

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022222\
 Data File : BM034041.D
 Acq On : 22 Feb 2022 16:15
 Operator : CG/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/23/2022
 Supervised By : Jagrut Upadhyay 02/23/2022

Quant Time: Feb 22 16:53:10 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM022222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 22 16:00:49 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.025	152	267394	20.000	ng	0.00
21) Naphthalene-d8	10.836	136	969151	20.000	ng	# 0.00
39) Acenaphthene-d10	14.654	164	633067	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	1319774	20.000	ng	# 0.00
76) Chrysene-d12	21.565	240	1372863	20.000	ng	0.00
86) Perylene-d12	24.030	264	1401076	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.566	112	1734204	105.014	ng	0.00
7) Phenol-d6	7.172	99	2204490	109.504	ng	0.00
23) Nitrobenzene-d5	9.189	82	2163516	104.858	ng	0.00
42) 2,4,6-Tribromophenol	16.136	330	989241	141.162	ng	0.00
45) 2-Fluorobiphenyl	13.289	172	5564797	115.410	ng	0.00
79) Terphenyl-d14	20.012	244	8564935	123.387	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.431	88	328313	49.164	ng	# 87
3) Pyridine	3.837	79	912577m	45.624	ng	
4) n-Nitrosodimethylamine	3.743	42	448961	62.228	ng	# 73
6) Aniline	7.342	93	1202555	52.780	ng	91
8) 2-Chlorophenol	7.584	128	945431	57.528	ng	84
10) Phenol	7.201	94	1077378	51.907	ng	91
11) bis(2-Chloroethyl)ether	7.437	93	845682	49.933	ng	91
12) 1,3-Dichlorobenzene	7.913	146	1116983	56.983	ng	# 93
13) 1,4-Dichlorobenzene	8.060	146	1138059	60.142	ng	93
14) 1,2-Dichlorobenzene	8.378	146	1096981	57.611	ng	95
15) Benzyl Alcohol	8.260	79	831258	53.601	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.560	45	1081849	59.278	ng	94
17) 2-Methylphenol	8.460	107	754769	55.735	ng	94
18) Hexachloroethane	9.119	117	398869	55.765	ng	# 88
19) n-Nitroso-di-n-propyla...	8.836	70	670813	51.641	ng	# 92
20) 3+4-Methylphenols	8.795	107	1028199	55.049	ng	93
22) Acetophenone	8.848	105	1360439	56.131	ng	# 90
24) Nitrobenzene	9.231	77	991957	53.820	ng	# 88
25) Isophorone	9.760	82	1681490	55.821	ng	# 94
26) 2-Nitrophenol	9.942	139	502945	64.419	ng	# 78
27) 2,4-Dimethylphenol	10.001	122	747017	54.705	ng	91
28) bis(2-Chloroethoxy)met...	10.242	93	1115688	56.037	ng	# 97
29) 2,4-Dichlorophenol	10.478	162	947350	66.196	ng	96
30) 1,2,4-Trichlorobenzene	10.701	180	1102617	63.414	ng	96
31) Naphthalene	10.889	128	2898460	57.533	ng	99
32) Benzoic acid	10.148	122	609030	66.916	ng	# 82
33) 4-Chloroaniline	10.989	127	1189755	60.300	ng	# 89
34) Hexachlorobutadiene	11.189	225	699701	63.219	ng	96
35) Caprolactam	11.772	113	198935	49.844	ng	# 70
36) 4-Chloro-3-methylphenol	12.107	107	903201	57.866	ng	86
37) 2-Methylnaphthalene	12.495	142	2103330	65.447	ng	97
38) 1-Methylnaphthalene	12.713	142	2040259	63.579	ng	98
40) 1,2,4,5-Tetrachloroben...	12.860	216	1203585	57.031	ng	# 97
41) Hexachlorocyclopentadiene	12.848	237	747872	67.898	ng	98
43) 2,4,6-Trichlorophenol	13.095	196	763536	62.879	ng	98
44) 2,4,5-Trichlorophenol	13.160	196	768655	56.268	ng	# 91

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46) 1,1'-Biphenyl	13.495	154	2827958	58.343	ng	98
47) 2-Chloronaphthalene	13.536	162	2221825	57.107	ng	95
48) 2-Nitroaniline	13.730	65	578008	54.345	ng	# 76
49) Acenaphthylene	14.377	152	3375555	63.431	ng	99
50) Dimethylphthalate	14.118	163	2768435	61.028	ng	99
51) 2,6-Dinitrotoluene	14.224	165	581458	61.810	ng	# 80
52) Acenaphthene	14.718	154	2233785	61.209	ng	98
53) 3-Nitroaniline	14.548	138	593694	63.125	ng	82
54) 2,4-Dinitrophenol	14.754	184	310078	61.799	ng	# 76
55) Dibenzofuran	15.054	168	3425601	64.906	ng	97
56) 4-Nitrophenol	14.842	139	502280	61.459	ng	# 80
57) 2,4-Dinitrotoluene	15.007	165	788486	67.517	ng	# 75
58) Fluorene	15.701	166	2800989	60.308	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.271	232	779000	69.950	ng	# 91
60) Diethylphthalate	15.477	149	2817290	57.996	ng	99
61) 4-Chlorophenyl-phenyle...	15.695	204	1488039	58.259	ng	94
62) 4-Nitroaniline	15.707	138	649818	66.925	ng	# 82
63) Azobenzene	15.989	77	2426927	53.202	ng	90
65) 4,6-Dinitro-2-methylph...	15.771	198	497385	65.818	ng	# 70
66) n-Nitrosodiphenylamine	15.907	169	2406704	58.556	ng	98
67) 4-Bromophenyl-phenylether	16.589	248	893704	64.380	ng	92
68) Hexachlorobenzene	16.707	284	1015137	58.516	ng	# 90
69) Atrazine	16.854	200	811539	65.342	ng	97
70) Pentachlorophenol	17.042	266	661719	62.026	ng	99
71) Phenanthrene	17.436	178	4304480	56.344	ng	99
72) Anthracene	17.530	178	4261833	57.007	ng	99
73) Carbazole	17.795	167	4076026	53.719	ng	99
74) Di-n-butylphthalate	18.371	149	4846916	59.281	ng	99
75) Fluoranthene	19.448	202	5010888	66.296	ng	97
77) Benzidine	19.624	184	1544482	62.977	ng	99
78) Pyrene	19.806	202	5160244	58.936	ng	98
80) Butylbenzylphthalate	20.706	149	2218928	44.137	ng	86
81) Benzo(a)anthracene	21.553	228	5196275	54.864	ng	97
82) 3,3'-Dichlorobenzidine	21.477	252	1877687	73.831	ng	98
83) Chrysene	21.606	228	4955975	48.133	ng	99
84) Bis(2-ethylhexyl)phtha...	21.489	149	3463749	85.916	ng	# 97
85) Di-n-octyl phthalate	22.436	149	5623935	50.552	ng	# 91
87) Indeno(1,2,3-cd)pyrene	26.618	276	6093032	59.793	ng	97
88) Benzo(b)fluoranthene	23.277	252	5228472	54.825	ng	98
89) Benzo(k)fluoranthene	23.330	252	5042413	63.828	ng	98
90) Benzo(a)pyrene	23.924	252	4345138	59.524	ng	99
91) Dibenzo(a,h)anthracene	26.641	278	5121810	61.170	ng	97
92) Benzo(g,h,i)perylene	27.412	276	4964794	59.268	ng	99

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

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