

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
 Data File : BF138988.D
 Acq On : 14 Aug 2024 12:03
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 14 12:36:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 17:50:01 2024
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	64	0.00
2	1,4-Dioxane	0.567	0.506	10.8	56	0.00
3	Pyridine	1.374	1.276	7.1	59	0.00
4	n-Nitrosodimethylamine	0.818	0.925	-13.1	73	0.02
5 S	2-Fluorophenol	1.296	1.279	1.3	64	0.00
6	Aniline	1.551	1.483	4.4	62	0.00
7 S	Phenol-d6	1.740	1.703	2.1	65	0.00
8	2-Chlorophenol	1.363	1.341	1.6	65	0.00
9	Benzaldehyde	1.043	0.855	18.0	61	0.00
10 C	Phenol	1.832	1.763	3.8	64	0.00
11	bis(2-Chloroethyl)ether	1.409	1.274	9.6	60	-0.01
12	1,3-Dichlorobenzene	1.526	1.474	3.4	64	-0.01
13 C	1,4-Dichlorobenzene	1.540	1.520	1.3	65	0.00
14	1,2-Dichlorobenzene	1.439	1.411	1.9	64	0.00
15	Benzyl Alcohol	1.254	1.292	-3.0	68	0.00
16	2,2'-oxybis(1-Chloropropane	2.426	2.134	12.0	58	-0.01
17	2-Methylphenol	1.126	1.081	4.0	63	0.00
18	Hexachloroethane	0.580	0.576	0.7	65	-0.01
19 P	n-Nitroso-di-n-propylamine	1.051	1.047	0.4	68	0.00
20	3+4-Methylphenols	1.444	1.458	-1.0	68	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	64	0.00
22	Acetophenone	0.490	0.526	-7.3	70	0.00
23 S	Nitrobenzene-d5	0.409	0.430	-5.1	68	0.00
24	Nitrobenzene	0.416	0.434	-4.3	67	0.00
25	Isophorone	0.699	0.701	-0.3	66	0.00
26 C	2-Nitrophenol	0.179	0.179	0.0	63	0.00
27	2,4-Dimethylphenol	0.214	0.214	0.0	64	0.00
28	bis(2-Chloroethoxy)methane	0.425	0.404	4.9	62	0.00
29 C	2,4-Dichlorophenol	0.275	0.280	-1.8	65	0.00
30	1,2,4-Trichlorobenzene	0.318	0.321	-0.9	65	0.00
31	Naphthalene	1.053	1.057	-0.4	65	-0.01
32	Benzoic acid	0.168	0.125	25.6#	48#	0.02
33	4-Chloroaniline	0.353	0.340	3.7	63	0.00
34 C	Hexachlorobutadiene	0.192	0.205	-6.8	69	-0.01
35	Caprolactam	0.082	0.089	-8.5	73	0.00
36 C	4-Chloro-3-methylphenol	0.315	0.334	-6.0	70	0.00
37	2-Methylnaphthalene	0.665	0.681	-2.4	67	-0.01
38	1-Methylnaphthalene	0.652	0.671	-2.9	68	-0.01
39 I	Acenaphthene-d10	1.000	1.000	0.0	69	-0.01
40	1,2,4,5-Tetrachlorobenzene	0.556	0.560	-0.7	70	-0.01
41 P	Hexachlorocyclopentadiene	0.120	0.104	13.3	56	-0.01
42 S	2,4,6-Tribromophenol	0.164	0.163	0.6	70	0.00
43 C	2,4,6-Trichlorophenol	0.339	0.331	2.4	67	0.00
44	2,4,5-Trichlorophenol	0.370	0.361	2.4	67	0.00
45 S	2-Fluorobiphenyl	1.331	1.376	-3.4	72	-0.01
46	1,1'-Biphenyl	1.566	1.563	0.2	70	-0.01
47	2-Chloronaphthalene	1.165	1.143	1.9	68	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
 Data File : BF138988.D
 Acq On : 14 Aug 2024 12:03
