Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102925\
Data File : BP026036.D
Acq On : 29 Oct 2025 16:09
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP102925

Quant Time: Oct 30 11:55:55 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 30 11:51:34 2025
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	-0.05
2		1,4-Dioxane	0.511	0.425	16.8	79	-0.02
3		Pyridine	1.453	1.186	18.4	77	-0.03
4		n-Nitrosodimethylamine	0.582	0.511	12.2	84	-0.03
5	S	2-Fluorophenol	1.212	1.037	14.4	80	-0.05
6		Aniline	2.059	1.845	10.4	84	-0.05
7	S	Phenol-d6	1.543	1.313	14.9	79	-0.05
8		2-Chlorophenol	1.342	1.161	13.5	80	-0.05
9		Benzaldehyde	0.802	0.923	-15.1	110	-0.05
10	C	Phenol	1.713	1.448	15.5	79	-0.05
11		bis(2-Chloroethyl)ether	1.352	1.145	15.3	80	-0.05
12		1,3-Dichlorobenzene	1.541	1.331	13.6	81	-0.05
13	C	1,4-Dichlorobenzene	1.558	1.367	12.3	83	-0.05
14		1,2-Dichlorobenzene	1.487	1.305	12.2	83	-0.04
15		Benzyl Alcohol	1.112	0.978	12.1	83	-0.05
16		2,2'-oxybis(1-Chloropropane	2.043	1.712	16.2	78	-0.03
17		2-Methylphenol	1.121	0.981	12.5	80	-0.04
18		Hexachloroethane	0.533	0.461	13.5	80	-0.05
19	P	n-Nitroso-di-n-propylamine	0.958	0.870	9.2	82	-0.05
20		3+4-Methylphenols	1.560	1.336	14.4	80	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	97	-0.04
22		Acetophenone	0.506	0.452	10.7	85	-0.05
23	S	Nitrobenzene-d5	0.321	0.275	14.3	79	-0.04
24		Nitrobenzene	0.339	0.295	13.0	80	-0.04
25		Isophorone	0.687	0.602	12.4	81	-0.05
26	C	2-Nitrophenol	0.104	0.110	-5.8	83	-0.04
27		2,4-Dimethylphenol	0.289	0.271	6.2	87	-0.05
28		bis(2-Chloroethoxy)methane	0.433	0.373	13.9	81	-0.03
29	C	2,4-Dichlorophenol	0.305	0.264	13.4	78	-0.03
30		1,2,4-Trichlorobenzene	0.331	0.294	11.2	84	-0.04
31		Naphthalene	1.091	0.924	15.3	80	-0.04
32		Benzoic acid	0.163	0.151	7.4	87	-0.09
33		4-Chloroaniline	0.419	0.407	2.9	90	-0.03
34	C	Hexachlorobutadiene	0.211	0.178	15.6	80	-0.04
35		Caprolactam	0.107	0.093	13.1	80	-0.07
36	C	4-Chloro-3-methylphenol	0.320	0.277	13.4	78	-0.04
37		2-Methylnaphthalene	0.763	0.591	22.5	72	-0.03
38		1-Methylnaphthalene	0.742	0.611	17.7	77	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
40		1,2,4,5-Tetrachlorobenzene	0.621	0.536	13.7	81	-0.03
41	P	Hexachlorocyclopentadiene	0.227	0.268	-18.1	109	-0.03
42	S	2,4,6-Tribromophenol	0.216	0.197	8.8	80	0.00
43	C	2,4,6-Trichlorophenol	0.367	0.333	9.3	80	-0.02
44		2,4,5-Trichlorophenol	0.419	0.370	11.7	78	-0.03
45	S	2-Fluorobiphenyl	1.392	1.166	16.2	79	-0.02
46		1,1'-Biphenyl	1.531	1.343	12.3	83	-0.02
47		2-Chloronaphthalene	1.205	1.002	16.8	78	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.272	0.255	6.3	81	-0.02
49		Acenaphthylene	1.756	1.709	2.7	90	0.00
50		Dimethylphthalate	1.455	1.277	12.2	81	-0.01
51		2,6-Dinitrotoluene	0.237	0.250	-5.5	91	-0.01
52	C	Acenaphthene	1.206	0.993	17.7	76	0.00
53		3-Nitroaniline	0.291	0.290	0.3	86	-0.01
54	P	2,4-Dinitrophenol	0.100	0.086	14.0	83	0.00
55		Dibenzofuran	1.822	1.545	15.2	79	0.00
56	P	4-Nitrophenol	0.290	0.237	18.3	80	-0.01
57		2,4-Dinitrotoluene	0.347	0.340	2.0	83	0.00
58		Fluorene	1.427	1.261	11.6	84	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.323	2.1	85	-0.01
60		Diethylphthalate	1.463	1.280	12.5	79	0.00
61		4-Chlorophenyl-phenylether	0.713	0.604	15.3	79	0.00
62		4-Nitroaniline	0.311	0.294	5.5	83	0.00
63		Azobenzene	1.325	1.161	12.4	81	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	0.077	0.065	15.6	84	0.00
66	c	n-Nitrosodiphenylamine	0.625	0.526	15.8	80	0.00
67		4-Bromophenyl-phenylether	0.222	0.182	18.0	78	0.00
68		Hexachlorobenzene	0.253	0.207	18.2	79	0.00
69		Atrazine	0.211	0.196	7.1	88	0.00
70	C	Pentachlorophenol	0.154	0.130	15.6	80	0.00
71		Phenanthrene	1.164	0.952	18.2	79	0.00
72		Anthracene	1.154	0.964	16.5	81	0.00
73		Carbazole	1.092	0.913	16.4	79	0.00
74		Di-n-butylphthalate	1.222	1.073	12.2	79	0.00
75	C	Fluoranthene	1.358	1.125	17.2	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	97	0.01
77		Benzidine	0.508	0.611	-20.3	115	0.00
78		Pyrene	1.353	1.113	17.7	76	0.00
79	S	Terphenyl-d14	0.984	0.831	15.5	80	0.00
80		Butylbenzylphthalate	0.432	0.411	4.9	81	0.00
81		Benzo(a)anthracene	1.374	1.136	17.3	78	0.01
82		3,3'-Dichlorobenzidine	0.443	0.412	7.0	87	0.00
83		Chrysene	1.302	1.059	18.7	77	0.00
84		Bis(2-ethylhexyl)phthalate	0.718	0.664	7.5	78	0.01
85	c	Di-n-octyl phthalate	1.191	1.002	15.9	76	0.02
86	I	Perylene-d12	1.000	1.000	0.0	96	0.02
87		Indeno(1,2,3-cd)pyrene	1.477	1.229	16.8	75	0.01
88		Benzo(b)fluoranthene	1.242	1.058	14.8	78	0.00
89		Benzo(k)fluoranthene	1.268	1.090	14.0	79	0.00
90	C	Benzo(a)pyrene	1.141	0.999	12.4	79	0.00
91		Dibenzo(a,h)anthracene	1.202	1.018	15.3	77	0.02
92		Benzo(g,h,i)perylene	1.156	0.979	15.3	77	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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