

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081123\
 Data File : BG058702.D
 Acq On : 11 Aug 2023 14:31
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 08/12/2023

Supervised By :mohammad ahmed 08/12/2023

Quant Time: Aug 11 15:05:03 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG072823.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Jul 28 22:29:07 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	50597	20.000	ng	0.00
21) Naphthalene-d8	10.614	136	213834	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	149495	20.000	ng	0.00
64) Phenanthrene-d10	17.207	188	362253	20.000	ng	0.00
76) Chrysene-d12	21.455	240	323384	20.000	ng	0.00
86) Perylene-d12	24.522	264	408113	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.379	112	247655	74.813	ng	0.00
7) Phenol-d6	6.983	99	360763	78.320	ng	0.00
23) Nitrobenzene-d5	8.975	82	338940	77.867	ng	0.00
42) 2,4,6-Tribromophenol	15.955	330	186891	76.272	ng	0.00
45) 2-Fluorobiphenyl	13.082	172	750561	75.239	ng	0.00
79) Terphenyl-d14	19.827	244	1290056	75.019	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.270	88	57648	35.938	ng	97
3) Pyridine	3.670	79	162518	37.161	ng	96
4) n-Nitrosodimethylamine	3.587	42	73931	39.898	ng	95
6) Aniline	7.136	93	225307	39.728	ng	97
8) 2-Chlorophenol	7.377	128	126920	39.082	ng	97
9) Benzaldehyde	6.948	77	92850m	37.102	ng	
10) Phenol	7.013	94	177809	39.517	ng	98
11) bis(2-Chloroethyl)ether	7.236	93	136887	38.589	ng	97
12) 1,3-Dichlorobenzene	7.700	146	128551	37.070	ng	97
13) 1,4-Dichlorobenzene	7.847	146	129244	37.119	ng	96
14) 1,2-Dichlorobenzene	8.164	146	125543	37.712	ng	97
15) Benzyl Alcohol	8.053	79	133623	45.743	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.335	45	233519	40.526	ng	98
17) 2-Methylphenol	8.258	107	132186	40.178	ng	99
18) Hexachloroethane	8.893	117	47965	37.900	ng	93
19) n-Nitroso-di-n-propyla...	8.617	70	120009	40.335	ng	98
20) 3+4-Methylphenols	8.587	107	184325	41.419	ng	98
22) Acetophenone	8.634	105	222509	38.862	ng	# 99
24) Nitrobenzene	9.022	77	171397	38.732	ng	98
25) Isophorone	9.539	82	350432	38.964	ng	99
26) 2-Nitrophenol	9.727	139	76666	39.800	ng	99
27) 2,4-Dimethylphenol	9.792	122	125951	39.419	ng	99
28) bis(2-Chloroethoxy)met...	10.021	93	191211	39.067	ng	100
29) 2,4-Dichlorophenol	10.268	162	128387	38.699	ng	100
30) 1,2,4-Trichlorobenzene	10.473	180	145344	37.989	ng	98
31) Naphthalene	10.661	128	423988	37.620	ng	100
32) Benzoic acid	9.956	122	71128	30.950	ng	96
33) 4-Chloroaniline	10.779	127	195066	40.965	ng	99
34) Hexachlorobutadiene	10.943	225	92713	37.629	ng	98
35) Caprolactam	11.560	113	47071	38.422	ng	91
36) 4-Chloro-3-methylphenol	11.913	107	161415	39.332	ng	99
37) 2-Methylnaphthalene	12.277	142	308703	38.292	ng	99
38) 1-Methylnaphthalene	12.500	142	289775	37.813	ng	99
40) 1,2,4,5-Tetrachloroben...	12.647	216	172553	37.596	ng	99
41) Hexachlorocyclopentadiene	12.624	237	86685	44.014	ng	97
43) 2,4,6-Trichlorophenol	12.894	196	117852	37.332	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.971	196	137948	37.100	ng	98
46) 1,1'-Biphenyl	13.294	154	389563	37.945	ng	100
47) 2-Chloronaphthalene	13.335	162	303915	37.966	ng	98
48) 2-Nitroaniline	13.546	65	122896	40.222	ng	98
49) Acenaphthylene	14.181	152	508054	37.909	ng	99
50) Dimethylphthalate	13.922	163	430703	38.362	ng	98
51) 2,6-Dinitrotoluene	14.046	165	94816	38.488	ng	98
52) Acenaphthene	14.528	154	290050	37.455	ng	98
53) 3-Nitroaniline	14.375	138	98886	40.121	ng	96
54) 2,4-Dinitrophenol	14.592	184	43736	36.032	ng	99
55) Dibenzofuran	14.862	168	502506	37.764	ng	98
56) 4-Nitrophenol	14.692	139	64702	37.280	ng	95
57) 2,4-Dinitrotoluene	14.833	165	136757	40.042	ng	95
58) Fluorene	15.509	166	397344	37.589	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.092	232	119154	37.554	ng	98
60) Diethylphthalate	15.280	149	438294	38.319	ng	99
61) 4-Chlorophenyl-phenyle...	15.503	204	224395	38.076	ng	97
62) 4-Nitroaniline	15.544	138	99500	42.616	ng	98
63) Azobenzene	15.797	77	429962	38.165	ng	98
65) 4,6-Dinitro-2-methylph...	15.603	198	80547	40.325	ng	99
66) n-Nitrosodiphenylamine	15.720	169	355021	37.654	ng	98
67) 4-Bromophenyl-phenylether	16.396	248	154114	37.509	ng	99
68) Hexachlorobenzene	16.513	284	165659	38.049	ng	99
69) Atrazine	16.666	200	120072	37.725	ng	99
70) Pentachlorophenol	16.866	266	96956	36.792	ng	97
71) Phenanthrene	17.248	178	643743	37.471	ng	99
72) Anthracene	17.342	178	661050	37.501	ng	99
73) Carbazole	17.612	167	605712	38.966	ng	100
74) Di-n-butylphthalate	18.164	149	754151	38.733	ng	99
75) Fluoranthene	19.263	202	795709	38.559	ng	100
77) Benzidine	19.451	184	239021	48.742	ng	98
78) Pyrene	19.627	202	816724	37.110	ng	99
80) Butylbenzylphthalate	20.515	149	344693	38.130	ng	99
81) Benzo(a)anthracene	21.431	228	852654	38.313	ng	98
82) 3,3'-Dichlorobenzidine	21.355	252	301886	40.120	ng	99
83) Chrysene	21.496	228	807964	37.865	ng	100
84) Bis(2-ethylhexyl)phtha...	21.337	149	498719	37.752	ng	# 99
85) Di-n-octyl phthalate	22.483	149	883002	38.018	ng	99
87) Indeno(1,2,3-cd)pyrene	28.029	276	1049580	38.658	ng	100
88) Benzo(b)fluoranthene	23.546	252	856611	38.699	ng	100
89) Benzo(k)fluoranthene	23.617	252	867228	38.999	ng	99
90) Benzo(a)pyrene	24.381	252	843668	38.811	ng	99
91) Dibenzo(a,h)anthracene	28.082	278	871243	38.809	ng	99
92) Benzo(g,h,i)perylene	29.128	276	865175	38.433	ng	99

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

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