

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080122\
 Data File : BM036035.D
 Acq On : 01 Aug 2022 17:29
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM080122

Quant Time: Aug 01 18:53:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 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	80	0.00
2	1,4-Dioxane	40.000	36.591	8.5	82	0.00
3	Pyridine	40.000	36.693	8.3	84	0.00
4	n-Nitrosodimethylamine	40.000	36.343	9.1	82	0.00
5 S	2-Fluorophenol	80.000	73.869	7.7	82	0.00
6	Aniline	40.000	36.919	7.7	82	0.00
7 S	Phenol-d6	80.000	73.552	8.1	82	0.00
8	2-Chlorophenol	40.000	36.885	7.8	81	0.00
9	Benzaldehyde	40.000	33.285	16.8	83	0.00
10 C	Phenol	40.000	36.913	7.7	82	0.00
11	bis(2-Chloroethyl)ether	40.000	36.276	9.3	79	0.00
12	1,3-Dichlorobenzene	40.000	36.658	8.4	79	0.00
13 C	1,4-Dichlorobenzene	40.000	36.580	8.6	79	0.00
14	1,2-Dichlorobenzene	40.000	36.865	7.8	80	0.00
15	Benzyl Alcohol	40.000	36.720	8.2	81	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.144	9.6	77	0.00
17	2-Methylphenol	40.000	36.904	7.7	82	0.00
18	Hexachloroethane	40.000	36.892	7.8	78	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.088	7.3	81	0.00
20	3+4-Methylphenols	40.000	36.919	7.7	82	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22	Acetophenone	40.000	35.247	11.9	80	0.00
23 S	Nitrobenzene-d5	80.000	71.134	11.1	81	0.00
24	Nitrobenzene	40.000	35.634	10.9	81	0.00
25	Isophorone	40.000	35.186	12.0	79	0.00
26 C	2-Nitrophenol	40.000	36.306	9.2	81	0.00
27	2,4-Dimethylphenol	40.000	35.492	11.3	81	0.00
28	bis(2-Chloroethoxy)methane	40.000	34.743	13.1	78	0.00
29 C	2,4-Dichlorophenol	40.000	35.920	10.2	81	0.00
30	1,2,4-Trichlorobenzene	40.000	35.190	12.0	79	0.00
31	Naphthalene	40.000	35.002	12.5	80	0.00
32	Benzoic acid	40.000	36.940	7.7	84	0.00
33	4-Chloroaniline	40.000	35.543	11.1	82	0.00
34 C	Hexachlorobutadiene	40.000	35.432	11.4	77	0.00
35	Caprolactam	40.000	37.164	7.1	89	0.00
36 C	4-Chloro-3-methylphenol	40.000	35.727	10.7	83	0.00
37	2-Methylnaphthalene	40.000	35.217	12.0	80	0.00
38	1-Methylnaphthalene	40.000	35.157	12.1	81	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	82	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	34.557	13.6	79	0.00
41 P	Hexachlorocyclopentadiene	40.000	35.397	11.5	77	0.00
42 S	2,4,6-Tribromophenol	80.000	72.291	9.6	85	0.00
43 C	2,4,6-Trichlorophenol	40.000	35.470	11.3	81	0.00
44	2,4,5-Trichlorophenol	40.000	35.715	10.7	82	0.00
45 S	2-Fluorobiphenyl	80.000	69.191	13.5	80	0.00
46	1,1'-Biphenyl	40.000	34.451	13.9	80	0.00
47	2-Chloronaphthalene	40.000	34.478	13.8	81	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080122\
 Data File : BM036035.D
 Acq On : 01 Aug 2022 17:29
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM080122

Quant Time: Aug 01 18:53:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 17:50:14 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	36.222	9.4	87	0.00
49	Acenaphthylene	40.000	35.193	12.0	83	0.00
50	Dimethylphthalate	40.000	34.839	12.9	83	0.00
51	2,6-Dinitrotoluene	40.000	36.322	9.2	84	0.00
52 C	Acenaphthene	40.000	34.946	12.6	83	0.00
53	3-Nitroaniline	40.000	36.685	8.3	88	0.00
54 P	2,4-Dinitrophenol	40.000	36.949	7.6	88	0.00
55	Dibenzofuran	40.000	34.837	12.9	83	0.00
56 P	4-Nitrophenol	40.000	37.490	6.3	91	0.00
57	2,4-Dinitrotoluene	40.000	36.933	7.7	87	0.00
58	Fluorene	40.000	34.807	13.0	84	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	35.718	10.7	85	0.00
60	Diethylphthalate	40.000	35.155	12.1	83	0.00
61	4-Chlorophenyl-phenylether	40.000	34.852	12.9	82	0.00
62	4-Nitroaniline	40.000	37.751	5.6	92	0.00
63	Azobenzene	40.000	35.401	11.5	82	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	87	0.00
65	4,6-Dinitro-2-methylphenol	40.000	36.681	8.3	87	0.00
66 c	n-Nitrosodiphenylamine	40.000	34.506	13.7	84	0.00
67	4-Bromophenyl-phenylether	40.000	34.444	13.9	84	0.00
68	Hexachlorobenzene	40.000	34.593	13.5	84	0.00
69	Atrazine	40.000	38.134	4.7	87	0.00
70 C	Pentachlorophenol	40.000	36.305	9.2	88	0.00
71	Phenanthrene	40.000	34.718	13.2	87	0.00
72	Anthracene	40.000	35.248	11.9	87	0.00
73	Carbazole	40.000	35.817	10.5	91	0.00
74	Di-n-butylphthalate	40.000	35.773	10.6	85	0.00
75 C	Fluoranthene	40.000	35.891	10.3	91	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	93	0.00
77	Benzydine	40.000	39.406	1.5	95	0.00
78	Pyrene	40.000	34.104	14.7	92	0.00
79 S	Terphenyl-d14	80.000	68.993	13.8	89	0.00
80	Butylbenzylphthalate	40.000	35.222	11.9	90	0.00
81	Benzo(a)anthracene	40.000	34.887	12.8	93	0.00
82	3,3'-Dichlorobenzidine	40.000	35.898	10.3	94	0.00
83	Chrysene	40.000	35.201	12.0	94	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	35.065	12.3	87	0.00
85 c	Di-n-octyl phthalate	40.000	35.861	10.3	88	0.00
86 I	Perylene-d12	20.000	20.000	0.0	94	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	34.899	12.8	94	0.00
88	Benzo(b)fluoranthene	40.000	34.949	12.6	93	0.00
89	Benzo(k)fluoranthene	40.000	34.821	12.9	93	0.00
90 C	Benzo(a)pyrene	40.000	35.039	12.4	94	0.00
91	Dibenzo(a,h)anthracene	40.000	35.098	12.3	94	0.00
92	Benzo(g,h,i)perylene	40.000	35.020	12.4	94	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080122\
Data File : BM036035.D
Acq On : 01 Aug 2022 17:29
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ICVBM080122

Quant Time: Aug 01 18:53:59 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:50:14 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