

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072617\
 Data File : BM010984.D
 Acq On : 26 Jul 2017 23:12
 Operator : SJ/JU
 Sample : PB100891BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB100891BS

Manual Integrations
 APPROVED

Sohil
 7/27/2017 6:21:37 PM

Quant Time: Jul 27 09:46:17 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 17:29:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	232741	20.00	ng	0.00
21) Naphthalene-d8	10.15	136	1042452	20.00	ng	0.00
38) Acenaphthene-d10	14.05	164	776808	20.00	ng	0.00
63) Phenanthrene-d10	16.81	188	1981422	20.00	ng	0.00
75) Chrysene-d12	21.03	240	1817914	20.00	ng	0.00
86) Perylene-d12	23.14	264	1224469	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.03	112	1840852	133.70	ng	0.00
7) Phenol-d6	6.60	99	2514305	128.53	ng	0.00
23) Nitrobenzene-d5	8.55	82	2109011	75.93	ng	0.00
41) 2,4,6-Tribromophenol	15.56	330	1512877	121.51	ng	0.00
44) 2-Fluorobiphenyl	12.66	172	4843645	78.20	ng	0.00
78) Terphenyl-d14	19.47	244	7490968	81.62	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.04	88	185917	30.075	ng	# 100
3) Pyridine	3.41	79	565369	33.483	ng	# 88
4) n-Nitrosodimethylamine	3.33	42	362901	42.174	ng	# 65
6) Aniline	6.74	93	850696	34.504	ng	# 90
8) 2-Chlorophenol	6.97	128	580380	41.520	ng	# 82
9) Benzaldehyde	6.55	77	282347	22.308	ng	# 88
10) Phenol	6.62	94	871344	41.010	ng	# 99
11) bis(2-Chloroethyl)ether	6.85	93	630777	38.101	ng	# 93
12) 1,3-Dichlorobenzene	7.28	146	637993	37.451	ng	# 83
13) 1,4-Dichlorobenzene	7.43	146	660884	37.847	ng	# 92
14) 1,2-Dichlorobenzene	7.74	146	630987	37.794	ng	# 92
15) Benzyl Alcohol	7.64	79	841485	44.841	ng	# 81
16) 2,2'-oxybis(1-Chloropropan	7.92	45	596910	40.068	ng	# 77
17) 2-Methylphenol	7.85	107	578545	42.604	ng	# 78
18) Hexachloroethane	8.45	117	291799	38.319	ng	# 81
19) n-Nitroso-di-n-propylamine	8.20	70	673446	40.512	ng	# 92
20) 3+4-Methylphenols	8.17	107	813695	43.106	ng	# 82
22) Acetophenone	8.21	105	1095249	35.050	ng	# 94
24) Nitrobenzene	8.58	77	1058127	36.764	ng	# 86
25) Isophorone	9.11	82	1712876	36.987	ng	# 86
26) 2-Nitrophenol	9.28	139	339715	37.095	ng	# 24
27) 2,4-Dimethylphenol	9.36	122	639446	39.664	ng	# 75
28) bis(2-Chloroethoxy)methane	9.59	93	983253	38.067	ng	# 98
29) 2,4-Dichlorophenol	9.81	162	731880	40.400	ng	# 92
30) 1,2,4-Trichlorobenzene	10.02	180	853684	38.112	ng	# 97
31) Naphthalene	10.20	128	1998736	38.113	ng	# 98
32) Benzoic acid	9.52	122	438900	33.870	ng	# 57
33) 4-Chloroaniline	10.33	127	377446	18.042	ng	# 74
34) Hexachlorobutadiene	10.49	225	676912	38.369	ng	# 97
35) Caprolactam	11.12	113	205502m	40.001	ng	#
36) 4-Chloro-3-methylphenol	11.46	107	896919	38.343	ng	# 75
