

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091718\
 Data File : BG036751.D
 Acq On : 17 Sep 2018 12:49
 Operator : JU/SJ
 Sample : PB112896BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112896BS

Manual Integrations
 APPROVED

Sohil
 9/18/2018 6:18:11 PM

Quant Time: Sep 17 18:12:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 31 13:03:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	64647	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	282007	20.00	ng	-0.01
38) Acenaphthene-d10	14.67	164	183143	20.00	ng	-0.01
63) Phenanthrene-d10	17.42	188	459690	20.00	ng	-0.01
75) Chrysene-d12	21.68	240	463203	20.00	ng	-0.01
86) Perylene-d12	24.89	264	447260	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	445997	109.44	ng	0.00
7) Phenol-d6	7.21	99	601561	103.10	ng	0.00
23) Nitrobenzene-d5	9.21	82	359769	69.22	ng	-0.01
41) 2,4,6-Tribromophenol	16.17	330	244248	111.77	ng	0.00
44) 2-Fluorobiphenyl	13.30	172	841649	65.62	ng	-0.01
78) Terphenyl-d14	20.03	244	1354217	68.33	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.53	88	64342	33.830	ng	99
3) Pyridine	3.93	79	172666	32.343	ng	100
4) n-Nitrosodimethylamine	3.85	42	96278	40.490	ng	95
6) Aniline	7.38	93	228707	30.880	ng	98
8) 2-Chlorophenol	7.62	128	178475	40.384	ng	97
9) Benzaldehyde	7.19	77	102378	25.479	ng	96
10) Phenol	7.24	94	245757	40.248	ng	98
11) bis(2-Chloroethyl)ether	7.47	93	173697	35.881	ng	98
12) 1,3-Dichlorobenzene	7.94	146	184635	36.587	ng	98
13) 1,4-Dichlorobenzene	8.09	146	187995	36.898	ng	98
14) 1,2-Dichlorobenzene	8.41	146	178239	36.631	ng	98
15) Benzyl Alcohol	8.29	79	179887	38.243	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.58	45	346525	35.496	ng	98
17) 2-Methylphenol	8.49	107	165869	40.910	ng	99
18) Hexachloroethane	9.13	117	69489	38.040	ng	94
19) n-Nitroso-di-n-propylamine	8.86	70	151774	34.804	ng	94
20) 3+4-Methylphenols	8.82	107	230830	40.778	ng	99
22) Acetophenone	8.87	105	275764	37.238	ng	98
24) Nitrobenzene	9.26	77	220943	39.420	ng	98
25) Isophorone	9.78	82	409189	39.630	ng	97
26) 2-Nitrophenol	9.96	139	96178	43.304	ng	98
27) 2,4-Dimethylphenol	10.02	122	176771	42.990	ng	99
28) bis(2-Chloroethoxy)methane	10.26	93	244140	38.526	ng	99
29) 2,4-Dichlorophenol	10.50	162	189999	42.162	ng	96
30) 1,2,4-Trichlorobenzene	10.71	180	193179	36.644	ng	97
31) Naphthalene	10.91	128	557176	39.524	ng	98
32) Benzoic acid	10.14	122	73038	26.202	ng	99
33) 4-Chloroaniline	11.01	127	173548	26.865	ng	97
34) Hexachlorobutadiene	11.19	225	133023	37.556	ng	98
35) Caprolactam	11.79	113	67700m	37.712	ng	
36) 4-Chloro-3-methylphenol	12.12	107	207673	39.660	ng	99
37) 2-Methylnaphthalene	12.51	142	429479	41.636	ng	99
39) 1,2,4,5-Tetrachlorobenzene	12.87	216	242546	37.423	ng	97
40) Hexachlorocyclopentadiene	12.85	237	294063	87.203	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.11	196	163206	41.339	ng	97
43) 2,4,5-Trichlorophenol	13.18	196	179493	42.625	ng	98
45) 1,1'-Biphenyl	13.51	154	553246	36.392	ng	99
46) 2-Chloronaphthalene	13.55	162	421428	36.312	ng	99
47) 2-Nitroaniline	13.75	65	152316	41.795	ng	97
48) Acenaphthylene	14.40	152	709923	39.140	ng	99
49) Dimethylphthalate	14.13	163	567803	37.968	ng	98
50) 2,6-Dinitrotoluene	14.25	165	125732	40.858	ng	98
51) Acenaphthene	14.74	154	420974	36.240	ng	99
52) 3-Nitroaniline	14.57	138	104578	31.536	ng	99
53) 2,4-Dinitrophenol	14.78	184	113106	97.037	ng	96
54) Dibenzofuran	15.07	168	677304	37.217	ng	99
55) 4-Nitrophenol	14.89	139	206194	76.048	ng	98
56) 2,4-Dinitrotoluene	15.03	165	176074	43.902	ng	94
57) Fluorene	15.72	166	541500	39.802	ng	96
58) 2,3,4,6-Tetrachlorophenol	15.30	232	176415	44.589	ng	97
59) Diethylphthalate	15.49	149	560001	37.157	ng	99
60) 4-Chlorophenyl-phenylether	15.71	204	309786	36.876	ng	100
61) 4-Nitroaniline	15.74	138	137343	39.579	ng	96
62) Azobenzene	16.01	77	562049	36.653	ng	97
64) 4,6-Dinitro-2-methylphenol	15.80	198	96314	47.696	ng	98
65) n-Nitrosodiphenylamine	15.93	169	498099	36.467	ng	98
66) 4-Bromophenyl-phenylether	16.61	248	204421	37.982	ng	99
67) Hexachlorobenzene	16.73	284	205988	37.430	ng	97
68) Atrazine	16.87	200	202556	39.509	ng	97
69) Pentachlorophenol	17.07	266	242690	64.886	ng	97
70) Phenanthrene	17.46	178	907367	38.846	ng	99
71) Anthracene	17.55	178	927850	39.849	ng	99
72) Carbazole	17.82	167	812990	34.962	ng	99
73) Di-n-butylphthalate	18.38	149	938753	36.253	ng	100
74) Fluoranthene	19.47	202	1113358	39.095	ng	100
76) Benzidine	19.65	184	402827	27.258	ng	98
77) Pyrene	19.83	202	1077162	37.816	ng	99
79) Butylbenzylphthalate	20.71	149	427224	37.688	ng	98
80) Benzo(a)anthracene	21.67	228	1081212	38.530	ng	100
81) 3,3'-Dichlorobenzidine	21.58	252	302634	27.060	ng	96
82) Chrysene	21.73	228	1000351	37.609	ng	99
83) Bis(2-ethylhexyl)phthalate	21.57	149	582102	36.696	ng	99
84) Di-n-octyl phthalate	22.78	149	990638	37.388	ng	98
85) Indeno(1,2,3-cd)pyrene	28.54	276	1197107	38.749	ng	99
87) Benzo(b)fluoranthene	23.88	252	1091875	43.963	ng	96
88) Benzo(k)fluoranthene	23.95	252	1008789	41.576	ng	98
89) Benzo(a)pyrene	24.75	252	1037800	43.484	ng	99
90) Dibenzo(a,h)anthracene	28.61	278	975350	42.333	ng	97
91) Benzo(g,h,i)perylene	29.68	276	1001618	43.260	ng	97

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

