

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031824\
 Data File : BF137622.D
 Acq On : 18 Mar 2024 10:36
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 18 11:35:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 14:45:20 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	81	0.00
2	1,4-Dioxane	0.722	0.668	7.5	73	0.00
3	Pyridine	1.741	1.673	3.9	71	0.00
4	n-Nitrosodimethylamine	0.652	0.593	9.0	69	0.00
5 S	2-Fluorophenol	1.321	1.285	2.7	75	0.00
6	Aniline	2.329	2.130	8.5	70	0.00
7 S	Phenol-d6	1.906	1.737	8.9	73	0.00
8	2-Chlorophenol	1.393	1.316	5.5	75	0.00
9	Benzaldehyde	0.934	0.868	7.1	70	0.00
10 C	Phenol	2.052	1.891	7.8	72	0.00
11	bis(2-Chloroethyl)ether	1.503	1.355	9.8	70	0.00
12	1,3-Dichlorobenzene	1.488	1.401	5.8	75	0.00
13 C	1,4-Dichlorobenzene	1.523	1.430	6.1	75	0.00
14	1,2-Dichlorobenzene	1.398	1.310	6.3	76	0.00
15	Benzyl Alcohol	1.187	1.119	5.7	71	0.00
16	2,2'-oxybis(1-Chloropropane	2.584	2.152	16.7	67	0.00
17	2-Methylphenol	1.303	1.213	6.9	74	0.00
18	Hexachloroethane	0.501	0.479	4.4	76	0.00
19 P	n-Nitroso-di-n-propylamine	1.177	0.995	15.5	69	0.00
20	3+4-Methylphenols	1.493	1.383	7.4	74	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	77	0.00
22	Acetophenone	0.484	0.463	4.3	74	0.00
23 S	Nitrobenzene-d5	0.430	0.408	5.1	71	0.00
24	Nitrobenzene	0.432	0.411	4.9	72	0.00
25	Isophorone	0.818	0.740	9.5	69	0.00
26 C	2-Nitrophenol	0.172	0.174	-1.2	75	0.00
27	2,4-Dimethylphenol	0.321	0.311	3.1	73	0.00
28	bis(2-Chloroethoxy)methane	0.480	0.453	5.6	71	0.00
29 C	2,4-Dichlorophenol	0.294	0.288	2.0	73	0.00
30	1,2,4-Trichlorobenzene	0.310	0.305	1.6	76	0.00
31	Naphthalene	1.073	1.018	5.1	74	0.00
32	Benzoic acid	0.167	0.138	17.4	64	0.00
33	4-Chloroaniline	0.421	0.407	3.3	70	0.00
34 C	Hexachlorobutadiene	0.183	0.180	1.6	75	0.00
35	Caprolactam	0.100	0.098	2.0	74	0.00
36 C	4-Chloro-3-methylphenol	0.329	0.308	6.4	71	0.00
37	2-Methylnaphthalene	0.684	0.646	5.6	73	0.00
38	1-Methylnaphthalene	0.644	0.604	6.2	73	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	76	0.00
40	1,2,4,5-Tetrachlorobenzene	0.558	0.551	1.3	74	0.00
41 P	Hexachlorocyclopentadiene	0.180	0.204	-13.3	81	0.00
42 S	2,4,6-Tribromophenol	0.176	0.172	2.3	74	0.00
43 C	2,4,6-Trichlorophenol	0.364	0.357	1.9	72	0.00
44	2,4,5-Trichlorophenol	0.419	0.398	5.0	70	0.00
45 S	2-Fluorobiphenyl	1.278	1.182	7.5	72	0.00
46	1,1'-Biphenyl	1.464	1.399	4.4	72	0.00
47	2-Chloronaphthalene	1.116	1.069	4.2	72	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031824\
 Data File : BF137622.D
 Acq On : 18 Mar 2024 10:36
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 18 11:35:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 14:45:20 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	0.379	0.368	2.9	69	0.00
49	Acenaphthylene	1.786	1.680	5.9	71	0.00
50	Dimethylphthalate	1.392	1.297	6.8	71	0.00
51	2,6-Dinitrotoluene	0.291	0.286	1.7	72	0.00
52 C	Acenaphthene	1.066	0.989	7.2	71	0.00
53	3-Nitroaniline	0.308	0.304	1.3	70	0.00
54 P	2,4-Dinitrophenol	0.111	0.116	-4.5	72	0.00
55	Dibenzofuran	1.624	1.501	7.6	70	0.00
56 P	4-Nitrophenol	0.190	0.195	-2.6	72	0.00
57	2,4-Dinitrotoluene	0.358	0.358	0.0	72	0.00
58	Fluorene	1.247	1.138	8.7	71	0.00
59	2,3,4,6-Tetrachlorophenol	0.336	0.314	6.5	70	0.00
60	Diethylphthalate	1.315	1.254	4.6	73	0.00
61	4-Chlorophenyl-phenylether	0.611	0.564	7.7	73	0.00
62	4-Nitroaniline	0.265	0.247	6.8	69	0.00
63	Azobenzene	1.514	1.332	12.0	66	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
65	4,6-Dinitro-2-methylphenol	0.111	0.113	-1.8	70	0.00
66 c	n-Nitrosodiphenylamine	0.644	0.621	3.6	71	0.00
67	4-Bromophenyl-phenylether	0.218	0.218	0.0	72	0.00
68	Hexachlorobenzene	0.220	0.218	0.9	73	0.00
69	Atrazine	0.196	0.198	-1.0	74	0.00
70 C	Pentachlorophenol	0.133	0.121	9.0	63	0.00
71	Phenanthrene	1.088	1.019	6.3	70	0.00
72	Anthracene	1.118	1.062	5.0	71	0.00
73	Carbazole	0.958	0.923	3.7	71	0.00
74	Di-n-butylphthalate	1.139	1.121	1.6	72	0.00
75 C	Fluoranthene	1.064	1.018	4.3	72	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	87	0.00
77	Benzidine	0.489	0.321	34.4#	49#	0.00
78	Pyrene	1.962	1.709	12.9	72	0.00
79 S	Terphenyl-d14	1.454	1.276	12.2	75	0.00
80	Butylbenzylphthalate	0.501	0.623	-24.4	103	0.00
81	Benzo(a)anthracene	1.402	1.303	7.1	80	0.00
82	3,3'-Dichlorobenzidine	0.373	0.394	-5.6	82	0.00
83	Chrysene	1.417	1.352	4.6	81	0.00
84	Bis(2-ethylhexyl)phthalate	0.637	0.800	-25.6#	103	0.00
85 c	Di-n-octyl phthalate	1.120	1.264	-12.9	92	0.00
86 I	Perylene-d12	1.000	1.000	0.0	82	0.00
87	Indeno(1,2,3-cd)pyrene	1.316	1.184	10.0	70	0.00
88	Benzo(b)fluoranthene	1.183	1.197	-1.2	82	0.00
89	Benzo(k)fluoranthene	1.271	1.139	10.4	76	0.00
90 C	Benzo(a)pyrene	1.166	1.109	4.9	75	0.00
91	Dibenzo(a,h)anthracene	1.068	0.942	11.8	68	0.00
92	Benzo(g,h,i)perylene	1.096	0.992	9.5	70	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031824\
Data File : BF137622.D
Acq On : 18 Mar 2024 10:36
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Mar 18 11:35:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 14 14:45:20 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