

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG113023\
 Data File : BG059919.D
 Acq On : 30 Nov 2023 21:19
 Operator : MA/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC060

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 12/01/2023

Supervised By :mohammad ahmed 12/01/2023

Quant Time: Dec 01 03:32:09 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG113023.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Dec 01 03:17:30 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.964	152	30064	20.000	ng	0.00
21) Naphthalene-d8	10.773	136	140661	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	101484	20.000	ng	0.00
64) Phenanthrene-d10	17.353	188	227914	20.000	ng	0.00
76) Chrysene-d12	21.625	240	210552	20.000	ng	0.00
86) Perylene-d12	24.821	264	251279	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.526	112	234378	122.116	ng	0.00
7) Phenol-d6	7.118	99	329326	118.872	ng	0.00
23) Nitrobenzene-d5	9.128	82	321853	133.073	ng	0.00
42) 2,4,6-Tribromophenol	16.096	330	185736	119.255	ng	0.00
45) 2-Fluorobiphenyl	13.229	172	844283	115.716	ng	0.00
79) Terphenyl-d14	19.980	244	1428524	112.819	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.440	88	29224m	48.832	ng	
3) Pyridine	3.834	79	94931m	61.556	ng	
4) n-Nitrosodimethylamine	3.746	42	50673	57.457	ng	99
6) Aniline	7.289	93	179780	60.567	ng	97
8) 2-Chlorophenol	7.524	128	130895	60.544	ng	96
10) Phenol	7.148	94	176058	60.703	ng	95
11) bis(2-Chloroethyl)ether	7.389	93	134617	60.203	ng	99
12) 1,3-Dichlorobenzene	7.853	146	140677	59.403	ng	98
13) 1,4-Dichlorobenzene	8.000	146	146182	59.901	ng	95
14) 1,2-Dichlorobenzene	8.317	146	143671	59.751	ng	98
15) Benzyl Alcohol	8.199	79	128587	60.045	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.499	45	255967	60.075	ng	97
17) 2-Methylphenol	8.399	107	120777	60.147	ng	94
18) Hexachloroethane	9.051	117	53040	60.498	ng	97
19) n-Nitroso-di-n-propyla...	8.775	70	116608	59.563	ng	97
20) 3+4-Methylphenols	8.728	107	171213	60.526	ng	99
22) Acetophenone	8.787	105	231145	59.357	ng	# 97
24) Nitrobenzene	9.169	77	170553	64.748	ng	97
25) Isophorone	9.692	82	347779	59.624	ng	99
27) 2,4-Dimethylphenol	9.933	122	100237	59.267	ng	98
28) bis(2-Chloroethoxy)met...	10.179	93	196638	59.549	ng	99
29) 2,4-Dichlorophenol	10.414	162	144143	62.431	ng	98
30) 1,2,4-Trichlorobenzene	10.632	180	162467	60.354	ng	94
31) Naphthalene	10.820	128	471444	59.411	ng	98
32) Benzoic acid	10.080	122	97482m	59.217	ng	
33) 4-Chloroaniline	10.926	127	194656	60.562	ng	99
34) Hexachlorobutadiene	11.108	225	107620	59.597	ng	97
35) Caprolactam	11.713	113	46378	56.530	ng	90
36) 4-Chloro-3-methylphenol	12.042	107	162587	59.075	ng	94
37) 2-Methylnaphthalene	12.430	142	332867	59.262	ng	98
38) 1-Methylnaphthalene	12.647	142	326362	57.424	ng	99
40) 1,2,4,5-Tetrachloroben...	12.794	216	191637	60.600	ng	100
41) Hexachlorocyclopentadiene	12.776	237	77686	64.486	ng	97
43) 2,4,6-Trichlorophenol	13.029	196	126663	61.731	ng	97
44) 2,4,5-Trichlorophenol	13.100	196	142211	61.152	ng	96
46) 1,1'-Biphenyl	13.440	154	447228	58.752	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.481	162	374712	59.532	ng	97
48) 2-Nitroaniline	13.681	65	110830	66.700	ng	99
49) Acenaphthylene	14.328	152	561426	58.071	ng	99
50) Dimethylphthalate	14.063	163	466559	57.301	ng	98
51) 2,6-Dinitrotoluene	14.175	165	94077	60.416	ng	97
52) Acenaphthene	14.668	154	353371	58.837	ng	100
53) 3-Nitroaniline	14.504	138	94402	66.409	ng	97
54) 2,4-Dinitrophenol	14.704	184	38101	59.789	ng	91
55) Dibenzofuran	15.003	168	541505	57.937	ng	99
56) 4-Nitrophenol	14.798	139	74757	64.665	ng	88
57) 2,4-Dinitrotoluene	14.962	165	126527	60.897	ng	# 96
58) Fluorene	15.655	166	425912	56.201	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.221	232	126838	60.003	ng	99
60) Diethylphthalate	15.432	149	464387	56.555	ng	98
61) 4-Chlorophenyl-phenyle...	15.650	204	230798	56.455	ng	96
62) 4-Nitroaniline	15.673	138	96787	63.612	ng	97
63) Azobenzene	15.943	77	448816	58.148	ng	97
65) 4,6-Dinitro-2-methylph...	15.726	198	58456	60.371	ng	98
66) n-Nitrosodiphenylamine	15.861	169	372572	60.679	ng	98
67) 4-Bromophenyl-phenylether	16.543	248	165252	60.706	ng	96
68) Hexachlorobenzene	16.654	284	192103	59.323	ng	98
69) Atrazine	16.813	200	122915	54.596	ng	97
70) Pentachlorophenol	16.995	266	119458	65.215	ng	98
71) Phenanthrene	17.395	178	675731	58.441	ng	99
72) Anthracene	17.489	178	680268	58.323	ng	99
73) Carbazole	17.753	167	625085	57.802	ng	99
74) Di-n-butylphthalate	18.323	149	763037	57.932	ng	98
75) Fluoranthene	19.410	202	844899	57.407	ng	98
77) Benzidine	19.592	184	313281	74.125	ng	98
78) Pyrene	19.774	202	877948	58.460	ng	98
80) Butylbenzylphthalate	20.673	149	328025	61.684	ng	99
81) Benzo(a)anthracene	21.607	228	868377	59.145	ng	99
82) 3,3'-Dichlorobenzidine	21.525	252	287989	60.207	ng	96
83) Chrysene	21.678	228	822773	59.089	ng	99
84) Bis(2-ethylhexyl)phtha...	21.537	149	469968	59.577	ng	99
85) Di-n-octyl phthalate	22.747	149	842461	62.211	ng	100
87) Indeno(1,2,3-cd)pyrene	28.487	276	1148345	61.068	ng	99
88) Benzo(b)fluoranthene	23.816	252	942579	60.002	ng	99
89) Benzo(k)fluoranthene	23.887	252	931847	60.637	ng	99
90) Benzo(a)pyrene	24.680	252	846861	60.649	ng	100
91) Dibenzo(a,h)anthracene	28.558	278	957433	61.064	ng	97
92) Benzo(g,h,i)perylene	29.633	276	961436	61.408	ng	97

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

