

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022023\
 Data File : BG056679.D
 Acq On : 20 Feb 2023 18:49
 Operator : CG/JU
 Sample : 01606-02MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 VNJ-210MSD

Quant Time: Feb 21 03:27:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG021523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 16 01:33:33 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.433	152	22602	20.000	ng	0.00
21) Naphthalene-d8	11.277	136	88928	20.000	ng	0.00
39) Acenaphthene-d10	15.037	164	64528	20.000	ng	0.00
64) Phenanthrene-d10	17.775	188	151489	20.000	ng	0.00
76) Chrysene-d12	22.093	240	133747	20.000	ng	0.00
86) Perylene-d12	25.636	264	148221	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.947	112	167539	120.608	ng	0.00
7) Phenol-d6	7.563	99	215614	105.078	ng	0.00
23) Nitrobenzene-d5	9.614	82	153742	88.519	ng	0.00
42) 2,4,6-Tribromophenol	16.517	330	109503	123.947	ng	0.00
45) 2-Fluorobiphenyl	13.668	172	359665	74.636	ng	0.00
79) Terphenyl-d14	20.342	244	691480	84.588	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.768	88	20130	33.054	ng	# 53
3) Pyridine	4.197	79	55583	28.854	ng	95
4) n-Nitrosodimethylamine	4.091	42	34578	38.509	ng	# 89
6) Aniline	7.745	93	78222	27.430	ng	95
8) 2-Chlorophenol	7.992	128	60195	43.797	ng	96
9) Benzaldehyde	7.557	77	40361	37.194	ng	90
10) Phenol	7.593	94	77829	37.079	ng	97
11) bis(2-Chloroethyl)ether	7.834	93	61209	38.865	ng	99
12) 1,3-Dichlorobenzene	8.327	146	65425	39.953	ng	99
13) 1,4-Dichlorobenzene	8.474	146	66064	40.059	ng	93
14) 1,2-Dichlorobenzene	8.797	146	63549	40.090	ng	99
15) Benzyl Alcohol	8.662	79	72242	39.319	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.956	45	81269	38.436	ng	99
17) 2-Methylphenol	8.862	107	55445	37.588	ng	98
18) Hexachloroethane	9.543	117	22581	43.225	ng	98
19) n-Nitroso-di-n-propyla...	9.238	70	57397	37.364	ng	100
20) 3+4-Methylphenols	9.185	107	76303	37.656	ng	99
22) Acetophenone	9.261	105	106095	39.046	ng	# 99
24) Nitrobenzene	9.655	77	84617	42.173	ng	98
25) Isophorone	10.178	82	154924	36.651	ng	99
26) 2-Nitrophenol	10.372	139	31101	43.489	ng	91
27) 2,4-Dimethylphenol	10.407	122	63708	38.983	ng	98
28) bis(2-Chloroethoxy)met...	10.648	93	84734	39.334	ng	97
29) 2,4-Dichlorophenol	10.906	162	68067	42.025	ng	97
30) 1,2,4-Trichlorobenzene	11.130	180	76067	39.425	ng	95
31) Naphthalene	11.324	128	189819	38.514	ng	99
32) Benzoic acid	10.483	122	12677m	12.036	ng	
33) 4-Chloroaniline	11.423	127	11429	5.076	ng	# 89
34) Hexachlorobutadiene	11.606	225	54902	39.818	ng	98
35) Caprolactam	12.176	113	17157m	32.934	ng	
36) 4-Chloro-3-methylphenol	12.499	107	70577	39.513	ng	98
37) 2-Methylnaphthalene	12.898	142	139514	35.851	ng	100
38) 1-Methylnaphthalene	13.116	142	130125	35.835	ng	97
40) 1,2,4,5-Tetrachloroben...	13.257	216	97635	40.702	ng	97
41) Hexachlorocyclopentadiene	13.233	237	85564	65.765	ng	98
43) 2,4,6-Trichlorophenol	13.480	196	63152	43.391	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.550	196	69847	40.042	ng	97
46) 1,1'-Biphenyl	13.879	154	194482	39.787	ng	97
47) 2-Chloronaphthalene	13.932	162	157014	40.592	ng	98
48) 2-Nitroaniline	14.114	65	45180	40.959	ng	89
49) Acenaphthylene	14.767	152	240956	38.337	ng	97
50) Dimethylphthalate	14.479	163	204662	38.367	ng	99
51) 2,6-Dinitrotoluene	14.596	165	40751	39.712	ng	88
52) Acenaphthene	15.107	154	151505	39.079	ng	98
53) 3-Nitroaniline	14.925	138	22460	21.691	ng #	97
54) 2,4-Dinitrophenol	15.125	184	25578	41.505	ng #	83
55) Dibenzofuran	15.431	168	248021	37.867	ng	98
56) 4-Nitrophenol	15.213	139	60858	76.711	ng	90
57) 2,4-Dinitrotoluene	15.378	165	55114	42.364	ng #	88
58) Fluorene	16.077	166	193520	37.369	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.648	232	64539	42.216	ng #	92
60) Diethylphthalate	15.824	149	199816	37.135	ng	99
61) 4-Chlorophenyl-phenyle...	16.059	204	117915	37.270	ng	100
62) 4-Nitroaniline	16.089	138	38544	35.717	ng	98
63) Azobenzene	16.353	77	200014	37.141	ng	98
65) 4,6-Dinitro-2-methylph...	16.136	198	21177	27.070	ng	90
66) n-Nitrosodiphenylamine	16.271	169	175440	39.809	ng	99
67) 4-Bromophenyl-phenylether	16.952	248	79404	40.825	ng	93
68) Hexachlorobenzene	17.076	284	86810	43.809	ng	99
69) Atrazine	17.205	200	72738	42.983	ng	99
70) Pentachlorophenol	17.416	266	87330	66.548	ng	97
71) Phenanthrene	17.816	178	330723	40.128	ng	99
72) Anthracene	17.910	178	335995	39.841	ng	99
73) Carbazole	18.168	167	292062	40.174	ng	99
74) Di-n-butylphthalate	18.697	149	327763	40.288	ng	99
75) Fluoranthene	19.802	202	412641	37.945	ng	100
77) Benzidine	19.960	184	134862	42.159	ng	98
78) Pyrene	20.160	202	415607	41.146	ng	99
80) Butylbenzylphthalate	21.024	149	135075	44.532	ng	96
81) Benzo(a)anthracene	22.070	228	403433	40.561	ng	98
82) 3,3'-Dichlorobenzidine	21.964	252	80108	24.824	ng	99
83) Chrysene	22.146	228	370798	39.878	ng	99
84) Bis(2-ethylhexyl)phtha...	21.941	149	191331	42.095	ng	100
85) Di-n-octyl phthalate	23.262	149	324269	42.924	ng	100
87) Indeno(1,2,3-cd)pyrene	29.737	276	471016	40.576	ng	99
88) Benzo(b)fluoranthene	24.502	252	408637	42.200	ng	100
89) Benzo(k)fluoranthene	24.579	252	386651	40.604	ng	99
90) Benzo(a)pyrene	25.472	252	348306	37.420	ng	99
91) Dibenzo(a,h)anthracene	29.808	278	391999	40.773	ng	97
92) Benzo(g,h,i)perylene	31.018	276	380851	40.541	ng	96

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

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