

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG043024\
 Data File : BG061063.D
 Acq On : 30 Apr 2024 14:17
 Operator : MA/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC050

Manual Integrations

APPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024

Supervised By :mohammad ahmed 05/03/2024

Quant Time: May 01 04:51:11 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG043024.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed May 01 04:45:52 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.959	152	20565	20.000	ng	0.00
21) Naphthalene-d8	10.761	136	92219	20.000	ng	# 0.00
39) Acenaphthene-d10	14.586	164	81615	20.000	ng	0.00
64) Phenanthrene-d10	17.330	188	211382	20.000	ng	0.00
76) Chrysene-d12	21.590	240	207712	20.000	ng	# 0.00
86) Perylene-d12	24.727	264	241672	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.550	112	100582	99.469	ng	0.00
7) Phenol-d6	7.136	99	147364	98.514	ng	0.00
23) Nitrobenzene-d5	9.122	82	185042	99.469	ng	0.00
42) 2,4,6-Tribromophenol	16.078	330	159545	97.800	ng	0.00
45) 2-Fluorobiphenyl	13.211	172	543422	98.458	ng	0.00
79) Terphenyl-d14	19.950	244	1226639	98.216	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.464	88	18860	46.128	ng	94
3) Pyridine	3.857	79	60761m	53.281	ng	
4) n-Nitrosodimethylamine	3.769	42	57294	49.566	ng	93
6) Aniline	7.289	93	74561	50.279	ng	96
8) 2-Chlorophenol	7.530	128	60565	48.820	ng	96
9) Benzaldehyde	7.101	77	36710	47.184	ng	91
10) Phenol	7.160	94	76161	50.066	ng	92
11) bis(2-Chloroethyl)ether	7.383	93	50968	48.499	ng	98
12) 1,3-Dichlorobenzene	7.853	146	70885	48.469	ng	98
13) 1,4-Dichlorobenzene	8.000	146	72519	47.977	ng	98
14) 1,2-Dichlorobenzene	8.317	146	71239	48.851	ng	95
15) Benzyl Alcohol	8.199	79	65895	48.833	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.493	45	114185	50.073	ng	97
17) 2-Methylphenol	8.405	107	54762	50.515	ng	98
18) Hexachloroethane	9.046	117	26532	49.193	ng	95
19) n-Nitroso-di-n-propyla...	8.769	70	56753	50.697	ng	# 94
20) 3+4-Methylphenols	8.734	107	78943	50.473	ng	97
22) Acetophenone	8.781	105	111924	48.734	ng	# 98
24) Nitrobenzene	9.163	77	90353	49.648	ng	97
25) Isophorone	9.686	82	163911	47.666	ng	99
26) 2-Nitrophenol	9.874	139	42360	49.128	ng	97
27) 2,4-Dimethylphenol	9.933	122	50747	51.049	ng	89
28) bis(2-Chloroethoxy)met...	10.168	93	81577	48.337	ng	96
29) 2,4-Dichlorophenol	10.409	162	77068	49.480	ng	96
30) 1,2,4-Trichlorobenzene	10.626	180	91215	48.603	ng	96
31) Naphthalene	10.808	128	229523	47.889	ng	99
32) Benzoic acid	10.085	122	60639	55.256	ng	# 84
33) 4-Chloroaniline	10.914	127	94992	49.696	ng	98
34) Hexachlorobutadiene	11.096	225	78512	48.756	ng	96
35) Caprolactam	11.695	113	26263	46.823	ng	# 90
36) 4-Chloro-3-methylphenol	12.042	107	89945	47.890	ng	93
37) 2-Methylnaphthalene	12.412	142	172663	48.075	ng	97
38) 1-Methylnaphthalene	12.635	142	172825	47.799	ng	95
40) 1,2,4,5-Tetrachloroben...	12.782	216	133784	49.118	ng	100
41) Hexachlorocyclopentadiene	12.765	237	55566	53.291	ng	98
43) 2,4,6-Trichlorophenol	13.023	196	87239	49.889	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.100	196	99126	50.087	ng	96
46) 1,1'-Biphenyl	13.423	154	256738	49.061	ng	96
47) 2-Chloronaphthalene	13.464	162	207423	49.500	ng	97
48) 2-Nitroaniline	13.664	65	77431	49.797	ng	98
49) Acenaphthylene	14.310	152	317947	49.402	ng	98
50) Dimethylphthalate	14.046	163	304352	48.280	ng	99
51) 2,6-Dinitrotoluene	14.163	165	66532	49.342	ng	97
52) Acenaphthene	14.651	154	198138	49.351	ng	99
53) 3-Nitroaniline	14.492	138	59014	47.833	ng	93
54) 2,4-Dinitrophenol	14.698	184	47830	50.868	ng	98
55) Dibenzofuran	14.986	168	334723	49.358	ng	98
56) 4-Nitrophenol	14.803	139	49363	49.508	ng	95
57) 2,4-Dinitrotoluene	14.950	165	99293	49.773	ng	92
58) Fluorene	15.638	166	285054	48.576	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.215	232	99646	48.911	ng	98
60) Diethylphthalate	15.409	149	303443	47.734	ng	98
61) 4-Chlorophenyl-phenyle...	15.632	204	172117	48.847	ng	97
62) 4-Nitroaniline	15.655	138	66244	49.506	ng	98
63) Azobenzene	15.920	77	238133	48.122	ng	98
65) 4,6-Dinitro-2-methylph...	15.714	198	68034	51.874	ng	97
66) n-Nitrosodiphenylamine	15.843	169	259101	49.429	ng	99
67) 4-Bromophenyl-phenylether	16.519	248	125143	50.256	ng	96
68) Hexachlorobenzene	16.642	284	151574	48.906	ng	99
69) Atrazine	16.795	200	103904	48.064	ng	97
70) Pentachlorophenol	16.983	266	96689	50.780	ng	95
71) Phenanthrene	17.377	178	474725	49.116	ng	99
72) Anthracene	17.465	178	470821	48.819	ng	99
73) Carbazole	17.735	167	407927	48.106	ng	98
74) Di-n-butylphthalate	18.299	149	483993	48.000	ng	99
75) Fluoranthene	19.386	202	638789	48.243	ng	100
77) Benzidine	19.563	184	200548	50.580	ng	98
78) Pyrene	19.745	202	656227	49.693	ng	99
80) Butylbenzylphthalate	20.638	149	206224	49.334	ng	97
81) Benzo(a)anthracene	21.572	228	665586	48.968	ng	99
82) 3,3'-Dichlorobenzidine	21.490	252	240334	49.133	ng	98
83) Chrysene	21.637	228	620837	48.752	ng	98
84) Bis(2-ethylhexyl)phtha...	21.496	149	267773	49.056	ng	# 99
85) Di-n-octyl phthalate	22.682	149	478951	49.733	ng	100
87) Indeno(1,2,3-cd)pyrene	28.288	276	783554	48.923	ng	99
88) Benzo(b)fluoranthene	23.734	252	659918	48.866	ng	100
89) Benzo(k)fluoranthene	23.805	252	625889	48.247	ng	99
90) Benzo(a)pyrene	24.580	252	570733	48.509	ng	99
91) Dibenzo(a,h)anthracene	28.352	278	646339	48.730	ng	# 98
92) Benzo(g,h,i)perylene	29.410	276	643788	48.844	ng	97

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

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