37) 2-Methylnaphthalene	11.83	142	1613948	41.894	ng	# 88
39) 1,2,4,5-Tetrachlorobenzene	12.21	216	1188581	35.614	ng	# 100
40) Hexachlorocyclopentadiene	12.19	237	1883828	96.871	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.46	196	761610	37.703	nq	97
43) 2,4,5-Trichlorophenol	12.53	196	811878	39.023	nq #	88
45) 1,1'-Biphenyl	12.87	154	2321474	34.562	nq	98
46) 2-Chloronaphthalene	12.91	162	1843715	37.255	nq	95
47) 2-Nitroaniline	13.13	65	710357	40.573	nq #	65
48) Acenaphthylene	13.77	152	2772763	37.542	nq	99
49) Dimethylphthalate	13.53	163	2432423	36.604	nq #	99
50) 2,6-Dinitrotoluene	13.64	165	508328	37.828	nq #	63
51) Acenaphthene	14.12	154	1721717	37.649	nq	98
52) 3-Nitroaniline	13.97	138	372384	31.934	nq #	47
53) 2,4-Dinitrophenol	14.18	184	723833	75.792	nq #	90
54) Dibenzofuran	14.46	168	3024591	40.818	nq #	91
55) 4-Nitrophenol	14.30	139	764504	74.621	nq	88
56) 2,4-Dinitrotoluene	14.44	165	746249	39.557	nq	94
57) Fluorene	15.11	166	2511093	38.923	nq	100
58) 2,3,4,6-Tetrachlorophenol	14.69	232	842166	43.186	nq #	100
59) Diethylphthalate	14.91	149	2592015	38.193	nq	100
60) 4-Chlorophenyl-phenylether	15.12	204	1529766	39.274	nq	98
61) 4-Nitroaniline	15.14	138	476841	38.498	nq #	1
62) Azobenzene	15.41	77	3170660	43.653	nq	89
64) 4,6-Dinitro-2-methylphenol	15.20	198	497468	37.086	nq	90
65) n-Nitrosodiphenylamine	15.33	169	2210977	38.991	nq	99
66) 4-Bromophenyl-phenylether	16.01	248	1021859	40.117	nq #	89
67) Hexachlorobenzene	16.12	284	1049451	36.460	nq #	88
68) Atrazine	16.30	200	906336	39.884	nq	96
69) Pentachlorophenol	16.46	266	1363407	74.610	nq	97
70) Phenanthrene	16.85	178	4125406	39.762	nq	99
71) Anthracene	16.94	178	4149659	40.104	nq	99
72) Carbazole	17.22	167	3420260	37.797	nq	100
73) Di-n-butylphthalate	17.80	149	4282098	38.859	nq #	97
74) Fluoranthene	18.88	202	4956890	38.553	nq	98
76) Benzidine	19.09	184	2132064	49.025	nq	100
77) Pyrene	19.24	202	4832909	44.742	nq	99
79) Butylbenzylphthalate	20.19	149	1704660	44.379	nq #	71
80) Benzo(a)anthracene	21.01	228	4157041	40.048	nq	100
81) 3,3'-Dichlorobenzidine	20.96	252	1084319	26.628	nq #	95
82) Chrysene	21.06	228	3824279	38.381	nq	99
83) Bis(2-ethylhexyl)phthalate	20.97	149	2338496	41.384	nq #	97
84) Di-n-octyl phthalate	21.82	149	3422624	37.319	nq	100
85) Indeno(1,2,3-cd)pyrene	25.21	276	2713106	26.292	nq #	97
87) Benzo(b)fluoranthene	22.51	252	3356718	42.713	nq #	93
88) Benzo(k)fluoranthene	22.56	252	3047155	40.454	nq #	95
89) Benzo(a)pyrene	23.05	252	2844844	40.005	nq #	96
90) Dibenzo(a,h)anthracene	25.23	278	2234175	33.891	nq #	91
91) Benzo(g,h,i)perylene	25.84	276	2184886	33.752	nq #	86

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

