Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF031816\
Data File : BF085544.D
Acq On : 19 Mar 2016 00:34
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF031816

Quant Time: Mar 19 03:35:54 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 18 23:31:39 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	99	0.00
2		1,4-Dioxane	0.579	0.568	1.9	94	-0.01
3		Pyridine	1.421	1.540	-8.4	106	-0.01
4		n-Nitrosodimethylamine	0.593	0.602	-1.5	102	0.00
5	S	2-Fluorophenol	1.191	1.220	-2.4	96	0.00
6		Aniline	2.051	2.032	0.9	98	0.00
7	S	Phenol-d6	1.527	1.563	-2.4	99	0.00
8		2-Chlorophenol	1.377	1.403	-1.9	99	0.00
9		Benzaldehyde	0.768	0.724	5.7	100	0.00
10	C	Phenol	1.739	1.836	-5.6	94	0.00
11		bis(2-Chloroethyl)ether	1.311	1.382	-5.4	100	0.00
12		1,3-Dichlorobenzene	1.501	1.481	1.3	99	-0.01
13	C	1,4-Dichlorobenzene	1.531	1.446	5.6	98	0.00
14		1,2-Dichlorobenzene	1.359	1.411	-3.8	96	0.00
15		Benzyl Alcohol	1.046	0.990	5.4	95	0.00
16		2,2'-oxybis(1-Chloropropane	1.725	1.766	-2.4	97	0.00
17		2-Methylphenol	1.142	1.098	3.9	95	0.00
18		Hexachloroethane	0.553	0.561	-1.4	100	0.00
19	P	n-Nitroso-di-n-propylamine	0.912	0.887	2.7	93	0.00
20		3+4-Methylphenols	1.323	1.313	0.8	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.475	0.448	5.7	100	0.00
23	S	Nitrobenzene-d5	0.345	0.356	-3.2	96	-0.01
24		Nitrobenzene	0.377	0.379	-0.5	104	0.00
25		Isophorone	0.690	0.678	1.7	98	0.00
26	C	2-Nitrophenol	0.186	0.214	-15.1	101	0.00
27		2,4-Dimethylphenol	0.330	0.315	4.5	99	0.00
28		bis(2-Chloroethoxy)methane	0.392	0.406	-3.6	95	0.00
29	C	2,4-Dichlorophenol	0.272	0.287	-5.5	100	0.00
30		1,2,4-Trichlorobenzene	0.305	0.293	3.9	97	0.00
31		Naphthalene	1.010	0.942	6.7	97	0.00
32		Benzoic acid	0.276	0.274	0.7	95	0.00
33		4-Chloroaniline	0.414	0.428	-3.4	95	0.00
34	C	Hexachlorobutadiene	0.162	0.168	-3.7	98	0.00
35		Caprolactam	0.094	0.100	-6.4	103	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.329	-3.5	98	0.00
37		2-Methylnaphthalene	0.615	0.657	-6.8	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
39		1,2,4,5-Tetrachlorobenzene	0.607	0.556	8.4	99	0.00
40	P	Hexachlorocyclopentadiene	0.273	0.287	-5.1	98	0.00
41	S	2,4,6-Tribromophenol	0.196	0.206	-5.1	102	0.00
42	C	2,4,6-Trichlorophenol	0.415	0.408	1.7	99	0.00
43		2,4,5-Trichlorophenol	0.423	0.410	3.1	93	0.00
44	S	2-Fluorobiphenyl	1.250	1.288	-3.0	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.725	1.606	6.9	95	0.00
46		2-Chloronaphthalene	1.260	1.250	0.8	95	0.00
47		2-Nitroaniline	0.381	0.408	-7.1	97	0.00
48		Acenaphthylene	2.040	2.144	-5.1	99	0.00
49		Dimethylphthalate	1.509	1.452	3.8	97	0.00
50		2,6-Dinitrotoluene	0.319	0.293	8.2	96	0.00
51	C	Acenaphthene	1.280	1.309	-2.3	98	0.00
52		3-Nitroaniline	0.399	0.400	-0.3	105	0.00
53	P	2,4-Dinitrophenol	0.195	0.185	5.1	97	0.00
54		Dibenzofuran	1.559	1.648	-5.7	100	0.00
55	P	4-Nitrophenol	0.308	0.324	-5.2	97	0.00
56		2,4-Dinitrotoluene	0.417	0.409	1.9	88	-0.01
57		Fluorene	1.299	1.361	-4.8	98	0.00
58		2,3,4,6-Tetrachlorophenol	0.304	0.299	1.6	99	0.00
59		Diethylphthalate	1.501	1.503	-0.1	100	0.00
60		4-Chlorophenyl-phenylether	0.634	0.602	5.0	101	0.00
61		4-Nitroaniline	0.393	0.414	-5.3	94	0.00
62		Azobenzene	1.440	1.428	0.8	97	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	95	-0.01
64		4,6-Dinitro-2-methylphenol	0.128	0.125	2.3	92	0.00
65	c	n-Nitrosodiphenylamine	0.623	0.629	-1.0	92	0.00
66		4-Bromophenyl-phenylether	0.200	0.214	-7.0	98	0.00
67		Hexachlorobenzene	0.212	0.227	-7.1	94	0.00
68		Atrazine	0.190	0.205	-7.9	95	0.00
69	C	Pentachlorophenol	0.129	0.130	-0.8	90	0.00
70		Phenanthrene	1.055	1.105	-4.7	95	0.00
71		Anthracene	1.106	0.995	10.0	95	0.00
72		Carbazole	0.954	1.014	-6.3	97	0.00
73		Di-n-butylphthalate	1.195	1.268	-6.1	97	0.00
74	C	Fluoranthene	1.104	1.055	4.4	95	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
76		Benzidine	0.673	0.649	3.6	92	0.00
77		Pyrene	1.436	1.348	6.1	88	-0.01
78	S	Terphenyl-d14	0.831	0.762	8.3	97	0.00
79		Butylbenzylphthalate	0.685	0.646	5.7	89	-0.01
80		Benzo(a)anthracene	1.161	1.111	4.3	95	0.00
81		3,3'-Dichlorobenzidine	0.426	0.400	6.1	93	0.00
82		Chrysene	1.117	1.085	2.9	86	0.00
83		Bis(2-ethylhexyl)phthalate	0.850	0.793	6.7	91	0.00
84	c	Di-n-octyl phthalate	1.386	1.391	-0.4	94	0.00
85		Indeno(1,2,3-cd)pyrene	0.933	0.835	10.5	93	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	91	0.00
87		Benzo(b)fluoranthene	1.305	1.178	9.7	91	0.00

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88	Benzo(k)fluoranthene	1.075	1.241	-15.4	88	0.00
89 C	Benzo(a)pyrene	1.106	1.116	-0.9	90	0.00
90	Dibenzo(a,h)anthracene	0.923	0.897	2.8	92	0.00
91	Benzo(g,h,i)perylene	0.893	0.855	4.3	92	0.00

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