

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100121\
 Data File : BM032304.D
 Acq On : 30 Sep 2021 20:41
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM100121

Quant Time: Oct 01 10:41:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 10:39:43 2021
 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	80	0.00
2	1,4-Dioxane	40.000	36.846	7.9	78	0.00
3	Pyridine	40.000	37.553	6.1	77	0.00
4	n-Nitrosodimethylamine	40.000	36.558	8.6	75	0.00
5 S	2-Fluorophenol	80.000	76.765	4.0	79	0.00
6	Aniline	40.000	38.758	3.1	78	0.00
7 S	Phenol-d6	80.000	77.171	3.5	78	0.00
8	2-Chlorophenol	40.000	39.076	2.3	80	0.00
9	Benzaldehyde	40.000	36.678	8.3	80	0.00
10 C	Phenol	40.000	38.595	3.5	78	0.00
11	bis(2-Chloroethyl)ether	40.000	37.900	5.3	78	0.00
12	1,3-Dichlorobenzene	40.000	38.326	4.2	80	0.00
13 C	1,4-Dichlorobenzene	40.000	38.357	4.1	80	0.00
14	1,2-Dichlorobenzene	40.000	38.375	4.1	80	0.00
15	Benzyl Alcohol	40.000	40.814	-2.0	81	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.303	9.2	74	0.00
17	2-Methylphenol	40.000	40.056	-0.1	80	0.00
18	Hexachloroethane	40.000	39.946	0.1	80	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	39.985	0.0	78	0.00
20	3+4-Methylphenols	40.000	40.309	-0.8	80	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22	Acetophenone	40.000	38.311	4.2	79	0.00
23 S	Nitrobenzene-d5	80.000	78.650	1.7	79	0.00
24	Nitrobenzene	40.000	38.898	2.8	79	0.00
25	Isophorone	40.000	40.816	-2.0	80	0.00
26 C	2-Nitrophenol	40.000	43.499	-8.7	84	0.00
27	2,4-Dimethylphenol	40.000	39.860	0.4	80	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.157	4.6	78	0.00
29 C	2,4-Dichlorophenol	40.000	40.631	-1.6	82	0.00
30	1,2,4-Trichlorobenzene	40.000	39.184	2.0	83	0.00
31	Naphthalene	40.000	38.391	4.0	80	0.00
32	Benzoic acid	40.000	41.861	-4.7	86	0.00
33	4-Chloroaniline	40.000	39.600	1.0	80	0.00
34 C	Hexachlorobutadiene	40.000	39.575	1.1	84	0.00
35	Caprolactam	40.000	41.036	-2.6	83	0.00
36 C	4-Chloro-3-methylphenol	40.000	41.086	-2.7	83	0.00
37	2-Methylnaphthalene	40.000	39.312	1.7	82	0.00
38	1-Methylnaphthalene	40.000	39.234	1.9	81	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	84	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.793	5.5	83	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.473	-3.7	84	0.00
42 S	2,4,6-Tribromophenol	80.000	85.424	-6.8	89	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.426	-1.1	85	0.00
44	2,4,5-Trichlorophenol	40.000	40.184	-0.5	85	0.00
45 S	2-Fluorobiphenyl	80.000	77.237	3.5	83	0.00
46	1,1'-Biphenyl	40.000	37.962	5.1	83	0.00
47	2-Chloronaphthalene	40.000	38.145	4.6	83	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100121\
 Data File : BM032304.D
 Acq On : 30 Sep 2021 20:41
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICBM100121

Quant Time: Oct 01 10:41:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 01 10:39:43 2021
 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	42.548	-6.4	83	0.00
49	Acenaphthylene	40.000	39.675	0.8	83	0.00
50	Dimethylphthalate	40.000	39.844	0.4	85	0.00
51	2,6-Dinitrotoluene	40.000	42.597	-6.5	86	0.00
52 C	Acenaphthene	40.000	39.083	2.3	84	0.00
53	3-Nitroaniline	40.000	42.843	-7.1	85	0.00
54 P	2,4-Dinitrophenol	40.000	41.096	-2.7	87	0.00
55	Dibenzofuran	40.000	38.556	3.6	84	0.00
56 P	4-Nitrophenol	40.000	43.428	-8.6	87	0.00
57	2,4-Dinitrotoluene	40.000	44.838	-12.1	88	0.00
58	Fluorene	40.000	37.662	5.8	83	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.957	-2.4	86	0.00
60	Diethylphthalate	40.000	40.621	-1.6	86	0.00
61	4-Chlorophenyl-phenylether	40.000	38.440	3.9	85	0.00
62	4-Nitroaniline	40.000	43.384	-8.5	85	0.00
63	Azobenzene	40.000	39.762	0.6	81	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	87	0.00
65	4,6-Dinitro-2-methylphenol	40.000	42.650	-6.6	88	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.850	5.4	84	0.00
67	4-Bromophenyl-phenylether	40.000	39.148	2.1	87	0.00
68	Hexachlorobenzene	40.000	39.096	2.3	89	0.00
69	Atrazine	40.000	40.606	-1.5	89	0.00
70 C	Pentachlorophenol	40.000	43.267	-8.2	90	0.00
71	Phenanthrene	40.000	37.664	5.8	85	0.00
72	Anthracene	40.000	38.821	2.9	86	0.00
73	Carbazole	40.000	38.629	3.4	85	0.00
74	Di-n-butylphthalate	40.000	42.179	-5.4	89	0.00
75 C	Fluoranthene	40.000	38.949	2.6	87	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	87	0.00
77	Benzydine	40.000	39.946	0.1	86	0.00
78	Pyrene	40.000	39.041	2.4	87	0.00
79 S	Terphenyl-d14	80.000	82.495	-3.1	90	0.00
80	Butylbenzylphthalate	40.000	44.806	-12.0	92	0.00
81	Benzo(a)anthracene	40.000	39.122	2.2	88	0.00
82	3,3'-Dichlorobenzidine	40.000	41.371	-3.4	88	0.00
83	Chrysene	40.000	38.458	3.9	88	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	44.234	-10.6	90	0.00
85 c	Di-n-octyl phthalate	40.000	45.660	-14.1	90	0.00
86 I	Perylene-d12	20.000	20.000	0.0	89	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.270	4.3	88	0.00
88	Benzo(b)fluoranthene	40.000	40.081	-0.2	93	0.00
89	Benzo(k)fluoranthene	40.000	35.981	10.0	84	0.00
90 C	Benzo(a)pyrene	40.000	39.562	1.1	89	0.00
91	Dibenzo(a,h)anthracene	40.000	38.238	4.4	88	0.00
92	Benzo(g,h,i)perylene	40.000	38.644	3.4	87	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100121\
Data File : BM032304.D
Acq On : 30 Sep 2021 20:41
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM100121

Quant Time: Oct 01 10:41:47 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 01 10:39:43 2021
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