

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081917\
 Data File : BM011313.D
 Acq On : 19 Aug 2017 12:01
 Operator : SJ/JU
 Sample : PB101645BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB101645BS

Manual Integrations
 APPROVED

Sohil
 8/21/2017 5:20:48 PM

Quant Time: Aug 21 01:58:49 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 02:25:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.27	152	179739	20.00	ng	-0.02
21) Naphthalene-d8	10.02	136	648448	20.00	ng	-0.02
38) Acenaphthene-d10	13.94	164	431048	20.00	ng	-0.02
63) Phenanthrene-d10	16.70	188	1053268	20.00	ng	-0.02
75) Chrysene-d12	20.93	240	1154291	20.00	ng	-0.02
86) Perylene-d12	22.99	264	1107098	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	4.91	112	1149069	108.07	ng	-0.02
7) Phenol-d6	6.48	99	1669580	110.52	ng	-0.02
23) Nitrobenzene-d5	8.42	82	1438734	83.27	ng	-0.02
41) 2,4,6-Tribromophenol	15.45	330	690869	100.00	ng	-0.02
44) 2-Fluorobiphenyl	12.55	172	2460405	71.59	ng	-0.02
78) Terphenyl-d14	19.37	244	3761326	64.55	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.90	88	132908	27.840	ng	# 100
3) Pyridine	3.28	79	403485	30.942	ng	# 87
4) n-Nitrosodimethylamine	3.20	42	256148	38.546	ng	# 71
6) Aniline	6.62	93	547845	28.773	ng	# 82
8) 2-Chlorophenol	6.85	128	338707	31.376	ng	# 72
9) Benzaldehyde	6.43	77	401817	41.108	ng	# 71
10) Phenol	6.50	94	564898	34.427	ng	# 93
11) bis(2-Chloroethyl)ether	6.72	93	414059	32.386	ng	# 83
12) 1,3-Dichlorobenzene	7.16	146	393227	29.890	ng	# 77
13) 1,4-Dichlorobenzene	7.30	146	410809	30.464	ng	# 90
14) 1,2-Dichlorobenzene	7.61	146	389695	30.224	ng	# 87
15) Benzyl Alcohol	7.52	79	530205	36.585	ng	# 77
16) 2,2'-oxybis(1-Chloropropan	7.81	45	448485	38.982	ng	# 77
17) 2-Methylphenol	7.73	107	359767	34.306	ng	# 70
18) Hexachloroethane	8.32	117	185164	31.486	ng	# 79
19) n-Nitroso-di-n-propylamine	8.08	70	473844	36.910	ng	# 90
20) 3+4-Methylphenols	8.05	107	478196	32.803	ng	# 80
22) Acetophenone	8.09	105	670847	34.513	ng	# 87
24) Nitrobenzene	8.46	77	726051	40.554	ng	# 81
25) Isophorone	8.98	82	1103427	38.304	ng	# 83
26) 2-Nitrophenol	9.16	139	191469	33.611	ng	# 7
27) 2,4-Dimethylphenol	9.24	122	337394	33.644	ng	# 70
28) bis(2-Chloroethoxy)methane	9.47	93	578368	35.998	ng	# 95
29) 2,4-Dichlorophenol	9.69	162	369545	32.794	ng	# 91
30) 1,2,4-Trichlorobenzene	9.89	180	469404	33.689	ng	# 96
31) Naphthalene	10.07	128	1079669	33.097	ng	# 99
32) Benzoic acid	9.37	122	212966	26.420	ng	# 36
33) 4-Chloroaniline	10.21	127	182044	13.989	ng	# 60
34) Hexachlorobutadiene	10.36	225	379262	34.560	ng	# 94
35) Caprolactam	10.99	113	110348m	34.530	ng	#
36) 4-Chloro-3-methylphenol	11.35	107	477178	32.794	ng	# 71
