

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG101722\
 Data File : BG055192.D
 Acq On : 17 Oct 2022 13:30
 Operator : CG/JU
 Sample : PB148386BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB148386BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/18/2022
 Supervised By : Jagrut Upadhyay 10/18/2022

Quant Time: Oct 17 16:33:23 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG101322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 13 05:14:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.299	152	21461	20.000	ng	0.00
21) Naphthalene-d8	11.131	136	102931	20.000	ng	0.00
39) Acenaphthene-d10	14.915	164	87326	20.000	ng	0.00
64) Phenanthrene-d10	17.647	188	226013	20.000	ng	0.00
76) Chrysene-d12	21.936	240	199949	20.000	ng	0.01
86) Perylene-d12	25.368	264	210087	20.000	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.814	112	169582	154.598	ng	0.00
7) Phenol-d6	7.436	99	239590	147.654	ng	0.00
23) Nitrobenzene-d5	9.469	82	147644	76.919	ng	0.00
42) 2,4,6-Tribromophenol	16.396	330	182715	145.616	ng	0.00
45) 2-Fluorobiphenyl	13.546	172	433258	81.047	ng	0.00
79) Terphenyl-d14	20.227	244	905535	91.211	ng	0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.628	88	13338	28.736	ng	96
3) Pyridine	4.045	79	39976	27.313	ng	95
4) n-Nitrosodimethylamine	3.951	42	26840	34.557	ng	94
6) Aniline	7.606	93	89399	42.905	ng	97
8) 2-Chlorophenol	7.853	128	66916	50.026	ng	98
9) Benzaldehyde	7.418	77	21804	19.980	ng	94
10) Phenol	7.465	94	89174	49.466	ng	98
11) bis(2-Chloroethyl)ether	7.700	93	55619	43.620	ng	97
12) 1,3-Dichlorobenzene	8.188	146	65565	42.172	ng	96
13) 1,4-Dichlorobenzene	8.335	146	66961	42.660	ng	96
14) 1,2-Dichlorobenzene	8.658	146	67350	44.805	ng	98
15) Benzyl Alcohol	8.528	79	69447	47.873	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.816	45	84751	39.760	ng	96
17) 2-Methylphenol	8.728	107	65009	50.121	ng	99
18) Hexachloroethane	9.398	117	22278	41.867	ng	93
19) n-Nitroso-di-n-propyla...	9.098	70	56926	43.470	ng	97
20) 3+4-Methylphenols	9.057	107	89822	48.266	ng	97
22) Acetophenone	9.128	105	108767	42.678	ng	# 98
24) Nitrobenzene	9.516	77	79737	40.152	ng	97
25) Isophorone	10.038	82	162331	43.548	ng	98
26) 2-Nitrophenol	10.232	139	42000	44.639	ng	94
27) 2,4-Dimethylphenol	10.279	122	76140	53.211	ng	97
28) bis(2-Chloroethoxy)met...	10.514	93	86378	46.757	ng	99
29) 2,4-Dichlorophenol	10.767	162	81900	47.890	ng	99
30) 1,2,4-Trichlorobenzene	10.990	180	76563	41.503	ng	97
31) Naphthalene	11.184	128	222893	41.175	ng	99
32) Benzoic acid	10.409	122	39176m	33.468	ng	
33) 4-Chloroaniline	11.278	127	94263	40.511	ng	96
34) Hexachlorobutadiene	11.460	225	48761	41.454	ng	95
35) Caprolactam	12.054	113	32598m	46.400	ng	
36) 4-Chloro-3-methylphenol	12.377	107	94263	48.878	ng	95
37) 2-Methylnaphthalene	12.765	142	174710	43.100	ng	99
38) 1-Methylnaphthalene	12.982	142	169870	42.951	ng	98
40) 1,2,4,5-Tetrachloroben...	13.123	216	105982	43.029	ng	99
41) Hexachlorocyclopentadiene	13.100	237	120922	98.255	ng	98
43) 2,4,6-Trichlorophenol	13.358	196	77194	43.120	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.429	196	88959	44.782	ng	98
46) 1,1'-Biphenyl	13.752	154	252044	43.473	ng	99
47) 2-Chloronaphthalene	13.805	162	198514	42.329	ng	97
48) 2-Nitroaniline	13.993	65	68025	40.658	ng	98
49) Acenaphthylene	14.645	152	336272	42.713	ng	99
50) Dimethylphthalate	14.357	163	301631	42.970	ng	99
51) 2,6-Dinitrotoluene	14.480	165	64552	44.687	ng	93
52) Acenaphthene	14.980	154	239505m	47.027	ng	
53) 3-Nitroaniline	14.809	138	61386	36.653	ng	99
54) 2,4-Dinitrophenol	15.015	184	78244	89.043	ng	89
55) Dibenzofuran	15.309	168	341835	43.020	ng	99
56) 4-Nitrophenol	15.109	139	110095	84.418	ng	98
57) 2,4-Dinitrotoluene	15.262	165	95334	45.059	ng	96
58) Fluorene	15.955	166	277746	42.571	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.532	232	80791	41.214	ng	99
60) Diethylphthalate	15.708	149	317566	42.950	ng	99
61) 4-Chlorophenyl-phenyle...	15.937	204	156253	44.728	ng	99
62) 4-Nitroaniline	15.967	138	73593	40.701	ng	98
63) Azobenzene	16.231	77	256803	40.911	ng	97
65) 4,6-Dinitro-2-methylph...	16.026	198	57246	43.092	ng	96
66) n-Nitrosodiphenylamine	16.149	169	259387	43.503	ng	99
67) 4-Bromophenyl-phenylether	16.831	248	109697	44.523	ng	98
68) Hexachlorobenzene	16.954	284	129925	45.880	ng	98
69) Atrazine	17.083	200	107646	45.794	ng	99
70) Pentachlorophenol	17.295	266	131382	77.869	ng	97
71) Phenanthrene	17.694	178	491605	41.934	ng	99
72) Anthracene	17.782	178	499343	43.586	ng	99
73) Carbazole	18.047	167	459942	43.118	ng	100
74) Di-n-butylphthalate	18.581	149	542112	41.282	ng	99
75) Fluoranthene	19.680	202	571053	38.968	ng	99
77) Benzidine	19.850	184	322480	71.920	ng	99
78) Pyrene	20.038	202	575102	43.146	ng	98
80) Butylbenzylphthalate	20.902	149	215250	42.068	ng	99
81) Benzo(a)anthracene	21.919	228	545534	42.643	ng	99
82) 3,3'-Dichlorobenzidine	21.819	252	194073	44.203	ng	96
83) Chrysene	21.989	228	512272	42.449	ng	98
84) Bis(2-ethylhexyl)phtha...	21.789	149	303959	41.703	ng	98
85) Di-n-octyl phthalate	23.070	149	502698	41.369	ng	97
87) Indeno(1,2,3-cd)pyrene	29.328	276	666921	42.856	ng	# 93
88) Benzo(b)fluoranthene	24.269	252	580314	46.345	ng	99
89) Benzo(k)fluoranthene	24.345	252	554053	44.041	ng	99
90) Benzo(a)pyrene	25.203	252	505987	47.256	ng	99
91) Dibenzo(a,h)anthracene	29.392	278	592694	46.025	ng	98
92) Benzo(g,h,i)perylene	30.561	276	575418	45.615	ng	98

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

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