

Data Path : Z:\HPCHEM1\BNA N\DATA\BN040918\
 Data File : BN000715.D
 Acq On : 09 Apr 2018 18:34
 Operator : JU/SJ
 Sample : PB108094BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB108094BS

Manual Integrations
 APPROVED

Sohil
 4/10/2018 3:38:25 PM

Quant Time: Apr 10 04:55:10 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN040918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 09 13:41:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	153176	20.00	ng	0.00
21) Naphthalene-d8	10.40	136	615574	20.00	ng	0.00
38) Acenaphthene-d10	14.27	164	289343	20.00	ng	0.00
63) Phenanthrene-d10	17.02	188	522039	20.00	ng	0.00
75) Chrysene-d12	21.22	240	417633	20.00	ng	0.00
86) Perylene-d12	23.41	264	389580	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.26	112	1020642	98.39	ng	0.00
7) Phenol-d6	6.82	99	1321258	93.56	ng	0.00
23) Nitrobenzene-d5	8.78	82	1138913	97.91	ng	0.00
41) 2,4,6-Tribromophenol	15.76	330	270228	94.99	ng	0.00
44) 2-Fluorobiphenyl	12.89	172	2057667	98.01	ng	0.00
78) Terphenyl-d14	19.66	244	1961153	102.57	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	189236	39.762	ng	96
3) Pyridine	3.63	79	542038	42.198	ng	98
4) n-Nitrosodimethylamine	3.55	42	182986	52.596	ng	# 82
6) Aniline	6.98	93	764772	40.569	ng	93
8) 2-Chlorophenol	7.21	128	520949	48.620	ng	90
9) Benzaldehyde	6.79	77	126969	17.725	ng	95
10) Phenol	6.85	94	722331	48.102	ng	87
11) bis(2-Chloroethyl)ether	7.08	93	553626	44.875	ng	84
12) 1,3-Dichlorobenzene	7.53	146	539486	43.443	ng	99
13) 1,4-Dichlorobenzene	7.68	146	547773	43.542	ng	98
14) 1,2-Dichlorobenzene	7.99	146	521019	43.131	ng	96
15) Benzyl Alcohol	7.88	79	464026	48.240	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.18	45	580058	43.340	ng	78
17) 2-Methylphenol	8.08	107	454452	45.389	ng	95
18) Hexachloroethane	8.70	117	210663	45.432	ng	# 82
19) n-Nitroso-di-n-propylamine	8.45	70	392053	44.413	ng	# 82
20) 3+4-Methylphenols	8.40	107	604993	44.024	ng	94
22) Acetophenone	8.45	105	794643	44.834	ng	# 88
24) Nitrobenzene	8.82	77	603072	47.744	ng	# 95
25) Isophorone	9.35	82	1064643	49.577	ng	99
26) 2-Nitrophenol	9.53	139	246777	49.355	ng	93
27) 2,4-Dimethylphenol	9.59	122	472561	55.484	ng	97
28) bis(2-Chloroethoxy)methane	9.83	93	701391	49.644	ng	97
29) 2,4-Dichlorophenol	10.06	162	420864	50.899	ng	97
30) 1,2,4-Trichlorobenzene	10.27	180	443104	44.485	ng	97
31) Naphthalene	10.46	128	1504953	45.246	ng	98
32) Benzoic acid	9.73	122	287400	38.257	ng	95
33) 4-Chloroaniline	10.56	127	494141	36.553	ng	97
34) Hexachlorobutadiene	10.75	225	235694	43.994	ng	97
35) Caprolactam	11.33	113	127730	47.333	ng	96
36) 4-Chloro-3-methylphenol	11.69	107	481457	48.939	ng	100
37) 2-Methylnaphthalene	12.07	142	1003543	46.106	ng	96
39) 1,2,4,5-Tetrachlorobenzene	12.45	216	430994	48.443	ng	# 96
40) Hexachlorocyclopentadiene	12.43	237	591148	119.142	ng	94

Data Path : Z:\HPCHEM1\BNA N\DATA\BN040918\
 Data File : BN000715.D
 Acq On : 09 Apr 2018 18:34
 Operator : JU/SJ
 Sample : PB108094BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB108094BS

