

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081723\
 Data File : BM041429.D
 Acq On : 17 Aug 2023 09:32
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 17 17:53:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 02:13:45 2023
 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	97	0.00
2	1,4-Dioxane	40.000	41.390	-3.5	106	0.00
3	Pyridine	40.000	41.442	-3.6	102	0.00
4	n-Nitrosodimethylamine	40.000	40.486	-1.2	102	0.00
5 S	2-Fluorophenol	80.000	83.767	-4.7	104	0.00
6	Aniline	40.000	40.198	-0.5	99	0.00
7 S	Phenol-d6	80.000	81.147	-1.4	101	0.00
8	2-Chlorophenol	40.000	40.859	-2.1	103	0.00
9	Benzaldehyde	40.000	37.539	6.2	100	0.00
10 C	Phenol	40.000	40.607	-1.5	101	0.00
11	bis(2-Chloroethyl)ether	40.000	40.459	-1.1	103	0.00
12	1,3-Dichlorobenzene	40.000	40.911	-2.3	104	0.00
13 C	1,4-Dichlorobenzene	40.000	41.120	-2.8	104	0.00
14	1,2-Dichlorobenzene	40.000	40.676	-1.7	102	0.00
15	Benzyl Alcohol	40.000	40.692	-1.7	97	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.345	1.6	102	-0.02
17	2-Methylphenol	40.000	40.385	-1.0	101	0.00
18	Hexachloroethane	40.000	40.823	-2.1	104	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.453	3.9	98	0.00
20	3+4-Methylphenols	40.000	40.083	-0.2	101	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	93	0.00
22	Acetophenone	40.000	40.650	-1.6	101	0.00
23 S	Nitrobenzene-d5	80.000	81.357	-1.7	100	0.00
24	Nitrobenzene	40.000	41.269	-3.2	101	0.00
25	Isophorone	40.000	39.231	1.9	93	0.00
26 C	2-Nitrophenol	40.000	41.930	-4.8	99	0.00
27	2,4-Dimethylphenol	40.000	40.214	-0.5	96	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.626	0.9	96	0.00
29 C	2,4-Dichlorophenol	40.000	41.117	-2.8	99	0.00
30	1,2,4-Trichlorobenzene	40.000	41.636	-4.1	102	0.00
31	Naphthalene	40.000	40.276	-0.7	98	0.00
32	Benzoic acid	40.000	38.342	4.1	89	0.00
33	4-Chloroaniline	40.000	40.487	-1.2	96	0.00
34 C	Hexachlorobutadiene	40.000	41.670	-4.2	103	0.00
35	Caprolactam	40.000	36.396	9.0	86	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.148	2.1	93	0.00
37	2-Methylnaphthalene	40.000	39.597	1.0	96	0.00
38	1-Methylnaphthalene	40.000	39.228	1.9	95	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	88	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	42.822	-7.1	98	0.00
41 P	Hexachlorocyclopentadiene	40.000	41.275	-3.2	98	0.00
42 S	2,4,6-Tribromophenol	80.000	77.350	3.3	90	0.00
43 C	2,4,6-Trichlorophenol	40.000	42.464	-6.2	96	0.00
44	2,4,5-Trichlorophenol	40.000	42.460	-6.2	96	0.00
45 S	2-Fluorobiphenyl	80.000	84.377	-5.5	96	0.00
46	1,1'-Biphenyl	40.000	41.368	-3.4	94	0.00
47	2-Chloronaphthalene	40.000	41.819	-4.5	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081723\
 Data File : BM041429.D
 Acq On : 17 Aug 2023 09:32
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 17 17:53:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 17 02:13:45 2023
 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	40.411	-1.0	87	0.00
49	Acenaphthylene	40.000	40.849	-2.1	92	0.00
50	Dimethylphthalate	40.000	39.404	1.5	90	0.00
51	2,6-Dinitrotoluene	40.000	39.942	0.1	90	0.00
52 C	Acenaphthene	40.000	40.675	-1.7	93	0.00
53	3-Nitroaniline	40.000	40.407	-1.0	89	0.00
54 P	2,4-Dinitrophenol	40.000	38.716	3.2	87	0.00
55	Dibenzofuran	40.000	40.384	-1.0	93	0.00
56 P	4-Nitrophenol	40.000	37.669	5.8	83	0.00
57	2,4-Dinitrotoluene	40.000	39.661	0.8	89	0.00
58	Fluorene	40.000	39.563	1.1	91	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.679	0.8	90	0.00
60	Diethylphthalate	40.000	38.050	4.9	87	0.00
61	4-Chlorophenyl-phenylether	40.000	40.139	-0.3	93	0.00
62	4-Nitroaniline	40.000	38.753	3.1	84	0.00
63	Azobenzene	40.000	38.948	2.6	86	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	83	0.00
65	4,6-Dinitro-2-methylphenol	40.000	43.262	-8.2	89	0.00
66 c	n-Nitrosodiphenylamine	40.000	41.986	-5.0	90	0.00
67	4-Bromophenyl-phenylether	40.000	42.212	-5.5	91	0.00
68	Hexachlorobenzene	40.000	41.785	-4.5	91	0.00
69	Atrazine	40.000	40.538	-1.3	86	0.00
70 C	Pentachlorophenol	40.000	41.434	-3.6	88	0.00
71	Phenanthrene	40.000	40.478	-1.2	87	0.00
72	Anthracene	40.000	40.756	-1.9	87	0.00
73	Carbazole	40.000	39.973	0.1	86	0.00
74	Di-n-butylphthalate	40.000	38.519	3.7	82	0.00
75 C	Fluoranthene	40.000	38.714	3.2	85	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	86	0.00
77	Benzydine	40.000	52.350	-30.9#	107	0.00
78	Pyrene	40.000	41.795	-4.5	87	0.00
79 S	Terphenyl-d14	80.000	83.459	-4.3	85	0.00
80	Butylbenzylphthalate	40.000	39.314	1.7	83	0.00
81	Benzo(a)anthracene	40.000	40.519	-1.3	91	0.00
82	3,3'-Dichlorobenzidine	40.000	42.164	-5.4	93	0.00
83	Chrysene	40.000	40.624	-1.6	92	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.504	3.7	85	0.00
85 c	Di-n-octyl phthalate	40.000	38.239	4.4	90	0.00
86 I	Perylene-d12	20.000	20.000	0.0	97	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	44.261	-10.7	113	-0.01
88	Benzo(b)fluoranthene	40.000	40.354	-0.9	99	0.00
89	Benzo(k)fluoranthene	40.000	40.446	-1.1	103	0.00
90 C	Benzo(a)pyrene	40.000	41.224	-3.1	104	0.00
91	Dibenzo(a,h)anthracene	40.000	44.075	-10.2	113	0.00
92	Benzo(g,h,i)perylene	40.000	44.846	-12.1	113	-0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081723\
Data File : BM041429.D
Acq On : 17 Aug 2023 09:32
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 17 17:53:23 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 17 02:13:45 2023
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