

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061119\
 Data File : BG041183.D
 Acq On : 11 Jun 2019 11:00
 Operator : HP/JU
 Sample : PB120413BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB120413BS

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:38:19 AM

Quant Time: Jun 12 02:14:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 23:08:34 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	33962	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	140612	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	110588	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	268858	20.00	ng	0.00
76) Chrysene-d12	21.76	240	265565	20.00	ng	0.00
87) Perylene-d12	25.09	264	279259	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	259269	137.66	ng	0.00
7) Phenol-d6	7.26	99	381027	130.99	ng	0.00
23) Nitrobenzene-d5	9.28	82	262391	82.84	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	218843	133.30	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	649123	84.53	ng	0.00
79) Terphenyl-d14	20.08	244	1157028	92.47	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.50	88	34667	40.563	ng	92
3) Pyridine	3.91	79	92877	36.726	ng	98
4) n-Nitrosodimethylamine	3.83	42	53513	45.594	ng	81
6) Aniline	7.43	93	93782	24.155	ng	98
8) 2-Chlorophenol	7.66	128	106120	47.261	ng	97
9) Benzaldehyde	7.24	77	29701	17.117	ng	90
10) Phenol	7.29	94	143747	46.091	ng	94
11) bis(2-Chloroethyl)ether	7.52	93	94839	41.918	ng	97
12) 1,3-Dichlorobenzene	7.99	146	114571	42.722	ng	99
13) 1,4-Dichlorobenzene	8.14	146	114288	42.102	ng	96
14) 1,2-Dichlorobenzene	8.46	146	114336	43.379	ng	96
15) Benzyl Alcohol	8.35	79	119471	44.401	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.63	45	141035	38.148	ng	95
17) 2-Methylphenol	8.54	107	100386	44.456	ng	98
18) Hexachloroethane	9.19	117	44481	42.515	ng	92
19) n-Nitroso-di-n-propylamine	8.91	70	87620	39.143	ng	97
20) 3+4-Methylphenols	8.87	107	138410	44.250	ng	97
22) Acetophenone	8.94	105	174520	44.245	ng	# 97
24) Nitrobenzene	9.33	77	147603	45.728	ng	98
25) Isophorone	9.84	82	255239	42.344	ng	99
26) 2-Nitrophenol	10.03	139	64947	48.807	ng	87
27) 2,4-Dimethylphenol	10.08	122	112581	51.227	ng	94
28) bis(2-Chloroethoxy)methane	10.32	93	138040	42.430	ng	98
29) 2,4-Dichlorophenol	10.57	162	133871	47.969	ng	98
30) 1,2,4-Trichlorobenzene	10.78	180	142466	44.874	ng	98
31) Naphthalene	10.98	128	347453	46.023	ng	99
32) Benzoic acid	10.22	122	68794	41.569	ng	96
33) 4-Chloroaniline	11.09	127	71236	20.336	ng	95
34) Hexachlorobutadiene	11.24	225	101421	41.751	ng	97
35) Caprolactam	11.89	113	40036m	43.442	ng	
36) 4-Chloro-3-methylphenol	12.20	107	148211	46.501	ng	96
37) 2-Methylnaphthalene	12.57	142	271613	46.713	ng	97
38) 1-Methylnaphthalene	12.80	142	258105	47.106	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	203802	45.723	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.90	237	220875	90.930	ng	98
43) 2,4,6-Trichlorophenol	13.18	196	132667	46.009	ng	98
44) 2,4,5-Trichlorophenol	13.25	196	140872	46.324	ng	96
46) 1,1'-Biphenyl	13.57	154	374325	45.728	ng	99
47) 2-Chloronaphthalene	13.62	162	308158	43.850	ng	98
48) 2-Nitroaniline	13.83	65	111074	44.058	ng	97
49) Acenaphthylene	14.46	152	482713	45.639	ng	98
50) Dimethylphthalate	14.19	163	396618	42.368	ng	99
51) 2,6-Dinitrotoluene	14.32	165	91931	45.547	ng	93
52) Acenaphthene	14.80	154	282081	44.024	ng	98
53) 3-Nitroaniline	14.65	138	63774	31.598	ng	99
54) 2,4-Dinitrophenol	14.86	184	87395	82.133	ng	95
55) Dibenzofuran	15.13	168	493286	45.754	ng	99
56) 4-Nitrophenol	14.95	139	125787	90.862	ng	92
57) 2,4-Dinitrotoluene	15.11	165	131420	46.094	ng	95
58) Fluorene	15.78	166	388137	45.205	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.36	232	142878	47.808	ng	99
60) Diethylphthalate	15.54	149	408621	42.772	ng	100
61) 4-Chlorophenyl-phenylether	15.77	204	244671	43.841	ng	94
62) 4-Nitroaniline	15.82	138	88966	41.636	ng	97
63) Azobenzene	16.06	77	379177	43.669	ng	96
65) 4,6-Dinitro-2-methylphenol	15.86	198	72979	50.180	ng	98
66) n-Nitrosodiphenylamine	15.99	169	349946	46.302	ng	99
67) 4-Bromophenyl-phenylether	16.66	248	161257	44.444	ng	98
68) Hexachlorobenzene	16.77	284	167437	43.337	ng	98
69) Atrazine	16.93	200	146051	50.082	ng	95
70) Pentachlorophenol	17.13	266	188823	90.633	ng	97
71) Phenanthrene	17.53	178	655636	46.816	ng	98
72) Anthracene	17.61	178	679365	49.186	ng	99
73) Carbazole	17.89	167	570567	48.144	ng	99
74) Di-n-butylphthalate	18.42	149	682321	46.322	ng	99
75) Fluoranthene	19.53	202	841006	48.619	ng	99
77) Benzidine	19.71	184	260928	41.187	ng	97
78) Pyrene	19.89	202	849902	49.231	ng	98
80) Butylbenzylphthalate	20.76	149	313226	46.623	ng	97
81) Benzo(a)anthracene	21.74	228	806042	45.597	ng	99
82) 3,3'-Dichlorobenzidine	21.66	252	187397	30.140	ng	98
83) Chrysene	21.81	228	798537	47.204	ng	98
84) Bis(2-ethylhexyl)phthalate	21.61	149	435414	46.040	ng	98
85) Di-n-octyl phthalate	22.85	149	738711	46.900	ng	99
86) Indeno(1,2,3-cd)pyrene	28.92	276	744806	35.942	ng	99
88) Benzo(b)fluoranthene	24.02	252	786812	46.452	ng	100
89) Benzo(k)fluoranthene	24.09	252	791012	47.193	ng	98
90) Benzo(a)pyrene	24.93	252	773490	47.864	ng	97
91) Dibenzo(a,h)anthracene	28.98	278	633219	39.215	ng	97
92) Benzo(g,h,i)perylene	30.13	276	585279	37.308	ng	96

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