Manual Integrations
 APPROVED

Sohil
 4/10/2018 3:38:25 PM

Quant Time: Apr 10 04:55:10 2018
 Quant Method : Z:\HPCHEM1\BNA N\METHODS\8270-BN040918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 09 13:41:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	12.69	196	282022	53.609	nq	99
43) 2,4,5-Trichlorophenol	12.76	196	302506	49.478	nq	# 93
45) 1,1'-Biphenyl	13.10	154	1216290	46.824	nq	97
46) 2-Chloronaphthalene	13.13	162	913774	46.810	nq	97
47) 2-Nitroaniline	13.34	65	272488	49.217	nq	89
48) Acenaphthylene	13.99	152	1426792	46.384	nq	98
49) Dimethylphthalate	13.74	163	1041208	45.205	nq	99
50) 2,6-Dinitrotoluene	13.85	165	227531	47.707	nq	93
51) Acenaphthene	14.33	154	843307	44.369	nq	99
52) 3-Nitroaniline	14.17	138	195098	35.718	nq	# 96
53) 2,4-Dinitrophenol	14.38	184	247576	97.989	nq	96
54) Dibenzofuran	14.67	168	1284706	47.018	nq	95
55) 4-Nitrophenol	14.49	139	421312	101.340	nq	# 86
56) 2,4-Dinitrotoluene	14.63	165	299303	48.333	nq	95
57) Fluorene	15.32	166	994077	46.065	nq	96
58) 2,3,4,6-Tetrachlorophenol	14.89	232	241940m	51.071	nq	
59) Diethylphthalate	15.12	149	1026241	45.337	nq	97
60) 4-Chlorophenyl-phenylether	15.32	204	458821	45.478	nq	91
61) 4-Nitroaniline	15.34	138	243804	45.794	nq	94
62) Azobenzene	15.62	77	1189417	47.612	nq	89
64) 4,6-Dinitro-2-methylphenol	15.40	198	142693	45.873	nq	89
65) n-Nitrosodiphenylamine	15.54	169	861239	49.437	nq	97
66) 4-Bromophenyl-phenylether	16.22	248	254849	48.822	nq	# 87
67) Hexachlorobenzene	16.32	284	281747	46.747	nq	# 91
68) Atrazine	16.50	200	256203	53.148	nq	94
69) Pentachlorophenol	16.66	266	351872	89.914	nq	98
70) Phenanthrene	17.06	178	1434075	47.263	nq	99
71) Anthracene	17.15	178	1455171	49.171	nq	98
72) Carbazole	17.42	167	1304891	46.540	nq	96
73) Di-n-butylphthalate	18.01	149	1531814	46.985	nq	99
74) Fluoranthene	19.08	202	1417507	46.674	nq	93
76) Benzidine	19.27	184	965280	84.553	nq	96
77) Pyrene	19.45	202	1444244	48.384	nq	100
79) Butylbenzylphthalate	20.37	149	617885	46.746	nq	# 94
80) Benzo(a)anthracene	21.20	228	1273232	48.450	nq	99
81) 3,3'-Dichlorobenzidine	21.15	252	337046	42.091	nq	# 98
82) Chrysene	21.26	228	1195246	46.630	nq	98
83) Bis(2-ethylhexyl)phthalate	21.17	149	920443	48.901	nq	# 93
84) Di-n-octyl phthalate	22.05	149	1438392	54.503	nq	# 94
85) Indeno(1,2,3-cd)pyrene	26.26	276	1070040	49.104	nq	# 91
87) Benzo(b)fluoranthene	22.76	252	1151701	48.713	nq	# 96
88) Benzo(k)fluoranthene	22.80	252	1157874	48.727	nq	# 95
89) Benzo(a)pyrene	23.32	252	1105045	51.165	nq	# 94
90) Dibenzo(a,h)anthracene	25.61	278	1086219	51.020	nq	# 92
91) Benzo(g,h,i)perylene	26.26	276	1070040	51.777	nq	# 89

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

