

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121719\
 Data File : BG043816.D
 Acq On : 17 Dec 2019 12:24
 Operator : JU
 Sample : PB125466BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB125466BS

Manual Integrations
 APPROVED

mohammad
 12/18/2019 2:40:35 PM

Quant Time: Dec 17 13:07:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG120919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 09 18:16:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	104973	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	476933	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	364882	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	886421	20.00	ng	0.00
76) Chrysene-d12	21.63	240	744900	20.00	ng	0.00
87) Perylene-d12	24.79	264	777061	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	748291	135.06	ng	0.00
7) Phenol-d6	7.17	99	1037409	129.84	ng	0.01
23) Nitrobenzene-d5	9.16	82	650548	76.59	ng	0.00
42) 2,4,6-Tribromophenol	16.12	330	458905	131.92	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	1545776	74.41	ng	0.00
79) Terphenyl-d14	19.99	244	2434938	78.64	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.48	88	100831	39.301	ng	96
3) Pyridine	3.88	79	254853	35.947	ng	99
4) n-Nitrosodimethylamine	3.79	42	133633	39.313	ng	94
6) Aniline	7.32	93	365105	34.534	ng	96
8) 2-Chlorophenol	7.57	128	320934	49.708	ng	95
9) Benzaldehyde	7.14	77	220746	56.462	ng	95
10) Phenol	7.20	94	415959	50.425	ng	99
11) bis(2-Chloroethyl)ether	7.42	93	292673	41.812	ng	98
12) 1,3-Dichlorobenzene	7.89	146	320918	41.477	ng	99
13) 1,4-Dichlorobenzene	8.04	146	320334	41.419	ng	99
14) 1,2-Dichlorobenzene	8.36	146	307261	41.235	ng	99
15) Benzyl Alcohol	8.24	79	295711	46.506	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.53	45	443408	39.525	ng	99
17) 2-Methylphenol	8.45	107	295838	50.240	ng	99
18) Hexachloroethane	9.09	117	115764	41.192	ng	97
19) n-Nitroso-di-n-propylamine	8.81	70	249665	40.872	ng	99
20) 3+4-Methylphenols	8.78	107	413759	50.462	ng	97
22) Acetophenone	8.82	105	450940	37.912	ng	97
24) Nitrobenzene	9.20	77	353941	40.908	ng	98
25) Isophorone	9.73	82	684688	40.360	ng	100
26) 2-Nitrophenol	9.91	139	174418	50.939	ng	98
27) 2,4-Dimethylphenol	9.98	122	332304	55.432	ng	98
28) bis(2-Chloroethoxy)methane	10.21	93	427935	39.408	ng	98
29) 2,4-Dichlorophenol	10.46	162	357468	48.558	ng	99
30) 1,2,4-Trichlorobenzene	10.67	180	334078	38.336	ng	99
31) Naphthalene	10.86	128	959054	41.908	ng	98
32) Benzoic acid	10.11	122	193299	46.462	ng	98
33) 4-Chloroaniline	10.96	127	272723	24.579	ng	99
34) Hexachlorobutadiene	11.15	225	191910	36.345	ng	97
35) Caprolactam	11.72	113	150988m	52.261	ng	
36) 4-Chloro-3-methylphenol	12.08	107	398285	50.254	ng	99
37) 2-Methylnaphthalene	12.46	142	763062	43.020	ng	99
38) 1-Methylnaphthalene	12.68	142	723404	42.889	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	389129	35.490	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	464047	90.170	nq	97
43) 2,4,6-Trichlorophenol	13.06	196	314605	45.144	nq	99
44) 2,4,5-Trichlorophenol	13.14	196	349483	47.867	nq	100
46) 1,1'-Biphenyl	13.47	154	986207	39.205	nq	100
47) 2-Chloronaphthalene	13.50	162	789207	38.377	nq	99
48) 2-Nitroaniline	13.70	65	275503	45.425	nq	97
49) Acenaphthylene	14.35	152	1304932	42.597	nq	99
50) Dimethylphthalate	14.09	163	1090412	41.869	nq	100
51) 2,6-Dinitrotoluene	14.20	165	261319	46.561	nq	99
52) Acenaphthene	14.70	154	929867	46.140	nq	96
53) 3-Nitroaniline	14.53	138	202383	31.511	nq	93
54) 2,4-Dinitrophenol	14.73	184	271695	104.851	nq	92
55) Dibenzofuran	15.03	168	1281318	42.952	nq	99
56) 4-Nitrophenol	14.84	139	502843	107.317	nq	99
57) 2,4-Dinitrotoluene	14.98	165	372310	49.094	nq	97
58) Fluorene	15.68	166	1053871	43.829	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.26	232	324650m	49.558	nq	
60) Diethylphthalate	15.46	149	1125632	43.270	nq	99
61) 4-Chlorophenyl-phenylether	15.67	204	566232	42.048	nq	98
62) 4-Nitroaniline	15.69	138	280273	42.117	nq	98
63) Azobenzene	15.97	77	1043300	43.656	nq	99
65) 4,6-Dinitro-2-methylphenol	15.75	198	192360	54.615	nq	96
66) n-Nitrosodiphenylamine	15.88	169	1010899	41.032	nq	98
67) 4-Bromophenyl-phenylether	16.57	248	365292	38.545	nq	98
68) Hexachlorobenzene	16.69	284	377166	37.524	nq	99
69) Atrazine	16.84	200	404737	51.532	nq	98
70) Pentachlorophenol	17.02	266	474787	88.800	nq	98
71) Phenanthrene	17.42	178	1710698	40.766	nq	99
72) Anthracene	17.51	178	1768653	42.578	nq	99
73) Carbazole	17.77	167	1614358	41.575	nq	100
74) Di-n-butylphthalate	18.35	149	1858623	41.754	nq	99
75) Fluoranthene	19.42	202	1969269	41.003	nq	99
77) Benzidine	19.60	184	846806	52.108	nq	98
78) Pyrene	19.78	202	1974782	40.566	nq	98
80) Butylbenzylphthalate	20.68	149	879937	43.401	nq	97
81) Benzo(a)anthracene	21.61	228	1915084	40.734	nq	99
82) 3,3'-Dichlorobenzidine	21.52	252	562200	32.371	nq	100
83) Chrysene	21.68	228	1816204	40.548	nq	100
84) Bis(2-ethylhexyl)phthalate	21.54	149	1217814	41.877	nq	99
85) Di-n-octyl phthalate	22.75	149	2108334	44.347	nq	100
86) Indeno(1,2,3-cd)pyrene	28.38	276	1901541	33.169	nq	# 96
88) Benzo(b)fluoranthene	23.79	252	1923473	42.992	nq	99
89) Benzo(k)fluoranthene	23.86	252	1901990	44.008	nq	100
90) Benzo(a)pyrene	24.64	252	1811605	42.489	nq	100
91) Dibenzo(a,h)anthracene	28.45	278	1700434	39.928	nq	100
92) Benzo(g,h,i)perylene	29.49	276	1509473	35.918	nq	99

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

