Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF071618\
Data File : BF107408.D
Acq On : 16 Jul 2018 17:06
Operator : JU/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF071618

Quant Time: Jul 16 17:33:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 16 17:24:26 2018
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	96	0.00
2		1,4-Dioxane	0.650	0.599	7.8	93	-0.02
3		Pyridine	1.754	1.606	8.4	93	-0.02
4		n-Nitrosodimethylamine	0.720	0.717	0.4	98	-0.02
5	S	2-Fluorophenol	1.317	1.254	4.8	97	-0.02
6		Aniline	2.251	2.156	4.2	95	0.00
7	S	Phenol-d6	1.703	1.595	6.3	96	-0.01
8		2-Chlorophenol	1.332	1.232	7.5	92	0.00
9		Benzaldehyde	1.370	1.282	6.4	93	-0.01
10	C	Phenol	1.825	1.704	6.6	94	-0.02
11		bis(2-Chloroethyl)ether	1.490	1.402	5.9	97	0.00
12		1,3-Dichlorobenzene	1.617	1.570	2.9	102	0.00
13	C	1,4-Dichlorobenzene	1.603	1.483	7.5	92	0.00
14		1,2-Dichlorobenzene	1.500	1.417	5.5	98	0.00
15		Benzyl Alcohol	1.710	1.671	2.3	97	-0.01
16		2,2'-oxybis(1-Chloropropane	1.590	1.551	2.5	95	0.00
17		2-Methylphenol	1.266	1.196	5.5	96	-0.01
18		Hexachloroethane	0.596	0.599	-0.5	99	0.00
19	P	n-Nitroso-di-n-propylamine	1.219	1.147	5.9	92	0.00
20		3+4-Methylphenols	1.608	1.555	3.3	97	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.550	0.518	5.8	97	0.00
23	S	Nitrobenzene-d5	0.375	0.406	-8.3	99	-0.01
24		Nitrobenzene	0.421	0.451	-7.1	99	0.00
25		Isophorone	0.743	0.701	5.7	95	0.00
26	C	2-Nitrophenol	0.143	0.158	-10.5	102	0.00
27		2,4-Dimethylphenol	0.299	0.291	2.7	99	-0.01
28		bis(2-Chloroethoxy)methane	0.466	0.450	3.4	98	0.00
29	C	2,4-Dichlorophenol	0.311	0.308	1.0	99	-0.01
30		1,2,4-Trichlorobenzene	0.377	0.351	6.9	95	0.00
31		Naphthalene	0.971	0.898	7.5	95	0.00
32		Benzoic acid	0.183	0.205	-12.0	112	-0.02
33		4-Chloroaniline	0.421	0.402	4.5	95	0.00
34	C	Hexachlorobutadiene	0.252	0.256	-1.6	101	0.00
35		Caprolactam	0.101	0.094	6.9	89	-0.02
36	C	4-Chloro-3-methylphenol	0.406	0.397	2.2	97	-0.01
37		2-Methylnaphthalene	0.640	0.623	2.7	98	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
39		1,2,4,5-Tetrachlorobenzene	0.778	0.687	11.7	97	0.00
40	P	Hexachlorocyclopentadiene	0.406	0.401	1.2	100	0.00
41	S	2,4,6-Tribromophenol	0.176	0.162	8.0	99	-0.01
42	C	2,4,6-Trichlorophenol	0.486	0.449	7.6	98	-0.01
43		2,4,5-Trichlorophenol	0.465	0.449	3.4	98	-0.02
44	S	2-Fluorobiphenyl	1.392	1.184	14.9	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.729	1.477	14.6	95	0.00
46		2-Chloronaphthalene	1.366	1.209	11.5	98	0.00
47		2-Nitroaniline	0.418	0.432	-3.3	104	-0.01
48		Acenaphthylene	1.906	1.746	8.4	99	0.00
49		Dimethylphthalate	1.589	1.445	9.1	99	0.00
50		2,6-Dinitrotoluene	0.303	0.305	-0.7	102	-0.01
51	C	Acenaphthene	1.264	1.113	11.9	97	0.00
52		3-Nitroaniline	0.319	0.307	3.8	105	-0.01
53	P	2,4-Dinitrophenol	0.107	0.117	-9.3	114	-0.01
54		Dibenzofuran	1.953	1.710	12.4	99	0.00
55	P	4-Nitrophenol	0.258	0.258	0.0	104	-0.02
56		2,4-Dinitrotoluene	0.395	0.411	-4.1	106	0.00
57		Fluorene	1.293	1.136	12.1	99	-0.01
58		2,3,4,6-Tetrachlorophenol	0.450	0.397	11.8	98	-0.01
59		Diethylphthalate	1.544	1.431	7.3	100	0.00
60		4-Chlorophenyl-phenylether	0.770	0.724	6.0	102	0.00
61		4-Nitroaniline	0.303	0.300	1.0	102	-0.01
62		Azobenzene	1.813	1.664	8.2	98	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	-0.01
64		4,6-Dinitro-2-methylphenol	0.095	0.114	-20.0	121	-0.01
65	c	n-Nitrosodiphenylamine	0.632	0.591	6.5	101	0.00
66		4-Bromophenyl-phenylether	0.229	0.220	3.9	103	0.00
67		Hexachlorobenzene	0.212	0.191	9.9	98	-0.01
68		Atrazine	0.239	0.217	9.2	98	0.00
69	C	Pentachlorophenol	0.157	0.158	-0.6	98	-0.01
70		Phenanthrene	0.997	0.891	10.6	97	0.00
71		Anthracene	0.993	0.894	10.0	98	0.00
72		Carbazole	0.940	0.847	9.9	98	-0.01
73		Di-n-butylphthalate	1.080	1.020	5.6	101	0.00
74	C	Fluoranthene	1.133	1.014	10.5	98	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76		Benzidine	0.776	0.728	6.2	96	-0.01
77		Pyrene	1.360	1.293	4.9	98	0.00
78	S	Terphenyl-d14	0.805	0.733	8.9	103	0.00
79		Butylbenzylphthalate	0.495	0.526	-6.3	103	0.00
80		Benzo(a)anthracene	1.224	1.126	8.0	97	-0.01
81		3,3'-Dichlorobenzidine	0.410	0.401	2.2	102	0.00
82		Chrysene	1.169	1.065	8.9	95	0.00
83		Bis(2-ethylhexyl)phthalate	0.710	0.699	1.5	101	0.00
84	c	Di-n-octyl phthalate	1.104	1.093	1.0	101	0.00
85		Indeno(1,2,3-cd)pyrene	0.832	0.691	16.9	96	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
87		Benzo(b)fluoranthene	1.171	1.171	0.0	101	0.00

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88	Benzo(k)fluoranthene	1.142	1.065	6.7	90	0.00
89 C	Benzo(a)pyrene	1.074	1.033	3.8	95	0.00
90	Dibenzo(a,h)anthracene	0.811	0.751	7.4	94	-0.02
91	Benzo(g,h,i)perylene	0.824	0.736	10.7	98	-0.01

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