

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112122\
 Data File : BP012689.D
 Acq On : 21 Nov 2022 16:49
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP112122

Quant Time: Nov 21 23:59:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 23:51:17 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	113	0.00
2	1,4-Dioxane	40.000	37.476	6.3	111	0.00
3	Pyridine	40.000	38.597	3.5	111	0.00
4	n-Nitrosodimethylamine	40.000	37.586	6.0	108	0.00
5 S	2-Fluorophenol	80.000	79.671	0.4	114	0.00
6	Aniline	40.000	40.674	-1.7	114	0.00
7 S	Phenol-d6	80.000	80.808	-1.0	114	0.00
8	2-Chlorophenol	40.000	40.436	-1.1	114	0.00
9	Benzaldehyde	40.000	35.443	11.4	114	0.00
10 C	Phenol	40.000	40.320	-0.8	114	0.00
11	bis(2-Chloroethyl)ether	40.000	38.512	3.7	111	0.00
12	1,3-Dichlorobenzene	40.000	38.594	3.5	112	0.00
13 C	1,4-Dichlorobenzene	40.000	38.341	4.1	112	0.00
14	1,2-Dichlorobenzene	40.000	38.387	4.0	112	0.00
15	Benzyl Alcohol	40.000	42.826	-7.1	117	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	37.962	5.1	109	0.00
17	2-Methylphenol	40.000	41.449	-3.6	115	0.00
18	Hexachloroethane	40.000	37.918	5.2	111	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	42.188	-5.5	113	0.00
20	3+4-Methylphenols	40.000	41.910	-4.8	115	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	111	0.00
22	Acetophenone	40.000	38.859	2.9	112	0.00
23 S	Nitrobenzene-d5	80.000	79.397	0.8	111	0.00
24	Nitrobenzene	40.000	39.169	2.1	110	0.00
25	Isophorone	40.000	39.020	2.4	112	0.00
26 C	2-Nitrophenol	40.000	39.469	1.3	125	0.00
27	2,4-Dimethylphenol	40.000	42.514	-6.3	116	0.00
28	bis(2-Chloroethoxy)methane	40.000	40.084	-0.2	112	0.00
29 C	2,4-Dichlorophenol	40.000	38.829	2.9	114	0.00
30	1,2,4-Trichlorobenzene	40.000	38.938	2.7	113	0.00
31	Naphthalene	40.000	38.459	3.9	111	0.00
32	Benzoic acid	40.000	40.611	-1.5	127	0.00
33	4-Chloroaniline	40.000	38.836	2.9	114	0.00
34 C	Hexachlorobutadiene	40.000	38.874	2.8	113	0.00
35	Caprolactam	40.000	40.064	-0.2	120	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.163	-2.9	112	0.00
37	2-Methylnaphthalene	40.000	38.891	2.8	110	0.00
38	1-Methylnaphthalene	40.000	38.552	3.6	111	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	110	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.687	0.8	112	0.00
41 P	Hexachlorocyclopentadiene	40.000	39.837	0.4	108	0.00
42 S	2,4,6-Tribromophenol	80.000	81.199	-1.5	126	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.214	2.0	114	0.00
44	2,4,5-Trichlorophenol	40.000	42.149	-5.4	113	0.00
45 S	2-Fluorobiphenyl	80.000	77.834	2.7	109	0.00
46	1,1'-Biphenyl	40.000	39.155	2.1	110	0.00
47	2-Chloronaphthalene	40.000	39.129	2.2	110	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112122\
 Data File : BP012689.D
 Acq On : 21 Nov 2022 16:49
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP112122

Quant Time: Nov 21 23:59:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 23:51:17 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	38.605	3.5	115	0.00
49	Acenaphthylene	40.000	38.881	2.8	109	0.00
50	Dimethylphthalate	40.000	39.675	0.8	111	0.00
51	2,6-Dinitrotoluene	40.000	41.459	-3.6	111	0.00
52 C	Acenaphthene	40.000	39.145	2.1	110	0.00
53	3-Nitroaniline	40.000	38.368	4.1	112	0.00
54 P	2,4-Dinitrophenol	40.000	40.353	-0.9	121	0.00
55	Dibenzofuran	40.000	38.762	3.1	110	0.00
56 P	4-Nitrophenol	40.000	38.846	2.9	114	0.00
57	2,4-Dinitrotoluene	40.000	38.229	4.4	111	0.00
58	Fluorene	40.000	38.501	3.7	109	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	42.134	-5.3	113	0.00
60	Diethylphthalate	40.000	40.818	-2.0	112	0.00
61	4-Chlorophenyl-phenylether	40.000	39.194	2.0	112	0.00
62	4-Nitroaniline	40.000	38.670	3.3	113	0.00
63	Azobenzene	40.000	38.608	3.5	109	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	109	0.00
65	4,6-Dinitro-2-methylphenol	40.000	38.653	3.4	119	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.732	0.7	110	0.00
67	4-Bromophenyl-phenylether	40.000	41.530	-3.8	120	0.00
68	Hexachlorobenzene	40.000	39.309	1.7	112	0.00
69	Atrazine	40.000	43.844	-9.6	116	0.00
70 C	Pentachlorophenol	40.000	42.071	-5.2	115	0.00
71	Phenanthrene	40.000	38.657	3.4	108	0.00
72	Anthracene	40.000	39.263	1.8	109	0.00
73	Carbazole	40.000	40.043	-0.1	110	0.00
74	Di-n-butylphthalate	40.000	40.116	-0.3	118	0.00
75 C	Fluoranthene	40.000	39.331	1.7	108	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	107	0.00
77	Benzydine	40.000	44.151	-10.4	147	0.00
78	Pyrene	40.000	39.334	1.7	108	0.00
79 S	Terphenyl-d14	80.000	80.271	-0.3	107	0.00
80	Butylbenzylphthalate	40.000	43.998	-10.0	136	0.00
81	Benzo(a)anthracene	40.000	39.410	1.5	107	0.00
82	3,3'-Dichlorobenzidine	40.000	39.231	1.9	119	0.00
83	Chrysene	40.000	38.939	2.7	107	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	43.902	-9.8	134	0.00
85 c	Di-n-octyl phthalate	40.000	42.549	-6.4	134	0.00
86 I	Perylene-d12	20.000	20.000	0.0	106	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.927	2.7	100	0.01
88	Benzo(b)fluoranthene	40.000	40.683	-1.7	107	0.00
89	Benzo(k)fluoranthene	40.000	39.102	2.2	106	0.00
90 C	Benzo(a)pyrene	40.000	40.257	-0.6	106	0.00
91	Dibenzo(a,h)anthracene	40.000	39.053	2.4	101	0.01
92	Benzo(g,h,i)perylene	40.000	38.095	4.8	99	0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112122\
Data File : BP012689.D
Acq On : 21 Nov 2022 16:49
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP112122

Quant Time: Nov 21 23:59:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP112122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 21 23:51:17 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