

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG082321\
 Data File : BG049939.D
 Acq On : 23 Aug 2021 18:52
 Operator : CG/JU
 Sample : M3505-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DRILL-CUTTINGMSD

Manual Integrations
 APPROVED

mohammad
 8/24/2021 11:18:58 AM

Quant Time: Aug 24 02:27:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 16:13:24 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.978	152	28244	20.000	ng	0.00
21) Naphthalene-d8	10.798	136	116550	20.000	ng	#-0.01
39) Acenaphthene-d10	14.640	164	82137	20.000	ng	0.00
64) Phenanthrene-d10	17.395	188	178535	20.000	ng	# 0.00
76) Chrysene-d12	21.672	240	145236	20.000	ng	0.00
86) Perylene-d12	24.897	264	148219	20.000	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.535	112	179869	114.160	ng	0.00
7) Phenol-d6	7.156	99	255238	107.017	ng	0.00
23) Nitrobenzene-d5	9.153	82	171212	65.005	ng	0.00
42) 2,4,6-Tribromophenol	16.138	330	132315	109.595	ng	0.00
45) 2-Fluorobiphenyl	13.254	172	369306	65.423	ng	0.00
79) Terphenyl-d14	20.003	244	521302	65.190	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.367	88	22131	32.143	ng	98
3) Pyridine	3.784	79	54323	28.783	ng	# 83
4) n-Nitrosodimethylamine	3.702	42	40495	39.903	ng	# 72
6) Aniline	7.297	93	101692	40.231	ng	99
8) 2-Chlorophenol	7.549	128	83043	48.342	ng	98
9) Benzaldehyde	7.115	77	59517	37.399	ng	# 85
10) Phenol	7.185	94	108357	46.888	ng	91
11) bis(2-Chloroethyl)ether	7.397	93	68783	44.082	ng	94
12) 1,3-Dichlorobenzene	7.867	146	85054	43.232	ng	96
13) 1,4-Dichlorobenzene	8.014	146	88054	44.216	ng	94
14) 1,2-Dichlorobenzene	8.337	146	83763	44.399	ng	96
15) Benzyl Alcohol	8.225	79	93590	46.725	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.507	45	75465	42.564	ng	97
17) 2-Methylphenol	8.437	107	73914	48.570	ng	97
18) Hexachloroethane	9.059	117	34117	44.566	ng	89
19) n-Nitroso-di-n-propyla...	8.789	70	65600	41.102	ng	90
20) 3+4-Methylphenols	8.771	107	102423	47.923	ng	97
22) Acetophenone	8.812	105	131707	41.595	ng	# 93
24) Nitrobenzene	9.194	77	106984	38.494	ng	95
25) Isophorone	9.717	82	174247	37.075	ng	98
26) 2-Nitrophenol	9.911	139	47177	44.480	ng	94
27) 2,4-Dimethylphenol	9.976	122	78233	49.905	ng	97
28) bis(2-Chloroethoxy)met...	10.199	93	96520	46.947	ng	# 91
29) 2,4-Dichlorophenol	10.457	162	87040	45.246	ng	98
30) 1,2,4-Trichlorobenzene	10.663	180	91840	43.359	ng	91
31) Naphthalene	10.851	128	252261	41.328	ng	99
32) Benzoic acid	10.146	122	40569m	37.818	ng	
33) 4-Chloroaniline	10.962	127	73907	30.669	ng	96
34) Hexachlorobutadiene	11.127	225	64060	41.618	ng	92
35) Caprolactam	11.756	113	26422m	44.453	ng	
36) 4-Chloro-3-methylphenol	12.102	107	101223	45.553	ng	# 92
37) 2-Methylnaphthalene	12.460	142	190073	41.337	ng	96
38) 1-Methylnaphthalene	12.678	142	178074	41.206	ng	92
40) 1,2,4,5-Tetrachloroben...	12.831	216	108757	44.938	ng	99
41) Hexachlorocyclopentadiene	12.801	237	75206	58.714	ng	93
43) 2,4,6-Trichlorophenol	13.077	196	71954	44.101	ng	95

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44) 2,4,5-Trichlorophenol	13.160	196	78286	44.228	ng	93
46) 1,1'-Biphenyl	13.471	154	242860	41.514	ng	98
47) 2-Chloronaphthalene	13.518	162	200654	42.944	ng	96
48) 2-Nitroaniline	13.723	65	71014	41.391	ng	92
49) Acenaphthylene	14.364	152	312433	42.188	ng	98
50) Dimethylphthalate	14.094	163	258790	41.760	ng	99
51) 2,6-Dinitrotoluene	14.217	165	57375	42.050	ng	86
52) Acenaphthene	14.704	154	181146	40.605	ng	94
53) 3-Nitroaniline	14.546	138	46152	34.685	ng	99
54) 2,4-Dinitrophenol	14.769	184	68442	88.356	ng	94
55) Dibenzofuran	15.039	168	297559	41.115	ng	96
56) 4-Nitrophenol	14.887	139	74466	76.603	ng	94
57) 2,4-Dinitrotoluene	15.010	165	80079	41.786	ng	# 93
58) Fluorene	15.691	166	247082	42.004	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.274	232	66236	40.290	ng	98
60) Diethylphthalate	15.451	149	261955	42.029	ng	96
61) 4-Chlorophenyl-phenyle...	15.674	204	133584	42.548	ng	93
62) 4-Nitroaniline	15.721	138	57992	42.370	ng	90
63) Azobenzene	15.973	77	251607	38.140	ng	88
65) 4,6-Dinitro-2-methylph...	15.785	198	45638	39.769	ng	94
66) n-Nitrosodiphenylamine	15.897	169	216468	42.911	ng	97
67) 4-Bromophenyl-phenylether	16.573	248	91270	43.407	ng	97
68) Hexachlorobenzene	16.702	284	99202	42.465	ng	96
69) Atrazine	16.849	200	93664	46.416	ng	98
70) Pentachlorophenol	17.054	266	83058	66.306	ng	98
71) Phenanthrene	17.436	178	394601	42.010	ng	98
72) Anthracene	17.530	178	396716	43.005	ng	98
73) Carbazole	17.800	167	354350	40.558	ng	98
74) Di-n-butylphthalate	18.347	149	424260	41.656	ng	98
75) Fluoranthene	19.451	202	449858	38.920	ng	98
77) Benzidine	19.627	184	238982	64.221	ng	97
78) Pyrene	19.809	202	439511	44.389	ng	98
80) Butylbenzylphthalate	20.685	149	165590	43.912	ng	96
81) Benzo(a)anthracene	21.648	228	394101	41.667	ng	97
82) 3,3'-Dichlorobenzidine	21.560	252	113724	41.180	ng	96
83) Chrysene	21.719	228	372654	41.217	ng	97
84) Bis(2-ethylhexyl)phtha...	21.536	149	226328	42.401	ng	96
85) Di-n-octyl phthalate	22.747	149	368703	40.888	ng	100
87) Indeno(1,2,3-cd)pyrene	28.615	276	458950	42.935	ng	98
88) Benzo(b)fluoranthene	23.874	252	399335	41.080	ng	95
89) Benzo(k)fluoranthene	23.945	252	382415	42.175	ng	99
90) Benzo(a)pyrene	24.756	252	382304	43.457	ng	# 98
91) Dibenzo(a,h)anthracene	28.674	278	390219	43.322	ng	99
92) Benzo(g,h,i)perylene	29.778	276	386333	43.779	ng	95

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

