

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042325\
 Data File : BF142225.D
 Acq On : 23 Apr 2025 14:40
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF042325

Quant Time: Apr 23 14:57:15 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 23 14:53:02 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	98	0.00
2	1,4-Dioxane	40.000	39.289	1.8	99	0.00
3	Pyridine	40.000	39.587	1.0	98	0.00
4	n-Nitrosodimethylamine	40.000	39.811	0.5	99	0.00
5 S	2-Fluorophenol	80.000	77.414	3.2	98	0.00
6	Aniline	40.000	40.178	-0.4	99	0.00
7 S	Phenol-d6	80.000	77.512	3.1	98	0.00
8	2-Chlorophenol	40.000	39.439	1.4	100	0.00
9	Benzaldehyde	40.000	39.919	0.2	102	0.00
10 C	Phenol	40.000	38.852	2.9	97	0.00
11	bis(2-Chloroethyl)ether	40.000	38.358	4.1	97	0.00
12	1,3-Dichlorobenzene	40.000	38.987	2.5	99	0.00
13 C	1,4-Dichlorobenzene	40.000	38.859	2.9	99	0.00
14	1,2-Dichlorobenzene	40.000	38.803	3.0	99	0.00
15	Benzyl Alcohol	40.000	39.539	1.2	97	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	38.197	4.5	97	0.00
17	2-Methylphenol	40.000	38.673	3.3	98	0.00
18	Hexachloroethane	40.000	38.980	2.6	98	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.993	5.0	97	0.00
20	3+4-Methylphenols	40.000	39.159	2.1	99	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	97	0.00
22	Acetophenone	40.000	39.555	1.1	99	0.00
23 S	Nitrobenzene-d5	80.000	79.426	0.7	97	0.00
24	Nitrobenzene	40.000	39.614	1.0	98	0.00
25	Isophorone	40.000	39.514	1.2	99	0.00
26 C	2-Nitrophenol	40.000	40.678	-1.7	99	0.00
27	2,4-Dimethylphenol	40.000	39.879	0.3	99	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.460	1.3	99	0.00
29 C	2,4-Dichlorophenol	40.000	40.332	-0.8	100	0.00
30	1,2,4-Trichlorobenzene	40.000	40.107	-0.3	99	0.00
31	Naphthalene	40.000	39.408	1.5	97	0.00
32	Benzoic acid	40.000	42.451	-6.1	103	0.00
33	4-Chloroaniline	40.000	41.103	-2.8	99	0.00
34 C	Hexachlorobutadiene	40.000	39.641	0.9	99	0.00
35	Caprolactam	40.000	39.934	0.2	98	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.143	-0.4	98	0.00
37	2-Methylnaphthalene	40.000	39.896	0.3	99	0.00
38	1-Methylnaphthalene	40.000	39.654	0.9	98	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.184	2.0	100	0.00
41 P	Hexachlorocyclopentadiene	40.000	38.370	4.1	94	0.00
42 S	2,4,6-Tribromophenol	80.000	79.901	0.1	97	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.697	0.8	97	0.00
44	2,4,5-Trichlorophenol	40.000	40.667	-1.7	102	0.00
45 S	2-Fluorobiphenyl	80.000	77.396	3.3	99	0.00
46	1,1'-Biphenyl	40.000	39.042	2.4	98	0.00
47	2-Chloronaphthalene	40.000	38.906	2.7	99	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042325\
 Data File : BF142225.D
 Acq On : 23 Apr 2025 14:40
 Operator : RC/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF042325

Quant Time: Apr 23 14:57:15 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 23 14:53:02 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	40.011	-0.0	98	0.00
49	Acenaphthylene	40.000	39.277	1.8	99	0.00
50	Dimethylphthalate	40.000	39.175	2.1	99	0.00
51	2,6-Dinitrotoluene	40.000	40.033	-0.1	99	0.00
52 C	Acenaphthene	40.000	39.694	0.8	99	0.00
53	3-Nitroaniline	40.000	39.499	1.3	98	0.00
54 P	2,4-Dinitrophenol	40.000	41.360	-3.4	97	0.00
55	Dibenzofuran	40.000	38.948	2.6	99	0.00
56 P	4-Nitrophenol	40.000	41.515	-3.8	97	0.00
57	2,4-Dinitrotoluene	40.000	40.019	-0.0	99	0.00
58	Fluorene	40.000	38.544	3.6	98	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.591	-1.5	99	0.00
60	Diethylphthalate	40.000	39.155	2.1	98	0.00
61	4-Chlorophenyl-phenylether	40.000	38.967	2.6	98	0.00
62	4-Nitroaniline	40.000	39.808	0.5	97	0.00
63	Azobenzene	40.000	39.183	2.0	98	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	97	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.809	-2.0	96	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.545	1.1	98	0.00
67	4-Bromophenyl-phenylether	40.000	40.173	-0.4	99	0.00
68	Hexachlorobenzene	40.000	40.223	-0.6	100	0.00
69	Atrazine	40.000	33.914	15.2	97	0.00
70 C	Pentachlorophenol	40.000	42.077	-5.2	98	0.00
71	Phenanthrene	40.000	39.272	1.8	98	0.00
72	Anthracene	40.000	39.408	1.5	98	0.00
73	Carbazole	40.000	38.956	2.6	96	0.00
74	Di-n-butylphthalate	40.000	39.325	1.7	96	0.00
75 C	Fluoranthene	40.000	38.044	4.9	95	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	92	0.00
77	Benzydine	40.000	42.821	-7.1	88	0.00
78	Pyrene	40.000	40.944	-2.4	94	0.00
79 S	Terphenyl-d14	80.000	81.750	-2.2	95	0.00
80	Butylbenzylphthalate	40.000	40.646	-1.6	92	0.00
81	Benzo(a)anthracene	40.000	38.927	2.7	94	0.00
82	3,3'-Dichlorobenzidine	40.000	38.740	3.1	91	0.00
83	Chrysene	40.000	40.389	-1.0	93	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	41.700	-4.3	95	0.00
85 c	Di-n-octyl phthalate	40.000	41.719	-4.3	97	0.00
86 I	Perylene-d12	20.000	20.000	0.0	99	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	39.392	1.5	101	0.00
88	Benzo(b)fluoranthene	40.000	39.285	1.8	101	0.00
89	Benzo(k)fluoranthene	40.000	38.757	3.1	96	0.00
90 C	Benzo(a)pyrene	40.000	39.548	1.1	99	0.00
91	Dibenzo(a,h)anthracene	40.000	39.582	1.0	100	0.00
92	Benzo(g,h,i)perylene	40.000	38.663	3.3	98	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF042325\
Data File : BF142225.D
Acq On : 23 Apr 2025 14:40
Operator : RC/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF042325

Quant Time: Apr 23 14:57:15 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF042325.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 23 14:53:02 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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

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