

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101124\
 Data File : BF139763.D
 Acq On : 11 Oct 2024 09:29
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 14 02:30:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 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	102	0.00
2	1,4-Dioxane	0.499	0.504	-1.0	106	-0.02
3	Pyridine	1.213	1.262	-4.0	110	-0.02
4	n-Nitrosodimethylamine	0.794	0.749	5.7	97	0.00
5 S	2-Fluorophenol	1.154	1.165	-1.0	106	0.00
6	Aniline	1.377	1.285	6.7	102	0.00
7 S	Phenol-d6	1.552	1.528	1.5	104	0.00
8	2-Chlorophenol	1.236	1.223	1.1	104	0.00
9	Benzaldehyde	0.822	0.795	3.3	103	0.00
10 C	Phenol	1.632	1.621	0.7	103	0.00
11	bis(2-Chloroethyl)ether	1.329	1.217	8.4	98	0.00
12	1,3-Dichlorobenzene	1.391	1.373	1.3	104	0.00
13 C	1,4-Dichlorobenzene	1.409	1.387	1.6	105	0.00
14	1,2-Dichlorobenzene	1.320	1.275	3.4	101	0.00
15	Benzyl Alcohol	1.186	1.116	5.9	97	0.00
16	2,2'-oxybis(1-Chloropropane	2.300	2.108	8.3	96	0.00
17	2-Methylphenol	1.032	1.003	2.8	101	0.00
18	Hexachloroethane	0.525	0.505	3.8	101	0.00
19 P	n-Nitroso-di-n-propylamine	0.998	0.895	10.3	95	0.00
20	3+4-Methylphenols	1.296	1.233	4.9	100	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22	Acetophenone	0.475	0.452	4.8	98	0.00
23 S	Nitrobenzene-d5	0.395	0.380	3.8	98	0.00
24	Nitrobenzene	0.409	0.396	3.2	98	0.00
25	Isophorone	0.702	0.663	5.6	96	0.00
26 C	2-Nitrophenol	0.176	0.178	-1.1	98	0.00
27	2,4-Dimethylphenol	0.215	0.210	2.3	100	0.00
28	bis(2-Chloroethoxy)methane	0.419	0.400	4.5	97	0.00
29 C	2,4-Dichlorophenol	0.281	0.277	1.4	99	0.00
30	1,2,4-Trichlorobenzene	0.318	0.315	0.9	100	0.00
31	Naphthalene	1.023	0.998	2.4	100	0.00
32	Benzoic acid	0.214	0.214	0.0	94	0.00
33	4-Chloroaniline	0.346	0.318	8.1	92	0.00
34 C	Hexachlorobutadiene	0.205	0.197	3.9	98	0.00
35	Caprolactam	0.092	0.088	4.3	98	0.00
36 C	4-Chloro-3-methylphenol	0.318	0.297	6.6	96	0.00
37	2-Methylnaphthalene	0.666	0.625	6.2	96	0.00
38	1-Methylnaphthalene	0.651	0.616	5.4	97	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
40	1,2,4,5-Tetrachlorobenzene	0.555	0.539	2.9	96	0.00
41 P	Hexachlorocyclopentadiene	0.170	0.214	-25.9#	112	0.00
42 S	2,4,6-Tribromophenol	0.193	0.171	11.4	88	0.00
43 C	2,4,6-Trichlorophenol	0.374	0.350	6.4	92	0.00
44	2,4,5-Trichlorophenol	0.404	0.382	5.4	94	0.00
45 S	2-Fluorobiphenyl	1.303	1.213	6.9	95	0.00
46	1,1'-Biphenyl	1.488	1.401	5.8	95	0.00
47	2-Chloronaphthalene	1.115	1.088	2.4	97	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101124\
 Data File : BF139763.D
 Acq On : 11 Oct 2024 09:29
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Oct 14 02:30:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 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	0.400	0.372	7.0	91	0.00
49	Acenaphthylene	1.696	1.602	5.5	94	0.00
50	Dimethylphthalate	1.309	1.220	6.8	94	0.00
51	2,6-Dinitrotoluene	0.296	0.278	6.1	93	0.00
52 C	Acenaphthene	1.164	1.090	6.4	93	0.00
53	3-Nitroaniline	0.302	0.274	9.3	90	0.00
54 P	2,4-Dinitrophenol	0.159	0.129	18.9	76	0.00
55	Dibenzofuran	1.630	1.511	7.3	93	0.00
56 P	4-Nitrophenol	0.230	0.192	16.5	81	0.00
57	2,4-Dinitrotoluene	0.387	0.357	7.8	90	0.00
58	Fluorene	1.298	1.202	7.4	94	0.00
59	2,3,4,6-Tetrachlorophenol	0.326	0.298	8.6	91	0.00
60	Diethylphthalate	1.352	1.213	10.3	90	0.00
61	4-Chlorophenyl-phenylether	0.650	0.597	8.2	94	0.00
62	4-Nitroaniline	0.309	0.276	10.7	89	0.00
63	Azobenzene	1.360	1.230	9.6	90	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
65	4,6-Dinitro-2-methylphenol	0.113	0.109	3.5	83	0.00
66 c	n-Nitrosodiphenylamine	0.590	0.584	1.0	93	0.00
67	4-Bromophenyl-phenylether	0.205	0.199	2.9	91	0.00
68	Hexachlorobenzene	0.228	0.223	2.2	92	0.00
69	Atrazine	0.158	0.140	11.4	96	0.00
70 C	Pentachlorophenol	0.135	0.115	14.8	75	0.00
71	Phenanthrene	0.943	0.899	4.7	90	0.00
72	Anthracene	0.933	0.898	3.8	91	0.00
73	Carbazole	0.896	0.844	5.8	88	0.00
74	Di-n-butylphthalate	1.090	0.987	9.4	84	0.00
75 C	Fluoranthene	1.031	0.928	10.0	85	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	83	0.00
77	Benzydine	0.353	0.334	5.4	80	0.00
78	Pyrene	1.789	1.838	-2.7	84	0.00
79 S	Terphenyl-d14	1.219	1.211	0.7	83	0.00
80	Butylbenzylphthalate	0.640	0.620	3.1	79	0.00
81	Benzo(a)anthracene	1.315	1.268	3.6	84	0.00
82	3,3'-Dichlorobenzidine	0.402	0.398	1.0	84	0.00
83	Chrysene	1.202	1.174	2.3	82	0.00
84	Bis(2-ethylhexyl)phthalate	0.799	0.810	-1.4	83	0.00
85 c	Di-n-octyl phthalate	1.174	1.213	-3.3	85	0.00
86 I	Perylene-d12	1.000	1.000	0.0	92	0.00
87	Indeno(1,2,3-cd)pyrene	1.177	1.332	-13.2	102	0.00
88	Benzo(b)fluoranthene	1.293	1.132	12.5	91	0.00
89	Benzo(k)fluoranthene	1.097	1.082	1.4	89	0.00
90 C	Benzo(a)pyrene	1.020	1.002	1.8	93	0.00
91	Dibenzo(a,h)anthracene	0.976	1.107	-13.4	101	0.00
92	Benzo(g,h,i)perylene	0.979	1.140	-16.4	103	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101124\
Data File : BF139763.D
Acq On : 11 Oct 2024 09:29
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Oct 14 02:30:57 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 02 04:58:47 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