

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040423\
 Data File : BG056973.D
 Acq On : 4 Apr 2023 20:17
 Operator : CG/JU
 Sample : 02161-02MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-6MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 04/05/2023
 Supervised By : Jagrut Upadhyay 04/05/2023

Quant Time: Apr 05 04:10:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 03:23:11 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.416	152	14145	20.000	ng	0.00
21) Naphthalene-d8	11.253	136	57298	20.000	ng	0.00
39) Acenaphthene-d10	15.025	164	45913	20.000	ng	0.00
64) Phenanthrene-d10	17.768	188	118329	20.000	ng	# 0.00
76) Chrysene-d12	22.086	240	107846	20.000	ng	#-0.01
86) Perylene-d12	25.628	264	112075	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.925	112	120411	136.240	ng	0.00
7) Phenol-d6	7.546	99	167339	126.594	ng	0.00
23) Nitrobenzene-d5	9.591	82	119747	85.514	ng	0.00
42) 2,4,6-Tribromophenol	16.505	330	107904	147.113	ng	0.00
45) 2-Fluorobiphenyl	13.656	172	290653	85.842	ng	0.00
79) Terphenyl-d14	20.335	244	613844	96.315	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.740	88	14506	37.441	ng	# 54
3) Pyridine	4.163	79	22157	17.756	ng	95
4) n-Nitrosodimethylamine	4.057	42	25122	43.787	ng	91
6) Aniline	7.729	93	22266	13.176	ng	99
8) 2-Chlorophenol	7.969	128	43339	49.231	ng	97
9) Benzaldehyde	7.535	77	22664	27.336	ng	93
10) Phenol	7.570	94	57392	43.818	ng	93
11) bis(2-Chloroethyl)ether	7.811	93	40374	43.274	ng	98
12) 1,3-Dichlorobenzene	8.304	146	44790m	46.046	ng	
13) 1,4-Dichlorobenzene	8.451	146	46356	47.218	ng	97
14) 1,2-Dichlorobenzene	8.774	146	44173	46.572	ng	100
15) Benzyl Alcohol	8.645	79	50484	45.963	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.933	45	52279m	46.494	ng	
17) 2-Methylphenol	8.845	107	42598	45.586	ng	89
18) Hexachloroethane	9.514	117	16405	45.718	ng	91
19) n-Nitroso-di-n-propyla...	9.215	70	40226	42.545	ng	# 92
20) 3+4-Methylphenols	9.168	107	58500	44.770	ng	89
22) Acetophenone	9.244	105	74047	43.430	ng	# 97
24) Nitrobenzene	9.638	77	62717	43.807	ng	99
25) Isophorone	10.155	82	114865	43.076	ng	98
26) 2-Nitrophenol	10.354	139	25058	45.345	ng	96
27) 2,4-Dimethylphenol	10.396	122	45255	46.383	ng	97
28) bis(2-Chloroethoxy)met...	10.631	93	56902	42.623	ng	98
29) 2,4-Dichlorophenol	10.889	162	49610	47.279	ng	98
30) 1,2,4-Trichlorobenzene	11.112	180	50043	44.885	ng	95
31) Naphthalene	11.306	128	134098	42.937	ng	97
32) Benzoic acid	10.519	122	31044	46.320	ng	88
34) Hexachlorobutadiene	11.576	225	36331	43.380	ng	96
35) Caprolactam	12.164	113	13094m	36.184	ng	
36) 4-Chloro-3-methylphenol	12.487	107	56323	47.997	ng	97
37) 2-Methylnaphthalene	12.880	142	104522	42.168	ng	98
38) 1-Methylnaphthalene	13.098	142	98635	42.378	ng	99
40) 1,2,4,5-Tetrachloroben...	13.239	216	71139	44.682	ng	99
41) Hexachlorocyclopentadiene	13.215	237	97787	111.473	ng	98
43) 2,4,6-Trichlorophenol	13.468	196	47267	46.523	ng	98
44) 2,4,5-Trichlorophenol	13.544	196	52489	43.610	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.867	154	154691	46.026	ng	96
47) 2-Chloronaphthalene	13.914	162	116149	46.726	ng	96
48) 2-Nitroaniline	14.108	65	41104	46.245	ng	96
49) Acenaphthylene	14.754	152	184999	45.297	ng	98
50) Dimethylphthalate	14.461	163	167158	46.207	ng	98
51) 2,6-Dinitrotoluene	14.590	165	36552	46.764	ng	93
52) Acenaphthene	15.089	154	154547m	54.867	ng	
53) 3-Nitroaniline	14.919	138	10198	13.181	ng	92
54) 2,4-Dinitrophenol	15.124	184	52465	119.253	ng	89
55) Dibenzofuran	15.424	168	192296	45.486	ng	97
56) 4-Nitrophenol	15.213	139	55970	97.833	ng	95
57) 2,4-Dinitrotoluene	15.371	165	51460	46.680	ng	87
58) Fluorene	16.064	166	164508	47.025	ng	95
59) 2,3,4,6-Tetrachlorophenol	15.641	232	53623	48.931	ng	# 98
60) Diethylphthalate	15.812	149	168079	46.346	ng	98
61) 4-Chlorophenyl-phenyle...	16.047	204	94812	46.161	ng	99
62) 4-Nitroaniline	16.076	138	30983	38.653	ng	94
63) Azobenzene	16.340	77	164878	45.619	ng	99
65) 4,6-Dinitro-2-methylph...	16.135	198	35360	49.341	ng	96
66) n-Nitrosodiphenylamine	16.258	169	135604	42.690	ng	98
67) 4-Bromophenyl-phenylether	16.940	248	64618	44.135	ng	97
68) Hexachlorobenzene	17.069	284	68754	47.499	ng	98
69) Atrazine	17.192	200	43509	32.997	ng	92
70) Pentachlorophenol	17.410	266	90150	86.812	ng	98
71) Phenanthrene	17.809	178	276159	45.775	ng	100
72) Anthracene	17.897	178	276302	44.676	ng	99
73) Carbazole	18.162	167	242987	45.063	ng	94
74) Di-n-butylphthalate	18.684	149	288550	46.212	ng	99
75) Fluoranthene	19.795	202	336964	44.061	ng	99
78) Pyrene	20.153	202	345276	46.289	ng	98
80) Butylbenzylphthalate	21.016	149	116035	45.205	ng	96
81) Benzo(a)anthracene	22.062	228	339478	45.489	ng	99
82) 3,3'-Dichlorobenzidine	21.956	252	19306	7.661	ng	96
83) Chrysene	22.138	228	310466	44.431	ng	99
84) Bis(2-ethylhexyl)phtha...	21.921	149	165964	44.652	ng	97
85) Di-n-octyl phthalate	23.237	149	274014	43.747	ng	99
87) Indeno(1,2,3-cd)pyrene	29.722	276	376191	45.418	ng	97
88) Benzo(b)fluoranthene	24.494	252	350102	49.955	ng	99
89) Benzo(k)fluoranthene	24.570	252	339965	48.589	ng	100
90) Benzo(a)pyrene	25.463	252	297961	43.265	ng	# 99
91) Dibenzo(a,h)anthracene	29.781	278	313047	45.980	ng	97
92) Benzo(g,h,i)perylene	31.009	276	294154	44.123	ng	98

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

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