

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013124\
 Data File : BP019500.D
 Acq On : 31 Jan 2024 21:37
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP013124

Quant Time: Feb 01 00:53:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 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	88	0.00
2	1,4-Dioxane	0.478	0.439	8.2	92	0.00
3	Pyridine	1.296	1.220	5.9	89	0.00
4	n-Nitrosodimethylamine	0.564	0.548	2.8	90	0.00
5 S	2-Fluorophenol	1.241	1.160	6.5	90	0.00
6	Aniline	1.626	1.549	4.7	87	0.00
7 S	Phenol-d6	1.673	1.606	4.0	89	0.00
8	2-Chlorophenol	1.274	1.213	4.8	89	0.00
9	Benzaldehyde	1.173	0.955	18.6	63	0.00
10 C	Phenol	1.667	1.596	4.3	88	0.00
11	bis(2-Chloroethyl)ether	1.297	1.232	5.0	86	0.00
12	1,3-Dichlorobenzene	1.559	1.461	6.3	89	0.00
13 C	1,4-Dichlorobenzene	1.579	1.500	5.0	90	0.00
14	1,2-Dichlorobenzene	1.528	1.416	7.3	87	0.00
15	Benzyl Alcohol	1.337	1.283	4.0	86	0.00
16	2,2'-oxybis(1-Chloropropane	1.297	1.230	5.2	85	0.00
17	2-Methylphenol	1.123	1.075	4.3	86	0.00
18	Hexachloroethane	0.612	0.571	6.7	86	0.00
19 P	n-Nitroso-di-n-propylamine	1.143	1.061	7.2	82	0.00
20	3+4-Methylphenols	1.532	1.479	3.5	88	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	84	0.00
22	Acetophenone	0.570	0.550	3.5	85	0.00
23 S	Nitrobenzene-d5	0.462	0.450	2.6	86	0.00
24	Nitrobenzene	0.465	0.454	2.4	87	0.00
25	Isophorone	0.804	0.755	6.1	81	0.00
26 C	2-Nitrophenol	0.177	0.182	-2.8	88	0.00
27	2,4-Dimethylphenol	0.227	0.219	3.5	84	0.00
28	bis(2-Chloroethoxy)methane	0.455	0.433	4.8	82	0.00
29 C	2,4-Dichlorophenol	0.343	0.329	4.1	84	0.00
30	1,2,4-Trichlorobenzene	0.420	0.400	4.8	85	0.00
31	Naphthalene	1.097	1.048	4.5	85	0.00
32	Benzoic acid	0.221	0.219	0.9	83	-0.01
33	4-Chloroaniline	0.406	0.381	6.2	82	0.00
34 C	Hexachlorobutadiene	0.311	0.302	2.9	86	0.00
35	Caprolactam	0.097	0.085	12.4	82	0.00
36 C	4-Chloro-3-methylphenol	0.373	0.350	6.2	82	0.00
37	2-Methylnaphthalene	0.759	0.712	6.2	82	0.00
38	1-Methylnaphthalene	0.746	0.697	6.6	82	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	80	0.00
40	1,2,4,5-Tetrachlorobenzene	0.771	0.762	1.2	83	0.00
41 P	Hexachlorocyclopentadiene	0.348	0.359	-3.2	81	0.00
42 S	2,4,6-Tribromophenol	0.292	0.268	8.2	80	0.00
43 C	2,4,6-Trichlorophenol	0.467	0.460	1.5	81	0.00
44	2,4,5-Trichlorophenol	0.513	0.504	1.8	81	0.00
45 S	2-Fluorobiphenyl	1.618	1.568	3.1	81	0.00
46	1,1'-Biphenyl	1.596	1.539	3.6	80	0.00
47	2-Chloronaphthalene	1.310	1.272	2.9	82	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013124\
 Data File : BP019500.D
 Acq On : 31 Jan 2024 21:37
 Operator : MA/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ICVBP013124

Quant Time: Feb 01 00:53:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 00:51:48 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	0.366	0.355	3.0	82	0.00
49	Acenaphthylene	1.838	1.752	4.7	81	0.00
50	Dimethylphthalate	1.531	1.395	8.9	79	0.00
51	2,6-Dinitrotoluene	0.336	0.316	6.0	80	0.00
52 C	Acenaphthene	1.141	1.086	4.8	80	0.00
53	3-Nitroaniline	0.309	0.290	6.1	83	0.00
54 P	2,4-Dinitrophenol	0.175	0.164	6.3	83	0.00
55	Dibenzofuran	1.858	1.731	6.8	80	0.00
56 P	4-Nitrophenol	0.253	0.234	7.5	86	0.00
57	2,4-Dinitrotoluene	0.441	0.415	5.9	82	0.00
58	Fluorene	1.480	1.361	8.0	80	0.00
59	2,3,4,6-Tetrachlorophenol	0.431	0.404	6.3	81	0.00
60	Diethylphthalate	1.523	1.336	12.3	78	0.00
61	4-Chlorophenyl-phenylether	0.832	0.759	8.8	78	0.00
62	4-Nitroaniline	0.324	0.288	11.1	84	0.00
63	Azobenzene	1.473	1.323	10.2	78	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	82	0.00
65	4,6-Dinitro-2-methylphenol	0.114	0.118	-3.5	84	0.00
66 c	n-Nitrosodiphenylamine	0.599	0.587	2.0	80	0.00
67	4-Bromophenyl-phenylether	0.250	0.246	1.6	79	0.00
68	Hexachlorobenzene	0.292	0.280	4.1	79	0.00
69	Atrazine	0.199	0.188	5.5	81	0.00
70 C	Pentachlorophenol	0.182	0.182	0.0	83	0.00
71	Phenanthrene	1.053	1.002	4.8	83	0.00
72	Anthracene	1.050	0.996	5.1	81	0.00
73	Carbazole	0.954	0.895	6.2	86	0.00
74	Di-n-butylphthalate	1.153	1.053	8.7	81	0.00
75 C	Fluoranthene	1.286	1.195	7.1	88	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
77	Benzydine	0.418	0.357	14.6	91	0.00
78	Pyrene	1.412	1.296	8.2	90	0.00
79 S	Terphenyl-d14	1.216	1.100	9.5	87	0.00
80	Butylbenzylphthalate	0.527	0.485	8.0	89	0.00
81	Benzo(a)anthracene	1.408	1.343	4.6	97	0.00
82	3,3'-Dichlorobenzidine	0.513	0.495	3.5	97	0.00
83	Chrysene	1.337	1.286	3.8	98	0.00
84	Bis(2-ethylhexyl)phthalate	0.803	0.732	8.8	89	0.00
85 c	Di-n-octyl phthalate	1.412	1.327	6.0	90	0.00
86 I	Perylene-d12	1.000	1.000	0.0	103	0.00
87	Indeno(1,2,3-cd)pyrene	1.531	1.463	4.4	102	0.00
88	Benzo(b)fluoranthene	1.261	1.187	5.9	99	0.00
89	Benzo(k)fluoranthene	1.204	1.111	7.7	101	0.00
90 C	Benzo(a)pyrene	1.122	1.066	5.0	102	0.00
91	Dibenzo(a,h)anthracene	1.259	1.201	4.6	101	0.00
92	Benzo(g,h,i)perylene	1.277	1.219	4.5	101	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP013124\
Data File : BP019500.D
Acq On : 31 Jan 2024 21:37
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP013124

Quant Time: Feb 01 00:53:56 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 01 00:51:48 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