

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040921\
 Data File : BP005251.D
 Acq On : 09 Apr 2021 14:46
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 4/12/2021 2:43:23 PM

Quant Time: Apr 09 16:47:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP040921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 09 16:43:24 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.681	152	23690	20.00	ng	# 0.00
21) Naphthalene-d8	10.469	136	79314	20.00	ng	# 0.00
39) Acenaphthene-d10	14.339	164	47953	20.00	ng	0.00
64) Phenanthrene-d10	17.104	188	99661	20.00	ng	# 0.00
76) Chrysene-d12	21.204	240	108529	20.00	ng	0.00
86) Perylene-d12	23.474	264	116683	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.281	112	46748	30.37	ng	0.00
7) Phenol-d6	6.863	99	53297	24.17	ng	0.00
23) Nitrobenzene-d5	8.840	82	49283	26.66	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	11389	18.23	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	70696	19.94	ng	0.00
79) Terphenyl-d14	19.680	244	108315	18.17	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.222	88	14013	20.90	ng	# 93
3) Pyridine	3.640	79	29014	17.41	ng	89
4) n-Nitrosodimethylamine	3.534	42	9536	16.00	ng	# 16
6) Aniline	7.016	93	28938	11.58	ng	# 81
8) 2-Chlorophenol	7.252	128	16841	10.37	ng	# 78
9) Benzaldehyde	6.834	77	19480	17.35	ng	# 62
10) Phenol	6.893	94	28375	12.56	ng	88
11) bis(2-Chloroethyl)ether	7.116	93	19467	11.77	ng	88
12) 1,3-Dichlorobenzene	7.569	146	18385	10.19	ng	# 82
13) 1,4-Dichlorobenzene	7.716	146	18581	10.15	ng	# 85
14) 1,2-Dichlorobenzene	8.028	146	17579	10.01	ng	# 91
15) Benzyl Alcohol	7.934	79	17360	10.20	ng	# 78
16) 2,2'-oxybis(1-Chloropr...	8.210	45	11439	9.43	ng	81
17) 2-Methylphenol	8.134	107	14822	10.28	ng	# 86
18) Hexachloroethane	8.746	117	7776	11.14	ng	# 82
19) n-Nitroso-di-n-propyla...	8.487	70	15235	10.65	ng	# 84
20) 3+4-Methylphenols	8.463	107	19484	9.89	ng	84
22) Acetophenone	8.505	105	28105	12.33	ng	# 89
24) Nitrobenzene	8.881	77	25139	12.90	ng	# 65
25) Isophorone	9.404	82	40907	11.99	ng	# 82
26) 2-Nitrophenol	9.593	139	7038	9.32	ng	# 18
27) 2,4-Dimethylphenol	9.657	122	11566	9.52	ng	# 86
28) bis(2-Chloroethoxy)met...	9.893	93	19201	11.12	ng	# 95
29) 2,4-Dichlorophenol	10.128	162	12511	9.20	ng	92
30) 1,2,4-Trichlorobenzene	10.334	180	15661	10.26	ng	95
31) Naphthalene	10.516	128	43745	9.76	ng	98
32) Benzoic acid	9.775	122	5725	6.35	ng	# 52
33) 4-Chloroaniline	10.646	127	15448	8.52	ng	# 73
34) Hexachlorobutadiene	10.799	225	10380	10.56	ng	97
35) Caprolactam	11.434	113	3748	8.40	ng	# 75
36) 4-Chloro-3-methylphenol	11.781	107	15709	9.57	ng	# 79
37) 2-Methylnaphthalene	12.140	142	28038	8.67	ng	# 90
38) 1-Methylnaphthalene	12.363	142	27812	9.17	ng	# 95
40) 1,2,4,5-Tetrachloroben...	12.516	216	15995	10.24	ng	# 96
41) Hexachlorocyclopentadiene	12.487	237	5550	6.53	ng	97
43) 2,4,6-Trichlorophenol	12.769	196	9343m	8.66	ng	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.845	196	10582	9.05	ng	# 82
46) 1,1'-Biphenyl	13.169	154	35279	9.65	ng	94
47) 2-Chloronaphthalene	13.204	162	28736	9.65	ng	98
48) 2-Nitroaniline	13.428	65	8920	9.40	ng	# 41
49) Acenaphthylene	14.057	152	44082	9.52	ng	94
50) Dimethylphthalate	13.804	163	34288	9.28	ng	98
51) 2,6-Dinitrotoluene	13.922	165	6935	7.97	ng	# 9
52) Acenaphthene	14.404	154	26241	9.21	ng	99
53) 3-Nitroaniline	14.263	138	5936	7.41	ng	# 51
54) 2,4-Dinitrophenol	14.481	184	3036	5.76	ng	# 59
55) Dibenzofuran	14.745	168	43653	9.35	ng	# 89
56) 4-Nitrophenol	14.592	139	4634	7.05	ng	# 1
57) 2,4-Dinitrotoluene	14.722	165	9314	8.10	ng	# 13
58) Fluorene	15.398	166	35091	9.31	ng	98
59) 2,3,4,6-Tetrachlorophenol	14.981	232	9628	8.93	ng	96
60) Diethylphthalate	15.175	149	33292	9.22	ng	98
61) 4-Chlorophenyl-phenyle...	15.392	204	18501	9.24	ng	94
62) 4-Nitroaniline	15.428	138	6263	7.61	ng	# 35
63) Azobenzene	15.686	77	51747	12.53	ng	# 78
65) 4,6-Dinitro-2-methylph...	15.492	198	5389	7.41	ng	# 65
66) n-Nitrosodiphenylamine	15.610	169	29574	9.23	ng	98
67) 4-Bromophenyl-phenylether	16.292	248	11514	9.97	ng	93
68) Hexachlorobenzene	16.404	284	12206	9.40	ng	# 90
69) Atrazine	16.575	200	9698	9.11	ng	95
70) Pentachlorophenol	16.757	266	5891	6.87	ng	96
71) Phenanthrene	17.145	178	54319	9.40	ng	99
72) Anthracene	17.239	178	52127	9.23	ng	99
73) Carbazole	17.516	167	49665	9.36	ng	98
74) Di-n-butylphthalate	18.075	149	51301	8.79	ng	# 94
75) Fluoranthene	19.127	202	65244	9.60	ng	98
77) Benzidine	19.316	184	15357	6.45	ng	# 96
78) Pyrene	19.480	202	67782	9.09	ng	98
80) Butylbenzylphthalate	20.357	149	22873	8.27	ng	# 58
81) Benzo(a)anthracene	21.186	228	72085	9.78	ng	99
82) 3,3'-Dichlorobenzidine	21.127	252	19609	7.91	ng	97
83) Chrysene	21.239	228	69060	9.54	ng	96
84) Bis(2-ethylhexyl)phtha...	21.116	149	36133	8.91	ng	98
85) Di-n-octyl phthalate	21.998	149	66164	9.67	ng	# 90
87) Indeno(1,2,3-cd)pyrene	25.804	276	90132	10.30	ng	# 95
88) Benzo(b)fluoranthene	22.786	252	75327	9.09	ng	98
89) Benzo(k)fluoranthene	22.833	252	72798m	9.17	ng	
90) Benzo(a)pyrene	23.374	252	70021	9.42	ng	98
91) Dibenzo(a,h)anthracene	25.815	278	77442	10.24	ng	96
92) Benzo(g,h,i)perylene	26.509	276	75855	10.39	ng	# 91

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

