

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082919\
 Data File : BG042576.D
 Acq On : 29 Aug 2019 22:06
 Operator : HP/JU
 Sample : K4549-02MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S700A-N-1.0-1.5-082619MSD

Manual Integrations
 APPROVED

Jagrut
 8/30/2019 7:58:18 PM

Quant Time: Aug 30 05:59:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	36623	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	140130	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	102251	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	237214	20.00	ng	0.00
76) Chrysene-d12	21.69	240	250211	20.00	ng	0.00
87) Perylene-d12	24.93	264	310263	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	238188	131.72	ng	0.00
7) Phenol-d6	7.17	99	314375	128.71	ng	0.00
23) Nitrobenzene-d5	9.17	82	179285	88.41	ng	0.00
42) 2,4,6-Tribromophenol	16.16	330	188982	130.03	ng	0.00
45) 2-Fluorobiphenyl	13.28	172	515507	85.27	ng	0.00
79) Terphenyl-d14	20.02	244	895477	77.37	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.42	88	25229	33.432 ng	98
3) Pyridine	3.82	79	56121	29.002 ng	99
4) n-Nitrosodimethylamine	3.73	42	31178	45.112 ng	92
6) Aniline	7.32	93	91500	28.086 ng	99
8) 2-Chlorophenol	7.57	128	99037	45.191 ng	99
9) Benzaldehyde	7.13	77	55696	45.546 ng	98
10) Phenol	7.20	94	111954	45.079 ng	95
11) bis(2-Chloroethyl)ether	7.41	93	77857	39.918 ng	98
12) 1,3-Dichlorobenzene	7.89	146	112724	40.434 ng	98
13) 1,4-Dichlorobenzene	8.04	146	115615	40.941 ng	98
14) 1,2-Dichlorobenzene	8.36	146	110679	40.927 ng	95
15) Benzyl Alcohol	8.24	79	75291	42.845 ng	99
16) 2,2'-oxybis(1-Chloropropan	8.52	45	100055	37.848 ng	98
17) 2-Methylphenol	8.45	107	78837	43.099 ng	97
18) Hexachloroethane	9.09	117	37832	39.133 ng	91
19) n-Nitroso-di-n-propylamine	8.81	70	61200	38.674 ng	# 92
20) 3+4-Methylphenols	8.78	107	107697	42.548 ng	99
22) Acetophenone	8.82	105	138451	46.194 ng	98
24) Nitrobenzene	9.21	77	91611	44.613 ng	99
25) Isophorone	9.73	82	170767	42.671 ng	98
26) 2-Nitrophenol	9.93	139	61026	47.424 ng	95
27) 2,4-Dimethylphenol	9.99	122	92995	52.462 ng	98
28) bis(2-Chloroethoxy)methane	10.22	93	106841	42.358 ng	99
29) 2,4-Dichlorophenol	10.47	162	113948	49.133 ng	99
30) 1,2,4-Trichlorobenzene	10.69	180	126968	44.602 ng	98
31) Naphthalene	10.87	128	294822	45.316 ng	99
32) Benzoic acid	10.15	122	52270m	41.885 ng	
33) 4-Chloroaniline	10.98	127	58396	19.444 ng	98
34) Hexachlorobutadiene	11.16	225	85847	44.872 ng	98
35) Caprolactam	11.76	113	32327m	41.774 ng	
36) 4-Chloro-3-methylphenol	12.11	107	100457	45.629 ng	97
37) 2-Methylnaphthalene	12.48	142	237637	45.490 ng	99
38) 1-Methylnaphthalene	12.70	142	220111	44.359 ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.85	216	160699	48.939 ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG082919\
 Data File : BG042576.D
 Acq On : 29 Aug 2019 22:06
 Operator : HP/JU
 Sample : K4549-02MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S700A-N-1.0-1.5-082619MSD

Manual Integrations
 APPROVED

Jagrut
 8/30/2019 7:58:18 PM

Quant Time: Aug 30 05:59:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.83	237	162302	94.096	nq	98
43) 2,4,6-Trichlorophenol	13.09	196	104006	46.859	nq	96
44) 2,4,5-Trichlorophenol	13.17	196	109801	47.738	nq	99
46) 1,1'-Biphenyl	13.49	154	312236	47.240	nq	99
47) 2-Chloronaphthalene	13.53	162	254342	44.628	nq	98
48) 2-Nitroaniline	13.73	65	57651	41.949	nq	99
49) Acenaphthylene	14.38	152	401787	46.860	nq	99
50) Dimethylphthalate	14.11	163	395724	53.637	nq	100
51) 2,6-Dinitrotoluene	14.23	165	78531	45.922	nq	98
52) Acenaphthene	14.72	154	232372	44.632	nq	98
53) 3-Nitroaniline	14.56	138	45546	26.631	nq	97
54) 2,4-Dinitrophenol	14.78	184	98304	84.863	nq	100
55) Dibenzofuran	15.06	168	402170	46.666	nq	99
56) 4-Nitrophenol	14.89	139	113317	93.243	nq	97
57) 2,4-Dinitrotoluene	15.03	165	110039	44.935	nq	98
58) Fluorene	15.71	166	325987	46.273	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.29	232	111325	48.943	nq	# 94
60) Diethylphthalate	15.47	149	316215	43.845	nq	99
61) 4-Chlorophenyl-phenylether	15.70	204	194792	45.869	nq	98
62) 4-Nitroaniline	15.73	138	78617	42.950	nq	95
63) Azobenzene	15.99	77	217100	43.930	nq	98
65) 4,6-Dinitro-2-methylphenol	15.80	198	76863	51.682	nq	96
66) n-Nitrosodiphenylamine	15.91	169	301362	46.196	nq	99
67) 4-Bromophenyl-phenylether	16.60	248	135259	44.953	nq	99
68) Hexachlorobenzene	16.72	284	142695	43.625	nq	98
69) Atrazine	16.87	200	114701	49.720	nq	98
70) Pentachlorophenol	17.07	266	172582	101.548	nq	98
71) Phenanthrene	17.45	178	551879	46.838	nq	99
72) Anthracene	17.55	178	570345	48.487	nq	99
73) Carbazole	17.81	167	498494	47.550	nq	99
74) Di-n-butylphthalate	18.37	149	566068	45.679	nq	100
75) Fluoranthene	19.47	202	719919	50.064	nq	97
77) Benzidine	19.64	184	372139	51.876	nq	99
78) Pyrene	19.83	202	724928	47.256	nq	100
80) Butylbenzylphthalate	20.71	149	262383	43.486	nq	99
81) Benzo(a)anthracene	21.67	228	712512	46.812	nq	98
82) 3,3'-Dichlorobenzidine	21.58	252	205172	33.516	nq	99
83) Chrysene	21.73	228	691420	47.102	nq	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	378879	45.226	nq	98
85) Di-n-octyl phthalate	22.78	149	655183	45.472	nq	100
86) Indeno(1,2,3-cd)pyrene	28.66	276	926737	46.456	nq	# 97
88) Benzo(b)fluoranthene	23.90	252	786164	46.090	nq	98
89) Benzo(k)fluoranthene	23.98	252	774693	47.250	nq	99
90) Benzo(a)pyrene	24.78	252	785190	48.511	nq	99
91) Dibenzo(a,h)anthracene	28.71	278	770431	47.011	nq	99
92) Benzo(g,h,i)perylene	29.81	276	773409	47.186	nq	98

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

