

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG051820\
 Data File : BG045415.D
 Acq On : 18 May 2020 16:28
 Operator : CG/JU
 Sample : PB128957BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB128957BS

Manual Integrations
 APPROVED

mohammad
 5/19/2020 3:05:57 PM

Quant Time: May 18 17:04:48 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG051520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 18 15:47:17 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	76624	20.00	ng	0.00
21) Naphthalene-d8	10.78	136	324708	20.00	ng	0.00
39) Acenaphthene-d10	14.61	164	233661	20.00	ng	0.00
64) Phenanthrene-d10	17.36	188	559693	20.00	ng	0.00
76) Chrysene-d12	21.62	240	514145	20.00	ng	0.00
87) Perylene-d12	24.78	264	549152	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	582892	148.67	ng	0.00
7) Phenol-d6	7.21	99	840386	150.59	ng	0.00
23) Nitrobenzene-d5	9.13	82	525763	88.30	ng	0.00
42) 2,4,6-Tribromophenol	16.11	330	392391	131.31	ng	0.00
45) 2-Fluorobiphenyl	13.23	172	1205962	87.54	ng	0.00
79) Terphenyl-d14	19.98	244	2037064	92.41	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.41	88	74580	38.359	ng	97
3) Pyridine	3.82	79	175376	32.387	ng	95
4) n-Nitrosodimethylamine	3.73	42	76774	43.328	ng	92
6) Aniline	7.30	93	284116	39.001	ng	98
8) 2-Chlorophenol	7.56	128	216062	50.251	ng	99
9) Benzaldehyde	7.10	77	152328	41.896	ng	97
10) Phenol	7.24	94	290977	50.592	ng	96
11) bis(2-Chloroethyl)ether	7.39	93	196057	43.703	ng	98
12) 1,3-Dichlorobenzene	7.86	146	226837	43.901	ng	97
13) 1,4-Dichlorobenzene	8.00	146	233663	45.825	ng	98
14) 1,2-Dichlorobenzene	8.32	146	217256	44.299	ng	98
15) Benzyl Alcohol	8.23	79	223019	48.377	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.49	45	193668	45.480	ng	92
17) 2-Methylphenol	8.47	107	202895	47.131	ng	95
18) Hexachloroethane	9.05	117	85021	43.535	ng	95
19) n-Nitroso-di-n-propylamine	8.77	70	184286	47.755	ng	98
20) 3+4-Methylphenols	8.80	107	268616	45.822	ng	# 60
22) Acetophenone	8.79	105	329362	42.797	ng	# 94
24) Nitrobenzene	9.17	77	270138	42.344	ng	95
25) Isophorone	9.70	82	508631	43.374	ng	100
26) 2-Nitrophenol	9.88	139	125235	45.889	ng	93
27) 2,4-Dimethylphenol	9.98	122	216258	53.428	ng	98
28) bis(2-Chloroethoxy)methane	10.18	93	277082	45.055	ng	# 96
29) 2,4-Dichlorophenol	10.46	162	225434	47.164	ng	94
30) 1,2,4-Trichlorobenzene	10.64	180	242677	42.967	ng	98
31) Naphthalene	10.82	128	672811	44.465	ng	99
32) Benzoic acid	10.19	122	156325	54.951	ng	95
33) 4-Chloroaniline	10.95	127	159112	25.157	ng	94
34) Hexachlorobutadiene	11.12	225	162253	40.640	ng	98
35) Caprolactam	11.72	113	82439m	46.989	ng	
36) 4-Chloro-3-methylphenol	12.12	107	265276	49.252	ng	96
37) 2-Methylnaphthalene	12.43	142	513390	45.522	ng	99
38) 1-Methylnaphthalene	12.66	142	490236	46.017	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.80	216	281686	43.160	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.79	237	392085	104.084	ng	99
43) 2,4,6-Trichlorophenol	13.06	196	206464	45.540	ng	99
44) 2,4,5-Trichlorophenol	13.16	196	220192	42.501	ng	98
46) 1,1'-Biphenyl	13.44	154	665676	46.365	ng	100
47) 2-Chloronaphthalene	13.49	162	504073	43.236	ng	98
48) 2-Nitroaniline	13.70	65	168721	43.994	ng	94
49) Acenaphthylene	14.33	152	852293	45.512	ng	99
50) Dimethylphthalate	14.07	163	706702	44.089	ng	99
51) 2,6-Dinitrotoluene	14.18	165	158940	43.943	ng	97
52) Acenaphthene	14.68	154	497446	39.683	ng	97
53) 3-Nitroaniline	14.53	138	103193	28.066	ng	91
54) 2,4-Dinitrophenol	14.73	184	177961	75.444	ng	97
55) Dibenzofuran	15.01	168	813968	43.726	ng	98
56) 4-Nitrophenol	14.91	139	251129	90.215	ng	92
57) 2,4-Dinitrotoluene	14.98	165	227641	43.457	ng	96
58) Fluorene	15.66	166	677789	44.319	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.26	232	197652	43.364	ng	99
60) Diethylphthalate	15.43	149	726869	43.568	ng	97
61) 4-Chlorophenyl-phenylether	15.65	204	370968	42.389	ng	99
62) 4-Nitroaniline	15.70	138	163027	42.472	ng	94
63) Azobenzene	15.95	77	690847	44.409	ng	98
65) 4,6-Dinitro-2-methylphenol	15.75	198	139297	42.734	ng	97
66) n-Nitrosodiphenylamine	15.87	169	614181	45.704	ng	97
67) 4-Bromophenyl-phenylether	16.55	248	257842	42.544	ng	99
68) Hexachlorobenzene	16.67	284	279083	44.027	ng	96
69) Atrazine	16.82	200	238850	43.152	ng	97
70) Pentachlorophenol	17.03	266	260210	79.058	ng	95
71) Phenanthrene	17.40	178	1102997	45.142	ng	100
72) Anthracene	17.49	178	1143797	46.615	ng	97
73) Carbazole	17.77	167	1051617	43.968	ng	100
74) Di-n-butylphthalate	18.33	149	1259486	43.873	ng	99
75) Fluoranthene	19.41	202	1379076	44.374	ng	98
77) Benzidine	19.60	184	712335	49.331	ng	98
78) Pyrene	19.77	202	1363525	46.523	ng	97
80) Butylbenzylphthalate	20.66	149	561774	44.407	ng	99
81) Benzo(a)anthracene	21.60	228	1284949	44.863	ng	97
82) 3,3'-Dichlorobenzidine	21.52	252	328553	29.244	ng	98
83) Chrysene	21.67	228	1241549	45.082	ng	99
84) Bis(2-ethylhexyl)phthalate	21.52	149	771490	44.365	ng	100
85) Di-n-octyl phthalate	22.71	149	1297993	43.459	ng	100
86) Indeno(1,2,3-cd)pyrene	28.37	276	1556938	43.233	ng	# 95
88) Benzo(b)fluoranthene	23.78	252	1322936	44.919	ng	99
89) Benzo(k)fluoranthene	23.85	252	1325840	47.917	ng	99
90) Benzo(a)pyrene	24.64	252	1233794	44.981	ng	# 96
91) Dibenzo(a,h)anthracene	28.43	278	1284525	46.602	ng	100
92) Benzo(g,h,i)perylene	29.49	276	1294468	46.957	ng	98

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

