

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101525\
 Data File : BP025980.D
 Acq On : 15 Oct 2025 22:00
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP101525

Quant Time: Oct 15 23:57:42 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP101525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 15 23:48:51 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	124	0.00
2	1,4-Dioxane	0.483	0.503	-4.1	115	0.00
3	Pyridine	1.348	1.494	-10.8	122	0.00
4	n-Nitrosodimethylamine	0.591	0.648	-9.6	122	0.00
5 S	2-Fluorophenol	1.220	1.366	-12.0	123	0.00
6	Aniline	2.279	2.517	-10.4	123	0.00
7 S	Phenol-d6	1.702	1.907	-12.0	124	0.00
8	2-Chlorophenol	1.414	1.558	-10.2	123	0.00
9	Benzaldehyde	0.908	1.759	-93.7#	223#	0.00
10 C	Phenol	1.719	1.960	-14.0	124	0.00
11	bis(2-Chloroethyl)ether	1.401	1.575	-12.4	124	0.00
12	1,3-Dichlorobenzene	1.567	1.701	-8.6	122	0.00
13 C	1,4-Dichlorobenzene	1.597	1.733	-8.5	122	0.00
14	1,2-Dichlorobenzene	1.562	1.700	-8.8	123	0.00
15	Benzyl Alcohol	1.449	1.609	-11.0	129	0.00
16	2,2'-oxybis(1-Chloropropane	2.158	2.338	-8.3	123	0.00
17	2-Methylphenol	1.232	1.364	-10.7	123	0.00
18	Hexachloroethane	0.609	0.652	-7.1	119	0.00
19 P	n-Nitroso-di-n-propylamine	1.214	1.370	-12.9	128	0.00
20	3+4-Methylphenols	1.717	1.869	-8.9	125	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
22	Acetophenone	0.544	0.610	-12.1	126	0.00
23 S	Nitrobenzene-d5	0.445	0.503	-13.0	126	0.00
24	Nitrobenzene	0.396	0.446	-12.6	125	0.00
25	Isophorone	0.799	0.888	-11.1	129	0.00
26 C	2-Nitrophenol	0.192	0.221	-15.1	127	0.00
27	2,4-Dimethylphenol	0.325	0.368	-13.2	126	0.00
28	bis(2-Chloroethoxy)methane	0.467	0.521	-11.6	126	0.00
29 C	2,4-Dichlorophenol	0.324	0.368	-13.6	128	-0.01
30	1,2,4-Trichlorobenzene	0.366	0.403	-10.1	124	0.00
31	Naphthalene	1.079	1.183	-9.6	125	0.00
32	Benzoic acid	0.235	0.271	-15.3	140	0.01
33	4-Chloroaniline	0.458	0.517	-12.9	127	0.00
34 C	Hexachlorobutadiene	0.242	0.270	-11.6	127	0.00
35	Caprolactam	0.115	0.127	-10.4	136	0.00
36 C	4-Chloro-3-methylphenol	0.380	0.431	-13.4	133	0.00
37	2-Methylnaphthalene	0.706	0.791	-12.0	132	0.00
38	1-Methylnaphthalene	0.745	0.810	-8.7	128	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	132	0.00
40	1,2,4,5-Tetrachlorobenzene	0.651	0.726	-11.5	128	0.00
41 P	Hexachlorocyclopentadiene	0.329	0.398	-21.0	139	0.00
42 S	2,4,6-Tribromophenol	0.269	0.312	-16.0	142	0.00
43 C	2,4,6-Trichlorophenol	0.413	0.478	-15.7	133	0.00
44	2,4,5-Trichlorophenol	0.450	0.516	-14.7	132	0.00
45 S	2-Fluorobiphenyl	1.560	1.686	-8.1	126	0.00
46	1,1'-Biphenyl	1.511	1.648	-9.1	128	0.00
47	2-Chloronaphthalene	1.181	1.299	-10.0	129	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.368	0.425	-15.5	135	0.00
49	Acenaphthylene	1.910	2.075	-8.6	129	0.00
50	Dimethylphthalate	1.512	1.677	-10.9	137	0.00
51	2,6-Dinitrotoluene	0.326	0.371	-13.8	136	0.00
52 C	Acenaphthene	1.103	1.198	-8.6	130	0.00
53	3-Nitroaniline	0.325	0.375	-15.4	137	0.00
54 P	2,4-Dinitrophenol	0.190	0.220	-15.8	145	-0.01
55	Dibenzofuran	1.766	1.944	-10.1	132	0.00
56 P	4-Nitrophenol	0.197	0.214	-8.6	170#	0.00
57	2,4-Dinitrotoluene	0.442	0.512	-15.8	140	0.00
58	Fluorene	1.419	1.544	-8.8	133	0.00
59	2,3,4,6-Tetrachlorophenol	0.404	0.449	-11.1	137	0.00
60	Diethylphthalate	1.491	1.629	-9.3	137	0.00
61	4-Chlorophenyl-phenylether	0.746	0.818	-9.7	135	0.00
62	4-Nitroaniline	0.301	0.342	-13.6	137	0.00
63	Azobenzene	1.389	1.526	-9.9	134	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	139	0.00
65	4,6-Dinitro-2-methylphenol	0.145	0.164	-13.1	142	0.00
66 c	n-Nitrosodiphenylamine	0.643	0.712	-10.7	135	0.00
67	4-Bromophenyl-phenylether	0.247	0.275	-11.3	137	0.00
68	Hexachlorobenzene	0.277	0.299	-7.9	134	0.00
69	Atrazine	0.232	0.254	-9.5	137	0.00
70 C	Pentachlorophenol	0.150	0.170	-13.3	145	0.00
71	Phenanthrene	1.134	1.230	-8.5	136	0.00
72	Anthracene	1.158	1.260	-8.8	134	0.00
73	Carbazole	1.021	1.118	-9.5	137	0.00
74	Di-n-butylphthalate	1.256	1.360	-8.3	137	0.00
75 C	Fluoranthene	1.279	1.371	-7.2	135	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	119	-0.02
77	Benzydine	0.641	1.217	-89.9#	205#	-0.01
78	Pyrene	1.475	1.701	-15.3	132	0.00
79 S	Terphenyl-d14	1.251	1.454	-16.2	133	0.00
80	Butylbenzylphthalate	0.587	0.657	-11.9	120	-0.01
81	Benzo(a)anthracene	1.368	1.538	-12.4	120	-0.02
82	3,3'-Dichlorobenzidine	0.487	0.548	-12.5	118	-0.01
83	Chrysene	1.308	1.444	-10.4	119	-0.02
84	Bis(2-ethylhexyl)phthalate	0.810	0.887	-9.5	114	-0.02
85 c	Di-n-octyl phthalate	1.384	1.538	-11.1	111	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	114	-0.02
87	Indeno(1,2,3-cd)pyrene	1.507	1.710	-13.5	113	-0.04
88	Benzo(b)fluoranthene	1.246	1.361	-9.2	112	-0.01
89	Benzo(k)fluoranthene	1.264	1.388	-9.8	113	-0.02
90 C	Benzo(a)pyrene	1.210	1.340	-10.7	113	-0.02
91	Dibenzo(a,h)anthracene	1.228	1.409	-14.7	114	-0.02
92	Benzo(g,h,i)perylene	1.203	1.367	-13.6	113	-0.02

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(#) = Out of Range

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