

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011119\
 Data File : BG039091.D
 Acq On : 11 Jan 2019 18:48
 Operator : JU/SJ
 Sample : PB116292BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116292BS

Manual Integrations
 APPROVED

Sohil
 1/14/2019 4:31:12 PM

Quant Time: Jan 12 00:34:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 09 14:23:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	24798	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	109485	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	74229	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	193903	20.00	ng	0.00
76) Chrysene-d12	21.70	240	188848	20.00	ng	0.00
87) Perylene-d12	24.96	264	205923	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	219772	168.12	ng	0.00
7) Phenol-d6	7.19	99	312714	166.22	ng	0.00
23) Nitrobenzene-d5	9.20	82	156660	78.30	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	154853	124.45	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	424013	78.18	ng	0.00
79) Terphenyl-d14	20.02	244	801623	78.52	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.45	88	20411	39.872 ng	97
3) Pyridine	3.86	79	56481	40.605 ng	90
4) n-Nitrosodimethylamine	3.78	42	27069	43.246 ng	88
6) Aniline	7.35	93	67856	30.033 ng	99
8) 2-Chlorophenol	7.59	128	71810	46.601 ng	96
9) Benzaldehyde	7.17	77	22241	20.500 ng	97
10) Phenol	7.21	94	96198	51.003 ng	96
11) bis(2-Chloroethyl)ether	7.44	93	58839	40.816 ng	98
12) 1,3-Dichlorobenzene	7.91	146	70060	37.947 ng	97
13) 1,4-Dichlorobenzene	8.06	146	76559	40.944 ng	98
14) 1,2-Dichlorobenzene	8.38	146	71181	39.926 ng	99
15) Benzyl Alcohol	8.27	79	64293	45.381 ng	95
16) 2,2'-oxybis(1-Chloropropan	8.55	45	108839	39.374 ng	99
17) 2-Methylphenol	8.47	107	63057	48.147 ng	99
18) Hexachloroethane	9.10	117	25569	40.251 ng	92
19) n-Nitroso-di-n-propylamine	8.83	70	50470	40.854 ng	92
20) 3+4-Methylphenols	8.80	107	87676	46.954 ng	98
22) Acetophenone	8.86	105	104277	36.471 ng	# 94
24) Nitrobenzene	9.25	77	76740	37.945 ng	99
25) Isophorone	9.76	82	145374	42.180 ng	98
26) 2-Nitrophenol	9.96	139	42759	39.089 ng	99
27) 2,4-Dimethylphenol	10.00	122	76843	49.931 ng	96
28) bis(2-Chloroethoxy)methane	10.24	93	90727	43.800 ng	99
29) 2,4-Dichlorophenol	10.49	162	79811	41.744 ng	94
30) 1,2,4-Trichlorobenzene	10.70	180	83463	36.332 ng	95
31) Naphthalene	10.89	128	222772	40.569 ng	99
32) Benzoic acid	10.16	122	29456m	34.443 ng	
33) 4-Chloroaniline	11.01	127	53862	21.921 ng	99
34) Hexachlorobutadiene	11.15	225	51778	32.757 ng	98
35) Caprolactam	11.80	113	27160	35.426 ng	96
36) 4-Chloro-3-methylphenol	12.12	107	90125	44.073 ng	89
37) 2-Methylnaphthalene	12.50	142	164634	39.989 ng	97
38) 1-Methylnaphthalene	12.71	142	158973	40.230 ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.86	216	101650	36.463 ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	113668	72.942	ng	98
43) 2,4,6-Trichlorophenol	13.10	196	72270	41.190	ng	97
44) 2,4,5-Trichlorophenol	13.18	196	79032	42.619	ng	98
46) 1,1'-Biphenyl	13.50	154	226464	38.503	ng	96
47) 2-Chloronaphthalene	13.55	162	180990	38.026	ng	99
48) 2-Nitroaniline	13.76	65	53498	40.120	ng	96
49) Acenaphthylene	14.39	152	300226	41.966	ng	98
50) Dimethylphthalate	14.12	163	239888	38.243	ng	99
51) 2,6-Dinitrotoluene	14.25	165	55960	39.903	ng	94
52) Acenaphthene	14.73	154	164496	38.270	ng	94
53) 3-Nitroaniline	14.59	138	34596	24.318	ng	97
54) 2,4-Dinitrophenol	14.80	184	37514	47.116	ng	97
55) Dibenzofuran	15.06	168	301772	39.752	ng	97
56) 4-Nitrophenol	14.90	139	83679	81.747	ng	88
57) 2,4-Dinitrotoluene	15.04	165	79244	39.386	ng	97
58) Fluorene	15.71	166	236835	41.447	ng	95
59) 2,3,4,6-Tetrachlorophenol	15.29	232	71909	41.080	ng	99
60) Diethylphthalate	15.47	149	247893	38.357	ng	98
61) 4-Chlorophenyl-phenylether	15.70	204	133187	36.997	ng	98
62) 4-Nitroaniline	15.76	138	57466	38.775	ng	90
63) Azobenzene	15.99	77	192731	38.134	ng	98
65) 4,6-Dinitro-2-methylphenol	15.80	198	39995	27.841	ng	97
66) n-Nitrosodiphenylamine	15.92	169	214009	37.230	ng	96
67) 4-Bromophenyl-phenylether	16.60	248	87842	35.242	ng	96
68) Hexachlorobenzene	16.71	284	93562	34.400	ng	98
69) Atrazine	16.86	200	88375	36.190	ng	99
70) Pentachlorophenol	17.07	266	86199	53.376	ng	96
71) Phenanthrene	17.46	178	401967	39.220	ng	100
72) Anthracene	17.55	178	398838	39.358	ng	99
73) Carbazole	17.82	167	359608	35.947	ng	99
74) Di-n-butylphthalate	18.35	149	404939	36.386	ng	99
75) Fluoranthene	19.47	202	485091	38.179	ng	98
77) Benzidine	19.65	184	190010	32.896	ng	99
78) Pyrene	19.83	202	496458	41.251	ng	99
80) Butylbenzylphthalate	20.70	149	186695	40.786	ng	99
81) Benzo(a)anthracene	21.67	228	478012	41.580	ng	99
82) 3,3'-Dichlorobenzidine	21.59	252	112827	24.088	ng	97
83) Chrysene	21.74	228	445213	40.719	ng	99
84) Bis(2-ethylhexyl)phthalate	21.54	149	261384	41.126	ng	96
85) Di-n-octyl phthalate	22.75	149	459269	42.735	ng	99
86) Indeno(1,2,3-cd)pyrene	28.72	276	532509	40.759	ng	99
88) Benzo(b)fluoranthene	23.92	252	529861	46.665	ng	# 99
89) Benzo(k)fluoranthene	23.99	252	517354	46.083	ng	# 98
90) Benzo(a)pyrene	24.81	252	467374	42.585	ng	97
91) Dibenzo(a,h)anthracene	28.78	278	442119	40.500	ng	96
92) Benzo(g,h,i)perylene	29.90	276	447535	41.411	ng	# 98

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

