

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110122\
 Data File : BG055404.D
 Acq On : 1 Nov 2022 22:41
 Operator : CG/JU
 Sample : PB148503BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

PB148503BS

Manual Integrations

APPROVED

Reviewed By :Christian

Giraldo

11/02/2022

Supervised By :Jagrut

Upadhyay

11/02/2022

Quant Time: Nov 02 05:41:38 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG110122.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Nov 02 05:05:12 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.241	152	24006	20.000	ng	0.00
21) Naphthalene-d8	11.073	136	117748	20.000	ng	0.00
39) Acenaphthene-d10	14.863	164	88693	20.000	ng	0.00
64) Phenanthrene-d10	17.595	188	209208	20.000	ng	0.00
76) Chrysene-d12	21.866	240	201689	20.000	ng	0.00
86) Perylene-d12	25.239	264	217261	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.767	112	222366	166.348	ng	0.00
7) Phenol-d6	7.395	99	304391	164.204	ng	0.00
23) Nitrobenzene-d5	9.422	82	180546	81.718	ng	0.00
42) 2,4,6-Tribromophenol	16.343	330	172589	142.660	ng	0.00
45) 2-Fluorobiphenyl	13.488	172	490397	81.978	ng	0.00
79) Terphenyl-d14	20.168	244	916770	88.388	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.564	88	19483	32.134	ng	96
3) Pyridine	3.987	79	55020	30.577	ng	98
4) n-Nitrosodimethylamine	3.899	42	29695	41.774	ng	99
6) Aniline	7.559	93	110113	45.798	ng	97
8) 2-Chlorophenol	7.806	128	84581	52.077	ng	99
9) Benzaldehyde	7.371	77	37631	34.529	ng	98
10) Phenol	7.418	94	111203	53.000	ng	99
11) bis(2-Chloroethyl)ether	7.648	93	73078	46.554	ng	98
12) 1,3-Dichlorobenzene	8.129	146	81835	44.453	ng	98
13) 1,4-Dichlorobenzene	8.276	146	83228	44.388	ng	100
14) 1,2-Dichlorobenzene	8.599	146	82644	46.049	ng	98
15) Benzyl Alcohol	8.482	79	75859m	51.309	ng	
16) 2,2'-oxybis(1-Chloropr...	8.770	45	96928	46.410	ng	98
17) 2-Methylphenol	8.682	107	79573	52.709	ng	98
18) Hexachloroethane	9.340	117	28489	44.540	ng	97
19) n-Nitroso-di-n-propyla...	9.046	70	67874	46.858	ng	99
20) 3+4-Methylphenols	9.017	107	109589	51.090	ng	98
22) Acetophenone	9.075	105	131696	42.671	ng	# 99
24) Nitrobenzene	9.463	77	95557	41.484	ng	97
25) Isophorone	9.980	82	191864	43.849	ng	100
26) 2-Nitrophenol	10.180	139	51724	43.609	ng	99
27) 2,4-Dimethylphenol	10.227	122	88133	52.994	ng	98
28) bis(2-Chloroethoxy)met...	10.456	93	110442	47.388	ng	98
29) 2,4-Dichlorophenol	10.720	162	93995	46.333	ng	98
30) 1,2,4-Trichlorobenzene	10.932	180	88686	40.370	ng	99
31) Naphthalene	11.126	128	270617	40.820	ng	99
32) Benzoic acid	10.391	122	39466	40.907	ng	95
33) 4-Chloroaniline	11.226	127	102612	37.250	ng	99
34) Hexachlorobutadiene	11.396	225	53770	39.377	ng	97
35) Caprolactam	12.001	113	34700m	46.197	ng	
36) 4-Chloro-3-methylphenol	12.336	107	105600	47.540	ng	98
37) 2-Methylnaphthalene	12.712	142	206152	42.334	ng	99
38) 1-Methylnaphthalene	12.924	142	198488	41.790	ng	98
40) 1,2,4,5-Tetrachloroben...	13.071	216	111551	42.577	ng	99
41) Hexachlorocyclopentadiene	13.047	237	122814	96.746	ng	95
43) 2,4,6-Trichlorophenol	13.306	196	80796	44.284	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.388	196	89994	44.283	ng	99
46) 1,1'-Biphenyl	13.699	154	288093	44.143	ng	99
47) 2-Chloronaphthalene	13.746	162	220908	42.125	ng	99
48) 2-Nitroaniline	13.946	65	70418	43.718	ng	98
49) Acenaphthylene	14.586	152	368452	42.789	ng	99
50) Dimethylphthalate	14.304	163	315766	42.652	ng	99
51) 2,6-Dinitrotoluene	14.428	165	68439	44.157	ng	97
52) Acenaphthene	14.921	154	217916	42.522	ng	99
53) 3-Nitroaniline	14.763	138	62445	35.847	ng	97
54) 2,4-Dinitrophenol	14.974	184	79454	84.505	ng	96
55) Dibenzofuran	15.256	168	365769	43.119	ng	99
56) 4-Nitrophenol	15.074	139	113945	95.398	ng	99
57) 2,4-Dinitrotoluene	15.215	165	99891	45.074	ng	98
58) Fluorene	15.903	166	292854	42.498	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.480	232	80640	44.130	ng	98
60) Diethylphthalate	15.650	149	327951	42.626	ng	100
61) 4-Chlorophenyl-phenyle...	15.879	204	156846	44.025	ng	97
62) 4-Nitroaniline	15.926	138	77542	42.074	ng	99
63) Azobenzene	16.173	77	272579	42.621	ng	98
65) 4,6-Dinitro-2-methylph...	15.979	198	57804	42.894	ng	96
66) n-Nitrosodiphenylamine	16.096	169	270500	43.489	ng	99
67) 4-Bromophenyl-phenylether	16.772	248	107125	43.317	ng	99
68) Hexachlorobenzene	16.901	284	122644	43.487	ng	96
69) Atrazine	17.031	200	111196	53.250	ng	98
70) Pentachlorophenol	17.248	266	117956	80.661	ng	97
71) Phenanthrene	17.636	178	498114	42.329	ng	100
72) Anthracene	17.730	178	510277	44.240	ng	99
73) Carbazole	17.994	167	477158	44.464	ng	100
74) Di-n-butylphthalate	18.517	149	578173	43.138	ng	99
75) Fluoranthene	19.628	202	602941	42.596	ng	99
77) Benzidine	19.792	184	306732	73.048	ng	99
78) Pyrene	19.986	202	603233	42.397	ng	99
80) Butylbenzylphthalate	20.844	149	256707	43.275	ng	97
81) Benzo(a)anthracene	21.849	228	586368	42.352	ng	99
82) 3,3'-Dichlorobenzidine	21.749	252	197891	41.681	ng	97
83) Chrysene	21.919	228	553567	41.909	ng	100
84) Bis(2-ethylhexyl)phtha...	21.708	149	376157	43.623	ng	99
85) Di-n-octyl phthalate	22.965	149	640611	43.359	ng	100
87) Indeno(1,2,3-cd)pyrene	29.140	276	749013	42.939	ng	100
88) Benzo(b)fluoranthene	24.163	252	640130	46.385	ng	99
89) Benzo(k)fluoranthene	24.234	252	623742	44.976	ng	99
90) Benzo(a)pyrene	25.086	252	569759	47.747	ng	99
91) Dibenzo(a,h)anthracene	29.199	278	651760	45.288	ng	99
92) Benzo(g,h,i)perylene	30.362	276	641811	44.841	ng	98

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

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