

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032918\
 Data File : BG033424.D
 Acq On : 29 Mar 2018 17:41
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/30/2018 6:04:46 PM

Quant Time: Mar 30 07:51:41 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 18:34:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.87	152	34766	20.00	ng	-0.01
21) Naphthalene-d8	11.75	136	155115	20.00	ng	0.00
38) Acenaphthene-d10	15.49	164	122682	20.00	ng	0.00
63) Phenanthrene-d10	18.21	188	331373	20.00	ng	0.00
75) Chrysene-d12	22.56	240	325707	20.00	ng	0.00
86) Perylene-d12	26.33	264	364695	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.30	112	151191	81.55	ng	0.00
7) Phenol-d6	7.98	99	250926	78.77	ng	0.00
23) Nitrobenzene-d5	10.06	82	257354	76.23	ng	-0.01
41) 2,4,6-Tribromophenol	16.96	330	157916	86.06	ng	0.00
44) 2-Fluorobiphenyl	14.11	172	657503	77.37	ng	0.00
78) Terphenyl-d14	20.73	244	1149207	75.46	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.93	88	34752	36.909	ng	88
3) Pyridine	4.39	79	107086	41.142	ng	90
4) n-Nitrosodimethylamine	4.30	42	42179	39.488	ng	# 95
6) Aniline	8.16	93	151478	40.434	ng	# 91
8) 2-Chlorophenol	8.41	128	83419	39.356	ng	96
9) Benzaldehyde	7.97	77	66213m	32.706	ng	
10) Phenol	8.01	94	118902	37.804	ng	89
11) bis(2-Chloroethyl)ether	8.25	93	94298	36.731	ng	90
12) 1,3-Dichlorobenzene	8.75	146	104590	37.743	ng	96
13) 1,4-Dichlorobenzene	8.91	146	104755	37.746	ng	95
14) 1,2-Dichlorobenzene	9.23	146	105465	38.936	ng	99
15) Benzyl Alcohol	9.10	79	104798	41.782	ng	92
16) 2,2'-oxybis(1-Chloropropan	9.39	45	158558	37.886	ng	87
17) 2-Methylphenol	9.31	107	86401m	40.325	ng	
18) Hexachloroethane	9.99	117	41637	38.872	ng	98
19) n-Nitroso-di-n-propylamine	9.67	70	95598	40.953	ng	98
20) 3+4-Methylphenols	9.65	107	124561	40.830	ng	89
22) Acetophenone	9.71	105	169351	39.851	ng	# 97
24) Nitrobenzene	10.11	77	140563	38.897	ng	91
25) Isophorone	10.63	82	277724	42.384	ng	99
26) 2-Nitrophenol	10.84	139	54397	39.760	ng	# 86
27) 2,4-Dimethylphenol	10.88	122	81067	40.788	ng	89
28) bis(2-Chloroethoxy)methane	11.11	93	126599	38.722	ng	97
29) 2,4-Dichlorophenol	11.39	162	101978	41.722	ng	94
30) 1,2,4-Trichlorobenzene	11.60	180	121148	38.149	ng	93
31) Naphthalene	11.80	128	318194	39.586	ng	98
32) Benzoic acid	11.03	122	10051m	5.440	ng	
33) 4-Chloroaniline	11.90	127	137846	38.277	ng	93
34) Hexachlorobutadiene	12.06	225	92501	38.518	ng	98
35) Caprolactam	12.64	113	45367	47.378	ng	97
36) 4-Chloro-3-methylphenol	12.97	107	135383	42.468	ng	94
37) 2-Methylnaphthalene	13.35	142	246394	39.493	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.71	216	171396	38.569	ng	98
40) Hexachlorocyclopentadiene	13.68	237	95674	36.726	ng	90

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.94	196	111713	43.268	nq	97
43) 2,4,5-Trichlorophenol	14.01	196	129755	42.518	nq	98
45) 1,1'-Biphenyl	14.32	154	367244	38.963	nq	98
46) 2-Chloronaphthalene	14.38	162	277574	38.194	nq	98
47) 2-Nitroaniline	14.57	65	115803	42.542	nq	92
48) Acenaphthylene	15.22	152	478162	39.417	nq	99
49) Dimethylphthalate	14.91	163	447646	41.928	nq	99
50) 2,6-Dinitrotoluene	15.05	165	92017	42.150	nq	94
51) Acenaphthene	15.55	154	298649	38.191	nq	94
52) 3-Nitroaniline	15.38	138	92471	40.403	nq	# 96
53) 2,4-Dinitrophenol	15.58	184	18293	22.278	nq	93
54) Dibenzofuran	15.88	168	457374	39.596	nq	93
55) 4-Nitrophenol	15.68	139	64035	39.789	nq	# 86
56) 2,4-Dinitrotoluene	15.82	165	137280	42.708	nq	92
57) Fluorene	16.52	166	394590	40.098	nq	97
58) 2,3,4,6-Tetrachlorophenol	16.09	232	130464m	45.410	nq	
59) Diethylphthalate	16.25	149	442432	42.676	nq	97
60) 4-Chlorophenyl-phenylether	16.49	204	224380	39.665	nq	98
61) 4-Nitroaniline	16.53	138	99273	42.832	nq	# 86
62) Azobenzene	16.79	77	400185	39.848	nq	96
64) 4,6-Dinitro-2-methylphenol	16.58	198	58971	29.748	nq	94
65) n-Nitrosodiphenylamine	16.71	169	362669	38.336	nq	99
66) 4-Bromophenyl-phenylether	17.39	248	148345	38.119	nq	96
67) Hexachlorobenzene	17.51	284	158240	38.528	nq	94
68) Atrazine	17.63	200	150545	39.271	nq	99
69) Pentachlorophenol	17.86	266	96957	34.395	nq	96
70) Phenanthrene	18.26	178	660682	38.283	nq	98
71) Anthracene	18.35	178	675159	38.638	nq	98
72) Carbazole	18.61	167	594961	38.423	nq	98
73) Di-n-butylphthalate	19.10	149	702704	38.989	nq	# 98
74) Fluoranthene	20.21	202	793093	37.136	nq	97
76) Benzidine	20.37	184	375513	42.514	nq	98
77) Pyrene	20.57	202	812435	37.400	nq	100
79) Butylbenzylphthalate	21.42	149	314621	39.248	nq	96
80) Benzo(a)anthracene	22.54	228	795413	38.352	nq	99
81) 3,3'-Dichlorobenzidine	22.43	252	310227	42.035	nq	97
82) Chrysene	22.62	228	766409	38.175	nq	98
83) Bis(2-ethylhexyl)phthalate	22.36	149	445702	40.334	nq	# 96
84) Di-n-octyl phthalate	23.75	149	718555	42.469	nq	98
85) Indeno(1,2,3-cd)pyrene	30.71	276	906249	41.751	nq	# 89
87) Benzo(b)fluoranthene	25.12	252	783121	37.167	nq	# 96
88) Benzo(k)fluoranthene	25.20	252	792056	37.168	nq	# 97
89) Benzo(a)pyrene	26.15	252	758895	37.505	nq	# 96
90) Dibenzo(a,h)anthracene	30.79	278	758252	39.782	nq	# 97
91) Benzo(g,h,i)perylene	32.09	276	736373	39.016	nq	# 94

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

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