

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072021\
 Data File : BM031016.D
 Acq On : 20 Jul 2021 10:10
 Operator : CG/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Quant Time: Jul 20 11:24:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM072021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 20 11:22:19 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	79603	20.000	ng	0.00
21) Naphthalene-d8	10.404	136	353183	20.000	ng	# 0.00
39) Acenaphthene-d10	14.269	164	259863	20.000	ng	0.00
64) Phenanthrene-d10	17.016	188	595601	20.000	ng	0.00
76) Chrysene-d12	21.203	240	702846	20.000	ng	0.00
86) Perylene-d12	23.386	264	741928	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.257	112	187098	41.889	ng	0.00
7) Phenol-d6	6.816	99	265705	41.369	ng	0.00
23) Nitrobenzene-d5	8.787	82	290454	41.178	ng	0.00
42) 2,4,6-Tribromophenol	15.763	330	136263	39.956	ng	0.00
45) 2-Fluorobiphenyl	12.886	172	708672	40.147	ng	0.00
79) Terphenyl-d14	19.657	244	1442035	39.936	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.246	88	38863	21.880	ng	94
3) Pyridine	3.622	79	104190	20.505	ng	# 87
4) n-Nitrosodimethylamine	3.546	42	60613	21.054	ng	91
6) Aniline	6.975	93	141588	20.501	ng	97
8) 2-Chlorophenol	7.210	128	110152	20.917	ng	95
9) Benzaldehyde	6.793	77	84489	22.315	ng	99
10) Phenol	6.846	94	129730	20.751	ng	92
11) bis(2-Chloroethyl)ether	7.081	93	96845	20.808	ng	96
12) 1,3-Dichlorobenzene	7.528	146	118493	20.863	ng	94
13) 1,4-Dichlorobenzene	7.675	146	120013	20.695	ng	96
14) 1,2-Dichlorobenzene	7.987	146	117861	20.846	ng	97
15) Benzyl Alcohol	7.881	79	112034	20.547	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.163	45	122680	20.981	ng	97
17) 2-Methylphenol	8.081	107	90549	20.359	ng	94
18) Hexachloroethane	8.704	117	46860	20.725	ng	92
19) n-Nitroso-di-n-propyla...	8.445	70	94103	20.322	ng	98
20) 3+4-Methylphenols	8.404	107	128621	20.729	ng	98
22) Acetophenone	8.457	105	180874	20.486	ng	# 95
24) Nitrobenzene	8.828	77	146841	20.582	ng	98
25) Isophorone	9.351	82	258835	20.369	ng	99
26) 2-Nitrophenol	9.534	139	60718	20.036	ng	96
27) 2,4-Dimethylphenol	9.598	122	98257	20.445	ng	98
28) bis(2-Chloroethoxy)met...	9.834	93	130069	20.093	ng	100
29) 2,4-Dichlorophenol	10.057	162	115940	20.585	ng	94
30) 1,2,4-Trichlorobenzene	10.275	180	126155	20.622	ng	97
31) Naphthalene	10.457	128	369995	20.297	ng	99
32) Benzoic acid	9.716	122	80874	20.722	ng	99
33) 4-Chloroaniline	10.569	127	160951	20.611	ng	98
34) Hexachlorobutadiene	10.739	225	81167	20.329	ng	98
35) Caprolactam	11.351	113	38546	20.705	ng	99
36) 4-Chloro-3-methylphenol	11.698	107	132077	20.561	ng	94
37) 2-Methylnaphthalene	12.069	142	285306	20.392	ng	95
38) 1-Methylnaphthalene	12.292	142	274114	20.514	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.445	216	147473	19.967	ng	# 97
41) Hexachlorocyclopentadiene	12.422	237	84155	20.168	ng	99
43) 2,4,6-Trichlorophenol	12.692	196	104734	20.297	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.757	196	114294	20.208	ng	# 92
46) 1,1'-Biphenyl	13.098	154	374406	20.242	ng	99
47) 2-Chloronaphthalene	13.139	162	294831	20.096	ng	96
48) 2-Nitroaniline	13.351	65	93649	20.361	ng	99
49) Acenaphthylene	13.986	152	476961	20.471	ng	99
50) Dimethylphthalate	13.739	163	398692	20.471	ng	99
51) 2,6-Dinitrotoluene	13.857	165	88562	20.440	ng	87
52) Acenaphthene	14.333	154	292199	20.093	ng	99
53) 3-Nitroaniline	14.180	138	90817	20.740	ng	100
54) 2,4-Dinitrophenol	14.386	184	50880	20.203	ng	94
55) Dibenzofuran	14.669	168	484581	20.285	ng	93
56) 4-Nitrophenol	14.492	139	80922	21.034	ng	87
57) 2,4-Dinitrotoluene	14.639	165	123915	20.395	ng	# 82
58) Fluorene	15.322	166	411195	20.419	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.898	232	106954	20.261	ng	# 91
60) Diethylphthalate	15.116	149	429936	20.549	ng	98
61) 4-Chlorophenyl-phenylether	15.322	204	201342	19.873	ng	# 84
62) 4-Nitroaniline	15.351	138	100731	21.045	ng	89
63) Azobenzene	15.616	77	434474	21.032	ng	94
65) 4,6-Dinitro-2-methylph...	15.404	198	76343	20.430	ng	92
66) n-Nitrosodiphenylamine	15.539	169	360369	20.284	ng	99
67) 4-Bromophenyl-phenylether	16.221	248	126797	20.202	ng	95
68) Hexachlorobenzene	16.321	284	142748	20.327	ng	95
69) Atrazine	16.498	200	133118	20.656	ng	96
70) Pentachlorophenol	16.668	266	93726	20.453	ng	97
71) Phenanthrene	17.057	178	677146	20.562	ng	100
72) Anthracene	17.151	178	665154	20.473	ng	100
73) Carbazole	17.421	167	664573	20.865	ng	98
74) Di-n-butylphthalate	17.998	149	780302	20.562	ng	99
75) Fluoranthene	19.074	202	847125	20.865	ng	100
77) Benzidine	19.274	184	331666	18.143	ng	100
78) Pyrene	19.439	202	888368	19.862	ng	98
80) Butylbenzylphthalate	20.362	149	389386	20.332	ng	98
81) Benzo(a)anthracene	21.192	228	947652	20.388	ng	98
82) 3,3'-Dichlorobenzidine	21.133	252	260817	20.591	ng	99
83) Chrysene	21.245	228	922008	20.391	ng	98
84) Bis(2-ethylhexyl)phtha...	21.139	149	599873	20.376	ng	94
85) Di-n-octyl phthalate	22.009	149	1017689	20.430	ng	98
87) Indeno(1,2,3-cd)pyrene	25.556	276	1031337	19.982	ng	99
88) Benzo(b)fluoranthene	22.739	252	977799	19.790	ng	99
89) Benzo(k)fluoranthene	22.780	252	945147	20.069	ng	99
90) Benzo(a)pyrene	23.292	252	894463	20.065	ng	100
91) Dibenzo(a,h)anthracene	25.568	278	883849	19.819	ng	100
92) Benzo(g,h,i)perylene	26.215	276	847293	20.027	ng	98

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

