

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071423\
 Data File : BF134292.D
 Acq On : 14 Jul 2023 15:56
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 15 00:42:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 14 22:51:27 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	62	0.00
2	1,4-Dioxane	0.493	0.477	3.2	60	0.00
3	Pyridine	1.284	1.225	4.6	59	0.00
4	n-Nitrosodimethylamine	0.733	0.694	5.3	58	0.00
5 S	2-Fluorophenol	1.196	1.195	0.1	63	0.00
6	Aniline	1.789	1.760	1.6	61	0.00
7 S	Phenol-d6	1.470	1.438	2.2	63	0.00
8	2-Chlorophenol	1.214	1.204	0.8	63	0.00
9	Benzaldehyde	0.886	0.771	13.0	59	0.00
10 C	Phenol	1.546	1.526	1.3	63	0.00
11	bis(2-Chloroethyl)ether	1.138	1.089	4.3	60	0.00
12	1,3-Dichlorobenzene	1.294	1.269	1.9	62	0.00
13 C	1,4-Dichlorobenzene	1.307	1.320	-1.0	63	0.00
14	1,2-Dichlorobenzene	1.224	1.223	0.1	64	0.00
15	Benzyl Alcohol	1.134	1.129	0.4	61	0.00
16	2,2'-oxybis(1-Chloropropane	2.014	1.764	12.4	55	0.00
17	2-Methylphenol	1.087	1.067	1.8	61	0.00
18	Hexachloroethane	0.485	0.486	-0.2	62	0.00
19 P	n-Nitroso-di-n-propylamine	0.932	0.914	1.9	64	0.00
20	3+4-Methylphenols	1.319	1.353	-2.6	65	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	61	0.00
22	Acetophenone	0.455	0.474	-4.2	66	0.00
23 S	Nitrobenzene-d5	0.371	0.384	-3.5	64	0.00
24	Nitrobenzene	0.375	0.384	-2.4	62	0.00
25	Isophorone	0.674	0.675	-0.1	63	0.00
26 C	2-Nitrophenol	0.180	0.187	-3.9	62	0.00
27	2,4-Dimethylphenol	0.285	0.283	0.7	61	0.00
28	bis(2-Chloroethoxy)methane	0.390	0.388	0.5	61	0.00
29 C	2,4-Dichlorophenol	0.282	0.290	-2.8	63	0.00
30	1,2,4-Trichlorobenzene	0.291	0.303	-4.1	64	0.00
31	Naphthalene	0.966	0.971	-0.5	63	0.00
32	Benzoic acid	0.213	0.205	3.8	55	0.00
33	4-Chloroaniline	0.413	0.418	-1.2	62	0.00
34 C	Hexachlorobutadiene	0.184	0.199	-8.2	67	0.00
35	Caprolactam	0.093	0.089	4.3	58	0.00
36 C	4-Chloro-3-methylphenol	0.317	0.326	-2.8	64	0.00
37	2-Methylnaphthalene	0.683	0.686	-0.4	63	0.00
38	1-Methylnaphthalene	0.635	0.628	1.1	63	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	62	0.00
40	1,2,4,5-Tetrachlorobenzene	0.593	0.618	-4.2	65	0.00
41 P	Hexachlorocyclopentadiene	0.281	0.360	-28.1#	76	0.00
42 S	2,4,6-Tribromophenol	0.251	0.271	-8.0	68	0.00
43 C	2,4,6-Trichlorophenol	0.403	0.400	0.7	61	0.00
44	2,4,5-Trichlorophenol	0.474	0.473	0.2	62	0.00
45 S	2-Fluorobiphenyl	1.376	1.437	-4.4	67	0.00
46	1,1'-Biphenyl	1.604	1.617	-0.8	64	0.00
47	2-Chloronaphthalene	1.174	1.172	0.2	63	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071423\
 Data File : BF134292.D
 Acq On : 14 Jul 2023 15:56
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jul 15 00:42:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 14 22:51:27 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48	2-Nitroaniline	0.428	0.432	-0.9	62	0.00
49	Acenaphthylene	2.005	1.994	0.5	62	0.00
50	Dimethylphthalate	1.452	1.443	0.6	63	0.00
51	2,6-Dinitrotoluene	0.313	0.322	-2.9	63	0.00
52 C	Acenaphthene	1.164	1.182	-1.5	64	0.00
53	3-Nitroaniline	0.375	0.354	5.6	58	0.00
54 P	2,4-Dinitrophenol	0.162	0.171	-5.6	63	0.00
55	Dibenzofuran	1.818	1.832	-0.8	63	0.00
56 P	4-Nitrophenol	0.262	0.229	12.6	53	0.00
57	2,4-Dinitrotoluene	0.413	0.430	-4.1	64	0.00
58	Fluorene	1.415	1.433	-1.3	65	0.00
59	2,3,4,6-Tetrachlorophenol	0.374	0.382	-2.1	64	0.00
60	Diethylphthalate	1.513	1.533	-1.3	64	0.00
61	4-Chlorophenyl-phenylether	0.682	0.711	-4.3	66	0.00
62	4-Nitroaniline	0.378	0.364	3.7	60	0.00
63	Azobenzene	1.431	1.409	1.5	62	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	62	0.00
65	4,6-Dinitro-2-methylphenol	0.109	0.116	-6.4	62	0.00
66 c	n-Nitrosodiphenylamine	0.639	0.635	0.6	63	0.00
67	4-Bromophenyl-phenylether	0.213	0.217	-1.9	64	0.00
68	Hexachlorobenzene	0.226	0.241	-6.6	68	0.00
69	Atrazine	0.177	0.182	-2.8	68	0.00
70 C	Pentachlorophenol	0.135	0.124	8.1	55	0.00
71	Phenanthrene	1.007	0.998	0.9	64	0.00
72	Anthracene	1.047	1.058	-1.1	65	0.00
73	Carbazole	0.959	0.961	-0.2	63	0.00
74	Di-n-butylphthalate	1.140	1.195	-4.8	65	0.00
75 C	Fluoranthene	1.088	1.131	-4.0	67	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	88	0.00
77	Benzydine	0.476	0.541	-13.7	93	0.00
78	Pyrene	1.861	1.508	19.0	68	0.00
79 S	Terphenyl-d14	1.243	1.099	11.6	76	0.00
80	Butylbenzylphthalate	0.698	0.665	4.7	81	0.00
81	Benzo(a)anthracene	1.387	1.374	0.9	88	0.00
82	3,3'-Dichlorobenzidine	0.439	0.454	-3.4	93	0.00
83	Chrysene	1.343	1.313	2.2	87	0.00
84	Bis(2-ethylhexyl)phthalate	0.802	0.904	-12.7	97	0.00
85 c	Di-n-octyl phthalate	1.179	1.461	-23.9#	107	0.00
86 I	Perylene-d12	1.000	1.000	0.0	89	0.00
87	Indeno(1,2,3-cd)pyrene	1.388	1.268	8.6	74	0.00
88	Benzo(b)fluoranthene	1.176	1.246	-6.0	94	0.00
89	Benzo(k)fluoranthene	1.197	1.231	-2.8	94	0.00
90 C	Benzo(a)pyrene	1.137	1.153	-1.4	90	0.00
91	Dibenzo(a,h)anthracene	1.134	1.047	7.7	75	0.00
92	Benzo(g,h,i)perylene	1.169	1.048	10.4	74	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071423\
Data File : BF134292.D
Acq On : 14 Jul 2023 15:56
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jul 15 00:42:59 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 14 22:51:27 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
----------	-------	------	------	-------	----------

(#) = Out of Range

SPCC's out = 0 CCC's out = 1