

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG042420\
 Data File : BG045159.D
 Acq On : 24 Apr 2020 13:42
 Operator : CG/JU
 Sample : PB128463BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB128463BS

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:05:38 PM

Quant Time: Apr 24 14:21:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 13:00:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	64117	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	256072	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	185516	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	449779	20.00	ng	0.00
76) Chrysene-d12	21.65	240	435446	20.00	ng	0.00
87) Perylene-d12	24.82	264	485758	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	498784	128.43	ng	0.00
7) Phenol-d6	7.17	99	696340	120.68	ng	0.00
23) Nitrobenzene-d5	9.16	82	437582	77.97	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	336247	117.32	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	1009859	79.82	ng	0.00
79) Terphenyl-d14	20.00	244	1827641	91.48	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.46	88	64759	37.521	ng	97
3) Pyridine	3.86	79	184168	34.656	ng	93
4) n-Nitrosodimethylamine	3.77	42	71445	43.887	ng	# 90
6) Aniline	7.32	93	238084	31.735	ng	98
8) 2-Chlorophenol	7.57	128	193602	46.384	ng	98
9) Benzaldehyde	7.13	77	93325	25.236	ng	95
10) Phenol	7.19	94	265416	45.967	ng	98
11) bis(2-Chloroethyl)ether	7.42	93	180437	39.249	ng	99
12) 1,3-Dichlorobenzene	7.89	146	210712	42.842	ng	98
13) 1,4-Dichlorobenzene	8.03	146	211131	43.081	ng	99
14) 1,2-Dichlorobenzene	8.36	146	202758	43.245	ng	96
15) Benzyl Alcohol	8.24	79	196371	40.591	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.53	45	166049	36.796	ng	96
17) 2-Methylphenol	8.45	107	174638	39.201	ng	94
18) Hexachloroethane	9.09	117	81690	44.339	ng	95
19) n-Nitroso-di-n-propylamine	8.81	70	159478	39.081	ng	98
20) 3+4-Methylphenols	8.77	107	243380	39.030	ng	97
22) Acetophenone	8.82	105	296302	42.788	ng	# 97
24) Nitrobenzene	9.20	77	241442	41.971	ng	95
25) Isophorone	9.73	82	448879	39.058	ng	98
26) 2-Nitrophenol	9.91	139	113431	45.777	ng	100
27) 2,4-Dimethylphenol	9.98	122	178637	45.435	ng	94
28) bis(2-Chloroethoxy)methane	10.21	93	254567	42.158	ng	97
29) 2,4-Dichlorophenol	10.45	162	202553	44.616	ng	97
30) 1,2,4-Trichlorobenzene	10.67	180	217473	42.309	ng	98
31) Naphthalene	10.86	128	599837	42.820	ng	97
32) Benzoic acid	10.12	122	97318	32.945	ng	93
33) 4-Chloroaniline	10.96	127	133745	20.472	ng	98
34) Hexachlorobutadiene	11.15	225	149298	42.364	ng	97
35) Caprolactam	11.73	113	74056m	40.986	ng	
36) 4-Chloro-3-methylphenol	12.09	107	230318	43.186	ng	95
37) 2-Methylnaphthalene	12.46	142	454140	41.767	ng	95
38) 1-Methylnaphthalene	12.68	142	433711	42.253	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	246560	41.278	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG042420\
 Data File : BG045159.D
 Acq On : 24 Apr 2020 13:42
 Operator : CG/JU
 Sample : PB128463BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB128463BS

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:05:38 PM

Quant Time: Apr 24 14:21:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 13:00:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	372642	99.816	nq	100
43) 2,4,6-Trichlorophenol	13.07	196	186162	44.160	nq	98
44) 2,4,5-Trichlorophenol	13.14	196	198755	41.420	nq	98
46) 1,1'-Biphenyl	13.47	154	574615	40.751	nq	100
47) 2-Chloronaphthalene	13.52	162	454189	42.155	nq	99
48) 2-Nitroaniline	13.71	65	153787	43.382	nq	95
49) Acenaphthylene	14.36	152	740139	42.174	nq	99
50) Dimethylphthalate	14.09	163	633296	41.925	nq	98
51) 2,6-Dinitrotoluene	14.21	165	146288	43.297	nq	96
52) Acenaphthene	14.70	154	570027m	48.006	nq	
53) 3-Nitroaniline	14.54	138	86352	23.303	nq	96
54) 2,4-Dinitrophenol	14.74	184	188416	93.783	nq	97
55) Dibenzofuran	15.04	168	726849	41.987	nq	98
56) 4-Nitrophenol	14.85	139	246154	85.672	nq	96
57) 2,4-Dinitrotoluene	15.00	165	208292	42.185	nq	99
58) Fluorene	15.69	166	603949	42.711	nq	98
59) 2,3,4,6-Tetrachlorophenol	15.27	232	195225	45.724	nq	98
60) Diethylphthalate	15.46	149	653919	41.521	nq	100
61) 4-Chlorophenyl-phenylether	15.68	204	331069	40.677	nq	98
62) 4-Nitroaniline	15.70	138	154350	40.159	nq	94
63) Azobenzene	15.97	77	618559	42.600	nq	97
65) 4,6-Dinitro-2-methylphenol	15.76	198	140646	47.573	nq	98
66) n-Nitrosodiphenylamine	15.89	169	551350	42.664	nq	99
67) 4-Bromophenyl-phenylether	16.58	248	229344	39.525	nq	97
68) Hexachlorobenzene	16.70	284	254977	42.370	nq	100
69) Atrazine	16.84	200	236030	50.678	nq	100
70) Pentachlorophenol	17.04	266	305208	82.673	nq	99
71) Phenanthrene	17.43	178	1000731	43.297	nq	100
72) Anthracene	17.52	178	1028450	44.359	nq	99
73) Carbazole	17.79	167	976762	41.515	nq	99
74) Di-n-butylphthalate	18.35	149	1163481	45.495	nq	99
75) Fluoranthene	19.44	202	1289367	47.853	nq	98
77) Benzidine	19.61	184	656748	53.151	nq	98
78) Pyrene	19.80	202	1279182	42.611	nq	99
80) Butylbenzylphthalate	20.69	149	530655	42.645	nq	98
81) Benzo(a)anthracene	21.63	228	1210256	42.882	nq	98
82) 3,3'-Dichlorobenzidine	21.55	252	273103	24.312	nq	99
83) Chrysene	21.70	228	1172133	43.225	nq	99
84) Bis(2-ethylhexyl)phthalate	21.55	149	728791	42.584	nq	98
85) Di-n-octyl phthalate	22.75	149	1237710	41.822	nq	99
86) Indeno(1,2,3-cd)pyrene	28.44	276	1544843	42.909	nq	# 97
88) Benzo(b)fluoranthene	23.82	252	1309536	43.038	nq	97
89) Benzo(k)fluoranthene	23.89	252	1287991	44.511	nq	99
90) Benzo(a)pyrene	24.68	252	1212349	42.200	nq	98
91) Dibenzo(a,h)anthracene	28.50	278	1260985	43.560	nq	97
92) Benzo(g,h,i)perylene	29.56	276	1247563	42.987	nq	98

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

