

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG093022\
 Data File : BG055041.D
 Acq On : 30 Sep 2022 15:02
 Operator : CG/JU
 Sample : PB148010BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB148010BS

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 10/01/2022
 Supervised By :mohammad ahmed 10/10/2022

Quant Time: Oct 01 02:12:17 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG092822.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Sep 28 15:43:39 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.346	152	34287	20.000	ng	0.00
21) Naphthalene-d8	11.178	136	147676	20.000	ng	0.00
39) Acenaphthene-d10	14.950	164	102737	20.000	ng	0.00
64) Phenanthrene-d10	17.688	188	254739	20.000	ng	0.00
76) Chrysene-d12	21.983	240	263629	20.000	ng	#-0.01
86) Perylene-d12	25.438	264	283351	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.867	112	273556	152.943	ng	0.00
7) Phenol-d6	7.477	99	352982	148.741	ng	0.00
23) Nitrobenzene-d5	9.516	82	197157	86.143	ng	0.00
42) 2,4,6-Tribromophenol	16.431	330	162598	139.478	ng	0.00
45) 2-Fluorobiphenyl	13.587	172	531277	81.400	ng	0.00
79) Terphenyl-d14	20.262	244	1108642	87.691	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.693	88	30589	34.528	ng	96
3) Pyridine	4.110	79	82552	32.791	ng	98
4) n-Nitrosodimethylamine	4.016	42	52217	42.657	ng	90
6) Aniline	7.659	93	141531	45.267	ng	98
8) 2-Chlorophenol	7.900	128	92400	46.264	ng	99
9) Benzaldehyde	7.471	77	62364	40.335	ng	97
10) Phenol	7.500	94	123827	46.007	ng	95
11) bis(2-Chloroethyl)ether	7.747	93	87958	43.543	ng	97
12) 1,3-Dichlorobenzene	8.235	146	106848	43.704	ng	98
13) 1,4-Dichlorobenzene	8.382	146	108044	43.575	ng	99
14) 1,2-Dichlorobenzene	8.705	146	105592	45.067	ng	99
15) Benzyl Alcohol	8.576	79	91254	44.928	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.869	45	143791	42.541	ng	97
17) 2-Methylphenol	8.769	107	86358	45.934	ng	96
18) Hexachloroethane	9.451	117	35023	42.999	ng	93
19) n-Nitroso-di-n-propyla...	9.151	70	77068	41.337	ng	96
20) 3+4-Methylphenols	9.098	107	113670	43.877	ng	96
22) Acetophenone	9.169	105	152023	41.964	ng	# 98
24) Nitrobenzene	9.563	77	104495	42.140	ng	98
25) Isophorone	10.086	82	213294	41.507	ng	99
26) 2-Nitrophenol	10.274	139	37087	35.843	ng	99
27) 2,4-Dimethylphenol	10.321	122	96172	47.876	ng	98
28) bis(2-Chloroethoxy)met...	10.561	93	119539	44.565	ng	100
29) 2,4-Dichlorophenol	10.808	162	96269	42.909	ng	95
30) 1,2,4-Trichlorobenzene	11.031	180	105624	40.106	ng	98
31) Naphthalene	11.231	128	304386	39.517	ng	100
32) Benzoic acid	10.432	122	48088	35.326	ng	97
33) 4-Chloroaniline	11.319	127	96740	30.876	ng	97
34) Hexachlorobutadiene	11.507	225	68887	40.092	ng	98
35) Caprolactam	12.077	113	32440m	44.030	ng	
36) 4-Chloro-3-methylphenol	12.406	107	106697	43.102	ng	94
37) 2-Methylnaphthalene	12.806	142	218698	39.570	ng	98
38) 1-Methylnaphthalene	13.023	142	208667	39.106	ng	97
40) 1,2,4,5-Tetrachloroben...	13.164	216	125391	41.587	ng	99
41) Hexachlorocyclopentadiene	13.147	237	146833	96.167	ng	99
43) 2,4,6-Trichlorophenol	13.393	196	79817	41.556	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.464	196	90798	41.566	ng	94
46) 1,1'-Biphenyl	13.793	154	296079	42.129	ng	99
47) 2-Chloronaphthalene	13.840	162	228691	40.264	ng	98
48) 2-Nitroaniline	14.028	65	59198	37.881	ng	97
49) Acenaphthylene	14.680	152	376327	40.833	ng	98
50) Dimethylphthalate	14.398	163	311353	40.552	ng	100
51) 2,6-Dinitrotoluene	14.516	165	57501	40.253	ng	95
52) Acenaphthene	15.015	154	247229	43.214	ng	98
53) 3-Nitroaniline	14.845	138	58368	35.535	ng #	93
54) 2,4-Dinitrophenol	15.044	184	55419	85.579	ng	92
55) Dibenzofuran	15.350	168	373039	40.343	ng	98
56) 4-Nitrophenol	15.127	139	113696	91.501	ng	91
57) 2,4-Dinitrotoluene	15.291	165	81882	41.848	ng	94
58) Fluorene	15.990	166	295746	40.059	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.561	232	80613	40.961	ng	99
60) Diethylphthalate	15.749	149	326079	41.162	ng	100
61) 4-Chlorophenyl-phenyle...	15.979	204	164569	41.673	ng	100
62) 4-Nitroaniline	16.002	138	73182	44.773	ng	91
63) Azobenzene	16.272	77	293282	39.455	ng	96
65) 4,6-Dinitro-2-methylph...	16.055	198	35215	35.908	ng	86
66) n-Nitrosodiphenylamine	16.190	169	277617	40.484	ng	97
67) 4-Bromophenyl-phenylether	16.872	248	109335	39.845	ng	99
68) Hexachlorobenzene	16.989	284	125690	40.554	ng	99
69) Atrazine	17.124	200	117138	46.089	ng	99
70) Pentachlorophenol	17.330	266	146684	84.052	ng	99
71) Phenanthrene	17.730	178	525664	39.971	ng	98
72) Anthracene	17.824	178	534591	41.209	ng	99
73) Carbazole	18.082	167	512609	42.156	ng	98
74) Di-n-butylphthalate	18.623	149	601314	41.154	ng	99
75) Fluoranthene	19.715	202	687089	40.768	ng	98
77) Benzidine	19.886	184	541886	83.007	ng	98
78) Pyrene	20.074	202	691603	41.002	ng	99
80) Butylbenzylphthalate	20.949	149	247185	40.566	ng	96
81) Benzo(a)anthracene	21.966	228	679677	40.248	ng	98
82) 3,3'-Dichlorobenzidine	21.866	252	211919	38.806	ng	98
83) Chrysene	22.036	228	630793	39.661	ng	99
84) Bis(2-ethylhexyl)phtha...	21.854	149	374115	41.481	ng	99
85) Di-n-octyl phthalate	23.158	149	611730	40.418	ng	98
87) Indeno(1,2,3-cd)pyrene	29.428	276	847384	40.934	ng	97
88) Benzo(b)fluoranthene	24.339	252	736361	43.647	ng #	98
89) Benzo(k)fluoranthene	24.416	252	715315	42.859	ng #	97
90) Benzo(a)pyrene	25.279	252	651844	45.716	ng #	98
91) Dibenzo(a,h)anthracene	29.504	278	731444	42.927	ng #	97
92) Benzo(g,h,i)perylene	30.685	276	723137	43.140	ng #	94

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

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