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 14 12:36:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 17:50:01 2024
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.395	0.418	-5.8	74	0.00
49	Acenaphthylene	1.652	1.626	1.6	68	-0.01
50	Dimethylphthalate	1.279	1.330	-4.0	74	-0.01
51	2,6-Dinitrotoluene	0.289	0.299	-3.5	71	0.00
52 C	Acenaphthene	1.111	1.101	0.9	70	-0.01
53	3-Nitroaniline	0.298	0.287	3.7	68	0.00
54 P	2,4-Dinitrophenol	0.133	0.120	9.8	65	0.00
55	Dibenzofuran	1.568	1.584	-1.0	72	-0.01
56 P	4-Nitrophenol	0.179	0.214	-19.6	84	0.01
57	2,4-Dinitrotoluene	0.368	0.393	-6.8	74	0.00
58	Fluorene	1.249	1.314	-5.2	75	0.00
59	2,3,4,6-Tetrachlorophenol	0.283	0.275	2.8	69	0.00
60	Diethylphthalate	1.213	1.324	-9.2	79	0.00
61	4-Chlorophenyl-phenylether	0.614	0.637	-3.7	74	-0.01
62	4-Nitroaniline	0.284	0.283	0.4	71	0.00
63	Azobenzene	1.345	1.362	-1.3	72	-0.01
64 I	Phenanthrene-d10	1.000	1.000	0.0	75	-0.01
65	4,6-Dinitro-2-methylphenol	0.122	0.113	7.4	68	0.00
66 c	n-Nitrosodiphenylamine	0.625	0.622	0.5	74	0.00
67	4-Bromophenyl-phenylether	0.217	0.212	2.3	73	-0.01
68	Hexachlorobenzene	0.224	0.221	1.3	76	0.00
69	Atrazine	0.161	0.152	5.6	70	0.00
70 C	Pentachlorophenol	0.101	0.094	6.9	69	0.00
71	Phenanthrene	1.030	1.031	-0.1	76	-0.01
72	Anthracene	1.015	1.008	0.7	76	-0.01
73	Carbazole	0.875	0.850	2.9	73	0.00
74	Di-n-butylphthalate	0.984	1.072	-8.9	81	-0.01
75 C	Fluoranthene	0.961	0.935	2.7	73	-0.01
76 I	Chrysene-d12	1.000	1.000	0.0	66	0.00
77	Benzidine	0.478	0.426	10.9	52	0.00
78	Pyrene	1.883	1.903	-1.1	72	-0.01
79 S	Terphenyl-d14	1.195	1.229	-2.8	73	0.00
80	Butylbenzylphthalate	0.603	0.627	-4.0	66	-0.01
81	Benzo(a)anthracene	1.377	1.316	4.4	62	0.00
82	3,3'-Dichlorobenzidine	0.352	0.363	-3.1	67	0.00
83	Chrysene	1.243	1.276	-2.7	69	0.00
84	Bis(2-ethylhexyl)phthalate	0.883	0.812	8.0	56	-0.02
85 c	Di-n-octyl phthalate	1.634	1.348	17.5	51	-0.02
86 I	Perylene-d12	1.000	1.000	0.0	58	0.00
87	Indeno(1,2,3-cd)pyrene	1.433	1.318	8.0	54	0.00
88	Benzo(b)fluoranthene	1.240	1.293	-4.3	61	0.00
89	Benzo(k)fluoranthene	1.073	1.068	0.5	61	0.00
90 C	Benzo(a)pyrene	1.043	1.067	-2.3	60	0.00
91	Dibenzo(a,h)anthracene	1.177	1.083	8.0	55	0.00
92	Benzo(g,h,i)perylene	1.221	1.091	10.6	52	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081424\
Data File : BF138988.D
Acq On : 14 Aug 2024 12:03
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 14 12:36:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 30 17:50:01 2024
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
----------	-------	------	------	-------	----------

(#) = Out of Range

SPCC's out = 0 CCC's out = 0