

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100121\
 Data File : BM032303.D
 Acq On : 30 Sep 2021 20:05
 Operator : CG/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

mohammad
 10/1/2021 11:31:45 AM

Quant Time: Oct 01 01:05:53 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM100121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 30 19:01:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.643	152	293669	20.000	ng	0.00
21) Naphthalene-d8	10.436	136	1196323	20.000	ng	# 0.00
39) Acenaphthene-d10	14.295	164	783108	20.000	ng	0.00
64) Phenanthrene-d10	17.018	188	1661658	20.000	ng	# 0.00
76) Chrysene-d12	21.189	240	1595205	20.000	ng	0.00
86) Perylene-d12	23.342	264	1849068	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.213	112	2394340	145.485	ng	0.00
7) Phenol-d6	6.837	99	3498218	139.720	ng	0.00
23) Nitrobenzene-d5	8.807	82	3167569	118.001	ng	0.00
42) 2,4,6-Tribromophenol	15.777	330	1296197	208.274	ng	0.00
45) 2-Fluorobiphenyl	12.913	172	6391630	107.016	ng	0.00
79) Terphenyl-d14	0.000	244	0d	0.000	ng	
Target Compounds						Qvalue
2) 1,4-Dioxane	3.008	88	486087	65.258	ng	90
3) Pyridine	3.419	79	1490769	76.760	ng	94
4) n-Nitrosodimethylamine	3.331	42	574332	56.653	ng	# 62
6) Aniline	6.972	93	2049590	73.995	ng	91
8) 2-Chlorophenol	7.219	128	1376585	74.887	ng	92
10) Phenol	6.866	94	1708154	67.270	ng	82
11) bis(2-Chloroethyl)ether	7.066	93	1332960	69.552	ng	93
12) 1,3-Dichlorobenzene	7.537	146	1483917	65.271	ng	# 91
13) 1,4-Dichlorobenzene	7.678	146	1514575	65.797	ng	98
14) 1,2-Dichlorobenzene	8.001	146	1472164	65.382	ng	97
15) Benzyl Alcohol	7.890	79	1292765	65.626	ng	87
16) 2,2'-oxybis(1-Chloropr...	8.178	45	1701403	61.666	ng	97
17) 2-Methylphenol	8.107	107	1196229	70.359	ng	# 90
18) Hexachloroethane	8.731	117	529937	61.165	ng	94
19) n-Nitroso-di-n-propyla...	8.460	70	998905	54.538	ng	# 82
20) 3+4-Methylphenols	8.442	107	1625266	72.153	ng	94
22) Acetophenone	8.466	105	2060715	59.780	ng	# 87
24) Nitrobenzene	8.842	77	1534264	53.631	ng	# 88
25) Isophorone	9.366	82	2821811	56.600	ng	95
26) 2-Nitrophenol	9.554	139	791804	144.212	ng	# 73
27) 2,4-Dimethylphenol	9.631	122	1213718	73.035	ng	94
28) bis(2-Chloroethoxy)met...	9.848	93	1870521	74.922	ng	99
29) 2,4-Dichlorophenol	10.095	162	1381475	76.573	ng	97
30) 1,2,4-Trichlorobenzene	10.307	180	1454745	59.334	ng	98
31) Naphthalene	10.484	128	4371679	65.332	ng	99
32) Benzoic acid	9.831	122	1092219	189.891	ng	# 81
33) 4-Chloroaniline	10.595	127	1948341	75.707	ng	# 93
34) Hexachlorobutadiene	10.789	225	852444	51.801	ng	98
35) Caprolactam	11.383	113	515029	103.928	ng	# 79
36) 4-Chloro-3-methylphenol	11.742	107	1545372	73.019	ng	95
37) 2-Methylnaphthalene	12.101	142	3056173m	63.042	ng	
38) 1-Methylnaphthalene	12.325	142	2991058	65.253	ng	# 98
40) 1,2,4,5-Tetrachloroben...	12.483	216	1509543	57.613	ng	# 96
41) Hexachlorocyclopentadiene	12.472	237	776684	96.136	ng	97
43) 2,4,6-Trichlorophenol	12.725	196	1160076	109.053	ng	96
44) 2,4,5-Trichlorophenol	12.801	196	1261972	97.440	ng	# 90

