32		Benzoic acid	0.158	0.110	30.4#	51	0.00
33		4-Chloroaniline	0.415	0.425	-2.4	83	0.00
34	C	Hexachlorobutadiene	0.140	0.148	-5.7	89	0.00
35		Caprolactam	0.083	0.083	0.0	80	-0.01
36	C	4-Chloro-3-methylphenol	0.290	0.286	1.4	77	0.02
37		2-Methylnaphthalene	0.630	0.630	0.0	83	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	81	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.566	0.607	-7.2	87	-0.01
40	P	Hexachlorocyclopentadiene	0.072	0.007#	90.3#	7#	-0.01
41	S	2,4,6-Tribromophenol	0.152	0.125	17.8	67	0.00
42	C	2,4,6-Trichlorophenol	0.373	0.392	-5.1	82	0.00
43		2,4,5-Trichlorophenol	0.352	0.328	6.8	69	0.02
44	S	2-Fluorobiphenyl	1.325	1.325	0.0	83	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.716	1.803	-5.1	86	0.00
46		2-Chloronaphthalene	1.317	1.353	-2.7	84	-0.01
47		2-Nitroaniline	0.516	0.540	-4.7	77	0.00
48		Acenaphthylene	2.094	2.093	0.0	81	-0.01
49		Dimethylphthalate	1.541	1.594	-3.4	85	-0.01
50		2,6-Dinitrotoluene	0.318	0.320	-0.6	80	0.00
51	C	Acenaphthene	1.255	1.226	2.3	81	0.00
52		3-Nitroaniline	0.417	0.420	-0.7	80	0.00
53	P	2,4-Dinitrophenol	0.131	0.010#	92.4#	7#	0.06
54		Dibenzofuran	1.757	1.730	1.5	81	-0.01
55	P	4-Nitrophenol	0.159	0.069	56.6#	35#	0.06
56		2,4-Dinitrotoluene	0.425	0.429	-0.9	79	0.00
57		Fluorene	1.445	1.398	3.3	79	-0.01
58		2,3,4,6-Tetrachlorophenol	0.242	0.239	1.2	75	0.00
59		Diethylphthalate	1.456	1.522	-4.5	85	-0.01
60		4-Chlorophenyl-phenylether	0.617	0.628	-1.8	84	0.00
61		4-Nitroaniline	0.398	0.362	9.0	70	0.00
62		Azobenzene	1.448	1.405	3.0	78	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	68	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.055	54.9#	29#	0.00
65	c	n-Nitrosodiphenylamine	0.739	0.848	-14.7	80	0.00
66		4-Bromophenyl-phenylether	0.207	0.240	-15.9	79	-0.01
67		Hexachlorobenzene	0.226	0.241	-6.6	73	-0.01
68		Atrazine	0.185	0.196	-5.9	71	0.00
69	C	Pentachlorophenol	0.037	0.223	-502.7#	348#	0.00
70		Phenanthrene	1.152	1.148	0.3	69	0.00
71		Anthracene	1.179	1.165	1.2	69	0.00
72		Carbazole	1.122	1.076	4.1	67	0.00
73		Di-n-butylphthalate	1.313	1.486	-13.2	77	-0.01
74	C	Fluoranthene	1.041	0.988	5.1	61	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	61	-0.01
76		Benzidine	0.659	0.549	16.7	51	0.00
77		Pyrene	1.654	1.505	9.0	60	-0.01
78	S	Terphenyl-d14	0.966	0.847	12.3	62	-0.01
79		Butylbenzylphthalate	0.691	0.742	-7.4	62	0.00
80		Benzo(a)anthracene	1.168	1.182	-1.2	63	-0.01
81		3,3'-Dichlorobenzidine	0.403	0.429	-6.5	65	0.00
82		Chrysene	1.162	1.155	0.6	60	-0.01
83		Bis(2-ethylhexyl)phthalate	0.983	1.008	-2.5	59	-0.01
84	c	Di-n-octyl phthalate	1.717	1.708	0.5	63	-0.01
85		Indeno(1,2,3-cd)pyrene	0.904	1.032	-14.2	88	0.00
86	I	Perylene-d12	1.000	1.000	0.0	73	0.00
87		Benzo(b)fluoranthene	1.201	1.212	-0.9	75	0.00

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88	Benzo(k)fluoranthene	1.210	1.255	-3.7	74	0.00
89 C	Benzo(a)pyrene	1.102	1.073	2.6	72	0.00
90	Dibenzo(a,h)anthracene	0.789	0.946	-19.9	90	0.00
91	Benzo(g,h,i)perylene	0.866	0.932	-7.6	85	0.00

(#) = Out of Range SPCC's out = 2 CCC's out = 1