

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081621\
 Data File : BM031748.D
 Acq On : 16 Aug 2021 13:06
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Manual Integrations
 APPROVED

mohammad
 8/17/2021 1:21:53 PM

Quant Time: Aug 16 14:22:02 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 14:20:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.551	152	36823	20.000	ng	0.00
21) Naphthalene-d8	10.316	136	165448	20.000	ng	# 0.00
39) Acenaphthene-d10	14.192	164	118775	20.000	ng	0.00
64) Phenanthrene-d10	16.945	188	267634	20.000	ng	0.00
76) Chrysene-d12	21.145	240	289156	20.000	ng	0.00
86) Perylene-d12	23.291	264	287756	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	92060	44.894	ng	0.00
7) Phenol-d6	6.745	99	132180	44.465	ng	0.00
23) Nitrobenzene-d5	8.704	82	157257	47.724	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	55935	36.064	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	343735	43.180	ng	0.00
79) Terphenyl-d14	19.586	244	638793	43.709	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	19544	23.718	ng	91
3) Pyridine	3.575	79	54693	23.463	ng	87
4) n-Nitrosodimethylamine	3.499	42	30086	22.709	ng	# 84
6) Aniline	6.898	93	73090	22.902	ng	97
8) 2-Chlorophenol	7.122	128	53815	22.295	ng	93
9) Benzaldehyde	6.716	77	47591m	26.868	ng	
10) Phenol	6.769	94	65912	22.886	ng	90
11) bis(2-Chloroethyl)ether	6.998	93	48768	22.874	ng	92
12) 1,3-Dichlorobenzene	7.440	146	58689	22.360	ng	# 92
13) 1,4-Dichlorobenzene	7.587	146	59375	22.229	ng	96
14) 1,2-Dichlorobenzene	7.898	146	58325	22.354	ng	94
15) Benzyl Alcohol	7.804	79	57893	23.134	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.087	45	71612	26.847	ng	97
17) 2-Methylphenol	7.998	107	46782	23.218	ng	95
18) Hexachloroethane	8.610	117	24409	23.675	ng	89
19) n-Nitroso-di-n-propyla...	8.357	70	52533	25.022	ng	96
20) 3+4-Methylphenols	8.328	107	65826	23.103	ng	93
22) Acetophenone	8.375	105	94233	23.008	ng	# 93
24) Nitrobenzene	8.745	77	82574	24.817	ng	95
25) Isophorone	9.269	82	143048	24.277	ng	98
26) 2-Nitrophenol	9.445	139	31811	22.705	ng	# 80
27) 2,4-Dimethylphenol	9.516	122	49117	22.251	ng	93
28) bis(2-Chloroethoxy)met...	9.751	93	70391	23.694	ng	99
29) 2,4-Dichlorophenol	9.969	162	54897	21.025	ng	93
30) 1,2,4-Trichlorobenzene	10.181	180	58665	20.461	ng	# 94
31) Naphthalene	10.363	128	189739	22.289	ng	99
32) Benzoic acid	9.645	122	44780	23.809	ng	96
33) 4-Chloroaniline	10.492	127	78762	21.447	ng	95
34) Hexachlorobutadiene	10.639	225	37156	20.181	ng	96
35) Caprolactam	11.280	113	20011	22.799	ng	92
36) 4-Chloro-3-methylphenol	11.616	107	69239	23.237	ng	89
37) 2-Methylnaphthalene	11.980	142	140547	21.419	ng	# 96
38) 1-Methylnaphthalene	12.204	142	133454	21.359	ng	# 98
40) 1,2,4,5-Tetrachloroben...	12.357	216	64808	19.584	ng	# 95
41) Hexachlorocyclopentadiene	12.327	237	33008	17.930	ng	97
43) 2,4,6-Trichlorophenol	12.610	196	47720	20.528	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.680	196	54026	20.991	ng	# 90
46) 1,1'-Biphenyl	13.016	154	182350	21.652	ng	100
47) 2-Chloronaphthalene	13.051	162	142108	21.543	ng	# 90
48) 2-Nitroaniline	13.274	65	53306	25.358	ng	86
49) Acenaphthylene	13.910	152	233792	22.173	ng	99
50) Dimethylphthalate	13.663	163	192364	21.593	ng	99
51) 2,6-Dinitrotoluene	13.786	165	42847	21.471	ng	# 73
52) Acenaphthene	14.257	154	141499	21.325	ng	99
53) 3-Nitroaniline	14.116	138	45355	22.496	ng	95
54) 2,4-Dinitrophenol	14.327	184	22661	19.254	ng	# 76
55) Dibenzofuran	14.592	168	238820	21.946	ng	92
56) 4-Nitrophenol	14.433	139	38149	21.397	ng	# 73
57) 2,4-Dinitrotoluene	14.580	165	61169	22.013	ng	# 60
58) Fluorene	15.245	166	193656	21.095	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.821	232	48042	20.031	ng	# 84
60) Diethylphthalate	15.039	149	212776	22.277	ng	97
61) 4-Chlorophenyl-phenyle...	15.245	204	94874	20.871	ng	# 84
62) 4-Nitroaniline	15.286	138	49355	22.192	ng	# 79
63) Azobenzene	15.539	77	238842	25.124	ng	91
65) 4,6-Dinitro-2-methylph...	15.345	198	34843	20.757	ng	# 67
66) n-Nitrosodiphenylamine	15.468	169	169442	21.539	ng	99
67) 4-Bromophenyl-phenylether	16.145	248	55802	20.152	ng	# 85
68) Hexachlorobenzene	16.245	284	61329	19.655	ng	# 88
69) Atrazine	16.433	200	61629	21.340	ng	99
70) Pentachlorophenol	16.598	266	41328	20.258	ng	96
71) Phenanthrene	16.986	178	314024	21.189	ng	99
72) Anthracene	17.074	178	313770	21.470	ng	99
73) Carbazole	17.357	167	313322	21.658	ng	98
74) Di-n-butylphthalate	17.927	149	390773	22.994	ng	# 98
75) Fluoranthene	19.009	202	382204	20.607	ng	99
77) Benzidine	19.209	184	171335	24.001	ng	99
78) Pyrene	19.374	202	397682	22.024	ng	100
80) Butylbenzylphthalate	20.292	149	190796	24.747	ng	89
81) Benzo(a)anthracene	21.127	228	410888	21.442	ng	98
82) 3,3'-Dichlorobenzidine	21.068	252	124820	23.761	ng	98
83) Chrysene	21.180	228	399063	21.435	ng	98
84) Bis(2-ethylhexyl)phtha...	21.074	149	294330	24.641	ng	94
85) Di-n-octyl phthalate	21.933	149	500786	24.484	ng	# 94
87) Indeno(1,2,3-cd)pyrene	25.421	276	408986	19.890	ng	98
88) Benzo(b)fluoranthene	22.656	252	404632	21.221	ng	99
89) Benzo(k)fluoranthene	22.697	252	397734m	22.303	ng	
90) Benzo(a)pyrene	23.203	252	369412	21.522	ng	98
91) Dibenzo(a,h)anthracene	25.432	278	359113	20.175	ng	99
92) Benzo(g,h,i)perylene	26.068	276	336692	19.853	ng	98

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

