

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122123\
 Data File : BG060128.D
 Acq On : 21 Dec 2023 18:49
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG122123

Quant Time: Dec 22 03:43:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 03:42:04 2023
 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	94	0.00
2	1,4-Dioxane	0.530	0.515	2.8	91	0.00
3	Pyridine	1.714	1.851	-8.0	92	0.00
4	n-Nitrosodimethylamine	0.779	0.813	-4.4	91	0.00
5 S	2-Fluorophenol	1.269	1.277	-0.6	93	0.00
6	Aniline	2.231	2.489	-11.6	94	0.00
7 S	Phenol-d6	2.136	2.296	-7.5	93	0.00
8	2-Chlorophenol	1.406	1.520	-8.1	95	0.00
9	Benzaldehyde	1.185	1.243	-4.9	97	0.00
10 C	Phenol	2.134	2.282	-6.9	92	0.00
11	bis(2-Chloroethyl)ether	1.676	1.739	-3.8	91	0.00
12	1,3-Dichlorobenzene	1.572	1.561	0.7	90	0.00
13 C	1,4-Dichlorobenzene	1.626	1.580	2.8	89	0.00
14	1,2-Dichlorobenzene	1.637	1.665	-1.7	92	0.00
15	Benzyl Alcohol	1.568	1.802	-14.9	94	0.00
16	2,2'-oxybis(1-Chloropropane	3.117	2.988	4.1	84	0.00
17	2-Methylphenol	1.481	1.616	-9.1	94	0.00
18	Hexachloroethane	0.540	0.532	1.5	91	0.00
19 P	n-Nitroso-di-n-propylamine	1.661	1.885	-13.5	92	0.00
20	3+4-Methylphenols	2.126	2.419	-13.8	95	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22	Acetophenone	0.570	0.603	-5.8	95	0.00
23 S	Nitrobenzene-d5	0.408	0.433	-6.1	95	0.00
24	Nitrobenzene	0.422	0.448	-6.2	95	0.00
25	Isophorone	0.995	1.073	-7.8	94	0.00
26 C	2-Nitrophenol	0.167	0.178	-6.6	92	0.00
27	2,4-Dimethylphenol	0.230	0.244	-6.1	91	0.00
28	bis(2-Chloroethoxy)methane	0.522	0.534	-2.3	90	0.00
29 C	2,4-Dichlorophenol	0.343	0.378	-10.2	94	0.00
30	1,2,4-Trichlorobenzene	0.368	0.377	-2.4	92	0.00
31	Naphthalene	1.065	1.083	-1.7	89	0.00
32	Benzoic acid	0.237	0.285	-20.3	98	0.00
33	4-Chloroaniline	0.460	0.539	-17.2	96	0.00
34 C	Hexachlorobutadiene	0.206	0.209	-1.5	92	0.00
35	Caprolactam	0.179	0.241	-34.6#	108	0.00
36 C	4-Chloro-3-methylphenol	0.410	0.494	-20.5#	99	0.00
37	2-Methylnaphthalene	0.826	0.899	-8.8	94	0.00
38	1-Methylnaphthalene	0.829	0.927	-11.8	97	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
40	1,2,4,5-Tetrachlorobenzene	0.524	0.506	3.4	97	0.00
41 P	Hexachlorocyclopentadiene	0.177	0.164	7.3	92	0.00
42 S	2,4,6-Tribromophenol	0.269	0.311	-15.6	105	0.00
43 C	2,4,6-Trichlorophenol	0.391	0.396	-1.3	96	0.00
44	2,4,5-Trichlorophenol	0.443	0.463	-4.5	98	0.00
45 S	2-Fluorobiphenyl	1.332	1.324	0.6	96	0.00
46	1,1'-Biphenyl	1.453	1.407	3.2	96	0.00
47	2-Chloronaphthalene	1.192	1.164	2.3	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122123\
 Data File : BG060128.D
 Acq On : 21 Dec 2023 18:49
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG122123

