

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042318\
 Data File : BG033918.D
 Acq On : 23 Apr 2018 13:22
 Operator : SJ/JU
 Sample : J2490-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDP-B-3-18-20MS

Manual Integrations
 APPROVED

Sohil
 4/24/2018 4:37:24 PM

Quant Time: Apr 24 03:47:38 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 19 13:56:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	70495	20.00	ng	0.00
21) Naphthalene-d8	10.62	136	325027	20.00	ng	-0.01
38) Acenaphthene-d10	14.46	164	214378	20.00	ng	-0.01
63) Phenanthrene-d10	17.21	188	547171	20.00	ng	-0.01
75) Chrysene-d12	21.47	240	554297	20.00	ng	-0.02
86) Perylene-d12	24.53	264	578567	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.44	112	657531	148.91	ng	0.00
7) Phenol-d6	7.01	99	899265	139.63	ng	0.00
23) Nitrobenzene-d5	8.98	82	557943	97.69	ng	0.00
41) 2,4,6-Tribromophenol	15.96	330	412918	165.10	ng	-0.01
44) 2-Fluorobiphenyl	13.09	172	1199411	82.19	ng	-0.01
78) Terphenyl-d14	19.85	244	1940495	78.65	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.42	88	69783	37.401	ng	93
3) Pyridine	3.80	79	204575	37.896	ng	92
4) n-Nitrosodimethylamine	3.72	42	112996	45.944	ng	# 90
6) Aniline	7.17	93	189624	22.027	ng	97
8) 2-Chlorophenol	7.40	128	236778	47.857	ng	97
9) Benzaldehyde	6.98	77	121709	34.459	ng	93
10) Phenol	7.03	94	338833	51.357	ng	98
11) bis(2-Chloroethyl)ether	7.27	93	211860	41.985	ng	95
12) 1,3-Dichlorobenzene	7.72	146	237663	41.728	ng	97
13) 1,4-Dichlorobenzene	7.87	146	242597	42.054	ng	99
14) 1,2-Dichlorobenzene	8.18	146	230063	41.010	ng	96
15) Benzyl Alcohol	8.07	79	237140	42.227	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.36	45	380692	38.681	ng	95
17) 2-Methylphenol	8.27	107	209708	42.892	ng	96
18) Hexachloroethane	8.91	117	88886	42.167	ng	98
19) n-Nitroso-di-n-propylamine	8.63	70	203905	37.143	ng	98
20) 3+4-Methylphenols	8.59	107	294234	41.417	ng	96
22) Acetophenone	8.65	105	352545	39.852	ng	# 97
24) Nitrobenzene	9.02	77	279340	44.884	ng	96
25) Isophorone	9.55	82	536541	41.785	ng	96
26) 2-Nitrophenol	9.73	139	136415	55.474	ng	95
27) 2,4-Dimethylphenol	9.79	122	240393	52.264	ng	96
28) bis(2-Chloroethoxy)methane	10.03	93	315615	46.508	ng	96
29) 2,4-Dichlorophenol	10.26	162	261063	50.914	ng	95
30) 1,2,4-Trichlorobenzene	10.48	180	243870	42.464	ng	96
31) Naphthalene	10.67	128	703672	42.100	ng	99
32) Benzoic acid	9.90	122	3257m	5.719	ng	
33) 4-Chloroaniline	10.77	127	154201	19.628	ng	98
34) Hexachlorobutadiene	10.96	225	151049	43.966	ng	98
35) Caprolactam	11.55	113	106716m	48.952	ng	
36) 4-Chloro-3-methylphenol	11.90	107	290532	46.424	ng	92
37) 2-Methylnaphthalene	12.28	142	533590	41.697	ng	94
39) 1,2,4,5-Tetrachlorobenzene	12.65	216	306399	43.042	ng	98
40) Hexachlorocyclopentadiene	12.63	237	345906	88.373	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.89	196	229742	52.933	nq	91
43) 2,4,5-Trichlorophenol	12.96	196	254395	50.983	nq	98
45) 1,1'-Biphenyl	13.29	154	725245	41.792	nq	97
46) 2-Chloronaphthalene	13.34	162	555011	42.179	nq	98
47) 2-Nitroaniline	13.54	65	207409	50.483	nq	93
48) Acenaphthylene	14.19	152	903619	41.313	nq	99
49) Dimethylphthalate	13.93	163	872945	46.291	nq	98
50) 2,6-Dinitrotoluene	14.04	165	180564	50.525	nq	87
51) Acenaphthene	14.53	154	555753	40.447	nq	99
52) 3-Nitroaniline	14.36	138	104327	26.885	nq	# 84
53) 2,4-Dinitrophenol	14.57	184	68503	53.599	nq	# 58
54) Dibenzofuran	14.86	168	892935	43.835	nq	95
55) 4-Nitrophenol	14.68	139	344560	113.214	nq	90
56) 2,4-Dinitrotoluene	14.83	165	257659	54.218	nq	86
57) Fluorene	15.52	166	749830	42.537	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.09	232	255346	55.890	nq	99
59) Diethylphthalate	15.30	149	830956	41.688	nq	98
60) 4-Chlorophenyl-phenylether	15.52	204	402826	43.690	nq	97
61) 4-Nitroaniline	15.54	138	205910	49.976	nq	92
62) Azobenzene	15.81	77	743426	41.840	nq	99
64) 4,6-Dinitro-2-methylphenol	15.59	198	112534	43.426	nq	91
65) n-Nitrosodiphenylamine	15.73	169	678382	42.644	nq	98
66) 4-Bromophenyl-phenylether	16.41	248	261878	43.580	nq	96
67) Hexachlorobenzene	16.52	284	269518	42.057	nq	# 73
68) Atrazine	16.69	200	292544	46.942	nq	98
69) Pentachlorophenol	16.87	266	361481	82.119	nq	94
70) Phenanthrene	17.26	178	1220027	41.563	nq	97
71) Anthracene	17.35	178	1260316	42.934	nq	96
72) Carbazole	17.62	167	1174041	41.457	nq	96
73) Di-n-butylphthalate	18.20	149	1418127	40.252	nq	98
74) Fluoranthene	19.28	202	1487699	41.898	nq	95
76) Benzidine	19.46	184	418396	27.953	nq	99
77) Pyrene	19.64	202	1479082	40.683	nq	93
79) Butylbenzylphthalate	20.56	149	675730	42.323	nq	92
80) Benzo(a)anthracene	21.46	228	1476052	42.230	nq	96
81) 3,3'-Dichlorobenzidine	21.38	252	364674	27.982	nq	97
82) Chrysene	21.52	228	1397836	41.880	nq	95
83) Bis(2-ethylhexyl)phthalate	21.40	149	932396	41.352	nq	97
84) Di-n-octyl phthalate	22.57	149	1623961	45.376	nq	100
85) Indeno(1,2,3-cd)pyrene	28.00	276	1736574	45.707	nq	# 93
87) Benzo(b)fluoranthene	23.57	252	1496527	42.400	nq	99
88) Benzo(k)fluoranthene	23.64	252	1478955	42.183	nq	98
89) Benzo(a)pyrene	24.39	252	1470669	43.507	nq	100
90) Dibenzo(a,h)anthracene	28.07	278	1440638	44.045	nq	96
91) Benzo(g,h,i)perylene	29.08	276	1426133	44.311	nq	97

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

