

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061223\
 Data File : BG057681.D
 Acq On : 12 Jun 2023 14:40
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 06/13/2023

Supervised By :mohammad ahmed 06/13/2023

Quant Time: Jun 13 01:46:10 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061223.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Jun 13 01:42:42 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.941	152	27546	20.000	ng	0.00
21) Naphthalene-d8	10.749	136	115891	20.000	ng	# 0.00
39) Acenaphthene-d10	14.580	164	79139	20.000	ng	0.00
64) Phenanthrene-d10	17.318	188	190621	20.000	ng	0.00
76) Chrysene-d12	21.578	240	174449	20.000	ng	0.00
86) Perylene-d12	24.750	264	204924	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.508	112	36480	20.767	ng	0.00
7) Phenol-d6	7.095	99	53635	21.109	ng	0.00
23) Nitrobenzene-d5	9.104	82	26986	14.270	ng	0.00
42) 2,4,6-Tribromophenol	16.060	330	17421	18.048	ng	0.00
45) 2-Fluorobiphenyl	13.205	172	116782	21.378	ng	0.00
79) Terphenyl-d14	19.938	244	203641	21.560	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.428	88	9002	10.764	ng	95
3) Pyridine	3.822	79	24105	10.286	ng	92
4) n-Nitrosodimethylamine	3.734	42	11464	10.102	ng	# 87
6) Aniline	7.265	93	33502	10.588	ng	92
8) 2-Chlorophenol	7.506	128	17930	10.485	ng	97
9) Benzaldehyde	7.083	77	18304	11.987	ng	97
10) Phenol	7.118	94	26277	10.535	ng	96
11) bis(2-Chloroethyl)ether	7.365	93	19276	10.246	ng	97
12) 1,3-Dichlorobenzene	7.835	146	20163	10.838	ng	98
13) 1,4-Dichlorobenzene	7.982	146	20300	10.935	ng	95
14) 1,2-Dichlorobenzene	8.299	146	19829	10.995	ng	96
15) Benzyl Alcohol	8.176	79	18690	9.924	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.475	45	44119m	10.757	ng	
17) 2-Methylphenol	8.375	107	18561	10.342	ng	98
18) Hexachloroethane	9.028	117	6329	10.021	ng	94
19) n-Nitroso-di-n-propyla...	8.745	70	17429	10.400	ng	90
20) 3+4-Methylphenols	8.698	107	25093	10.323	ng	98
22) Acetophenone	8.763	105	33538	10.335	ng	# 97
24) Nitrobenzene	9.145	77	15612	8.010	ng	99
25) Isophorone	9.662	82	50249	10.189	ng	98
26) 2-Nitrophenol	9.856	139	3279	9.397	ng	# 91
27) 2,4-Dimethylphenol	9.909	122	18107	10.059	ng	95
28) bis(2-Chloroethoxy)met...	10.156	93	26067	9.944	ng	# 95
29) 2,4-Dichlorophenol	10.391	162	15777	9.465	ng	91
30) 1,2,4-Trichlorobenzene	10.608	180	20662	10.229	ng	100
31) Naphthalene	10.796	128	64565	10.446	ng	99
32) Benzoic acid	9.973	122	4831	10.112	ng	# 79
33) 4-Chloroaniline	10.902	127	27789	10.160	ng	98
34) Hexachlorobutadiene	11.084	225	13361	10.276	ng	96
35) Caprolactam	11.642	113	4960	8.772	ng	96
36) 4-Chloro-3-methylphenol	12.018	107	20776	9.619	ng	98
37) 2-Methylnaphthalene	12.406	142	45656	10.336	ng	90
38) 1-Methylnaphthalene	12.623	142	44012	10.578	ng	96
40) 1,2,4,5-Tetrachloroben...	12.770	216	23615	10.092	ng	97
41) Hexachlorocyclopentadiene	12.753	237	9845	8.735	ng	96
43) 2,4,6-Trichlorophenol	13.005	196	13103	9.014	ng	90

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.076	196	16439	9.486	ng	97
46) 1,1'-Biphenyl	13.411	154	57913	10.352	ng	98
47) 2-Chloronaphthalene	13.452	162	44269	10.376	ng	97
48) 2-Nitroaniline	13.652	65	6918	10.572	ng	95
49) Acenaphthylene	14.298	152	74705	10.414	ng	99
50) Dimethylphthalate	14.028	163	59889	10.252	ng	98
51) 2,6-Dinitrotoluene	14.145	165	5666	10.290	ng	90
52) Acenaphthene	14.644	154	44624	10.339	ng	95
53) 3-Nitroaniline	14.474	138	6772	9.931	ng #	92
54) 2,4-Dinitrophenol	14.674	184	1882	10.952	ng #	86
55) Dibenzofuran	14.974	168	76300	10.634	ng	99
56) 4-Nitrophenol	14.768	139	5304	10.831	ng	97
57) 2,4-Dinitrotoluene	14.927	165	6906	10.714	ng #	96
58) Fluorene	15.626	166	60536	10.746	ng	96
59) 2,3,4,6-Tetrachlorophenol	15.197	232	12572	9.056	ng #	97
60) Diethylphthalate	15.397	149	62330	10.384	ng	96
61) 4-Chlorophenyl-phenyle...	15.620	204	32454	10.669	ng	96
62) 4-Nitroaniline	15.632	138	7640	9.688	ng #	84
63) Azobenzene	15.914	77	65851	10.459	ng	98
65) 4,6-Dinitro-2-methylph...	15.690	198	3075	10.772	ng	84
66) n-Nitrosodiphenylamine	15.831	169	53766	10.616	ng	99
67) 4-Bromophenyl-phenylether	16.513	248	20806	10.311	ng	98
68) Hexachlorobenzene	16.625	284	20599	10.196	ng	90
69) Atrazine	16.777	200	18145	10.342	ng	98
70) Pentachlorophenol	16.965	266	10981	8.383	ng	94
71) Phenanthrene	17.365	178	99685	10.702	ng	100
72) Anthracene	17.453	178	100854	10.618	ng	99
73) Carbazole	17.717	167	93769	10.624	ng	98
74) Di-n-butylphthalate	18.293	149	107077	10.156	ng	98
75) Fluoranthene	19.374	202	120982	10.623	ng	97
77) Benzidine	19.556	184	43172	9.749	ng	99
78) Pyrene	19.733	202	127270	10.308	ng	98
80) Butylbenzylphthalate	20.632	149	37670	8.716	ng	96
81) Benzo(a)anthracene	21.560	228	125165	10.328	ng	98
82) 3,3'-Dichlorobenzidine	21.478	252	39363	9.809	ng	98
83) Chrysene	21.625	228	117611	10.231	ng	98
84) Bis(2-ethylhexyl)phtha...	21.489	149	63529	9.607	ng	99
85) Di-n-octyl phthalate	22.688	149	97177	8.701	ng	98
87) Indeno(1,2,3-cd)pyrene	28.352	276	132470	9.982	ng	99
88) Benzo(b)fluoranthene	23.740	252	113435	10.172	ng	99
89) Benzo(k)fluoranthene	23.804	252	116659	10.267	ng	99
90) Benzo(a)pyrene	24.598	252	111135	10.142	ng	99
91) Dibenzo(a,h)anthracene	28.422	278	111211	10.046	ng	100
92) Benzo(g,h,i)perylene	29.474	276	110247	10.087	ng	98

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

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