

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG111918\
 Data File : BG038078.D
 Acq On : 20 Nov 2018 1:54
 Operator : JU/SJ
 Sample : PB114916BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB114916BS

Manual Integrations
 APPROVED

Sohil
 11/20/2018 9:41:50 AM

Quant Time: Nov 20 03:05:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG111318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 13 16:39:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.09	152	42401	20.00	ng	0.00
21) Naphthalene-d8	10.91	136	179591	20.00	ng	0.00
39) Acenaphthene-d10	14.72	164	111515	20.00	ng	0.00
64) Phenanthrene-d10	17.47	188	255867	20.00	ng	0.00
76) Chrysene-d12	21.75	240	235623	20.00	ng	0.00
87) Perylene-d12	25.07	264	250937	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	349556	142.34	ng	0.00
7) Phenol-d6	7.24	99	482334	137.59	ng	0.00
23) Nitrobenzene-d5	9.26	82	293660	87.53	ng	0.00
42) 2,4,6-Tribromophenol	16.21	330	164894	129.65	ng	0.00
45) 2-Fluorobiphenyl	13.34	172	659263	86.13	ng	0.00
79) Terphenyl-d14	20.07	244	984474	98.31	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	38369	33.077	ng	96
3) Pyridine	3.92	79	114075	34.475	ng	95
4) n-Nitrosodimethylamine	3.84	42	58590	48.806	ng	97
6) Aniline	7.41	93	70707	15.628	ng	97
8) 2-Chlorophenol	7.65	128	134427	47.175	ng	97
9) Benzaldehyde	7.22	77	69274	27.326	ng	97
10) Phenol	7.27	94	175323	47.217	ng	97
11) bis(2-Chloroethyl)ether	7.51	93	125218	41.308	ng	97
12) 1,3-Dichlorobenzene	7.97	146	134533	40.982	ng	98
13) 1,4-Dichlorobenzene	8.12	146	136862	41.272	ng	98
14) 1,2-Dichlorobenzene	8.44	146	129952	41.046	ng	98
15) Benzyl Alcohol	8.33	79	117933	46.493	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.61	45	244914	43.510	ng	98
17) 2-Methylphenol	8.52	107	119397	47.562	ng	98
18) Hexachloroethane	9.17	117	49176	40.747	ng	93
19) n-Nitroso-di-n-propylamine	8.89	70	102602	40.450	ng	98
20) 3+4-Methylphenols	8.85	107	165176	46.415	ng	98
22) Acetophenone	8.92	105	191630	42.886	ng	# 98
24) Nitrobenzene	9.30	77	152958	44.569	ng	99
25) Isophorone	9.82	82	286545	47.453	ng	97
26) 2-Nitrophenol	10.02	139	74755	43.631	ng	96
27) 2,4-Dimethylphenol	10.06	122	137651	53.323	ng	96
28) bis(2-Chloroethoxy)methane	10.30	93	174509	45.194	ng	94
29) 2,4-Dichlorophenol	10.54	162	139715	48.117	ng	99
30) 1,2,4-Trichlorobenzene	10.76	180	138195	42.472	ng	95
31) Naphthalene	10.96	128	407452	46.127	ng	100
32) Benzoic acid	10.19	122	70438	43.822	ng	99
33) 4-Chloroaniline	11.07	127	42321	10.300	ng	95
34) Hexachlorobutadiene	11.22	225	81746	40.821	ng	98
35) Caprolactam	11.85	113	47392m	40.777	ng	
36) 4-Chloro-3-methylphenol	12.17	107	147307	46.436	ng	96
37) 2-Methylnaphthalene	12.55	142	302603	47.692	ng	99
38) 1-Methylnaphthalene	12.78	142	285641	46.769	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.92	216	153770	41.266	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	200183	100.324	nq	97
43) 2,4,6-Trichlorophenol	13.15	196	112777	44.414	nq	96
44) 2,4,5-Trichlorophenol	13.23	196	123124	46.083	nq	97
46) 1,1'-Biphenyl	13.56	154	381663	40.739	nq	99
47) 2-Chloronaphthalene	13.61	162	303434	42.445	nq	99
48) 2-Nitroaniline	13.81	65	104531	46.459	nq	99
49) Acenaphthylene	14.45	152	507673	47.223	nq	99
50) Dimethylphthalate	14.17	163	391819	44.320	nq	100
51) 2,6-Dinitrotoluene	14.30	165	91092	45.526	nq	94
52) Acenaphthene	14.79	154	287793	44.467	nq	99
53) 3-Nitroaniline	14.63	138	32942	14.920	nq #	94
54) 2,4-Dinitrophenol	14.84	184	92912	85.338	nq	92
55) Dibenzofuran	15.12	168	465592	43.506	nq	98
56) 4-Nitrophenol	14.93	139	148367	87.129	nq	95
57) 2,4-Dinitrotoluene	15.09	165	127701	46.908	nq	96
58) Fluorene	15.77	166	375886	47.075	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.34	232	106997	47.859	nq	99
60) Diethylphthalate	15.53	149	406923	43.335	nq	100
61) 4-Chlorophenyl-phenylether	15.76	204	195381	41.818	nq	99
62) 4-Nitroaniline	15.80	138	96127	43.827	nq	93
63) Azobenzene	16.05	77	377793	44.998	nq	99
65) 4,6-Dinitro-2-methylphenol	15.84	198	68563	42.366	nq	96
66) n-Nitrosodiphenylamine	15.97	169	341209	44.080	nq	99
67) 4-Bromophenyl-phenylether	16.65	248	122004	43.443	nq	96
68) Hexachlorobenzene	16.76	284	125640	43.342	nq	98
69) Atrazine	16.91	200	127726	44.761	nq	99
70) Pentachlorophenol	17.11	266	148690	75.525	nq	96
71) Phenanthrene	17.51	178	628607	47.985	nq	98
72) Anthracene	17.61	178	634706	49.152	nq	99
73) Carbazole	17.88	167	579134	44.701	nq	100
74) Di-n-butylphthalate	18.41	149	690077	47.450	nq	99
75) Fluoranthene	19.52	202	736763	49.294	nq	99
77) Benzidine	19.70	184	184770	22.348	nq	99
78) Pyrene	19.88	202	744991	48.360	nq	99
80) Butylbenzylphthalate	20.75	149	323063	47.954	nq	98
81) Benzo(a)anthracene	21.73	228	696283	48.900	nq	99
82) 3,3'-Dichlorobenzidine	21.65	252	124392	22.427	nq	97
83) Chrysene	21.80	228	652030	47.860	nq	98
84) Bis(2-ethylhexyl)phthalate	21.60	149	451946	47.041	nq	99
85) Di-n-octyl phthalate	22.84	149	769120	49.483	nq	100
86) Indeno(1,2,3-cd)pyrene	28.89	276	773921	51.148	nq	97
88) Benzo(b)fluoranthene	24.02	252	679441	49.202	nq	99
89) Benzo(k)fluoranthene	24.08	252	676360	49.637	nq	99
90) Benzo(a)pyrene	24.92	252	673107	51.004	nq	98
91) Dibenzo(a,h)anthracene	28.96	278	634869	49.998	nq	98
92) Benzo(g,h,i)perylene	30.09	276	618157	48.960	nq	99

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

