

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072723\
 Data File : BP016401.D
 Acq On : 27 Jul 2023 15:38
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP072723

Quant Time: Jul 28 01:50:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 01:48:25 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	115	0.00
2	1,4-Dioxane	40.000	39.860	0.4	119	0.00
3	Pyridine	40.000	38.603	3.5	112	0.00
4	n-Nitrosodimethylamine	40.000	38.253	4.4	113	0.00
5 S	2-Fluorophenol	80.000	80.557	-0.7	116	0.00
6	Aniline	40.000	38.198	4.5	109	0.00
7 S	Phenol-d6	80.000	76.973	3.8	110	0.00
8	2-Chlorophenol	40.000	39.615	1.0	112	0.00
9	Benzaldehyde	40.000	39.105	2.2	111	0.00
10 C	Phenol	40.000	38.260	4.4	110	0.00
11	bis(2-Chloroethyl)ether	40.000	38.023	4.9	110	0.00
12	1,3-Dichlorobenzene	40.000	40.196	-0.5	116	0.00
13 C	1,4-Dichlorobenzene	40.000	39.895	0.3	115	0.00
14	1,2-Dichlorobenzene	40.000	39.494	1.3	113	0.00
15	Benzyl Alcohol	40.000	37.992	5.0	107	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.555	8.6	107	0.00
17	2-Methylphenol	40.000	38.446	3.9	109	0.00
18	Hexachloroethane	40.000	39.349	1.6	114	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	36.602	8.5	105	0.00
20	3+4-Methylphenols	40.000	38.299	4.3	107	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22	Acetophenone	40.000	39.561	1.1	106	0.00
23 S	Nitrobenzene-d5	80.000	81.217	-1.5	108	0.00
24	Nitrobenzene	40.000	40.212	-0.5	108	0.00
25	Isophorone	40.000	39.229	1.9	104	0.00
26 C	2-Nitrophenol	40.000	41.378	-3.4	114	0.00
27	2,4-Dimethylphenol	40.000	40.241	-0.6	107	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.036	2.4	105	0.00
29 C	2,4-Dichlorophenol	40.000	42.110	-5.3	107	0.00
30	1,2,4-Trichlorobenzene	40.000	41.865	-4.7	109	0.00
31	Naphthalene	40.000	39.939	0.2	107	0.00
32	Benzoic acid	40.000	40.374	-0.9	114	0.00
33	4-Chloroaniline	40.000	39.711	0.7	105	0.00
34 C	Hexachlorobutadiene	40.000	43.348	-8.4	111	0.00
35	Caprolactam	40.000	39.038	2.4	102	0.00
36 C	4-Chloro-3-methylphenol	40.000	39.936	0.2	104	0.00
37	2-Methylnaphthalene	40.000	39.972	0.1	105	0.00
38	1-Methylnaphthalene	40.000	39.703	0.7	105	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	104	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	42.664	-6.7	107	0.00
41 P	Hexachlorocyclopentadiene	40.000	45.884	-14.7	109	0.00
42 S	2,4,6-Tribromophenol	80.000	88.205	-10.3	106	0.00
43 C	2,4,6-Trichlorophenol	40.000	43.378	-8.4	105	0.00
44	2,4,5-Trichlorophenol	40.000	43.019	-7.5	106	0.00
45 S	2-Fluorobiphenyl	80.000	81.313	-1.6	104	0.00
46	1,1'-Biphenyl	40.000	40.246	-0.6	104	0.00
47	2-Chloronaphthalene	40.000	40.371	-0.9	104	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072723\
 Data File : BP016401.D
 Acq On : 27 Jul 2023 15:38
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP072723

Quant Time: Jul 28 01:50:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 28 01:48:25 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	41.373	-3.4	103	0.00
49	Acenaphthylene	40.000	40.487	-1.2	104	0.00
50	Dimethylphthalate	40.000	40.074	-0.2	104	0.00
51	2,6-Dinitrotoluene	40.000	42.120	-5.3	104	0.00
52 C	Acenaphthene	40.000	40.130	-0.3	103	0.00
53	3-Nitroaniline	40.000	41.051	-2.6	103	0.00
54 P	2,4-Dinitrophenol	40.000	40.125	-0.3	114	0.00
55	Dibenzofuran	40.000	39.919	0.2	103	0.00
56 P	4-Nitrophenol	40.000	41.827	-4.6	104	0.00
57	2,4-Dinitrotoluene	40.000	43.018	-7.5	103	0.00
58	Fluorene	40.000	40.008	-0.0	102	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	43.640	-9.1	106	0.00
60	Diethylphthalate	40.000	39.527	1.2	102	0.00
61	4-Chlorophenyl-phenylether	40.000	41.212	-3.0	103	0.00
62	4-Nitroaniline	40.000	41.534	-3.8	103	0.00
63	Azobenzene	40.000	38.441	3.9	102	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	103	0.00
65	4,6-Dinitro-2-methylphenol	40.000	41.151	-2.9	111	0.00
66 c	n-Nitrosodiphenylamine	40.000	40.026	-0.1	102	0.00
67	4-Bromophenyl-phenylether	40.000	42.889	-7.2	104	0.00
68	Hexachlorobenzene	40.000	43.036	-7.6	105	0.00
69	Atrazine	40.000	41.612	-4.0	104	0.00
70 C	Pentachlorophenol	40.000	43.835	-9.6	107	0.00
71	Phenanthrene	40.000	40.125	-0.3	103	0.00
72	Anthracene	40.000	40.441	-1.1	102	0.00
73	Carbazole	40.000	39.933	0.2	102	0.00
74	Di-n-butylphthalate	40.000	41.199	-3.0	103	0.00
75 C	Fluoranthene	40.000	41.172	-2.9	103	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	101	0.00
77	Benzydine	40.000	45.012	-12.5	104	0.00
78	Pyrene	40.000	40.819	-2.0	102	0.00
79 S	Terphenyl-d14	80.000	85.764	-7.2	103	0.00
80	Butylbenzylphthalate	40.000	41.997	-5.0	103	0.00
81	Benzo(a)anthracene	40.000	40.823	-2.1	101	0.00
82	3,3'-Dichlorobenzidine	40.000	42.382	-6.0	101	0.00
83	Chrysene	40.000	40.273	-0.7	101	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	41.234	-3.1	102	0.00
85 c	Di-n-octyl phthalate	40.000	40.541	-1.4	100	0.00
86 I	Perylene-d12	20.000	20.000	0.0	97	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.243	4.4	92	0.00
88	Benzo(b)fluoranthene	40.000	42.329	-5.8	98	0.00
89	Benzo(k)fluoranthene	40.000	42.666	-6.7	100	0.00
90 C	Benzo(a)pyrene	40.000	41.850	-4.6	97	0.00
91	Dibenzo(a,h)anthracene	40.000	38.538	3.7	92	0.00
92	Benzo(g,h,i)perylene	40.000	37.324	6.7	91	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072723\
Data File : BP016401.D
Acq On : 27 Jul 2023 15:38
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP072723

Quant Time: Jul 28 01:50:34 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP072723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 28 01:48:25 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