

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG071519\
 Data File : BG041643.D
 Acq On : 15 Jul 2019 9:48
 Operator : HP/JU
 Sample : SSTDICC040
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICC040

Quant Time: Jul 15 11:47:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 15 11:42:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	154114	20.00	ng	0.00
21) Naphthalene-d8	10.86	136	661466	20.00	ng	0.00
39) Acenaphthene-d10	14.67	164	409798	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	952514	20.00	ng	0.00
76) Chrysene-d12	21.66	240	870427	20.00	ng	0.00
87) Perylene-d12	24.86	264	1027355	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.61	112	691570	75.30	ng	0.00
7) Phenol-d6	7.20	99	955036	76.76	ng	0.00
23) Nitrobenzene-d5	9.21	82	833732	76.82	ng	0.00
42) 2,4,6-Tribromophenol	16.15	330	366868	78.29	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	1903299	73.02	ng	0.00
79) Terphenyl-d14	20.01	244	2607486	73.21	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.49	88	156983	37.672	ng	100
3) Pyridine	3.89	79	446575	41.348	ng	100
4) n-Nitrosodimethylamine	3.81	42	211644	38.481	ng	100
6) Aniline	7.37	93	656694	38.515	ng	100
8) 2-Chlorophenol	7.61	128	407211	39.632	ng	100
9) Benzaldehyde	7.18	77	279788	40.750	ng	100
10) Phenol	7.23	94	508752	39.381	ng	100
11) bis(2-Chloroethyl)ether	7.47	93	413502	39.180	ng	100
12) 1,3-Dichlorobenzene	7.94	146	477161	38.035	ng	100
13) 1,4-Dichlorobenzene	8.09	146	476214	38.047	ng	100
14) 1,2-Dichlorobenzene	8.41	146	449523	37.766	ng	100
15) Benzyl Alcohol	8.28	79	387682	38.444	ng	100
16) 2,2'-oxybis(1-Chloropropan	8.58	45	850987	37.537	ng	100
17) 2-Methylphenol	8.48	107	355429	38.621	ng	100
18) Hexachloroethane	9.15	117	176777	37.991	ng	100
19) n-Nitroso-di-n-propylamine	8.87	70	354227	38.765	ng	100
20) 3+4-Methylphenols	8.81	107	484168	38.215	ng	100
22) Acetophenone	8.87	105	612419	37.416	ng	100
24) Nitrobenzene	9.25	77	460376	40.615	ng	100
25) Isophorone	9.78	82	909783	38.493	ng	100
26) 2-Nitrophenol	9.96	139	201693	40.418	ng	100
27) 2,4-Dimethylphenol	10.02	122	356729	37.939	ng	100
28) bis(2-Chloroethoxy)methane	10.26	93	557140	37.989	ng	100
29) 2,4-Dichlorophenol	10.50	162	415969	38.239	ng	100
30) 1,2,4-Trichlorobenzene	10.72	180	481806	38.007	ng	100
31) Naphthalene	10.91	128	1236903	38.292	ng	100
32) Benzoic acid	10.17	122	262608	38.013	ng	100
33) 4-Chloroaniline	11.00	127	584458	37.647	ng	100
34) Hexachlorobutadiene	11.20	225	311051	37.915	ng	100
35) Caprolactam	11.79	113	154078	37.153	ng	100
36) 4-Chloro-3-methylphenol	12.12	107	437310	38.681	ng	100
37) 2-Methylnaphthalene	12.51	142	955857	38.731	ng	100
38) 1-Methylnaphthalene	12.73	142	901886	38.502	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.88	216	525542	37.849	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.87	237	307442	36.136	ng	100
43) 2,4,6-Trichlorophenol	13.11	196	349444	38.333	ng	100
44) 2,4,5-Trichlorophenol	13.18	196	371607	39.916	ng	100
46) 1,1'-Biphenyl	13.51	154	1151275	38.019	ng	100
47) 2-Chloronaphthalene	13.55	162	968660	38.409	ng	100
48) 2-Nitroaniline	13.74	65	298191	42.721	ng	100
49) Acenaphthylene	14.39	152	1484380	38.648	ng	100
50) Dimethylphthalate	14.13	163	1221649	38.046	ng	100
51) 2,6-Dinitrotoluene	14.24	165	267828	41.814	ng	100
52) Acenaphthene	14.73	154	947600	38.486	ng	100
53) 3-Nitroaniline	14.56	138	291412	40.582	ng	100
54) 2,4-Dinitrophenol	14.76	184	116311	40.223	ng	100
55) Dibenzofuran	15.07	168	1424236	38.615	ng	100
56) 4-Nitrophenol	14.86	139	243271	42.580	ng	100
57) 2,4-Dinitrotoluene	15.02	165	358526	42.554	ng	100
58) Fluorene	15.72	166	1139234	38.094	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.29	232	349095	39.825	ng	100
60) Diethylphthalate	15.49	149	1232464	37.964	ng	100
61) 4-Chlorophenyl-phenylether	15.71	204	656951	38.336	ng	100
62) 4-Nitroaniline	15.73	138	313462	41.135	ng	100
63) Azobenzene	16.00	77	1107555	38.753	ng	100
65) 4,6-Dinitro-2-methylphenol	15.78	198	167643	40.544	ng	100
66) n-Nitrosodiphenylamine	15.92	169	1064564	36.923	ng	100
67) 4-Bromophenyl-phenylether	16.60	248	442546	37.585	ng	100
68) Hexachlorobenzene	16.72	284	439515	36.905	ng	100
69) Atrazine	16.87	200	387353	44.908	ng	100
70) Pentachlorophenol	17.06	266	290664	39.297	ng	100
71) Phenanthrene	17.45	178	1797554	36.113	ng	100
72) Anthracene	17.54	178	1805680	34.686	ng	100
73) Carbazole	17.80	167	1704125	35.511	ng	100
74) Di-n-butylphthalate	18.37	149	1969765	33.459	ng	100
75) Fluoranthene	19.45	202	2117807	35.339	ng	100
77) Benzidine	19.62	184	1186641	40.822	ng	100
78) Pyrene	19.81	202	2124268	37.930	ng	100
80) Butylbenzylphthalate	20.70	149	958898	36.224	ng	100
81) Benzo(a)anthracene	21.64	228	2139140	35.519	ng	100
82) 3,3'-Dichlorobenzidine	21.56	252	873025	34.861	ng	100
83) Chrysene	21.71	228	2030203	37.870	ng	100
84) Bis(2-ethylhexyl)phthalate	21.57	149	1315186	36.327	ng	100
85) Di-n-octyl phthalate	22.78	149	2267540	36.087	ng	100
86) Indeno(1,2,3-cd)pyrene	28.52	276	2772622	38.072	ng	100
88) Benzo(b)fluoranthene	23.85	252	2346602	38.473	ng	100
89) Benzo(k)fluoranthene	23.92	252	2197476	37.805	ng	100
90) Benzo(a)pyrene	24.72	252	2261958	38.698	ng	100
91) Dibenzo(a,h)anthracene	28.59	278	2211721	38.965	ng	100
92) Benzo(g,h,i)perylene	29.66	276	2220211	39.213	ng	100

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

