

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG032420\
 Data File : BG044939.D
 Acq On : 24 Mar 2020 20:46
 Operator : CG/JU
 Sample : PB127773BS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB127773BS

Manual Integrations
 APPROVED

mohammad
 3/25/2020 12:27:59 PM

Quant Time: Mar 25 06:35:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG031020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 23 07:18:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	73556	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	283797	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	187559	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	443944	20.00	ng	0.00
76) Chrysene-d12	21.69	240	448273	20.00	ng	0.00
87) Perylene-d12	24.89	264	509668	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	454100	96.10	ng	0.00
7) Phenol-d6	7.20	99	616717	89.83	ng	0.00
23) Nitrobenzene-d5	9.20	82	564081	87.22	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	256854	104.59	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	1203769	91.91	ng	0.00
79) Terphenyl-d14	20.03	244	2119222	88.43	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.51	88	86459	37.997	ng	98
3) Pyridine	3.91	79	248685	36.919	ng	98
4) n-Nitrosodimethylamine	3.82	42	121455	47.522	ng	99
6) Aniline	7.36	93	267976	31.369	ng	95
8) 2-Chlorophenol	7.61	128	242604	51.433	ng	98
9) Benzaldehyde	7.17	77	177615	43.559	ng	94
10) Phenol	7.23	94	319180	46.960	ng	96
11) bis(2-Chloroethyl)ether	7.45	93	231636	42.702	ng	92
12) 1,3-Dichlorobenzene	7.93	146	265154	45.633	ng	95
13) 1,4-Dichlorobenzene	8.08	146	260176	45.292	ng	97
14) 1,2-Dichlorobenzene	8.40	146	245656	45.090	ng	99
15) Benzyl Alcohol	8.28	79	241748	43.888	ng	93
16) 2,2'-oxybis(1-Chloropropan	8.57	45	349974	36.560	ng	97
17) 2-Methylphenol	8.48	107	214669	46.626	ng	97
18) Hexachloroethane	9.13	117	101006	44.662	ng	99
19) n-Nitroso-di-n-propylamine	8.85	70	194893	40.204	ng	99
20) 3+4-Methylphenols	8.81	107	300449	47.340	ng	94
22) Acetophenone	8.86	105	351680	43.741	ng	99
24) Nitrobenzene	9.24	77	300982	44.852	ng	99
25) Isophorone	9.77	82	532431	42.182	ng	97
26) 2-Nitrophenol	9.95	139	126502	53.576	ng	96
27) 2,4-Dimethylphenol	10.02	122	220226	53.576	ng	90
28) bis(2-Chloroethoxy)methane	10.25	93	308526	40.958	ng	96
29) 2,4-Dichlorophenol	10.49	162	240067	50.002	ng	98
30) 1,2,4-Trichlorobenzene	10.72	180	261964	45.632	ng	98
31) Naphthalene	10.90	128	726256	46.196	ng	99
32) Benzoic acid	10.16	122	123232	40.401	ng	98
33) 4-Chloroaniline	11.00	127	141693	20.005	ng	95
34) Hexachlorobutadiene	11.20	225	173327	45.911	ng	94
35) Caprolactam	11.77	113	83995m	46.207	ng	
36) 4-Chloro-3-methylphenol	12.12	107	271175	47.358	ng	98
37) 2-Methylnaphthalene	12.50	142	541236	46.024	ng	96
38) 1-Methylnaphthalene	12.72	142	519173	46.959	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.87	216	288238	46.664	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.86	237	399902	118.465	nq	98
43) 2,4,6-Trichlorophenol	13.11	196	213040	51.970	nq	98
44) 2,4,5-Trichlorophenol	13.18	196	229456	53.974	nq	96
46) 1,1'-Biphenyl	13.51	154	676055	45.106	nq	98
47) 2-Chloronaphthalene	13.55	162	527688	46.087	nq	97
48) 2-Nitroaniline	13.74	65	186520	46.531	nq	96
49) Acenaphthylene	14.39	152	857190	47.787	nq	99
50) Dimethylphthalate	14.13	163	701978	46.992	nq	99
51) 2,6-Dinitrotoluene	14.24	165	161263	53.557	nq	96
52) Acenaphthene	14.74	154	630883	53.703	nq	98
53) 3-Nitroaniline	14.56	138	100596	29.059	nq	96
54) 2,4-Dinitrophenol	14.78	184	171661	106.071	nq	96
55) Dibenzofuran	15.08	168	824428	48.394	nq	97
56) 4-Nitrophenol	14.88	139	276724	104.740	nq	94
57) 2,4-Dinitrotoluene	15.03	165	221967	53.750	nq	95
58) Fluorene	15.72	166	693910	49.517	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.30	232	213680m	54.594	nq	
60) Diethylphthalate	15.49	149	732104	46.989	nq	98
61) 4-Chlorophenyl-phenylether	15.72	204	374108	49.579	nq	98
62) 4-Nitroaniline	15.73	138	167825	45.464	nq	91
63) Azobenzene	16.01	77	731404	45.205	nq	99
65) 4,6-Dinitro-2-methylphenol	15.79	198	132545	60.959	nq	94
66) n-Nitrosodiphenylamine	15.93	169	631231	47.227	nq	99
67) 4-Bromophenyl-phenylether	16.61	248	254284	45.818	nq	99
68) Hexachlorobenzene	16.73	284	284648	47.275	nq	98
69) Atrazine	16.88	200	258840	52.858	nq	98
70) Pentachlorophenol	17.07	266	353045	106.567	nq	97
71) Phenanthrene	17.46	178	1142678	48.166	nq	99
72) Anthracene	17.55	178	1167334	49.901	nq	98
73) Carbazole	17.82	167	1104644	49.160	nq	100
74) Di-n-butylphthalate	18.39	149	1306754	47.839	nq	99
75) Fluoranthene	19.47	202	1469679	52.030	nq	99
77) Benzidine	19.64	184	674271	46.333	nq	98
78) Pyrene	19.83	202	1455343	44.251	nq	99
80) Butylbenzylphthalate	20.72	149	624021	42.650	nq	97
81) Benzo(a)anthracene	21.67	228	1434873	44.452	nq	100
82) 3,3'-Dichlorobenzidine	21.58	252	341255	27.327	nq	99
83) Chrysene	21.74	228	1394286	45.220	nq	100
84) Bis(2-ethylhexyl)phthalate	21.59	149	864462	42.810	nq	96
85) Di-n-octyl phthalate	22.80	149	1478357	43.750	nq	99
86) Indeno(1,2,3-cd)pyrene	28.54	276	1889272	46.737	nq	# 98
88) Benzo(b)fluoranthene	23.88	252	1584701	47.098	nq	99
89) Benzo(k)fluoranthene	23.95	252	1538276	48.310	nq	99
90) Benzo(a)pyrene	24.75	252	1475851	46.876	nq	97
91) Dibenzo(a,h)anthracene	28.61	278	1520718	48.960	nq	96
92) Benzo(g,h,i)perylene	29.68	276	1543449	49.184	nq	97

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

