Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF012315\
Data File : BF077045.D
Acq On : 23 Jan 2015 19:29
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Jan 26 11:07:24 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 23 22:30:04 2015
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	95	0.00
2		1,4-Dioxane	0.521	0.420	19.4	81	0.00
3		Pyridine	1.348	1.443	-7.0	84	0.00
4		n-Nitrosodimethylamine	0.668	0.634	5.1	89	0.00
5	S	2-Fluorophenol	1.216	1.229	-1.1	91	0.01
6		Aniline	2.124	2.167	-2.0	91	0.01
7	S	Phenol-d6	1.535	1.528	0.5	90	0.00
8		2-Chlorophenol	1.402	1.419	-1.2	90	0.00
9		Benzaldehyde	0.894	0.885	1.0	107	0.00
10	C	Phenol	1.629	1.619	0.6	86	0.00
11		bis(2-Chloroethyl)ether	1.275	1.321	-3.6	95	0.00
12		1,3-Dichlorobenzene	1.595	1.602	-0.4	92	0.01
13	C	1,4-Dichlorobenzene	1.692	1.682	0.6	92	0.01
14		1,2-Dichlorobenzene	1.559	1.568	-0.6	91	0.00
15		Benzyl Alcohol	1.242	1.259	-1.4	89	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.730	-0.9	92	0.00
17		2-Methylphenol	1.146	1.152	-0.5	89	0.00
18		Hexachloroethane	0.588	0.621	-5.6	95	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	1.078	-0.4	89	0.00
20		3+4-Methylphenols	1.517	1.475	2.8	88	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
22		Acetophenone	0.490	0.486	0.8	90	0.01
23	S	Nitrobenzene-d5	0.361	0.366	-1.4	91	0.00
24		Nitrobenzene	0.368	0.377	-2.4	91	0.00
25		Isophorone	0.657	0.655	0.3	89	0.00
26	C	2-Nitrophenol	0.172	0.183	-6.4	90	0.00
27		2,4-Dimethylphenol	0.284	0.294	-3.5	91	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.379	-1.3	89	0.00
29	C	2,4-Dichlorophenol	0.265	0.276	-4.2	91	-0.01
30		1,2,4-Trichlorobenzene	0.294	0.312	-6.1	92	0.00
31		Naphthalene	1.030	1.024	0.6	89	0.00
32		Benzoic acid	0.158	0.134	15.2	75	0.00
33		4-Chloroaniline	0.416	0.406	2.4	88	0.00
34	C	Hexachlorobutadiene	0.175	0.178	-1.7	91	0.00
35		Caprolactam	0.078	0.082	-5.1	90	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.303	0.0	89	0.00
37		2-Methylnaphthalene	0.703	0.702	0.1	91	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.497	-0.6	90	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.267	-32.2#	98	0.00
41	S	2,4,6-Tribromophenol	0.162	0.171	-5.6	90	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.369	-2.5	87	0.00
43		2,4,5-Trichlorophenol	0.367	0.395	-7.6	92	0.00
44	S	2-Fluorobiphenyl	1.314	1.354	-3.0	91	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.483	1.549	-4.5	90	0.00
46		2-Chloronaphthalene	1.220	1.232	-1.0	90	0.00
47		2-Nitroaniline	0.370	0.377	-1.9	86	0.00
48		Acenaphthylene	2.058	2.096	-1.8	89	0.00
49		Dimethylphthalate	1.418	1.417	0.1	89	0.00
50		2,6-Dinitrotoluene	0.283	0.290	-2.5	85	0.00
51	C	Acenaphthene	1.168	1.179	-0.9	89	0.00
52		3-Nitroaniline	0.339	0.360	-6.2	88	-0.01
53	P	2,4-Dinitrophenol	0.095	0.117	-23.2	101	0.00
54		Dibenzofuran	1.701	1.704	-0.2	89	0.00
55	P	4-Nitrophenol	0.222	0.207	6.8	80	-0.01
56		2,4-Dinitrotoluene	0.384	0.420	-9.4	90	0.00
57		Fluorene	1.441	1.459	-1.2	90	-0.01
58		2,3,4,6-Tetrachlorophenol	0.268	0.309	-15.3	98	0.00
59		Diethylphthalate	1.433	1.475	-2.9	91	0.00
60		4-Chlorophenyl-phenylether	0.590	0.618	-4.7	93	0.00
61		4-Nitroaniline	0.328	0.336	-2.4	85	0.00
62		Azobenzene	1.494	1.511	-1.1	89	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.122	-17.3	90	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.706	1.0	88	0.00
66		4-Bromophenyl-phenylether	0.203	0.206	-1.5	89	0.00
67		Hexachlorobenzene	0.214	0.215	-0.5	92	0.00
68		Atrazine	0.216	0.224	-3.7	88	0.00
69	C	Pentachlorophenol	0.082	0.104	-26.8#	111	0.00
70		Phenanthrene	1.247	1.219	2.2	90	0.00
71		Anthracene	1.216	1.218	-0.2	93	0.00
72		Carbazole	1.180	1.154	2.2	88	0.00
73		Di-n-butylphthalate	1.365	1.442	-5.6	94	-0.01
74	C	Fluoranthene	1.278	1.240	3.0	86	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.640	0.593	7.3	93	-0.01
77		Pyrene	1.371	1.308	4.6	87	0.00
78	S	Terphenyl-d14	0.869	0.833	4.1	89	0.00
79		Butylbenzylphthalate	0.621	0.664	-6.9	94	0.00
80		Benzo(a)anthracene	1.255	1.263	-0.6	93	0.00
81		3,3'-Dichlorobenzidine	0.399	0.423	-6.0	99	0.00
82		Chrysene	1.144	1.104	3.5	89	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.964	-12.2	104	-0.01
84	c	Di-n-octyl phthalate	1.491	1.625	-9.0	100	-0.01
85		Indeno(1,2,3-cd)pyrene	1.213	1.276	-5.2	96	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	98	0.00
87		Benzo(b)fluoranthene	1.347	1.263	6.2	96	-0.01

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88	Benzo(k)fluoranthene	1.177	1.298	-10.3	95	0.00
89 C	Benzo(a)pyrene	1.188	1.193	-0.4	93	0.00
90	Dibenzo(a,h)anthracene	1.090	1.146	-5.1	97	-0.01
91	Benzo(g,h,i)perylene	1.084	1.146	-5.7	98	-0.01

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