

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081823\
 Data File : BG058771.D
 Acq On : 18 Aug 2023 11:20
 Operator : MA/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 08/22/2023
 Supervised By :mohammad ahmed 08/22/2023

Quant Time: Aug 21 01:25:25 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG081823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 21 01:23:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.813	152	25070	20.000	ng	0.00
21) Naphthalene-d8	10.615	136	99130	20.000	ng	0.00
39) Acenaphthene-d10	14.458	164	65151	20.000	ng	0.00
64) Phenanthrene-d10	17.202	188	164034	20.000	ng	0.00
76) Chrysene-d12	21.450	240	163374	20.000	ng	0.00
86) Perylene-d12	24.522	264	195063	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.386	112	15177	9.742	ng	0.00
7) Phenol-d6	6.984	99	21144	9.674	ng	0.00
23) Nitrobenzene-d5	8.976	82	19441	9.429	ng	0.00
42) 2,4,6-Tribromophenol	15.956	330	9389	9.191	ng	0.00
45) 2-Fluorobiphenyl	13.077	172	49059	10.952	ng	0.00
79) Terphenyl-d14	19.828	244	96497	10.606	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.271	88	3920	5.270	ng	91
3) Pyridine	3.676	79	8855m	4.559	ng	
4) n-Nitrosodimethylamine	3.594	42	5901m	5.608	ng	
6) Aniline	7.137	93	13129	4.914	ng	97
8) 2-Chlorophenol	7.372	128	7893	5.011	ng	95
9) Benzaldehyde	6.955	77	7101	5.795	ng	# 80
10) Phenol	7.014	94	10114	4.799	ng	96
11) bis(2-Chloroethyl)ether	7.231	93	8948	5.441	ng	92
12) 1,3-Dichlorobenzene	7.701	146	8901	5.320	ng	93
13) 1,4-Dichlorobenzene	7.842	146	8773	5.185	ng	100
14) 1,2-Dichlorobenzene	8.165	146	8764	5.359	ng	94
15) Benzyl Alcohol	8.053	79	6488	4.152	ng	92
16) 2,2'-oxybis(1-Chloropr...	8.335	45	17810	5.082	ng	95
17) 2-Methylphenol	8.265	107	7488	4.837	ng	94
18) Hexachloroethane	8.888	117	3038	4.928	ng	96
19) n-Nitroso-di-n-propyla...	8.618	70	6596	4.669	ng	# 96
20) 3+4-Methylphenols	8.600	107	9606	4.616	ng	89
22) Acetophenone	8.635	105	12567	4.683	ng	# 91
24) Nitrobenzene	9.023	77	9504	4.634	ng	99
25) Isophorone	9.534	82	19492	4.754	ng	96
26) 2-Nitrophenol	9.728	139	3142	3.663	ng	97
27) 2,4-Dimethylphenol	9.798	122	6452	4.439	ng	93
28) bis(2-Chloroethoxy)met...	10.028	93	10312	4.748	ng	# 89
29) 2,4-Dichlorophenol	10.292	162	6161	4.117	ng	92
30) 1,2,4-Trichlorobenzene	10.480	180	8791	4.874	ng	96
31) Naphthalene	10.662	128	26574	5.106	ng	97
33) 4-Chloroaniline	10.797	127	9389m	4.277	ng	
34) Hexachlorobutadiene	10.938	225	5959	4.895	ng	96
35) Caprolactam	11.538	113	2202	4.114	ng	# 91
36) 4-Chloro-3-methylphenol	11.920	107	8684	4.675	ng	# 85
37) 2-Methylnaphthalene	12.284	142	18604	4.977	ng	95
38) 1-Methylnaphthalene	12.501	142	18241	5.140	ng	98
40) 1,2,4,5-Tetrachloroben...	12.654	216	10283	5.120	ng	99
43) 2,4,6-Trichlorophenol	12.907	196	5407	4.231	ng	96
44) 2,4,5-Trichlorophenol	12.983	196	7140	4.665	ng	# 83
46) 1,1'-Biphenyl	13.294	154	23393	5.212	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Chloronaphthalene	13.336	162	17663	5.050	ng	97
48) 2-Nitroaniline	13.559	65	5673	4.271	ng	86
49) Acenaphthylene	14.182	152	30277	5.220	ng	95
50) Dimethylphthalate	13.917	163	26601	5.273	ng	97
51) 2,6-Dinitrotoluene	14.041	165	4552	4.295	ng	# 86
52) Acenaphthene	14.522	154	18172	5.361	ng	98
53) 3-Nitroaniline	14.387	138	4134	4.044	ng	# 89
55) Dibenzofuran	14.857	168	31164	5.330	ng	98
57) 2,4-Dinitrotoluene	14.846	165	6191	4.188	ng	# 85
58) Fluorene	15.509	166	24844	5.339	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.098	232	5848	4.683	ng	# 94
60) Diethylphthalate	15.274	149	27602	5.307	ng	98
61) 4-Chlorophenyl-phenyle...	15.498	204	13282	5.150	ng	96
62) 4-Nitroaniline	15.568	138	3788	3.652	ng	# 82
63) Azobenzene	15.791	77	24318	5.092	ng	98
66) n-Nitrosodiphenylamine	15.721	169	21439	5.071	ng	95
67) 4-Bromophenyl-phenylether	16.397	248	8709	4.838	ng	94
68) Hexachlorobenzene	16.514	284	9826	5.072	ng	94
69) Atrazine	16.673	200	8423	5.630	ng	92
71) Phenanthrene	17.249	178	42723	5.425	ng	98
72) Anthracene	17.343	178	39663	5.022	ng	97
73) Carbazole	17.619	167	37561	5.193	ng	96
74) Di-n-butylphthalate	18.165	149	47932	5.254	ng	99
75) Fluoranthene	19.264	202	53895	5.519	ng	99
77) Benzidine	19.470	184	16499m	5.635	ng	
78) Pyrene	19.628	202	57354	5.038	ng	99
80) Butylbenzylphthalate	20.515	149	21683	4.697	ng	96
81) Benzo(a)anthracene	21.432	228	57290	5.038	ng	97
82) 3,3'-Dichlorobenzidine	21.361	252	18748	4.775	ng	96
83) Chrysene	21.491	228	58559	5.342	ng	100
84) Bis(2-ethylhexyl)phtha...	21.332	149	32590	4.813	ng	99
85) Di-n-octyl phthalate	22.478	149	53496	4.593	ng	# 99
87) Indeno(1,2,3-cd)pyrene	28.030	276	66414	5.032	ng	98
88) Benzo(b)fluoranthene	23.547	252	53741	4.965	ng	93
89) Benzo(k)fluoranthene	23.606	252	58914	5.278	ng	98
90) Benzo(a)pyrene	24.381	252	52848	4.959	ng	96
91) Dibenzo(a,h)anthracene	28.089	278	53479m	4.934	ng	
92) Benzo(g,h,i)perylene	29.123	276	53937m	4.967	ng	

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

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