

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG090619\
 Data File : BG042751.D
 Acq On : 6 Sep 2019 16:55
 Operator : HP/JU
 Sample : PB122813BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB122813BS

Manual Integrations
 APPROVED

mohammad
 9/10/2019 9:59:45 AM

Quant Time: Sep 06 17:59:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	27011	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	124184	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	103985	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	263930	20.00	ng	0.00
76) Chrysene-d12	21.69	240	216000	20.00	ng	0.00
87) Perylene-d12	24.93	264	205645	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	181371	135.99	ng	0.00
7) Phenol-d6	7.17	99	235461	130.70	ng	0.00
23) Nitrobenzene-d5	9.16	82	121168	67.43	ng	-0.01
42) 2,4,6-Tribromophenol	16.16	330	193198	130.72	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	438244	71.28	ng	0.00
79) Terphenyl-d14	20.02	244	873655	87.44	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.41	88	17667	31.743	ng	97
3) Pyridine	3.81	79	39418	27.620	ng	91
4) n-Nitrosodimethylamine	3.73	42	22632	44.399	ng	# 88
6) Aniline	7.31	93	79455	33.067	ng	99
8) 2-Chlorophenol	7.56	128	84064	52.009	ng	98
9) Benzaldehyde	7.12	77	29940	33.196	ng	97
10) Phenol	7.19	94	92602	50.556	ng	93
11) bis(2-Chloroethyl)ether	7.41	93	61069	42.453	ng	99
12) 1,3-Dichlorobenzene	7.88	146	83689	40.701	ng	98
13) 1,4-Dichlorobenzene	8.03	146	88118	42.308	ng	100
14) 1,2-Dichlorobenzene	8.35	146	84978	42.605	ng	98
15) Benzyl Alcohol	8.23	79	63744	49.182	ng	92
16) 2,2'-oxybis(1-Chloropropan	8.52	45	75876	38.916	ng	97
17) 2-Methylphenol	8.45	107	68917	51.083	ng	95
18) Hexachloroethane	9.08	117	28374	39.794	ng	90
19) n-Nitroso-di-n-propylamine	8.80	70	50697	43.438	ng	92
20) 3+4-Methylphenols	8.78	107	93741	50.213	ng	96
22) Acetophenone	8.82	105	114748	43.202	ng	# 98
24) Nitrobenzene	9.20	77	74792	41.100	ng	98
25) Isophorone	9.73	82	143748	40.532	ng	98
26) 2-Nitrophenol	9.93	139	51464	45.128	ng	97
27) 2,4-Dimethylphenol	9.99	122	84035	53.495	ng	98
28) bis(2-Chloroethoxy)methane	10.21	93	87838	39.296	ng	97
29) 2,4-Dichlorophenol	10.47	162	102945	50.088	ng	100
30) 1,2,4-Trichlorobenzene	10.68	180	108334	42.943	ng	97
31) Naphthalene	10.87	128	247661	42.955	ng	99
32) Benzoic acid	10.17	122	33121	32.277	ng	94
33) 4-Chloroaniline	10.98	127	70889	26.635	ng	97
34) Hexachlorobutadiene	11.16	225	72160	42.561	ng	96
35) Caprolactam	11.76	113	31461m	45.875	ng	
36) 4-Chloro-3-methylphenol	12.11	107	96158	49.285	ng	97
37) 2-Methylnaphthalene	12.48	142	213522	46.122	ng	100
38) 1-Methylnaphthalene	12.70	142	197637	44.944	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.85	216	148121	44.356	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	93534	53.323	nq	98
43) 2,4,6-Trichlorophenol	13.09	196	99208	43.952	nq	99
44) 2,4,5-Trichlorophenol	13.17	196	106362	45.472	nq	96
46) 1,1'-Biphenyl	13.49	154	290264	43.183	nq	97
47) 2-Chloronaphthalene	13.53	162	238393	41.132	nq	99
48) 2-Nitroaniline	13.73	65	53789	38.486	nq	98
49) Acenaphthylene	14.38	152	379648	43.539	nq	99
50) Dimethylphthalate	14.11	163	327507	43.651	nq	99
51) 2,6-Dinitrotoluene	14.23	165	78318	45.033	nq	98
52) Acenaphthene	14.72	154	222368	41.998	nq	99
53) 3-Nitroaniline	14.56	138	52761	30.335	nq	99
54) 2,4-Dinitrophenol	14.78	184	94079	80.230	nq	99
55) Dibenzofuran	15.06	168	392760	44.814	nq	99
56) 4-Nitrophenol	14.89	139	105352	85.244	nq	94
57) 2,4-Dinitrotoluene	15.03	165	114329	45.908	nq	98
58) Fluorene	15.71	166	325385	45.417	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.29	232	107785	46.597	nq	95
60) Diethylphthalate	15.47	149	319972	43.626	nq	99
61) 4-Chlorophenyl-phenylether	15.70	204	195541	45.277	nq	98
62) 4-Nitroaniline	15.74	138	78612	42.231	nq	94
63) Azobenzene	15.99	77	204630	40.716	nq	96
65) 4,6-Dinitro-2-methylphenol	15.80	198	78321	47.332	nq	98
66) n-Nitrosodiphenylamine	15.91	169	303421	41.803	nq	100
67) 4-Bromophenyl-phenylether	16.60	248	140417	41.943	nq	96
68) Hexachlorobenzene	16.72	284	150247	41.285	nq	99
69) Atrazine	16.86	200	117165	45.647	nq	98
70) Pentachlorophenol	17.07	266	160980	85.133	nq	99
71) Phenanthrene	17.45	178	570890	43.547	nq	99
72) Anthracene	17.55	178	574904	43.928	nq	99
73) Carbazole	17.81	167	505032	43.298	nq	99
74) Di-n-butylphthalate	18.37	149	584474	42.390	nq	99
75) Fluoranthene	19.47	202	715133	44.697	nq	97
77) Benzidine	19.65	184	131702	21.267	nq	98
78) Pyrene	19.83	202	709667	53.589	nq	98
80) Butylbenzylphthalate	20.71	149	201496	38.684	nq	94
81) Benzo(a)anthracene	21.67	228	575492	43.798	nq	97
82) 3,3'-Dichlorobenzidine	21.58	252	191293	36.199	nq	98
83) Chrysene	21.74	228	555781	43.859	nq	98
84) Bis(2-ethylhexyl)phthalate	21.57	149	280431	38.777	nq	99
85) Di-n-octyl phthalate	22.78	149	499656	40.170	nq	99
86) Indeno(1,2,3-cd)pyrene	28.65	276	719791	41.797	nq	# 97
88) Benzo(b)fluoranthene	23.90	252	613164	54.235	nq	# 97
89) Benzo(k)fluoranthene	23.98	252	615286	56.619	nq	# 98
90) Benzo(a)pyrene	24.79	252	608016	56.675	nq	# 98
91) Dibenzo(a,h)anthracene	28.71	278	606391	55.826	nq	98
92) Benzo(g,h,i)perylene	29.81	276	604306	55.625	nq	# 95

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

