

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090618\
 Data File : BG036590.D
 Acq On : 7 Sep 2018 10:32
 Operator : JU/SJ
 Sample : PB112623BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB112623BS

Quant Time: Sep 07 11:59:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 23 12:07:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	56328	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	241633	20.00	ng	-0.02
38) Acenaphthene-d10	14.68	164	166801	20.00	ng	-0.02
63) Phenanthrene-d10	17.42	188	428537	20.00	ng	-0.02
75) Chrysene-d12	21.69	240	427386	20.00	ng	-0.02
86) Perylene-d12	24.90	264	437349	20.00	ng	-0.04

System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	421550	118.72	ng	0.00
7) Phenol-d6	7.21	99	585189	115.10	ng	0.00
23) Nitrobenzene-d5	9.22	82	348829	78.33	ng	-0.02
41) 2,4,6-Tribromophenol	16.17	330	254323	127.78	ng	-0.02
44) 2-Fluorobiphenyl	13.30	172	843376	72.19	ng	-0.02
78) Terphenyl-d14	20.03	244	1408726	77.04	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.53	88	45862	27.675	ng	99
3) Pyridine	3.92	79	145495	31.278	ng	95
4) n-Nitrosodimethylamine	3.84	42	73656	35.551	ng	82
6) Aniline	7.38	93	185323	28.718	ng	98
8) 2-Chlorophenol	7.62	128	157152	40.811	ng	96
9) Benzaldehyde	7.19	77	100706	28.765	ng	99
10) Phenol	7.24	94	225676	42.417	ng	99
11) bis(2-Chloroethyl)ether	7.48	93	148613	35.234	ng	96
12) 1,3-Dichlorobenzene	7.95	146	159884	36.362	ng	98
13) 1,4-Dichlorobenzene	8.09	146	164271	37.003	ng	98
14) 1,2-Dichlorobenzene	8.41	146	157273	37.096	ng	97
15) Benzyl Alcohol	8.29	79	161207	39.334	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.58	45	278705	32.765	ng	98
17) 2-Methylphenol	8.49	107	148835	42.130	ng	99
18) Hexachloroethane	9.14	117	58318	36.640	ng	93
19) n-Nitroso-di-n-propylamine	8.86	70	129117	33.982	ng	96
20) 3+4-Methylphenols	8.82	107	206781	41.925	ng	99
22) Acetophenone	8.87	105	244057	38.463	ng	98
24) Nitrobenzene	9.26	77	194048	40.407	ng	97
25) Isophorone	9.78	82	363174	41.050	ng	98
26) 2-Nitrophenol	9.97	139	90171	47.383	ng	94
27) 2,4-Dimethylphenol	10.03	122	163181	46.316	ng	99
28) bis(2-Chloroethoxy)methane	10.26	93	214057	39.423	ng	97
29) 2,4-Dichlorophenol	10.51	162	176911	45.817	ng	96
30) 1,2,4-Trichlorobenzene	10.73	180	175885	38.938	ng	100
31) Naphthalene	10.91	128	504367	41.756	ng	99
32) Benzoic acid	10.14	122	94637	37.329	ng	94
33) 4-Chloroaniline	11.01	127	137386	24.821	ng	97
34) Hexachlorobutadiene	11.20	225	117356	38.669	ng	96
35) Caprolactam	11.78	113	67887	44.135	ng	94
36) 4-Chloro-3-methylphenol	12.13	107	196398	43.774	ng	99
37) 2-Methylnaphthalene	12.51	142	398568	45.096	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.88	216	228054	38.634	ng	100
40) Hexachlorocyclopentadiene	12.86	237	272625	88.766	ng	98

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42) 2,4,6-Trichlorophenol	13.11	196	155276	43.183	ng	99
43) 2,4,5-Trichlorophenol	13.19	196	171146	44.624	ng	95
45) 1,1'-Biphenyl	13.52	154	510386	36.862	ng	98
46) 2-Chloronaphthalene	13.56	162	404669	38.284	ng	100
47) 2-Nitroaniline	13.76	65	145235	43.756	ng	96
48) Acenaphthylene	14.40	152	679778	41.150	ng	99
49) Dimethylphthalate	14.13	163	547295	40.182	ng	99
50) 2,6-Dinitrotoluene	14.25	165	121258	43.265	ng	94
51) Acenaphthene	14.74	154	408195	38.583	ng	98
52) 3-Nitroaniline	14.58	138	101108	33.477	ng	96
53) 2,4-Dinitrophenol	14.79	184	82118	77.354	ng	97
54) Dibenzofuran	15.08	168	654249	39.472	ng	99
55) 4-Nitrophenol	14.88	139	191090	77.382	ng	98
56) 2,4-Dinitrotoluene	15.04	165	173689	47.551	ng	87
57) Fluorene	15.73	166	526148	42.463	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.30	232	172512	47.875	ng	94
59) Diethylphthalate	15.50	149	538336	39.219	ng	98
60) 4-Chlorophenyl-phenylether	15.72	204	301677	39.429	ng	95
61) 4-Nitroaniline	15.74	138	134881	42.677	ng	92
62) Azobenzene	16.01	77	526014	37.664	ng	96
64) 4,6-Dinitro-2-methylphenol	15.80	198	81563	43.327	ng	95
65) n-Nitrosodiphenylamine	15.93	169	484700	38.066	ng	98
66) 4-Bromophenyl-phenylether	16.61	248	198950	39.653	ng	99
67) Hexachlorobenzene	16.73	284	203367	39.640	ng	97
68) Atrazine	16.88	200	210554	44.054	ng	98
69) Pentachlorophenol	17.07	266	224534	64.396	ng	99
70) Phenanthrene	17.46	178	902208	41.433	ng	99
71) Anthracene	17.56	178	912260	42.028	ng	99
72) Carbazole	17.82	167	803042	37.045	ng	99
73) Di-n-butylphthalate	18.38	149	911149	37.745	ng	100
74) Fluoranthene	19.47	202	1088468	40.999	ng	99
76) Benzidine	19.65	184	423111	31.030	ng	100
77) Pyrene	19.83	202	1066771	40.590	ng	98
79) Butylbenzylphthalate	20.72	149	422671	40.411	ng	94
80) Benzo(a)anthracene	21.67	228	1063399	41.071	ng	99
81) 3,3'-Dichlorobenzidine	21.58	252	280560	27.189	ng	99
82) Chrysene	21.74	228	987132	40.222	ng	98
83) Bis(2-ethylhexyl)phthalate	21.58	149	567486	38.772	ng	99
84) Di-n-octyl phthalate	22.79	149	970036	39.679	ng	96
85) Indeno(1,2,3-cd)pyrene	28.55	276	1094855	38.409	ng	# 95
87) Benzo(b)fluoranthene	23.88	252	1058051	43.566	ng	98
88) Benzo(k)fluoranthene	23.95	252	1006009	42.400	ng	99
89) Benzo(a)pyrene	24.75	252	1010696	43.308	ng	99
90) Dibenzo(a,h)anthracene	28.61	278	890374	39.520	ng	98
91) Benzo(g,h,i)perylene	29.70	276	895187	39.539	ng	97

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

