

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG100418\
 Data File : BG037133.D
 Acq On : 3 Oct 2018 14:48
 Operator : JU/SJ
 Sample : PB113454BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB113454BS

Manual Integrations
 APPROVED

Sohil
 10/4/2018 12:59:16 PM

Quant Time: Oct 03 18:03:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 24 10:41:23 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	56204	20.00	ng	0.00
21) Naphthalene-d8	10.84	136	250354	20.00	ng	-0.01
38) Acenaphthene-d10	14.66	164	153804	20.00	ng	-0.01
63) Phenanthrene-d10	17.41	188	363325	20.00	ng	0.00
75) Chrysene-d12	21.68	240	360774	20.00	ng	-0.01
86) Perylene-d12	24.88	264	338588	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	393538	134.28	ng	0.00
7) Phenol-d6	7.21	99	532285	117.39	ng	0.00
23) Nitrobenzene-d5	9.20	82	329271	70.43	ng	-0.01
41) 2,4,6-Tribromophenol	16.16	330	187877	102.10	ng	-0.01
44) 2-Fluorobiphenyl	13.29	172	731911	70.88	ng	-0.01
78) Terphenyl-d14	20.02	244	1076952	78.49	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.52	88	55416	41.324	ng	99
3) Pyridine	3.92	79	144083	37.210	ng	99
4) n-Nitrosodimethylamine	3.84	42	76665	44.547	ng	# 94
6) Aniline	7.37	93	197322	33.538	ng	97
8) 2-Chlorophenol	7.61	128	156309	45.934	ng	97
9) Benzaldehyde	7.18	77	72702	24.265	ng	98
10) Phenol	7.23	94	212844	45.483	ng	96
11) bis(2-Chloroethyl)ether	7.46	93	159827	36.923	ng	99
12) 1,3-Dichlorobenzene	7.92	146	164984	39.547	ng	97
13) 1,4-Dichlorobenzene	8.08	146	170222	39.871	ng	98
14) 1,2-Dichlorobenzene	8.39	146	162045	39.167	ng	97
15) Benzyl Alcohol	8.28	79	143074	38.888	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.56	45	327440	39.401	ng	98
17) 2-Methylphenol	8.48	107	139326	42.743	ng	99
18) Hexachloroethane	9.12	117	66647	42.420	ng	97
19) n-Nitroso-di-n-propylamine	8.84	70	137274	36.393	ng	94
20) 3+4-Methylphenols	8.81	107	190757	42.066	ng	99
22) Acetophenone	8.86	105	247559	42.079	ng	# 96
24) Nitrobenzene	9.25	77	202222	40.505	ng	97
25) Isophorone	9.76	82	382904	44.824	ng	100
26) 2-Nitrophenol	9.96	139	86995	43.716	ng	96
27) 2,4-Dimethylphenol	10.01	122	147005	47.424	ng	93
28) bis(2-Chloroethoxy)methane	10.24	93	219491	41.641	ng	96
29) 2,4-Dichlorophenol	10.49	162	158039	43.980	ng	98
30) 1,2,4-Trichlorobenzene	10.70	180	174225	38.743	ng	98
31) Naphthalene	10.89	128	497284	41.746	ng	99
32) Benzoic acid	10.16	122	49546	24.151	ng	93
33) 4-Chloroaniline	11.00	127	135177	24.959	ng	97
34) Hexachlorobutadiene	11.17	225	117592	37.813	ng	97
35) Caprolactam	11.78	113	54519m	36.862	ng	
36) 4-Chloro-3-methylphenol	12.12	107	175021	40.820	ng	99
37) 2-Methylnaphthalene	12.49	142	378445	41.714	ng	97
39) 1,2,4,5-Tetrachlorobenzene	12.86	216	207018	38.591	ng	98
40) Hexachlorocyclopentadiene	12.84	237	230883	83.686	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.10	196	124744	41.876	nq	99
43) 2,4,5-Trichlorophenol	13.17	196	134973	42.492	nq	96
45) 1,1'-Biphenyl	13.49	154	479813	40.734	nq	97
46) 2-Chloronaphthalene	13.54	162	362978	39.184	nq	99
47) 2-Nitroaniline	13.75	65	130656	40.779	nq	98
48) Acenaphthylene	14.39	152	612218	42.176	nq	99
49) Dimethylphthalate	14.12	163	470326	37.990	nq	99
50) 2,6-Dinitrotoluene	14.24	165	106705	39.504	nq	96
51) Acenaphthene	14.73	154	346339	39.478	nq	98
52) 3-Nitroaniline	14.57	138	84109	29.550	nq	97
53) 2,4-Dinitrophenol	14.78	184	70831	58.253	nq	93
54) Dibenzofuran	15.06	168	571630	38.521	nq	99
55) 4-Nitrophenol	14.88	139	146712	80.684	nq	98
56) 2,4-Dinitrotoluene	15.03	165	150301	39.877	nq	89
57) Fluorene	15.71	166	455126	41.034	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.29	232	130952	40.640	nq	99
59) Diethylphthalate	15.48	149	471585	37.767	nq	98
60) 4-Chlorophenyl-phenylether	15.70	204	255160	36.327	nq	99
61) 4-Nitroaniline	15.74	138	112261	37.496	nq	95
62) Azobenzene	16.00	77	486835	40.577	nq	98
64) 4,6-Dinitro-2-methylphenol	15.79	198	73506	38.697	nq	96
65) n-Nitrosodiphenylamine	15.91	169	407961	40.673	nq	97
66) 4-Bromophenyl-phenylether	16.60	248	157986	38.229	nq	97
67) Hexachlorobenzene	16.71	284	160920	37.807	nq	98
68) Atrazine	16.87	200	159194	39.197	nq	98
69) Pentachlorophenol	17.06	266	168852	63.009	nq	97
70) Phenanthrene	17.45	178	741048	42.325	nq	99
71) Anthracene	17.54	178	745782	43.078	nq	99
72) Carbazole	17.81	167	663167	39.288	nq	99
73) Di-n-butylphthalate	18.36	149	785612	41.909	nq	100
74) Fluoranthene	19.46	202	889581	42.210	nq	99
76) Benzidine	19.64	184	317429	29.106	nq	97
77) Pyrene	19.82	202	872193	40.287	nq	100
79) Butylbenzylphthalate	20.70	149	360022	41.721	nq	96
80) Benzo(a)anthracene	21.65	228	848351	40.282	nq	100
81) 3,3'-Dichlorobenzidine	21.57	252	243689	30.399	nq	98
82) Chrysene	21.73	228	812820	39.945	nq	99
83) Bis(2-ethylhexyl)phthalate	21.55	149	505552	41.937	nq	99
84) Di-n-octyl phthalate	22.76	149	849659	42.808	nq	98
85) Indeno(1,2,3-cd)pyrene	28.53	276	928125	42.476	nq	# 97
87) Benzo(b)fluoranthene	23.86	252	842925	46.714	nq	97
88) Benzo(k)fluoranthene	23.93	252	810048	44.746	nq	99
89) Benzo(a)pyrene	24.73	252	802681	46.013	nq	99
90) Dibenzo(a,h)anthracene	28.59	278	746377	45.652	nq	99
91) Benzo(g,h,i)perylene	29.66	276	760311	46.337	nq	99

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

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