

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030818\
 Data File : BG033042.D
 Acq On : 8 Mar 2018 15:59
 Operator : SJ/JU
 Sample : J1748-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 AU-06-030618-AMS

Manual Integrations
 APPROVED

Sohil
 3/9/2018 11:41:23 AM

Quant Time: Mar 08 16:45:46 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 06 17:38:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.91	152	56592	20.00	ng	0.00
21) Naphthalene-d8	11.80	136	244728	20.00	ng	0.00
38) Acenaphthene-d10	15.53	164	136645	20.00	ng	0.00
63) Phenanthrene-d10	18.26	188	307582	20.00	ng	0.00
75) Chrysene-d12	22.62	240	287278	20.00	ng	0.00
86) Perylene-d12	26.41	264	312927	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.33	112	481008	135.98	ng	0.00
7) Phenol-d6	8.01	99	581388	117.92	ng	0.00
23) Nitrobenzene-d5	10.11	82	438197	119.68	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	244170	169.94	ng	0.00
44) 2-Fluorobiphenyl	14.16	172	901816	95.18	ng	0.00
78) Terphenyl-d14	20.77	244	1305676	102.11	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.98	88	57781	37.638	ng	# 89
3) Pyridine	4.44	79	111950	26.458	ng	92
4) n-Nitrosodimethylamine	4.33	42	94889	55.934	ng	# 85
6) Aniline	8.21	93	109061	16.341	ng	97
8) 2-Chlorophenol	8.46	128	185193	47.831	ng	96
9) Benzaldehyde	8.01	77	57581	20.263	ng	89
10) Phenol	8.04	94	200827	40.406	ng	97
11) bis(2-Chloroethyl)ether	8.29	93	176161	43.345	ng	98
12) 1,3-Dichlorobenzene	8.80	146	185267	40.854	ng	99
13) 1,4-Dichlorobenzene	8.95	146	187707	41.121	ng	98
14) 1,2-Dichlorobenzene	9.29	146	180778	41.328	ng	99
15) Benzyl Alcohol	9.15	79	169058	46.640	ng	95
16) 2,2'-oxybis(1-Chloropropan	9.44	45	349376	50.006	ng	99
17) 2-Methylphenol	9.34	107	158026	43.117	ng	97
18) Hexachloroethane	10.04	117	69880	44.962	ng	# 87
19) n-Nitroso-di-n-propylamine	9.73	70	151602	43.553	ng	# 92
20) 3+4-Methylphenols	9.67	107	215818	42.350	ng	95
22) Acetophenone	9.76	105	278268	45.023	ng	# 98
24) Nitrobenzene	10.16	77	212900	53.680	ng	95
25) Isophorone	10.68	82	408498	47.556	ng	96
26) 2-Nitrophenol	10.88	139	98045	63.683	ng	95
27) 2,4-Dimethylphenol	10.91	122	195124	52.291	ng	91
28) bis(2-Chloroethoxy)methane	11.15	93	243725	48.705	ng	96
29) 2,4-Dichlorophenol	11.42	162	173230	51.150	ng	93
30) 1,2,4-Trichlorobenzene	11.65	180	171664	43.241	ng	98
31) Naphthalene	11.85	128	599047	46.845	ng	99
32) Benzoic acid	11.03	122	104296	45.826	ng	# 83
33) 4-Chloroaniline	11.94	127	74299	12.994	ng	94
34) Hexachlorobutadiene	12.11	225	97467	43.770	ng	96
35) Caprolactam	12.68	113	47888	35.314	ng	96
36) 4-Chloro-3-methylphenol	12.99	107	205217	52.071	ng	92
37) 2-Methylnaphthalene	13.40	142	421165	45.522	ng	92
39) 1,2,4,5-Tetrachlorobenzene	13.75	216	181235	44.679	ng	# 96
40) Hexachlorocyclopentadiene	13.73	237	221819	107.301	ng	91

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.97	196	134492	52.575	nq	97
43) 2,4,5-Trichlorophenol	14.04	196	149298	50.427	nq #	95
45) 1,1'-Biphenyl	14.37	154	528920	44.296	nq	96
46) 2-Chloronaphthalene	14.43	162	400401	44.890	nq	97
47) 2-Nitroaniline	14.61	65	150026	62.863	nq	94
48) Acenaphthylene	15.26	152	651616	44.621	nq	98
49) Dimethylphthalate	14.95	163	526159	45.349	nq	99
50) 2,6-Dinitrotoluene	15.08	165	116379	58.266	nq	93
51) Acenaphthene	15.60	154	456912	50.239	nq	98
52) 3-Nitroaniline	15.41	138	72735	33.347	nq #	86
53) 2,4-Dinitrophenol	15.61	184	103557	135.329	nq	95
54) Dibenzofuran	15.92	168	611903	47.252	nq	94
55) 4-Nitrophenol	15.68	139	173956	90.534	nq	87
56) 2,4-Dinitrotoluene	15.86	165	166635	66.820	nq #	86
57) Fluorene	16.56	166	494744	46.288	nq	99
58) 2,3,4,6-Tetrachlorophenol	16.13	232	128849m	56.054	nq	
59) Diethylphthalate	16.29	149	569029	46.499	nq	97
60) 4-Chlorophenyl-phenylether	16.54	204	247323	47.302	nq	90
61) 4-Nitroaniline	16.56	138	123154	48.917	nq	88
62) Azobenzene	16.83	77	524392	47.899	nq	96
64) 4,6-Dinitro-2-methylphenol	16.61	198	76197	71.281	nq	90
65) n-Nitrosodiphenylamine	16.75	169	461193	46.091	nq	99
66) 4-Bromophenyl-phenylether	17.43	248	149736	46.039	nq #	85
67) Hexachlorobenzene	17.56	284	156318	44.476	nq #	91
68) Atrazine	17.66	200	163251	49.566	nq	96
69) Pentachlorophenol	17.89	266	208783	94.308	nq	98
70) Phenanthrene	18.30	178	798780	46.549	nq	98
71) Anthracene	18.39	178	816204	47.377	nq	99
72) Carbazole	18.64	167	761653	44.854	nq	98
73) Di-n-butylphthalate	19.14	149	979703	47.029	nq	99
74) Fluoranthene	20.25	202	913224	48.177	nq	94
76) Benzidine	20.40	184	57891	5.912	nq	98
77) Pyrene	20.61	202	914497	45.140	nq	98
79) Butylbenzylphthalate	21.46	149	462505	51.492	nq	95
80) Benzo(a)anthracene	22.59	228	871656	47.317	nq	100
81) 3,3'-Dichlorobenzidine	22.47	252	150602	22.073	nq #	99
82) Chrysene	22.67	228	826294	46.956	nq	97
83) Bis(2-ethylhexyl)phthalate	22.42	149	653759	49.171	nq #	96
84) Di-n-octyl phthalate	23.83	149	1122025	55.365	nq	99
85) Indeno(1,2,3-cd)pyrene	30.84	276	924960	45.070	nq	96
87) Benzo(b)fluoranthene	25.19	252	868588	46.945	nq #	97
88) Benzo(k)fluoranthene	25.27	252	873552	47.506	nq #	96
89) Benzo(a)pyrene	26.23	252	848210	48.198	nq #	97
90) Dibenzo(a,h)anthracene	30.91	278	746705	42.154	nq #	94
91) Benzo(g,h,i)perylene	32.23	276	777630	44.533	nq #	92

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

