

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081024\
 Data File : BM047142.D
 Acq On : 10 Aug 2024 14:31
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: Aug 11 23:32:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:24:37 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.375	152	292955	20.000	ng	0.00
21) Naphthalene-d8	10.128	136	1186627	20.000	ng	0.00
39) Acenaphthene-d10	14.033	164	788651	20.000	ng	0.00
64) Phenanthrene-d10	16.792	188	1690400	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1593166	20.000	ng	0.00
86) Perylene-d12	23.727	264	1901592	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.022	112	346900	21.159	ng	0.00
7) Phenol-d6	6.581	99	468308	20.730	ng	0.00
23) Nitrobenzene-d5	8.522	82	494443	20.981	ng	0.00
42) 2,4,6-Tribromophenol	15.539	330	183466	20.063	ng	0.00
45) 2-Fluorobiphenyl	12.639	172	1057713	20.688	ng	0.00
79) Terphenyl-d14	19.456	244	1459931	19.636	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.040	88	71740	10.561	ng	99
3) Pyridine	3.416	79	203646	10.230	ng	97
4) n-Nitrosodimethylamine	3.334	42	89658	10.302	ng	# 99
6) Aniline	6.722	93	222003	10.397	ng	99
8) 2-Chlorophenol	6.951	128	186177	10.412	ng	98
9) Benzaldehyde	6.540	77	153438	9.573	ng	99
10) Phenol	6.604	94	234318	10.395	ng	99
11) bis(2-Chloroethyl)ether	6.822	93	199316	10.325	ng	99
12) 1,3-Dichlorobenzene	7.263	146	220661	10.384	ng	99
13) 1,4-Dichlorobenzene	7.410	146	224423	10.431	ng	99
14) 1,2-Dichlorobenzene	7.716	146	215097	10.596	ng	98
15) Benzyl Alcohol	7.628	79	157264	9.781	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.910	45	260329	10.497	ng	98
17) 2-Methylphenol	7.834	107	154501	10.274	ng	99
18) Hexachloroethane	8.428	117	84578	10.149	ng	97
19) n-Nitroso-di-n-propyla...	8.181	70	160456	10.544	ng	97
20) 3+4-Methylphenols	8.157	107	211260	10.048	ng	98
22) Acetophenone	8.192	105	306786	10.711	ng	98
24) Nitrobenzene	8.563	77	253056	10.664	ng	99
25) Isophorone	9.086	82	418965	10.211	ng	100
26) 2-Nitrophenol	9.263	139	98575	9.380	ng	98
27) 2,4-Dimethylphenol	9.345	122	123790	10.056	ng	99
28) bis(2-Chloroethoxy)met...	9.575	93	273307	10.586	ng	99
29) 2,4-Dichlorophenol	9.792	162	178052	9.748	ng	99
30) 1,2,4-Trichlorobenzene	9.998	180	208733	10.098	ng	98
31) Naphthalene	10.181	128	617578	10.447	ng	99
32) Benzoic acid	9.457	122	125871	8.752	ng	96
33) 4-Chloroaniline	10.304	127	233198	10.258	ng	98
34) Hexachlorobutadiene	10.469	225	118512	9.795	ng	99
35) Caprolactam	11.086	113	61384	9.677	ng	98
36) 4-Chloro-3-methylphenol	11.451	107	192243	9.883	ng	94
37) 2-Methylnaphthalene	11.804	142	433416	10.298	ng	99
38) 1-Methylnaphthalene	12.027	142	430690	10.354	ng	100
40) 1,2,4,5-Tetrachloroben...	12.186	216	209468	9.738	ng	99
41) Hexachlorocyclopentadiene	12.163	237	70782	9.423	ng	96
43) 2,4,6-Trichlorophenol	12.445	196	146662	9.593	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.522	196	161605	9.520	ng	98
46) 1,1'-Biphenyl	12.851	154	584326	10.407	ng	99
47) 2-Chloronaphthalene	12.886	162	460927	10.503	ng	99
48) 2-Nitroaniline	13.110	65	136444	9.925	ng	95
49) Acenaphthylene	13.745	152	669998	10.358	ng	100
50) Dimethylphthalate	13.510	163	565056	10.259	ng	100
51) 2,6-Dinitrotoluene	13.621	165	128162	9.954	ng	95
52) Acenaphthene	14.098	154	425366	10.367	ng	98
53) 3-Nitroaniline	13.957	138	124808	9.674	ng	96
54) 2,4-Dinitrophenol	14.174	184	60036	7.780	ng	96
55) Dibenzofuran	14.439	168	682447	10.530	ng	99
56) 4-Nitrophenol	14.292	139	105813	9.511	ng	94
57) 2,4-Dinitrotoluene	14.421	165	171542	9.982	ng	92
58) Fluorene	15.092	166	562937	10.527	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.674	232	135880	9.739	ng	96
60) Diethylphthalate	14.892	149	586005	10.418	ng	98
61) 4-Chlorophenyl-phenyle...	15.098	204	261449	10.269	ng	98
62) 4-Nitroaniline	15.127	138	127938m	10.065	ng	
63) Azobenzene	15.392	77	569823	10.685	ng	99
65) 4,6-Dinitro-2-methylph...	15.192	198	90017	9.095	ng	95
66) n-Nitrosodiphenylamine	15.315	169	473749	10.289	ng	98
67) 4-Bromophenyl-phenylether	15.992	248	160370	9.845	ng	95
68) Hexachlorobenzene	16.098	284	198141	10.157	ng	97
69) Atrazine	16.286	200	157170	11.344	ng	98
70) Pentachlorophenol	16.451	266	128691	9.411	ng	99
71) Phenanthrene	16.833	178	874873	10.313	ng	99
72) Anthracene	16.927	178	845520	10.243	ng	99
73) Carbazole	17.204	167	879612	10.437	ng	99
74) Di-n-butylphthalate	17.792	149	1003294	10.231	ng	100
75) Fluoranthene	18.868	202	1047590	10.152	ng	99
77) Benzidine	19.074	184	151348	5.602	ng	100
78) Pyrene	19.233	202	1113925	9.792	ng	99
80) Butylbenzylphthalate	20.174	149	439513	9.467	ng	99
81) Benzo(a)anthracene	21.009	228	1060395	10.026	ng	100
82) 3,3'-Dichlorobenzidine	20.950	252	311004	9.398	ng	99
83) Chrysene	21.068	228	1022315	10.188	ng	99
84) Bis(2-ethylhexyl)phtha...	20.974	149	643719	9.777	ng	100
85) Di-n-octyl phthalate	22.003	149	1048294	9.368	ng	98
87) Indeno(1,2,3-cd)pyrene	26.715	276	1216673	9.899	ng	99
88) Benzo(b)fluoranthene	22.880	252	1140578	10.100	ng	98
89) Benzo(k)fluoranthene	22.933	252	1078126	10.108	ng	99
90) Benzo(a)pyrene	23.597	252	965036	9.865	ng	99
91) Dibenzo(a,h)anthracene	26.756	278	999745	10.051	ng	99
92) Benzo(g,h,i)perylene	27.644	276	1029079	9.964	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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