

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083122\
 Data File : BF130057.D
 Acq On : 31 Aug 2022 17:49
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF083122

Quant Time: Aug 31 18:18:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 31 18:15:32 2022
 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	95	0.00
2	1,4-Dioxane	40.000	37.811	5.5	93	0.00
3	Pyridine	40.000	38.842	2.9	95	-0.01
4	n-Nitrosodimethylamine	40.000	38.753	3.1	93	-0.01
5 S	2-Fluorophenol	80.000	77.195	3.5	97	0.00
6	Aniline	40.000	38.916	2.7	94	0.00
7 S	Phenol-d6	80.000	77.567	3.0	97	0.00
8	2-Chlorophenol	40.000	39.015	2.5	96	0.00
9	Benzaldehyde	40.000	37.167	7.1	101	0.00
10 C	Phenol	40.000	39.018	2.5	97	0.00
11	bis(2-Chloroethyl)ether	40.000	38.307	4.2	94	0.00
12	1,3-Dichlorobenzene	40.000	38.416	4.0	96	0.00
13 C	1,4-Dichlorobenzene	40.000	38.233	4.4	95	0.00
14	1,2-Dichlorobenzene	40.000	38.124	4.7	97	0.00
15	Benzyl Alcohol	40.000	39.455	1.4	97	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	38.557	3.6	94	0.00
17	2-Methylphenol	40.000	38.970	2.6	96	0.00
18	Hexachloroethane	40.000	40.033	-0.1	96	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.489	3.8	96	0.00
20	3+4-Methylphenols	40.000	37.943	5.1	97	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	94	0.00
22	Acetophenone	40.000	37.367	6.6	95	0.00
23 S	Nitrobenzene-d5	80.000	83.973	-5.0	99	0.00
24	Nitrobenzene	40.000	41.033	-2.6	98	0.00
25	Isophorone	40.000	38.638	3.4	95	0.00
26 C	2-Nitrophenol	40.000	42.551	-6.4	105	0.00
27	2,4-Dimethylphenol	40.000	39.327	1.7	97	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.293	4.3	95	0.00
29 C	2,4-Dichlorophenol	40.000	39.577	1.1	96	0.00
30	1,2,4-Trichlorobenzene	40.000	37.987	5.0	96	0.00
31	Naphthalene	40.000	37.471	6.3	94	0.00
32	Benzoic acid	40.000	41.861	-4.7	104	0.00
33	4-Chloroaniline	40.000	38.114	4.7	94	0.00
34 C	Hexachlorobutadiene	40.000	38.397	4.0	95	0.00
35	Caprolactam	40.000	39.831	0.4	96	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.071	2.3	96	0.00
37	2-Methylnaphthalene	40.000	37.812	5.5	97	0.00
38	1-Methylnaphthalene	40.000	37.679	5.8	96	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	94	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.784	5.5	95	0.00
41 P	Hexachlorocyclopentadiene	40.000	39.604	1.0	92	0.00
42 S	2,4,6-Tribromophenol	80.000	81.072	-1.3	97	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.100	-0.3	96	0.00
44	2,4,5-Trichlorophenol	40.000	39.388	1.5	95	0.00
45 S	2-Fluorobiphenyl	80.000	71.976	10.0	96	0.00
46	1,1'-Biphenyl	40.000	36.996	7.5	95	0.00
47	2-Chloronaphthalene	40.000	37.672	5.8	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083122\
 Data File : BF130057.D
 Acq On : 31 Aug 2022 17:49
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF083122

Quant Time: Aug 31 18:18:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 31 18:15:32 2022
 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	45.056	-12.6	100	0.00
49	Acenaphthylene	40.000	37.810	5.5	95	0.00
50	Dimethylphthalate	40.000	37.926	5.2	96	0.00
51	2,6-Dinitrotoluene	40.000	42.479	-6.2	98	0.00
52 C	Acenaphthene	40.000	37.932	5.2	95	0.00
53	3-Nitroaniline	40.000	43.168	-7.9	102	0.00
54 P	2,4-Dinitrophenol	40.000	41.229	-3.1	107	0.00
55	Dibenzofuran	40.000	38.057	4.9	96	0.00
56 P	4-Nitrophenol	40.000	42.631	-6.6	99	0.00
57	2,4-Dinitrotoluene	40.000	45.284	-13.2	100	0.00
58	Fluorene	40.000	36.416	9.0	96	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.156	2.1	95	0.00
60	Diethylphthalate	40.000	39.132	2.2	97	0.00
61	4-Chlorophenyl-phenylether	40.000	36.685	8.3	96	0.00
62	4-Nitroaniline	40.000	42.105	-5.3	97	0.00
63	Azobenzene	40.000	37.591	6.0	94	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	94	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.662	-4.2	104	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.971	5.1	96	0.00
67	4-Bromophenyl-phenylether	40.000	38.348	4.1	95	0.00
68	Hexachlorobenzene	40.000	38.419	4.0	95	0.00
69	Atrazine	40.000	39.579	1.1	96	0.00
70 C	Pentachlorophenol	40.000	41.577	-3.9	96	0.00
71	Phenanthrene	40.000	37.001	7.5	96	0.00
72	Anthracene	40.000	37.033	7.4	95	0.00
73	Carbazole	40.000	37.818	5.5	96	0.00
74	Di-n-butylphthalate	40.000	39.418	1.5	97	0.00
75 C	Fluoranthene	40.000	37.093	7.3	96	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	93	0.00
77	Benzydine	40.000	41.157	-2.9	95	0.00
78	Pyrene	40.000	39.641	0.9	95	0.00
79 S	Terphenyl-d14	80.000	79.562	0.5	96	0.00
80	Butylbenzylphthalate	40.000	44.351	-10.9	98	0.00
81	Benzo(a)anthracene	40.000	38.212	4.5	94	0.00
82	3,3'-Dichlorobenzidine	40.000	40.093	-0.2	94	0.00
83	Chrysene	40.000	38.695	3.3	96	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	41.788	-4.5	95	0.00
85 c	Di-n-octyl phthalate	40.000	43.874	-9.7	98	0.00
86 I	Perylene-d12	20.000	20.000	0.0	93	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	41.104	-2.8	93	0.00
88	Benzo(b)fluoranthene	40.000	38.832	2.9	99	0.00
89	Benzo(k)fluoranthene	40.000	38.063	4.8	90	0.00
90 C	Benzo(a)pyrene	40.000	39.290	1.8	94	0.00
91	Dibenzo(a,h)anthracene	40.000	41.145	-2.9	93	0.00
92	Benzo(g,h,i)perylene	40.000	41.569	-3.9	94	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF083122\
Data File : BF130057.D
Acq On : 31 Aug 2022 17:49
Operator : CG\JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF083122

Quant Time: Aug 31 18:18:58 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF083122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 18:15:32 2022
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