

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
 Data File : BF134544.D
 Acq On : 01 Aug 2023 12:50
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF080123

Quant Time: Aug 02 02:22:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 02:20:34 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	96	0.00
2	1,4-Dioxane	40.000	38.148	4.6	91	0.00
3	Pyridine	40.000	38.406	4.0	93	0.00
4	n-Nitrosodimethylamine	40.000	38.673	3.3	92	0.00
5 S	2-Fluorophenol	80.000	75.722	5.3	91	0.00
6	Aniline	40.000	38.117	4.7	92	0.00
7 S	Phenol-d6	80.000	76.467	4.4	93	0.00
8	2-Chlorophenol	40.000	38.929	2.7	93	0.00
9	Benzaldehyde	40.000	38.105	4.7	93	0.00
10 C	Phenol	40.000	38.983	2.5	92	0.00
11	bis(2-Chloroethyl)ether	40.000	38.714	3.2	93	0.00
12	1,3-Dichlorobenzene	40.000	38.142	4.6	92	0.00
13 C	1,4-Dichlorobenzene	40.000	38.755	3.1	93	0.00
14	1,2-Dichlorobenzene	40.000	38.479	3.8	93	0.00
15	Benzyl Alcohol	40.000	39.629	0.9	94	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	38.530	3.7	93	0.00
17	2-Methylphenol	40.000	38.910	2.7	93	0.00
18	Hexachloroethane	40.000	40.128	-0.3	97	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.716	5.7	93	0.00
20	3+4-Methylphenols	40.000	39.066	2.3	93	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	99	0.00
22	Acetophenone	40.000	37.043	7.4	93	0.00
23 S	Nitrobenzene-d5	80.000	84.473	-5.6	98	0.00
24	Nitrobenzene	40.000	40.832	-2.1	96	0.00
25	Isophorone	40.000	37.564	6.1	94	0.00
26 C	2-Nitrophenol	40.000	40.986	-2.5	106	0.00
27	2,4-Dimethylphenol	40.000	38.699	3.3	97	0.00
28	bis(2-Chloroethoxy)methane	40.000	37.995	5.0	95	0.00
29 C	2,4-Dichlorophenol	40.000	38.858	2.9	94	0.00
30	1,2,4-Trichlorobenzene	40.000	38.712	3.2	96	0.00
31	Naphthalene	40.000	37.728	5.7	93	0.00
32	Benzoic acid	40.000	40.424	-1.1	110	0.00
33	4-Chloroaniline	40.000	37.512	6.2	94	0.00
34 C	Hexachlorobutadiene	40.000	38.213	4.5	94	0.00
35	Caprolactam	40.000	38.700	3.2	92	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.337	4.2	94	0.00
37	2-Methylnaphthalene	40.000	37.519	6.2	94	0.00
38	1-Methylnaphthalene	40.000	37.754	5.6	94	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	101	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.878	5.3	95	0.00
41 P	Hexachlorocyclopentadiene	40.000	40.567	-1.4	95	0.00
42 S	2,4,6-Tribromophenol	80.000	79.927	0.1	95	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.802	0.5	98	0.00
44	2,4,5-Trichlorophenol	40.000	39.669	0.8	97	0.00
45 S	2-Fluorobiphenyl	80.000	76.977	3.8	96	0.00
46	1,1'-Biphenyl	40.000	37.830	5.4	95	0.00
47	2-Chloronaphthalene	40.000	37.782	5.5	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
 Data File : BF134544.D
 Acq On : 01 Aug 2023 12:50
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF080123

Quant Time: Aug 02 02:22:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 02:20:34 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	40.053	-0.1	100	0.00
49	Acenaphthylene	40.000	37.708	5.7	95	0.00
50	Dimethylphthalate	40.000	38.079	4.8	96	0.00
51	2,6-Dinitrotoluene	40.000	40.100	-0.3	100	0.00
52 C	Acenaphthene	40.000	37.648	5.9	95	0.00
53	3-Nitroaniline	40.000	43.450	-8.6	98	0.00
54 P	2,4-Dinitrophenol	40.000	41.497	-3.7	115	0.00
55	Dibenzofuran	40.000	37.710	5.7	93	0.00
56 P	4-Nitrophenol	40.000	40.955	-2.4	100	0.00
57	2,4-Dinitrotoluene	40.000	41.313	-3.3	102	0.00
58	Fluorene	40.000	36.552	8.6	95	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.840	0.4	97	0.00
60	Diethylphthalate	40.000	38.232	4.4	95	0.00
61	4-Chlorophenyl-phenylether	40.000	36.774	8.1	95	0.00
62	4-Nitroaniline	40.000	42.842	-7.1	97	0.00
63	Azobenzene	40.000	38.370	4.1	96	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	102	0.00
65	4,6-Dinitro-2-methylphenol	40.000	39.539	1.2	110	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.346	6.6	95	0.00
67	4-Bromophenyl-phenylether	40.000	37.577	6.1	95	0.00
68	Hexachlorobenzene	40.000	38.292	4.3	96	0.00
69	Atrazine	40.000	39.911	0.2	102	0.00
70 C	Pentachlorophenol	40.000	42.156	-5.4	99	0.00
71	Phenanthrene	40.000	37.048	7.4	94	0.00
72	Anthracene	40.000	37.550	6.1	95	0.00
73	Carbazole	40.000	37.490	6.3	95	0.00
74	Di-n-butylphthalate	40.000	39.352	1.6	97	0.00
75 C	Fluoranthene	40.000	37.371	6.6	93	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	99	0.00
77	Benzydine	40.000	38.156	4.6	85	0.00
78	Pyrene	40.000	38.209	4.5	94	0.00
79 S	Terphenyl-d14	80.000	77.255	3.4	96	0.00
80	Butylbenzylphthalate	40.000	41.344	-3.4	95	0.00
81	Benzo(a)anthracene	40.000	37.580	6.1	92	0.00
82	3,3'-Dichlorobenzidine	40.000	37.479	6.3	92	0.00
83	Chrysene	40.000	37.952	5.1	94	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	39.915	0.2	94	0.00
85 c	Di-n-octyl phthalate	40.000	39.252	1.9	92	0.00
86 I	Perylene-d12	20.000	20.000	0.0	97	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	39.633	0.9	90	0.00
88	Benzo(b)fluoranthene	40.000	38.756	3.1	95	0.00
89	Benzo(k)fluoranthene	40.000	38.834	2.9	89	0.00
90 C	Benzo(a)pyrene	40.000	38.454	3.9	91	0.00
91	Dibenzo(a,h)anthracene	40.000	40.305	-0.8	92	0.00
92	Benzo(g,h,i)perylene	40.000	40.140	-0.4	91	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080123\
Data File : BF134544.D
Acq On : 01 Aug 2023 12:50
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF080123

Quant Time: Aug 02 02:22:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 02 02:20:34 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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

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