

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011223\
 Data File : BG056267.D
 Acq On : 12 Jan 2023 13:51
 Operator : CG/JU
 Sample : 01094-01MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Jan 13 04:25:21 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 12 03:38:40 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut
 Upadhyay
 01/13/2023

Supervised By :mohammad
 ahmed
 01/13/2023

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.313	152	42638	20.000	ng	0.00
21) Naphthalene-d8	11.151	136	164236	20.000	ng	0.00
39) Acenaphthene-d10	14.928	164	136387	20.000	ng	0.00
64) Phenanthrene-d10	17.666	188	373101	20.000	ng	0.00
76) Chrysene-d12	21.961	240	354506	20.000	ng	0.00
86) Perylene-d12	25.399	264	390226	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.833	112	327704	137.228	ng	0.00
7) Phenol-d6	7.455	99	461750	125.909	ng	0.00
23) Nitrobenzene-d5	9.494	82	243521	75.845	ng	0.00
42) 2,4,6-Tribromophenol	16.409	330	225618	130.617	ng	0.00
45) 2-Fluorobiphenyl	13.559	172	756000	76.544	ng	0.00
79) Terphenyl-d14	20.240	244	1520054	76.797	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.642	88	44981	41.239	ng	# 95
3) Pyridine	4.065	79	126939	38.210	ng	96
4) n-Nitrosodimethylamine	3.971	42	29398	43.502	ng	# 96
6) Aniline	7.631	93	93486	19.439	ng	96
8) 2-Chlorophenol	7.878	128	130671	54.989	ng	94
9) Benzaldehyde	7.437	77	67340	30.882	ng	93
10) Phenol	7.478	94	191776	51.564	ng	93
11) bis(2-Chloroethyl)ether	7.719	93	116503	46.224	ng	96
12) 1,3-Dichlorobenzene	8.207	146	148008	51.176	ng	95
13) 1,4-Dichlorobenzene	8.354	146	150342	51.766	ng	99
14) 1,2-Dichlorobenzene	8.677	146	145959	52.086	ng	97
15) Benzyl Alcohol	8.548	79	122905	46.048	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.841	45	23365	52.460	ng	91
17) 2-Methylphenol	8.747	107	129737	47.540	ng	98
18) Hexachloroethane	9.417	117	45872	51.993	ng	97
19) n-Nitroso-di-n-propyla...	9.118	70	92962	41.686	ng	98
20) 3+4-Methylphenols	9.076	107	179462	46.773	ng	97
22) Acetophenone	9.141	105	206770	44.526	ng	# 98
24) Nitrobenzene	9.535	77	140716	44.224	ng	99
25) Isophorone	10.052	82	276827	40.978	ng	98
26) 2-Nitrophenol	10.246	139	80679	49.557	ng	97
27) 2,4-Dimethylphenol	10.293	122	146087	49.198	ng	98
28) bis(2-Chloroethoxy)met...	10.534	93	159729	45.492	ng	99
29) 2,4-Dichlorophenol	10.786	162	170939	52.071	ng	96
30) 1,2,4-Trichlorobenzene	11.004	180	181564	48.738	ng	100
31) Naphthalene	11.198	128	397746	46.017	ng	99
32) Benzoic acid	10.422	122	110248	50.041	ng	97
33) 4-Chloroaniline	11.297	127	45671	11.067	ng	94
34) Hexachlorobutadiene	11.474	225	126773	47.315	ng	97
35) Caprolactam	12.067	113	53436m	48.078	ng	
36) 4-Chloro-3-methylphenol	12.390	107	161614	48.516	ng	99
37) 2-Methylnaphthalene	12.778	142	314629	43.110	ng	98
38) 1-Methylnaphthalene	12.995	142	294821	43.501	ng	98
40) 1,2,4,5-Tetrachloroben...	13.142	216	243133	47.564	ng	99
41) Hexachlorocyclopentadiene	13.119	237	313140	107.448	ng	94
43) 2,4,6-Trichlorophenol	13.371	196	171682	50.792	ng	98

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44) 2,4,5-Trichlorophenol	13.442	196	187426	46.875	ng	98
46) 1,1'-Biphenyl	13.771	154	468854	48.001	ng	98
47) 2-Chloronaphthalene	13.818	162	375304	48.679	ng	99
48) 2-Nitroaniline	14.012	65	81573	49.659	ng	99
49) Acenaphthylene	14.658	152	581738	46.368	ng	99
50) Dimethylphthalate	14.370	163	504835	45.852	ng	99
51) 2,6-Dinitrotoluene	14.494	165	111523	47.537	ng	97
52) Acenaphthene	14.993	154	451553m	55.338	ng	
53) 3-Nitroaniline	14.823	138	41610	18.636	ng	95
54) 2,4-Dinitrophenol	15.028	184	188295	112.349	ng	95
55) Dibenzofuran	15.322	168	621491	46.025	ng	97
56) 4-Nitrophenol	15.122	139	185084	107.784	ng	94
57) 2,4-Dinitrotoluene	15.281	165	160937	49.052	ng	91
58) Fluorene	15.968	166	512611	47.198	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.545	232	191682m	51.057	ng	
60) Diethylphthalate	15.722	149	480783	46.564	ng	98
61) 4-Chlorophenyl-phenyle...	15.951	204	315571	46.559	ng	98
62) 4-Nitroaniline	15.986	138	107565	47.032	ng	97
63) Azobenzene	16.245	77	372781	46.786	ng	97
65) 4,6-Dinitro-2-methylph...	16.039	198	125469	49.913	ng	95
66) n-Nitrosodiphenylamine	16.168	169	475146	45.900	ng	97
67) 4-Bromophenyl-phenylether	16.844	248	217298	45.737	ng	95
68) Hexachlorobenzene	16.973	284	219283	49.470	ng	100
69) Atrazine	17.096	200	217073	50.597	ng	98
70) Pentachlorophenol	17.308	266	309103	90.779	ng	98
71) Phenanthrene	17.708	178	911576	47.146	ng	99
72) Anthracene	17.802	178	931353	46.695	ng	99
73) Carbazole	18.060	167	807440	48.160	ng	98
74) Di-n-butylphthalate	18.589	149	814710	48.747	ng	100
75) Fluoranthene	19.699	202	1167476	46.289	ng	100
77) Benzidine	19.864	184	353277	28.436	ng	99
78) Pyrene	20.058	202	1177143	47.110	ng	99
80) Butylbenzylphthalate	20.916	149	347786	48.515	ng	96
81) Benzo(a)anthracene	21.938	228	1174683	47.177	ng	99
82) 3,3'-Dichlorobenzidine	21.838	252	162401	18.118	ng	98
83) Chrysene	22.008	228	1095892	46.987	ng	98
84) Bis(2-ethylhexyl)phtha...	21.803	149	507885	49.176	ng	99
85) Di-n-octyl phthalate	23.089	149	839115	48.128	ng	100
87) Indeno(1,2,3-cd)pyrene	29.364	276	1401037	49.368	ng	# 99
88) Benzo(b)fluoranthene	24.300	252	1260461	50.838	ng	100
89) Benzo(k)fluoranthene	24.376	252	1172127	47.475	ng	98
90) Benzo(a)pyrene	25.240	252	1087321	44.963	ng	99
91) Dibenzo(a,h)anthracene	29.435	278	1167734	49.078	ng	99
92) Benzo(g,h,i)perylene	30.616	276	1134351	49.162	ng	98

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

Manual Integrations APPROVED

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ahmed
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[illegible]