

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030722\
 Data File : BP009280.D
 Acq On : 07 Mar 2022 10:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 07 16:13:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	107	0.00
2	1,4-Dioxane	40.000	35.977	10.1	107	0.00
3	Pyridine	40.000	36.667	8.3	106	0.00
4	n-Nitrosodimethylamine	40.000	36.480	8.8	103	0.00
5 S	2-Fluorophenol	80.000	73.084	8.6	107	0.00
6	Aniline	40.000	36.309	9.2	106	0.00
7 S	Phenol-d6	80.000	72.646	9.2	107	0.00
8	2-Chlorophenol	40.000	37.011	7.5	107	0.00
9	Benzaldehyde	40.000	36.642	8.4	106	0.00
10 C	Phenol	40.000	35.966	10.1	106	0.00
11	bis(2-Chloroethyl)ether	40.000	36.548	8.6	105	0.00
12	1,3-Dichlorobenzene	40.000	36.072	9.8	108	0.00
13 C	1,4-Dichlorobenzene	40.000	36.167	9.6	108	0.00
14	1,2-Dichlorobenzene	40.000	36.114	9.7	107	0.00
15	Benzyl Alcohol	40.000	36.991	7.5	107	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	35.784	10.5	106	0.00
17	2-Methylphenol	40.000	36.971	7.6	107	0.00
18	Hexachloroethane	40.000	36.594	8.5	108	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.563	6.1	108	0.00
20	3+4-Methylphenols	40.000	37.028	7.4	107	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	108	0.00
22	Acetophenone	40.000	36.679	8.3	107	0.00
23 S	Nitrobenzene-d5	80.000	76.492	4.4	108	0.00
24	Nitrobenzene	40.000	37.323	6.7	107	0.00
25	Isophorone	40.000	38.334	4.2	109	0.00
26 C	2-Nitrophenol	40.000	40.389	-1.0	110	0.00
27	2,4-Dimethylphenol	40.000	37.008	7.5	108	0.00
28	bis(2-Chloroethoxy)methane	40.000	37.597	6.0	109	0.00
29 C	2,4-Dichlorophenol	40.000	38.588	3.5	109	0.00
30	1,2,4-Trichlorobenzene	40.000	36.836	7.9	109	0.00
31	Naphthalene	40.000	36.460	8.8	107	0.00
32	Benzoic acid	40.000	42.626	-6.6	118	0.00
33	4-Chloroaniline	40.000	37.542	6.1	109	0.00
34 C	Hexachlorobutadiene	40.000	37.466	6.3	111	0.00
35	Caprolactam	40.000	40.995	-2.5	112	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.498	3.8	110	0.00
37	2-Methylnaphthalene	40.000	37.146	7.1	109	0.00
38	1-Methylnaphthalene	40.000	37.206	7.0	110	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	112	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	36.997	7.5	111	0.00
41 P	Hexachlorocyclopentadiene	40.000	37.978	5.1	113	0.00
42 S	2,4,6-Tribromophenol	80.000	81.257	-1.6	118	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.074	2.3	113	0.00
44	2,4,5-Trichlorophenol	40.000	38.569	3.6	112	0.00
45 S	2-Fluorobiphenyl	80.000	72.758	9.1	111	0.00
46	1,1'-Biphenyl	40.000	36.567	8.6	111	0.00
47	2-Chloronaphthalene	40.000	36.346	9.1	110	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030722\
 Data File : BP009280.D
 Acq On : 07 Mar 2022 10:53
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 07 16:13:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 05 01:55:59 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	41.242	-3.1	114	0.00
49	Acenaphthylene	40.000	37.190	7.0	111	0.00
50	Dimethylphthalate	40.000	38.024	4.9	114	0.00
51	2,6-Dinitrotoluene	40.000	40.555	-1.4	114	0.00
52 C	Acenaphthene	40.000	37.204	7.0	111	0.00
53	3-Nitroaniline	40.000	40.461	-1.2	114	0.00
54 P	2,4-Dinitrophenol	40.000	43.588	-9.0	120	0.00
55	Dibenzofuran	40.000	36.920	7.7	112	0.00
56 P	4-Nitrophenol	40.000	41.066	-2.7	114	0.00
57	2,4-Dinitrotoluene	40.000	42.452	-6.1	116	0.00
58	Fluorene	40.000	37.274	6.8	114	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.751	0.6	117	0.00
60	Diethylphthalate	40.000	38.946	2.6	116	0.00
61	4-Chlorophenyl-phenylether	40.000	37.789	5.5	115	0.00
62	4-Nitroaniline	40.000	41.014	-2.5	115	0.00
63	Azobenzene	40.000	38.354	4.1	114	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	114	0.00
65	4,6-Dinitro-2-methylphenol	40.000	42.803	-7.0	119	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.347	6.6	114	0.00
67	4-Bromophenyl-phenylether	40.000	37.864	5.3	116	0.00
68	Hexachlorobenzene	40.000	38.294	4.3	117	0.00
69	Atrazine	40.000	39.730	0.7	116	0.00
70 C	Pentachlorophenol	40.000	40.511	-1.3	116	0.00
71	Phenanthrene	40.000	36.824	7.9	113	0.00
72	Anthracene	40.000	37.565	6.1	114	0.00
73	Carbazole	40.000	37.789	5.5	113	0.00
74	Di-n-butylphthalate	40.000	40.414	-1.0	117	0.00
75 C	Fluoranthene	40.000	37.926	5.2	113	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	112	0.00
77	Benzydine	40.000	41.751	-4.4	110	0.00
78	Pyrene	40.000	37.951	5.1	113	0.00
79 S	Terphenyl-d14	80.000	75.779	5.3	114	0.00
80	Butylbenzylphthalate	40.000	41.580	-3.9	115	0.00
81	Benzo(a)anthracene	40.000	37.219	7.0	112	0.00
82	3,3'-Dichlorobenzidine	40.000	39.433	1.4	111	0.00
83	Chrysene	40.000	37.068	7.3	112	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.467	-1.2	114	0.00
85 c	Di-n-octyl phthalate	40.000	40.860	-2.1	112	0.00
86 I	Perylene-d12	20.000	20.000	0.0	109	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	37.601	6.0	108	-0.01
88	Benzo(b)fluoranthene	40.000	37.799	5.5	107	0.00
89	Benzo(k)fluoranthene	40.000	38.442	3.9	113	0.00
90 C	Benzo(a)pyrene	40.000	38.384	4.0	110	0.00
91	Dibenzo(a,h)anthracene	40.000	37.600	6.0	108	0.00
92	Benzo(g,h,i)perylene	40.000	37.833	5.4	109	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030722\
Data File : BP009280.D
Acq On : 07 Mar 2022 10:53
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Mar 07 16:13:04 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP030522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 05 01:55:59 2022
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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

(#) = Out of Range SPCC's out = 0 CCC's out = 0