

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_N\DATA\BN020918\
 Data File : BN000052.D
 Acq On : 09 Feb 2018 11:06
 Operator : JU/SJ
 Sample : MDL-S-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MDL-S-02

Manual Integrations
 APPROVED

Sohil
 2/13/2018 11:33:30 AM

Quant Time: Feb 12 14:58:52 2018
 Quant Method : Z:\HPCHEM1\BNA_N\METHODS\8270-BN020718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 18:24:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	406307	20.00	ng	0.00
21) Naphthalene-d8	11.30	136	2015954	20.00	ng	0.00
38) Acenaphthene-d10	15.07	164	1211148	20.00	ng	0.00
63) Phenanthrene-d10	17.80	188	2657541	20.00	ng	0.00
75) Chrysene-d12	21.91	240	2815155	20.00	ng	-0.01
86) Perylene-d12	24.40	264	2876861	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.94	112	3479577	139.57	ng	0.00
7) Phenol-d6	7.59	99	4961301	131.38	ng	0.00
23) Nitrobenzene-d5	9.65	82	2955163	87.16	ng	0.00
41) 2,4,6-Tribromophenol	16.55	330	1691213	141.07	ng	0.00
44) 2-Fluorobiphenyl	13.71	172	7443603	86.62	ng	0.00
78) Terphenyl-d14	20.36	244	9057839	86.88	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.66	88	36553	3.644	ng	98
3) Pyridine	4.12	79	97047	3.108	ng	96
4) n-Nitrosodimethylamine	4.02	42	28255	3.200	ng	# 88
6) Aniline	7.77	93	160831	3.091	ng	90
8) 2-Chlorophenol	8.02	128	100311	3.441	ng	90
9) Benzaldehyde	7.58	77	69223	2.947	ng	97
10) Phenol	7.62	94	136017	3.412	ng	91
11) bis(2-Chloroethyl)ether	7.87	93	109042	3.326	ng	84
12) 1,3-Dichlorobenzene	8.35	146	116830	3.527	ng	98
13) 1,4-Dichlorobenzene	8.50	146	120329	3.557	ng	96
14) 1,2-Dichlorobenzene	8.83	146	117962	3.542	ng	94
15) Benzyl Alcohol	8.70	79	75417	2.779	ng	86
16) 2,2'-oxybis(1-Chloropropan	9.00	45	120263	3.466	ng	80
17) 2-Methylphenol	8.89	107	84381	2.979	ng	93
18) Hexachloroethane	9.57	117	38137	3.238	ng	# 75
19) n-Nitroso-di-n-propylamine	9.27	70	72487	2.865	ng	# 85
20) 3+4-Methylphenols	9.22	107	114127	2.881	ng	93
22) Acetophenone	9.30	105	184957	3.274	ng	# 87
24) Nitrobenzene	9.69	77	139479	3.763	ng	# 90
25) Isophorone	10.20	82	204057	2.911	ng	98
26) 2-Nitrophenol	10.40	139	30009	8.872	ng	# 86
27) 2,4-Dimethylphenol	10.44	122	84972m	2.796	ng	
28) bis(2-Chloroethoxy)methane	10.69	93	145407	3.235	ng	96
29) 2,4-Dichlorophenol	10.94	162	69300	2.464	ng	96
30) 1,2,4-Trichlorobenzene	11.16	180	114549	3.537	ng	97
31) Naphthalene	11.36	128	389314	3.523	ng	97
32) Benzoic acid	10.48	122	22812m	11.332	ng	
33) 4-Chloroaniline	11.45	127	141445	2.913	ng	95
34) Hexachlorobutadiene	11.63	225	59246	3.528	ng	97
35) Caprolactam	12.17	113	20202	5.524	ng	92
36) 4-Chloro-3-methylphenol	12.53	107	85048	2.382	ng	99
37) 2-Methylnaphthalene	12.93	142	258998	3.368	ng	95
39) 1,2,4,5-Tetrachlorobenzene	13.29	216	124719	3.489	ng	# 97
40) Hexachlorocyclopentadiene	13.27	237	39454	2.648	ng	89

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.52	196	49177	2.295	ng	96
43) 2,4,5-Trichlorophenol	13.59	196	60564	2.351	ng	# 94
45) 1,1'-Biphenyl	13.92	154	397664	3.728	ng	98
46) 2-Chloronaphthalene	13.97	162	280334	3.503	ng	95
47) 2-Nitroaniline	14.15	65	42480	6.504	ng	# 82
48) Acenaphthylene	14.80	152	406832	3.155	ng	98
49) Dimethylphthalate	14.51	163	328422	3.244	ng	99
50) 2,6-Dinitrotoluene	14.63	165	46179	2.209	ng	90
51) Acenaphthene	15.14	154	289920	3.454	ng	99
52) 3-Nitroaniline	14.96	138	56441	2.222	ng	# 87
53) 2,4-Dinitrophenol	15.16	184	9864	10.405	ng	85
54) Dibenzofuran	15.47	168	425199	3.577	ng	95
55) 4-Nitrophenol	15.23	139	31117	6.378	ng	# 78
56) 2,4-Dinitrotoluene	15.41	165	52304	5.779	ng	# 97
57) Fluorene	16.11	166	344370	3.548	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.68	232	47135	2.319	ng	# 85
59) Diethylphthalate	15.86	149	324539	3.154	ng	97
60) 4-Chlorophenyl-phenylether	16.09	204	156906	3.547	ng	# 87
61) 4-Nitroaniline	16.11	138	55790	2.137	ng	90
62) Azobenzene	16.39	77	310307	3.030	ng	90
64) 4,6-Dinitro-2-methylphenol	16.16	198	18401	9.904	ng	89
65) n-Nitrosodiphenylamine	16.30	169	273613	3.164	ng	98
66) 4-Bromophenyl-phenylether	16.98	248	83324	3.297	ng	# 83
67) Hexachlorobenzene	17.10	284	101433	3.437	ng	# 85
68) Atrazine	17.23	200	70546	2.507	ng	96
69) Pentachlorophenol	17.43	266	35963	7.472	ng	96
70) Phenanthrene	17.83	178	570011	3.631	ng	98
71) Anthracene	17.93	178	498652	3.281	ng	99
72) Carbazole	18.19	167	494956	3.232	ng	97
73) Di-n-butylphthalate	18.72	149	397425	2.351	ng	# 97
74) Fluoranthene	19.82	202	537067	3.261	ng	95
76) Benzidine	19.98	184	129578	1.444	ng	97
77) Pyrene	20.17	202	603892	3.375	ng	99
79) Butylbenzylphthalate	21.03	149	132386	6.761	ng	96
80) Benzo(a)anthracene	21.90	228	583453	3.372	ng	97
81) 3,3'-Dichlorobenzidine	21.82	252	142468	2.242	ng	# 97
82) Chrysene	21.95	228	625924	3.630	ng	98
83) Bis(2-ethylhexyl)phthalate	21.80	149	210299	4.472	ng	92
84) Di-n-octyl phthalate	22.76	149	282436	7.737	ng	99
85) Indeno(1,2,3-cd)pyrene	26.96	276	595334	2.811	ng	# 86
87) Benzo(b)fluoranthene	23.64	252	535770	3.182	ng	# 96
88) Benzo(k)fluoranthene	23.69	252	576028	3.396	ng	# 95
89) Benzo(a)pyrene	24.29	252	489249	3.030	ng	# 95
90) Dibenzo(a,h)anthracene	26.97	278	509985	3.020	ng	# 92
91) Benzo(g,h,i)perylene	27.74	276	499799	2.927	ng	# 86

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