37) 2-Methylnaphthalene	11.70	142	838160	34.976	ng	# 84
39) 1,2,4,5-Tetrachlorobenzene	12.09	216	614163	33.164	ng	# 100
40) Hexachlorocyclopentadiene	12.06	237	905109	83.876	ng	# 96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.35	196	372445	33.227	ng	94
43) 2,4,5-Trichlorophenol	12.42	196	396053	34.306	ng #	86
45) 1,1'-Biphenyl	12.76	154	1121927	30.101	ng	95
46) 2-Chloronaphthalene	12.79	162	905928	32.989	ng #	92
47) 2-Nitroaniline	13.02	65	404954	41.682	ng #	60
48) Acenaphthylene	13.65	152	1366941	33.354	ng	99
49) Dimethylphthalate	13.42	163	1187441	32.202	ng #	98
50) 2,6-Dinitrotoluene	13.53	165	249390	33.445	ng #	35
51) Acenaphthene	14.00	154	821131	32.358	ng	99
52) 3-Nitroaniline	13.86	138	198955	30.747	ng #	20
53) 2,4-Dinitrophenol	14.07	184	360662	68.057	ng #	77
54) Dibenzofuran	14.35	168	1432404	34.837	ng #	90
55) 4-Nitrophenol	14.20	139	300023	52.774	ng #	78
56) 2,4-Dinitrotoluene	14.33	165	345494	33.004	ng #	67
57) Fluorene	15.00	166	1170268	32.690	ng	98
58) 2,3,4,6-Tetrachlorophenol	14.58	232	417853	38.615	ng #	100
59) Diethylphthalate	14.80	149	1166824	30.984	ng	97
60) 4-Chlorophenyl-phenylether	15.00	204	703551	32.551	ng	95
61) 4-Nitroaniline	15.04	138	169702	24.691	ng #	1
62) Azobenzene	15.30	77	1728788	42.894	ng	84
64) 4,6-Dinitro-2-methylphenol	15.10	198	250765	35.168	ng	90
65) n-Nitrosodiphenylamine	15.23	169	1042842	34.597	ng	99
66) 4-Bromophenyl-phenylether	15.90	248	456771	33.734	ng #	91
67) Hexachlorobenzene	16.00	284	469456	30.682	ng #	82
68) Atrazine	16.20	200	414556	34.319	ng	98
69) Pentachlorophenol	16.36	266	629424	64.797	ng	97
70) Phenanthrene	16.74	178	1962119	35.577	ng	99
71) Anthracene	16.83	178	1959952	35.633	ng	99
72) Carbazole	17.12	167	1594665	33.152	ng	99
73) Di-n-butylphthalate	17.70	149	1907116	32.557	ng #	95
74) Fluoranthene	18.77	202	2454644	35.915	ng	99
76) Benzidine	18.99	184	1706253	61.790	ng	98
77) Pyrene	19.14	202	2430744	35.441	ng	99
79) Butylbenzylphthalate	20.09	149	822529	33.725	ng #	52
80) Benzo(a)anthracene	20.91	228	2347625	35.619	ng	99
81) 3,3'-Dichlorobenzidine	20.86	252	824716	31.897	ng #	96
82) Chrysene	20.96	228	2197930	34.741	ng	100
83) Bis(2-ethylhexyl)phthalate	20.88	149	1144883	31.909	ng #	99
84) Di-n-octyl phthalate	21.72	149	1824533	31.332	ng #	92
85) Indeno(1,2,3-cd)pyrene	25.00	276	2697850	41.174	ng #	96
87) Benzo(b)fluoranthene	22.39	252	2351994	33.101	ng #	91
88) Benzo(k)fluoranthene	22.43	252	2249649	33.033	ng #	93
89) Benzo(a)pyrene	22.90	252	2294386	35.685	ng #	96
90) Dibenzo(a,h)anthracene	25.02	278	2240059	37.582	ng #	92
91) Benzo(g,h,i)perylene	25.61	276	2304323	39.371	ng #	84

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

