

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080321\
 Data File : BP006476.D
 Acq On : 03 Aug 2021 12:29
 Operator : CG/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICCC040

Quant Time: Aug 03 13:20:45 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 03 13:16:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	176396	20.000	ng	0.00
21) Naphthalene-d8	10.546	136	761172	20.000	ng	# 0.00
39) Acenaphthene-d10	14.392	164	454713	20.000	ng	0.00
64) Phenanthrene-d10	17.139	188	921640	20.000	ng	# 0.00
76) Chrysene-d12	21.239	240	920772	20.000	ng	# 0.00
86) Perylene-d12	23.568	264	939503	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	945487	83.588	ng	0.00
7) Phenol-d6	6.934	99	1398114	87.550	ng	0.00
23) Nitrobenzene-d5	8.916	82	1371600	84.956	ng	0.00
42) 2,4,6-Tribromophenol	15.881	330	404150	86.923	ng	0.00
45) 2-Fluorobiphenyl	13.016	172	2445787	80.622	ng	0.00
79) Terphenyl-d14	19.716	244	3706828	81.062	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.305	88	197658	39.432	ng	90
3) Pyridine	3.693	79	615682	42.532	ng	94
4) n-Nitrosodimethylamine	3.611	42	230591	42.172	ng	# 81
6) Aniline	7.099	93	775046	44.400	ng	96
8) 2-Chlorophenol	7.322	128	523567	42.791	ng	90
9) Benzaldehyde	6.910	77	400798	40.049	ng	91
10) Phenol	6.963	94	703879	43.744	ng	97
11) bis(2-Chloroethyl)ether	7.199	93	524742	42.604	ng	89
12) 1,3-Dichlorobenzene	7.646	146	534365	41.382	ng	96
13) 1,4-Dichlorobenzene	7.793	146	541535	41.453	ng	96
14) 1,2-Dichlorobenzene	8.110	146	527920	42.055	ng	98
15) Benzyl Alcohol	8.004	79	537557	44.953	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.299	45	635978	43.754	ng	93
17) 2-Methylphenol	8.204	107	450567	44.756	ng	99
18) Hexachloroethane	8.828	117	213063	41.619	ng	87
19) n-Nitroso-di-n-propyla...	8.569	70	487781	46.068	ng	# 96
20) 3+4-Methylphenols	8.534	107	623714	45.427	ng	98
22) Acetophenone	8.581	105	842039	41.957	ng	# 96
24) Nitrobenzene	8.957	77	708178	42.259	ng	90
25) Isophorone	9.487	82	1262686	43.804	ng	94
26) 2-Nitrophenol	9.663	139	266025	44.345	ng	92
27) 2,4-Dimethylphenol	9.728	122	438958	42.760	ng	97
28) bis(2-Chloroethoxy)met...	9.969	93	658019	41.747	ng	98
29) 2,4-Dichlorophenol	10.193	162	451432	43.499	ng	94
30) 1,2,4-Trichlorobenzene	10.410	180	457819	40.619	ng	96
31) Naphthalene	10.593	128	1682874	41.502	ng	99
32) Benzoic acid	9.857	122	338267	43.398	ng	91
33) 4-Chloroaniline	10.710	127	704973	43.489	ng	95
34) Hexachlorobutadiene	10.881	225	264098	40.083	ng	98
35) Caprolactam	11.504	113	174758	45.930	ng	# 74
36) 4-Chloro-3-methylphenol	11.834	107	585598	45.235	ng	91
37) 2-Methylnaphthalene	12.204	142	1162765	42.721	ng	99
38) 1-Methylnaphthalene	12.428	142	1100497	42.570	ng	97
40) 1,2,4,5-Tetrachloroben...	12.575	216	475644	40.134	ng	# 96
41) Hexachlorocyclopentadiene	12.551	237	258796	41.202	ng	96
43) 2,4,6-Trichlorophenol	12.816	196	339624	42.729	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.887	196	376910	42.836	ng	# 88
46) 1,1'-Biphenyl	13.228	154	1392198	40.502	ng	96
47) 2-Chloronaphthalene	13.263	162	1075878	40.643	ng	95
48) 2-Nitroaniline	13.475	65	406350	45.772	ng	86
49) Acenaphthylene	14.110	152	1771386	42.121	ng	100
50) Dimethylphthalate	13.863	163	1375640	42.082	ng	99
51) 2,6-Dinitrotoluene	13.975	165	316093	44.440	ng	# 79
52) Acenaphthene	14.457	154	1076954	41.530	ng	99
53) 3-Nitroaniline	14.304	138	355171	45.053	ng	84
54) 2,4-Dinitrophenol	14.504	184	162000	41.959	ng	# 78
55) Dibenzofuran	14.792	168	1681120	41.389	ng	95
56) 4-Nitrophenol	14.610	139	316418	44.193	ng	85
57) 2,4-Dinitrotoluene	14.763	165	430842	45.377	ng	# 82
58) Fluorene	15.439	166	1360981	42.177	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.016	232	318533	43.160	ng	# 70
60) Diethylphthalate	15.228	149	1455518	42.822	ng	97
61) 4-Chlorophenyl-phenyle...	15.439	204	608907	41.528	ng	# 89
62) 4-Nitroaniline	15.469	138	385228	46.183	ng	89
63) Azobenzene	15.733	77	1724614	43.940	ng	92
65) 4,6-Dinitro-2-methylph...	15.522	198	233623	42.367	ng	73
66) n-Nitrosodiphenylamine	15.657	169	1188678	41.655	ng	99
67) 4-Bromophenyl-phenylether	16.339	248	357397	40.444	ng	# 88
68) Hexachlorobenzene	16.439	284	401770	40.347	ng	# 84
69) Atrazine	16.616	200	381663	43.474	ng	94
70) Pentachlorophenol	16.786	266	272543	44.029	ng	99
71) Phenanthrene	17.186	178	2120015	41.014	ng	99
72) Anthracene	17.275	178	2075056	41.899	ng	99
73) Carbazole	17.551	167	2154912	42.839	ng	98
74) Di-n-butylphthalate	18.116	149	2441613	42.985	ng	99
75) Fluoranthene	19.157	202	2430307	42.242	ng	94
77) Benzidine	19.345	184	1074357	47.405	ng	99
78) Pyrene	19.510	202	2529041	41.662	ng	99
80) Butylbenzylphthalate	20.404	149	1148750	43.800	ng	87
81) Benzo(a)anthracene	21.221	228	2471231	41.688	ng	100
82) 3,3'-Dichlorobenzidine	21.163	252	662760	42.183	ng	# 96
83) Chrysene	21.274	228	2387136	41.331	ng	99
84) Bis(2-ethylhexyl)phtha...	21.169	149	1693212	41.793	ng	# 97
85) Di-n-octyl phthalate	22.074	149	2812759	41.000	ng	# 93
87) Indeno(1,2,3-cd)pyrene	25.980	276	2837971	42.103	ng	# 88
88) Benzo(b)fluoranthene	22.863	252	2525183	41.869	ng	# 96
89) Benzo(k)fluoranthene	22.910	252	2456106	41.980	ng	# 95
90) Benzo(a)pyrene	23.468	252	2304977	42.690	ng	# 95
91) Dibenzo(a,h)anthracene	26.009	278	2448121	41.935	ng	# 94
92) Benzo(g,h,i)perylene	26.727	276	2348123	41.953	ng	# 89

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

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