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.119	154	3931868	65.820	ng	99
47) 2-Chloronaphthalene	13.160	162	3108826	64.170	ng	98
48) 2-Nitroaniline	13.366	65	1001599	81.421	ng #	78
49) Acenaphthylene	14.013	152	5000192	68.835	ng	99
50) Dimethylphthalate	13.748	163	4001994	70.379	ng	99
51) 2,6-Dinitrotoluene	13.860	165	892170	77.241	ng #	63
52) Acenaphthene	14.360	154	3271707	72.477	ng	98
53) 3-Nitroaniline	14.195	138	1052412	104.069	ng	85
54) 2,4-Dinitrophenol	14.389	184	586787	179.354	ng #	65
55) Dibenzofuran	14.689	168	4798433	63.760	ng	94
56) 4-Nitrophenol	14.519	139	906126	160.318	ng #	65
57) 2,4-Dinitrotoluene	14.648	165	1240175	87.481	ng #	61
58) Fluorene	15.336	166	3380718	57.175	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.924	232	1097056	115.889	ng #	83
60) Diethylphthalate	15.113	149	4005227	72.542	ng	95
61) 4-Chlorophenyl-phenyle...	15.324	204	1735953	54.659	ng	94
62) 4-Nitroaniline	15.366	138	1100408	124.049	ng #	69
63) Azobenzene	15.618	77	3810122	59.020	ng	85
65) 4,6-Dinitro-2-methylph...	15.419	198	819648	147.573	ng #	72
66) n-Nitrosodiphenylamine	15.548	169	3350773	61.616	ng	98
67) 4-Bromophenyl-phenylether	16.213	248	1201279	61.484	ng	94
68) Hexachlorobenzene	16.354	284	1314563	59.502	ng #	81
69) Atrazine	16.489	200	1236790	74.136	ng	95
70) Pentachlorophenol	16.683	266	925542	127.769	ng	98
71) Phenanthrene	17.060	178	5876718	61.822	ng	99
72) Anthracene	17.154	178	5854174	62.999	ng	98
73) Carbazole	17.418	167	5963385	69.873	ng	98
74) Di-n-butylphthalate	17.971	149	6586224	74.245	ng #	97
75) Fluoranthene	19.071	202	6758358	63.803	ng	96
77) Benzidine	19.248	184	3143506	106.988	ng	97
78) Pyrene	19.430	202	6954307	57.223	ng	96
80) Butylbenzylphthalate	20.318	149	3196796	89.764	ng	91
81) Benzo(a)anthracene	21.171	228	6800203	63.641	ng	96
82) 3,3'-Dichlorobenzidine	21.106	252	2296458	78.061	ng	99
83) Chrysene	21.224	228	6429117	56.683	ng	97
84) Bis(2-ethylhexyl)phtha...	21.089	149	4017990	73.941	ng #	94
85) Di-n-octyl phthalate	21.942	149	7577255	82.634	ng #	90
87) Indeno(1,2,3-cd)pyrene	25.506	276	8540789	68.111	ng	97
88) Benzo(b)fluoranthene	22.706	252	7125533	57.601	ng	99
89) Benzo(k)fluoranthene	22.753	252	7188303	53.744	ng	97
90) Benzo(a)pyrene	23.253	252	6443019	56.709	ng	97
91) Dibenzo(a,h)anthracene	25.512	278	6906729	63.882	ng #	95
92) Benzo(g,h,i)perylene	26.171	276	7738805	72.286	ng #	95

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

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