

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122722\
 Data File : BG056128.D
 Acq On : 27 Dec 2022 16:17
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDICC010

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 12/28/2022
 Supervised By :Jagrut Upadhyay 12/28/2022

Quant Time: Dec 28 04:53:30 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG122722.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Dec 28 04:51:15 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.336	152	42891	20.000	ng	0.00
21) Naphthalene-d8	11.168	136	167757	20.000	ng	# 0.00
39) Acenaphthene-d10	14.945	164	151521	20.000	ng	0.00
64) Phenanthrene-d10	17.678	188	411709	20.000	ng	0.00
76) Chrysene-d12	21.973	240	400467	20.000	ng	0.00
86) Perylene-d12	25.427	264	440912	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.856	112	50141	27.416	ng	0.00
7) Phenol-d6	7.466	99	74566	31.700	ng	0.00
23) Nitrobenzene-d5	9.505	82	64939	27.051	ng	0.00
42) 2,4,6-Tribromophenol	16.420	330	37767	10.337	ng	0.00
45) 2-Fluorobiphenyl	13.571	172	217093	20.094	ng	0.00
79) Terphenyl-d14	20.251	244	458622	20.240	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.676	88	11850m	22.264	ng	
3) Pyridine	4.099	79	33336	19.032	ng	# 92
4) n-Nitrosodimethylamine	4.005	42	6670	4.236	ng	# 91
6) Aniline	7.648	93	50560	17.953	ng	96
8) 2-Chlorophenol	7.889	128	24107	10.217	ng	93
9) Benzaldehyde	7.460	77	24676	16.704	ng	84
10) Phenol	7.495	94	38773	15.061	ng	99
11) bis(2-Chloroethyl)ether	7.736	93	26768	15.963	ng	90
12) 1,3-Dichlorobenzene	8.224	146	31855	10.437	ng	97
13) 1,4-Dichlorobenzene	8.371	146	30625	9.818	ng	92
14) 1,2-Dichlorobenzene	8.700	146	29252	9.813	ng	98
15) Benzyl Alcohol	8.565	79	26398	10.801	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.847	45	4861m	2.227	ng	
17) 2-Methylphenol	8.759	107	27325	13.490	ng	99
18) Hexachloroethane	9.440	117	9164	9.678	ng	93
19) n-Nitroso-di-n-propyla...	9.135	70	23623	14.011	ng	96
20) 3+4-Methylphenols	9.088	107	37792	13.075	ng	93
22) Acetophenone	9.164	105	48952	11.771	ng	# 97
24) Nitrobenzene	9.546	77	32429	13.421	ng	94
25) Isophorone	10.069	82	70616	14.028	ng	97
26) 2-Nitrophenol	10.269	139	16181	11.312	ng	# 87
27) 2,4-Dimethylphenol	10.310	122	31979	13.312	ng	96
28) bis(2-Chloroethoxy)met...	10.545	93	35824	16.103	ng	96
29) 2,4-Dichlorophenol	10.803	162	32289	9.470	ng	99
30) 1,2,4-Trichlorobenzene	11.027	180	38035	8.581	ng	96
31) Naphthalene	11.220	128	87332	9.866	ng	97
32) Benzoic acid	10.374	122	19692m	10.626	ng	
33) 4-Chloroaniline	11.314	127	41139	10.841	ng	98
34) Hexachlorobutadiene	11.497	225	27372	6.569	ng	94
35) Caprolactam	12.061	113	11316	11.912	ng	98
36) 4-Chloro-3-methylphenol	12.407	107	32248	10.086	ng	98
37) 2-Methylnaphthalene	12.795	142	74947	9.584	ng	95
38) 1-Methylnaphthalene	13.012	142	69133	9.074	ng	94
40) 1,2,4,5-Tetrachloroben...	13.153	216	54500	8.525	ng	99
41) Hexachlorocyclopentadiene	13.136	237	28110	7.751	ng	99
43) 2,4,6-Trichlorophenol	13.383	196	35016	9.253	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.453	196	43189	10.124	ng	95
46) 1,1'-Biphenyl	13.782	154	105921	10.313	ng	97
47) 2-Chloronaphthalene	13.835	162	83754	10.036	ng	95
48) 2-Nitroaniline	14.023	65	17159	12.292	ng	97
49) Acenaphthylene	14.669	152	137004	10.285	ng	96
50) Dimethylphthalate	14.381	163	123025	9.391	ng	99
51) 2,6-Dinitrotoluene	14.505	165	25939	12.796	ng	94
52) Acenaphthene	15.010	154	85336	9.543	ng	93
53) 3-Nitroaniline	14.834	138	23334	11.316	ng	# 76
54) 2,4-Dinitrophenol	15.039	184	17349	11.512	ng	# 85
55) Dibenzofuran	15.339	168	146888	9.768	ng	94
56) 4-Nitrophenol	15.122	139	18218	11.025	ng	98
57) 2,4-Dinitrotoluene	15.286	165	34942	11.419	ng	89
58) Fluorene	15.985	166	120341	9.790	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.557	232	40204m	8.250	ng	
60) Diethylphthalate	15.727	149	115070	8.987	ng	98
61) 4-Chlorophenyl-phenyle...	15.968	204	73853	8.819	ng	95
62) 4-Nitroaniline	15.985	138	25481	11.347	ng	88
63) Azobenzene	16.262	77	90864	11.659	ng	96
65) 4,6-Dinitro-2-methylph...	16.044	198	24763	12.591	ng	96
66) n-Nitrosodiphenylamine	16.179	169	112233	10.403	ng	95
67) 4-Bromophenyl-phenylether	16.861	248	50994	8.453	ng	96
68) Hexachlorobenzene	16.984	284	47192	6.351	ng	98
69) Atrazine	17.108	200	48390	9.906	ng	97
70) Pentachlorophenol	17.325	266	34788	8.189	ng	95
71) Phenanthrene	17.719	178	213851	9.912	ng	96
72) Anthracene	17.813	178	219066	10.263	ng	99
73) Carbazole	18.077	167	185738	10.551	ng	98
74) Di-n-butylphthalate	18.606	149	181879	8.884	ng	100
75) Fluoranthene	19.710	202	281952	9.332	ng	98
77) Benzidine	19.875	184	140168	19.930	ng	99
78) Pyrene	20.069	202	286541	11.825	ng	98
80) Butylbenzylphthalate	20.933	149	77622	11.304	ng	98
81) Benzo(a)anthracene	21.955	228	279946	9.934	ng	97
82) 3,3'-Dichlorobenzidine	21.855	252	99837	9.847	ng	97
83) Chrysene	22.025	228	262534	9.940	ng	98
84) Bis(2-ethylhexyl)phtha...	21.826	149	113450	11.121	ng	98
85) Di-n-octyl phthalate	23.112	149	194744	11.254	ng	100
87) Indeno(1,2,3-cd)pyrene	29.399	276	315023	8.885	ng	# 96
88) Benzo(b)fluoranthene	24.317	252	272349	9.544	ng	98
89) Benzo(k)fluoranthene	24.387	252	283075	9.943	ng	99
90) Benzo(a)pyrene	25.257	252	271843	11.270	ng	98
91) Dibenzo(a,h)anthracene	29.470	278	269355	9.077	ng	97
92) Benzo(g,h,i)perylene	30.645	276	256170	8.928	ng	95

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

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