

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG041423\
 Data File : BG057100.D
 Acq On : 14 Apr 2023 14:57
 Operator : CG/JU
 Sample : 02320-01MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E9074-BAYS-001MSD

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo

Quant Time: Apr 14 15:42:24 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 05:15:23 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.402	152	19583	20.000	ng	0.00
21) Naphthalene-d8	11.245	136	78587	20.000	ng	# 0.00
39) Acenaphthene-d10	15.017	164	50903	20.000	ng	0.00
64) Phenanthrene-d10	17.754	188	103634	20.000	ng	# 0.00
76) Chrysene-d12	22.072	240	89201	20.000	ng	# 0.00
86) Perylene-d12	25.608	264	98464	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.917	112	158254	129.335	ng	0.00
7) Phenol-d6	7.539	99	220800	120.653	ng	0.00
23) Nitrobenzene-d5	9.583	82	143419	74.673	ng	0.00
42) 2,4,6-Tribromophenol	16.497	330	95037	116.869	ng	0.00
45) 2-Fluorobiphenyl	13.642	172	288403	76.828	ng	0.00
79) Terphenyl-d14	20.321	244	424150	80.462	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.720	88	20674	38.543	ng	100
3) Pyridine	4.143	79	59289	34.318	ng	96
4) n-Nitrosodimethylamine	4.049	42	31659	39.858	ng	87
6) Aniline	7.715	93	84890	36.285	ng	97
8) 2-Chlorophenol	7.962	128	67004	54.977	ng	92
9) Benzaldehyde	7.521	77	50554	44.043	ng	92
10) Phenol	7.568	94	93990	51.833	ng	96
11) bis(2-Chloroethyl)ether	7.803	93	65038	50.352	ng	95
12) 1,3-Dichlorobenzene	8.291	146	70688	52.491	ng	98
13) 1,4-Dichlorobenzene	8.437	146	72632	53.439	ng	100
14) 1,2-Dichlorobenzene	8.761	146	67488	51.395	ng	96
15) Benzyl Alcohol	8.637	79	69718	45.849	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.919	45	82169	52.784	ng	99
17) 2-Methylphenol	8.837	107	63824	49.334	ng	89
18) Hexachloroethane	9.501	117	26221	52.782	ng	# 91
19) n-Nitroso-di-n-propyla...	9.207	70	56932	43.494	ng	97
20) 3+4-Methylphenols	9.166	107	85152	47.071	ng	93
22) Acetophenone	9.230	105	109132	46.668	ng	# 97
24) Nitrobenzene	9.624	77	88976	45.312	ng	95
25) Isophorone	10.147	82	151124	41.321	ng	99
26) 2-Nitrophenol	10.341	139	37944	50.063	ng	90
27) 2,4-Dimethylphenol	10.382	122	65252	48.761	ng	96
28) bis(2-Chloroethoxy)met...	10.617	93	83584	45.649	ng	93
29) 2,4-Dichlorophenol	10.881	162	66892	46.479	ng	94
30) 1,2,4-Trichlorobenzene	11.099	180	76112	49.774	ng	98
31) Naphthalene	11.292	128	202237	47.213	ng	98
32) Benzoic acid	10.535	122	45248m	49.224	ng	
33) 4-Chloroaniline	11.392	127	66628	34.440	ng	97
34) Hexachlorobutadiene	11.568	225	52319	45.547	ng	89
35) Caprolactam	12.162	113	20289m	40.879	ng	
36) 4-Chloro-3-methylphenol	12.485	107	70893	44.048	ng	97
37) 2-Methylnaphthalene	12.867	142	144912	42.625	ng	97
38) 1-Methylnaphthalene	13.084	142	134899	42.258	ng	99
40) 1,2,4,5-Tetrachloroben...	13.231	216	93091	52.738	ng	99
41) Hexachlorocyclopentadiene	13.202	237	68466	70.397	ng	98
43) 2,4,6-Trichlorophenol	13.460	196	58472	51.910	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.536	196	62809	47.068	ng	96	04/17/2023
46) 1,1'-Biphenyl	13.854	154	196680	52.783	ng	99	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.906	162	147816	53.636	ng	97	
48) 2-Nitroaniline	14.094	65	47833	48.540	ng	96	
49) Acenaphthylene	14.741	152	225715	49.848	ng	98	
50) Dimethylphthalate	14.453	163	184480	45.996	ng	99	
51) 2,6-Dinitrotoluene	14.582	165	41090	47.417	ng	93	04/17/2023
52) Acenaphthene	15.081	154	163521m	52.362	ng		
53) 3-Nitroaniline	14.911	138	35402	41.273	ng	91	
54) 2,4-Dinitrophenol	15.117	184	36422	74.672	ng	95	
55) Dibenzofuran	15.410	168	226731	48.374	ng	97	
56) 4-Nitrophenol	15.211	139	65253	102.878	ng	# 83	
57) 2,4-Dinitrotoluene	15.363	165	56994	46.632	ng	95	
58) Fluorene	16.057	166	177556	45.780	ng	99	
59) 2,3,4,6-Tetrachlorophenol	15.634	232	56662	46.636	ng	# 97	
60) Diethylphthalate	15.798	149	183787	45.709	ng	99	
61) 4-Chlorophenyl-phenyle...	16.033	204	103263	45.347	ng	97	
62) 4-Nitroaniline	16.074	138	39413	44.349	ng	# 84	
63) Azobenzene	16.327	77	188434	47.025	ng	96	
65) 4,6-Dinitro-2-methylph...	16.127	198	27821	44.326	ng	95	
66) n-Nitrosodiphenylamine	16.250	169	154619	55.579	ng	98	
67) 4-Bromophenyl-phenylether	16.932	248	64502	50.303	ng	93	
68) Hexachlorobenzene	17.061	284	69674	54.959	ng	97	
69) Atrazine	17.184	200	63509	54.995	ng	96	
70) Pentachlorophenol	17.402	266	85639	94.161	ng	94	
71) Phenanthrene	17.801	178	267844	50.692	ng	99	
72) Anthracene	17.895	178	286019	52.805	ng	100	
73) Carbazole	18.154	167	237017	50.189	ng	95	
74) Di-n-butylphthalate	18.671	149	284945	52.106	ng	100	
75) Fluoranthene	19.787	202	313139	46.751	ng	97	
77) Benzidine	19.951	184	55964	22.953	ng	96	
78) Pyrene	20.145	202	314454	50.968	ng	98	
80) Butylbenzylphthalate	21.003	149	120681	56.843	ng	93	
81) Benzo(a)anthracene	22.048	228	310887	50.365	ng	100	
82) 3,3'-Dichlorobenzidine	21.949	252	98184	47.104	ng	97	
83) Chrysene	22.125	228	284455	49.217	ng	98	
84) Bis(2-ethylhexyl)phtha...	21.902	149	225230	73.263	ng	97	
85) Di-n-octyl phthalate	23.212	149	284661	54.947	ng	99	
87) Indeno(1,2,3-cd)pyrene	29.691	276	349686m	48.054	ng		
88) Benzo(b)fluoranthene	24.475	252	316846	51.459	ng	98	
89) Benzo(k)fluoranthene	24.551	252	316245	51.447	ng	100	
90) Benzo(a)pyrene	25.444	252	284006	46.939	ng	99	
91) Dibenzo(a,h)anthracene	29.756	278	288819	48.285	ng	98	
92) Benzo(g,h,i)perylene	30.966	276	262705m	44.853	ng		

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

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