Quant Time: Dec 22 03:43:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 03:42:04 2023
 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)
48	2-Nitroaniline	0.389	0.437	-12.3	104	0.00
49	Acenaphthylene	1.863	1.933	-3.8	97	0.00
50	Dimethylphthalate	1.759	1.884	-7.1	101	0.00
51	2,6-Dinitrotoluene	0.352	0.400	-13.6	104	0.00
52 C	Acenaphthene	1.145	1.189	-3.8	98	0.00
53	3-Nitroaniline	0.374	0.444	-18.7	107	0.00
54 P	2,4-Dinitrophenol	0.194	0.233	-20.1	105	0.00
55	Dibenzofuran	1.929	2.028	-5.1	99	0.00
56 P	4-Nitrophenol	0.310	0.385	-24.2	112	0.00
57	2,4-Dinitrotoluene	0.517	0.622	-20.3	107	0.00
58	Fluorene	1.632	1.778	-8.9	102	0.00
59	2,3,4,6-Tetrachlorophenol	0.398	0.449	-12.8	104	0.00
60	Diethylphthalate	1.900	2.135	-12.4	104	0.00
61	4-Chlorophenyl-phenylether	0.778	0.839	-7.8	101	0.00
62	4-Nitroaniline	0.439	0.524	-19.4	106	0.00
63	Azobenzene	1.696	1.818	-7.2	100	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
65	4,6-Dinitro-2-methylphenol	0.108	0.119	-10.2	103	0.00
66 c	n-Nitrosodiphenylamine	0.528	0.529	-0.2	102	0.00
67	4-Bromophenyl-phenylether	0.188	0.187	0.5	103	0.00
68	Hexachlorobenzene	0.216	0.214	0.9	104	0.00
69	Atrazine	0.196	0.176	10.2	107	0.00
70 C	Pentachlorophenol	0.134	0.148	-10.4	106	0.00
71	Phenanthrene	1.008	1.047	-3.9	105	0.00
72	Anthracene	1.014	1.059	-4.4	106	0.00
73	Carbazole	1.077	1.125	-4.5	107	0.00
74	Di-n-butylphthalate	1.277	1.320	-3.4	105	0.00
75 C	Fluoranthene	1.323	1.381	-4.4	107	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	116	0.00
77	Benzidine	0.541	0.476	12.0	113	0.00
78	Pyrene	1.527	1.497	2.0	108	0.00
79 S	Terphenyl-d14	1.105	0.961	13.0	106	0.00
80	Butylbenzylphthalate	0.512	0.569	-11.1	119	0.00
81	Benzo(a)anthracene	1.329	1.362	-2.5	116	0.00
82	3,3'-Dichlorobenzidine	0.394	0.400	-1.5	114	0.00
83	Chrysene	1.248	1.286	-3.0	115	0.00
84	Bis(2-ethylhexyl)phthalate	0.682	0.710	-4.1	116	0.00
85 c	Di-n-octyl phthalate	1.136	1.169	-2.9	113	0.00
86 I	Perylene-d12	1.000	1.000	0.0	115	0.00
87	Indeno(1,2,3-cd)pyrene	1.396	1.447	-3.7	115	0.00
88	Benzo(b)fluoranthene	1.265	1.294	-2.3	114	0.00
89	Benzo(k)fluoranthene	1.239	1.253	-1.1	117	0.00
90 C	Benzo(a)pyrene	1.114	1.131	-1.5	114	0.00
91	Dibenzo(a,h)anthracene	1.153	1.196	-3.7	115	0.00
92	Benzo(g,h,i)perylene	1.165	1.199	-2.9	117	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122123\
Data File : BG060128.D
Acq On : 21 Dec 2023 18:49
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
ICVBG122123

Quant Time: Dec 22 03:43:29 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 22 03:42:04 2023
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)
----------	-------	------	-----------	------------

(#) = Out of Range

SPCC's out = 0 CCC's out = 1