

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072719\
 Data File : BG041779.D
 Acq On : 27 Jul 2019 16:24
 Operator : HP/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

mohammad
 7/29/2019 5:30:19 PM

Quant Time: Jul 29 06:44:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 29 05:48:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	92440	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	393863	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	273040	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	603431	20.00	ng	0.00
76) Chrysene-d12	21.74	240	590326	20.00	ng	0.00
87) Perylene-d12	25.02	264	695393	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	589967	104.33	ng	0.00
7) Phenol-d6	7.20	99	782907	102.94	ng	0.00
23) Nitrobenzene-d5	9.22	82	719073	111.03	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	398325	129.45	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	1836000	104.13	ng	0.00
79) Terphenyl-d14	20.07	244	2634868	104.30	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.48	88	135634	50.503	ng	92
3) Pyridine	3.88	79	381287	53.186	ng	91
4) n-Nitrosodimethylamine	3.79	42	153437	46.346	ng	90
6) Aniline	7.37	93	540831	52.908	ng	99
8) 2-Chlorophenol	7.61	128	337633	52.742	ng	95
9) Benzaldehyde	7.17	77	270429	61.499	ng	95
10) Phenol	7.22	94	405405	50.591	ng	98
11) bis(2-Chloroethyl)ether	7.46	93	329330	50.887	ng	96
12) 1,3-Dichlorobenzene	7.94	146	429606	56.242	ng	98
13) 1,4-Dichlorobenzene	8.08	146	431120	56.359	ng	99
14) 1,2-Dichlorobenzene	8.41	146	405402	56.211	ng	96
15) Benzyl Alcohol	8.28	79	314988	52.020	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.58	45	580829	43.312	ng	95
17) 2-Methylphenol	8.49	107	295331	53.293	ng	98
18) Hexachloroethane	9.14	117	150463	53.702	ng	96
19) n-Nitroso-di-n-propylamine	8.86	70	279438	50.266	ng	91
20) 3+4-Methylphenols	8.82	107	407697	53.686	ng	100
22) Acetophenone	8.87	105	526780	54.914	ng	96
24) Nitrobenzene	9.26	77	379382	53.699	ng	96
25) Isophorone	9.79	82	749235	52.832	ng	98
26) 2-Nitrophenol	9.97	139	183449	60.023	ng	93
27) 2,4-Dimethylphenol	10.03	122	305682	54.478	ng	99
28) bis(2-Chloroethoxy)methane	10.27	93	464456	53.138	ng	99
29) 2,4-Dichlorophenol	10.51	162	376948	58.922	ng	95
30) 1,2,4-Trichlorobenzene	10.73	180	460640	61.324	ng	99
31) Naphthalene	10.92	128	1070510	55.208	ng	99
32) Benzoic acid	10.19	122	222573m	56.479	ng	
33) 4-Chloroaniline	11.02	127	509934	57.012	ng	98
34) Hexachlorobutadiene	11.22	225	311343	64.488	ng	98
35) Caprolactam	11.80	113	128914	54.824	ng	# 83
36) 4-Chloro-3-methylphenol	12.15	107	367196	54.789	ng	94
37) 2-Methylnaphthalene	12.53	142	844072	56.976	ng	98
38) 1-Methylnaphthalene	12.75	142	803553	56.895	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.90	216	531733	56.854	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.88	237	322171	60.895	nq	99
43) 2,4,6-Trichlorophenol	13.13	196	353385	58.152	nq	98
44) 2,4,5-Trichlorophenol	13.21	196	363466	56.767	nq	97
46) 1,1'-Biphenyl	13.53	154	1065015	52.064	nq	98
47) 2-Chloronaphthalene	13.58	162	915828	53.720	nq	97
48) 2-Nitroaniline	13.77	65	250752	50.517	nq	94
49) Acenaphthylene	14.43	152	1349662	50.713	nq	99
50) Dimethylphthalate	14.16	163	1134390	52.899	nq	100
51) 2,6-Dinitrotoluene	14.27	165	254418	55.934	nq	89
52) Acenaphthene	14.77	154	888113	52.953	nq	98
53) 3-Nitroaniline	14.60	138	271514	54.682	nq	92
54) 2,4-Dinitrophenol	14.80	184	134131	68.227	nq	87
55) Dibenzofuran	15.10	168	1323791	52.599	nq	99
56) 4-Nitrophenol	14.90	139	209365	52.080	nq	93
57) 2,4-Dinitrotoluene	15.06	165	354202	58.909	nq	95
58) Fluorene	15.76	166	1063966	51.956	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.33	232	358530	60.655	nq	94
60) Diethylphthalate	15.52	149	1114372	51.583	nq	99
61) 4-Chlorophenyl-phenylether	15.74	204	657453	56.796	nq	96
62) 4-Nitroaniline	15.77	138	285450	54.205	nq	97
63) Azobenzene	16.04	77	897541	45.954	nq	97
65) 4,6-Dinitro-2-methylphenol	15.82	198	187039	70.371	nq	88
66) n-Nitrosodiphenylamine	15.96	169	991296	55.940	nq	99
67) 4-Bromophenyl-phenylether	16.64	248	454846	63.349	nq	97
68) Hexachlorobenzene	16.77	284	472727	65.382	nq	95
69) Atrazine	16.91	200	394002	61.268	nq	98
70) Pentachlorophenol	17.10	266	291961	64.582	nq	99
71) Phenanthrene	17.50	178	1709416	56.251	nq	98
72) Anthracene	17.59	178	1714833	56.611	nq	98
73) Carbazole	17.85	167	1529660	54.449	nq	99
74) Di-n-butylphthalate	18.42	149	1779102	51.923	nq	98
75) Fluoranthene	19.51	202	2040425	54.681	nq	96
77) Benzidine	19.68	184	1165749	58.625	nq	98
78) Pyrene	19.87	202	2030204	50.850	nq	99
80) Butylbenzylphthalate	20.76	149	852134	51.362	nq	89
81) Benzo(a)anthracene	21.72	228	2061468	54.048	nq	99
82) 3,3'-Dichlorobenzidine	21.63	252	858880	57.225	nq	99
83) Chrysene	21.79	228	1960818	53.801	nq	99
84) Bis(2-ethylhexyl)phthalate	21.64	149	1142746	48.783	nq	99
85) Di-n-octyl phthalate	22.88	149	1952488	48.944	nq	# 92
86) Indeno(1,2,3-cd)pyrene	28.77	276	2714118	56.440	nq	# 89
88) Benzo(b)fluoranthene	23.98	252	2271045m	53.841	nq	
89) Benzo(k)fluoranthene	24.05	252	2227656	56.995	nq	# 98
90) Benzo(a)pyrene	24.86	252	2201650	55.167	nq	98
91) Dibenzo(a,h)anthracene	28.84	278	2174001	56.183	nq	97
92) Benzo(g,h,i)perylene	29.94	276	2165781	55.852	nq	95

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

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