

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070220\
 Data File : BG045930.D
 Acq On : 3 Jul 2020 2:47
 Operator : CG/JU
 Sample : PB129945BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB129945BS

Manual Integrations
 APPROVED

Sohil
 7/7/2020 8:19:39 AM

Quant Time: Jul 03 08:22:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 14:02:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	47524	20.00	ng	0.00
21) Naphthalene-d8	10.66	136	180470	20.00	ng	0.00
39) Acenaphthene-d10	14.51	164	127566	20.00	ng	0.00
64) Phenanthrene-d10	17.27	188	323337	20.00	ng	0.00
76) Chrysene-d12	21.51	240	323446	20.00	ng	0.00
86) Perylene-d12	24.59	264	345022	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	393359	139.82	ng	0.00
7) Phenol-d6	7.07	99	556843	130.12	ng	0.00
23) Nitrobenzene-d5	9.02	82	357826	90.58	ng	0.00
42) 2,4,6-Tribromophenol	16.01	330	251740	130.71	ng	0.00
45) 2-Fluorobiphenyl	13.13	172	771070	88.83	ng	0.00
79) Terphenyl-d14	19.89	244	1508742	94.53	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.27	88	48723	37.813	ng	93
3) Pyridine	3.68	79	138468	36.264	ng	96
4) n-Nitrosodimethylamine	3.60	42	57541	44.895	ng	97
6) Aniline	7.18	93	176224	34.974	ng	99
8) 2-Chlorophenol	7.43	128	142239	47.234	ng	96
9) Benzaldehyde	6.99	77	60593	23.574	ng	96
10) Phenol	7.10	94	192738	46.483	ng	99
11) bis(2-Chloroethyl)ether	7.27	93	134466	42.695	ng	99
12) 1,3-Dichlorobenzene	7.74	146	153194	42.688	ng	97
13) 1,4-Dichlorobenzene	7.89	146	154327	43.054	ng	98
14) 1,2-Dichlorobenzene	8.21	146	144360	42.067	ng	99
15) Benzyl Alcohol	8.11	79	142145	42.923	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.38	45	126535	42.630	ng	98
17) 2-Methylphenol	8.34	107	127446	41.301	ng	97
18) Hexachloroethane	8.94	117	59827	43.945	ng	95
19) n-Nitroso-di-n-propylamine	8.67	70	121418	43.157	ng	98
20) 3+4-Methylphenols	8.67	107	171996	42.382	ng	# 88
22) Acetophenone	8.68	105	211861	42.771	ng	# 98
24) Nitrobenzene	9.06	77	184282	43.750	ng	99
25) Isophorone	9.58	82	327609	42.838	ng	99
26) 2-Nitrophenol	9.78	139	78768	44.886	ng	94
27) 2,4-Dimethylphenol	9.86	122	129418	48.305	ng	97
28) bis(2-Chloroethoxy)methane	10.07	93	177494	44.431	ng	98
29) 2,4-Dichlorophenol	10.33	162	144046	46.083	ng	98
30) 1,2,4-Trichlorobenzene	10.53	180	158087	42.721	ng	99
31) Naphthalene	10.71	128	428744	43.454	ng	98
32) Benzoic acid	10.06	122	64576	39.191	ng	88
33) 4-Chloroaniline	10.83	127	114153	27.628	ng	98
34) Hexachlorobutadiene	11.00	225	107689	40.919	ng	97
35) Caprolactam	11.60	113	51143m	50.561	ng	
36) 4-Chloro-3-methylphenol	12.00	107	165726	45.919	ng	91
37) 2-Methylnaphthalene	12.33	142	323223	43.993	ng	98
38) 1-Methylnaphthalene	12.55	142	309748	44.551	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.70	216	181765	43.273	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG070220\
 Data File : BG045930.D
 Acq On : 3 Jul 2020 2:47
 Operator : CG/JU
 Sample : PB129945BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleID :
 PB129945BS

Manual Integrations
 APPROVED

Sohil
 7/7/2020 8:19:39 AM

Quant Time: Jul 03 08:22:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 02 14:02:29 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.68	237	174309	75.640	ng	97
43) 2,4,6-Trichlorophenol	12.95	196	120831	42.428	ng	97
44) 2,4,5-Trichlorophenol	13.05	196	125791	38.755	ng	97
46) 1,1'-Biphenyl	13.34	154	401772	43.991	ng	99
47) 2-Chloronaphthalene	13.38	162	315093	42.710	ng	97
48) 2-Nitroaniline	13.59	65	109759	44.726	ng	96
49) Acenaphthylene	14.23	152	518770	43.901	ng	99
50) Dimethylphthalate	13.97	163	443528	44.382	ng	100
51) 2,6-Dinitrotoluene	14.09	165	102172	45.520	ng	97
52) Acenaphthene	14.58	154	311419	43.477	ng	99
53) 3-Nitroaniline	14.43	138	72015	32.121	ng	97
54) 2,4-Dinitrophenol	14.64	184	100646	80.542	ng	95
55) Dibenzofuran	14.92	168	512409	44.036	ng	98
56) 4-Nitrophenol	14.79	139	128668	80.098	ng	95
57) 2,4-Dinitrotoluene	14.89	165	149281	46.419	ng	99
58) Fluorene	15.56	166	437614	45.228	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.16	232	119345	42.364	ng	95
60) Diethylphthalate	15.34	149	463423	45.362	ng	99
61) 4-Chlorophenyl-phenylether	15.56	204	236946	42.489	ng	98
62) 4-Nitroaniline	15.60	138	104822	45.556	ng	97
63) Azobenzene	15.85	77	458526	46.744	ng	99
65) 4,6-Dinitro-2-methylphenol	15.66	198	88595	44.337	ng	98
66) n-Nitrosodiphenylamine	15.77	169	389359	42.515	ng	99
67) 4-Bromophenyl-phenylether	16.45	248	168159	40.025	ng	96
68) Hexachlorobenzene	16.58	284	183959	42.226	ng	98
69) Atrazine	16.73	200	159835	44.976	ng	96
70) Pentachlorophenol	16.93	266	113925	55.598	ng	97
71) Phenanthrene	17.31	178	734466	44.044	ng	97
72) Anthracene	17.39	178	758045	45.650	ng	100
73) Carbazole	17.67	167	713792	45.144	ng	99
74) Di-n-butylphthalate	18.23	149	875795	46.780	ng	98
75) Fluoranthene	19.32	202	973726	47.486	ng	99
77) Benzidine	19.50	184	423043	50.311	ng	99
78) Pyrene	19.68	202	958209	44.240	ng	100
80) Butylbenzylphthalate	20.57	149	404750	44.645	ng	98
81) Benzo(a)anthracene	21.50	228	931355	44.413	ng	98
82) 3,3'-Dichlorobenzidine	21.41	252	256968	32.396	ng	98
83) Chrysene	21.56	228	898538	44.300	ng	100
84) Bis(2-ethylhexyl)phthalate	21.41	149	576377	45.707	ng	100
85) Di-n-octyl phthalate	22.57	149	988104	45.426	ng	100
87) Indeno(1,2,3-cd)pyrene	28.10	276	1173266	46.907	ng	99
88) Benzo(b)fluoranthene	23.63	252	993262	45.915	ng	99
89) Benzo(k)fluoranthene	23.69	252	988714	47.754	ng	99
90) Benzo(a)pyrene	24.45	252	930216	46.033	ng	99
91) Dibenzo(a,h)anthracene	28.15	278	951732	47.450	ng	100
92) Benzo(g,h,i)perylene	29.18	276	951610	47.400	ng	97

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

