

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100124\
 Data File : BF139592.D
 Acq On : 01 Oct 2024 17:50
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF100124

Quant Time: Oct 02 05:05:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	99	0.00
2	1,4-Dioxane	40.000	39.145	2.1	100	0.00
3	Pyridine	40.000	39.389	1.5	102	0.00
4	n-Nitrosodimethylamine	40.000	39.212	2.0	98	0.00
5 S	2-Fluorophenol	80.000	77.086	3.6	99	0.00
6	Aniline	40.000	37.668	5.8	101	0.00
7 S	Phenol-d6	80.000	77.200	3.5	99	0.00
8	2-Chlorophenol	40.000	39.042	2.4	100	0.00
9	Benzaldehyde	40.000	41.611	-4.0	108	0.00
10 C	Phenol	40.000	38.908	2.7	98	0.00
11	bis(2-Chloroethyl)ether	40.000	37.003	7.5	97	0.00
12	1,3-Dichlorobenzene	40.000	39.113	2.2	100	0.00
13 C	1,4-Dichlorobenzene	40.000	38.390	4.0	99	0.00
14	1,2-Dichlorobenzene	40.000	38.358	4.1	98	0.00
15	Benzyl Alcohol	40.000	39.165	2.1	98	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.242	1.9	100	0.00
17	2-Methylphenol	40.000	39.320	1.7	99	0.00
18	Hexachloroethane	40.000	39.154	2.1	100	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.536	3.7	99	0.00
20	3+4-Methylphenols	40.000	39.282	1.8	100	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	99	0.00
22	Acetophenone	40.000	38.356	4.1	100	0.00
23 S	Nitrobenzene-d5	80.000	77.670	2.9	100	0.00
24	Nitrobenzene	40.000	39.301	1.7	101	0.00
25	Isophorone	40.000	38.591	3.5	100	0.00
26 C	2-Nitrophenol	40.000	40.302	-0.8	99	0.00
27	2,4-Dimethylphenol	40.000	39.187	2.0	102	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.562	3.6	100	0.00
29 C	2,4-Dichlorophenol	40.000	39.142	2.1	99	0.00
30	1,2,4-Trichlorobenzene	40.000	38.237	4.4	98	0.00
31	Naphthalene	40.000	38.165	4.6	99	0.00
32	Benzoic acid	40.000	42.446	-6.1	101	0.00
33	4-Chloroaniline	40.000	36.080	9.8	91	0.00
34 C	Hexachlorobutadiene	40.000	38.931	2.7	101	0.00
35	Caprolactam	40.000	38.243	4.4	99	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.002	2.5	102	0.00
37	2-Methylnaphthalene	40.000	38.108	4.7	99	0.00
38	1-Methylnaphthalene	40.000	38.135	4.7	99	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	100	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	38.649	3.4	100	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.862	-4.7	97	0.00
42 S	2,4,6-Tribromophenol	80.000	75.239	6.0	97	0.00
43 C	2,4,6-Trichlorophenol	40.000	38.808	3.0	99	0.00
44	2,4,5-Trichlorophenol	40.000	38.800	3.0	100	0.00
45 S	2-Fluorobiphenyl	80.000	75.237	6.0	100	0.00
46	1,1'-Biphenyl	40.000	38.153	4.6	100	0.00
47	2-Chloronaphthalene	40.000	38.613	3.5	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100124\
 Data File : BF139592.D
 Acq On : 01 Oct 2024 17:50
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF100124

Quant Time: Oct 02 05:05:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	38.762	3.1	98	0.00
49	Acenaphthylene	40.000	38.013	5.0	99	0.00
50	Dimethylphthalate	40.000	38.624	3.4	101	0.00
51	2,6-Dinitrotoluene	40.000	39.296	1.8	101	0.00
52 C	Acenaphthene	40.000	38.189	4.5	99	0.00
53	3-Nitroaniline	40.000	38.035	4.9	98	0.00
54 P	2,4-Dinitrophenol	40.000	42.326	-5.8	103	0.00
55	Dibenzofuran	40.000	38.200	4.5	100	0.00
56 P	4-Nitrophenol	40.000	38.896	2.8	98	0.00
57	2,4-Dinitrotoluene	40.000	39.333	1.7	100	0.00
58	Fluorene	40.000	37.321	6.7	99	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	37.849	5.4	98	0.00
60	Diethylphthalate	40.000	38.204	4.5	100	0.00
61	4-Chlorophenyl-phenylether	40.000	37.483	6.3	100	0.00
62	4-Nitroaniline	40.000	37.635	5.9	97	0.00
63	Azobenzene	40.000	38.460	3.8	99	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.324	-3.3	98	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.192	4.5	100	0.00
67	4-Bromophenyl-phenylether	40.000	38.339	4.2	100	0.00
68	Hexachlorobenzene	40.000	38.513	3.7	100	0.00
69	Atrazine	40.000	32.554	18.6	98	0.00
70 C	Pentachlorophenol	40.000	39.355	1.6	95	0.00
71	Phenanthrene	40.000	37.575	6.1	98	0.00
72	Anthracene	40.000	37.863	5.3	99	0.00
73	Carbazole	40.000	37.627	5.9	98	0.00
74	Di-n-butylphthalate	40.000	38.808	3.0	99	0.00
75 C	Fluoranthene	40.000	36.956	7.6	96	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	96	0.00
77	Benzydine	40.000	32.751	18.1	80	0.00
78	Pyrene	40.000	40.566	-1.4	95	0.00
79 S	Terphenyl-d14	80.000	81.327	-1.7	98	0.00
80	Butylbenzylphthalate	40.000	40.748	-1.9	96	0.00
81	Benzo(a)anthracene	40.000	38.038	4.9	95	0.00
82	3,3'-Dichlorobenzidine	40.000	38.860	2.9	94	0.00
83	Chrysene	40.000	38.983	2.5	94	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	41.366	-3.4	97	0.00
85 c	Di-n-octyl phthalate	40.000	42.335	-5.8	100	0.00
86 I	Perylene-d12	20.000	20.000	0.0	100	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	42.613	-6.5	104	0.00
88	Benzo(b)fluoranthene	40.000	36.343	9.1	102	0.00
89	Benzo(k)fluoranthene	40.000	38.929	2.7	95	0.00
90 C	Benzo(a)pyrene	40.000	38.942	2.6	101	0.00
91	Dibenzo(a,h)anthracene	40.000	42.543	-6.4	103	0.00
92	Benzo(g,h,i)perylene	40.000	42.806	-7.0	103	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100124\
Data File : BF139592.D
Acq On : 01 Oct 2024 17:50
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF100124

Quant Time: Oct 02 05:05:20 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 02 04:58:47 2024
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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

(#) = Out of Range

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