

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021221\
 Data File : BP004732.D
 Acq On : 12 Feb 2021 16:24
 Operator : CG/JU
 Sample : M1396-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 FK-BPM-04-013-ABCDMSD

Manual Integrations
 APPROVED

mohammad
 2/15/2021 11:26:37 AM

Quant Time: Feb 12 17:40:03 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 18:04:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	49366	20.00	ng	0.00
21) Naphthalene-d8	10.563	136	203243	20.00	ng	# 0.00
39) Acenaphthene-d10	14.416	164	135066	20.00	ng	0.00
64) Phenanthrene-d10	17.175	188	292448	20.00	ng	0.00
76) Chrysene-d12	21.257	240	297998	20.00	ng	0.00
86) Perylene-d12	23.568	264	289172	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.358	112	391690	122.51	ng	0.00
7) Phenol-d6	6.940	99	525879	119.60	ng	0.00
23) Nitrobenzene-d5	8.928	82	328133	70.34	ng	0.00
42) 2,4,6-Tribromophenol	15.916	330	213493	118.97	ng	0.00
45) 2-Fluorobiphenyl	13.040	172	741844	67.13	ng	0.00
79) Terphenyl-d14	19.739	244	1119017	63.08	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.281	88	49552	36.22	ng	# 89
3) Pyridine	3.669	79	139997	40.41	ng	# 95
4) n-Nitrosodimethylamine	3.587	42	69385	48.13	ng	# 61
6) Aniline	7.099	93	170055	34.68	ng	# 85
8) 2-Chlorophenol	7.334	128	167822	50.24	ng	# 85
9) Benzaldehyde	6.910	77	67602	36.61	ng	# 80
10) Phenol	6.969	94	243292	57.16	ng	# 78
11) bis(2-Chloroethyl)ether	7.199	93	149081	45.80	ng	# 94
12) 1,3-Dichlorobenzene	7.657	146	175736	43.23	ng	# 91
13) 1,4-Dichlorobenzene	7.805	146	178887	43.32	ng	# 94
14) 1,2-Dichlorobenzene	8.122	146	169828	42.88	ng	# 96
15) Benzyl Alcohol	8.010	79	164677	49.88	ng	# 85
16) 2,2'-oxybis(1-Chloropr...	8.304	45	120925	45.93	ng	# 95
17) 2-Methylphenol	8.210	107	150988	52.50	ng	# 89
18) Hexachloroethane	8.846	117	67790	43.81	ng	# 94
19) n-Nitroso-di-n-propyla...	8.575	70	128561	48.39	ng	# 88
20) 3+4-Methylphenols	8.540	107	203706	53.37	ng	# 92
22) Acetophenone	8.593	105	257607	43.61	ng	# 93
24) Nitrobenzene	8.969	77	193842	42.04	ng	# 79
25) Isophorone	9.493	82	352279	47.26	ng	# 91
26) 2-Nitrophenol	9.681	139	92093	50.42	ng	# 66
27) 2,4-Dimethylphenol	9.740	122	158950	55.95	ng	# 89
28) bis(2-Chloroethoxy)met...	9.981	93	201278	41.91	ng	# 99
29) 2,4-Dichlorophenol	10.210	162	167766	48.88	ng	# 94
30) 1,2,4-Trichlorobenzene	10.428	180	170590	40.23	ng	# 98
31) Naphthalene	10.616	128	504078	42.49	ng	# 100
32) Benzoic acid	9.887	122	130435	53.33	ng	# 75
33) 4-Chloroaniline	10.722	127	82324	16.91	ng	# 89
34) Hexachlorobutadiene	10.898	225	110685	38.87	ng	# 100
35) Caprolactam	11.504	113	59666m	55.01	ng	# 83
36) 4-Chloro-3-methylphenol	11.851	107	204196	50.10	ng	# 95
37) 2-Methylnaphthalene	12.228	142	375520	44.50	ng	# 99
38) 1-Methylnaphthalene	12.451	142	351003	43.95	ng	# 98
40) 1,2,4,5-Tetrachloroben...	12.604	216	203625	41.96	ng	# 99
41) Hexachlorocyclopentadiene	12.581	237	256349	92.88	ng	# 99
43) 2,4,6-Trichlorophenol	12.840	196	144520	47.27	ng	# 99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.910	196	155138	48.29	ng	# 93
46) 1,1'-Biphenyl	13.251	154	485934	42.98	ng	98
47) 2-Chloronaphthalene	13.287	162	377130	40.69	ng	96
48) 2-Nitroaniline	13.492	65	127118	46.47	ng	# 61
49) Acenaphthylene	14.139	152	605650	44.24	ng	98
50) Dimethylphthalate	13.881	163	547172	46.06	ng	99
51) 2,6-Dinitrotoluene	13.992	165	112558	47.01	ng	# 58
52) Acenaphthene	14.481	154	367965	42.39	ng	98
53) 3-Nitroaniline	14.322	138	63371	25.45	ng	# 68
54) 2,4-Dinitrophenol	14.528	184	139818	101.39	ng	# 72
55) Dibenzofuran	14.822	168	591948	43.37	ng	97
56) 4-Nitrophenol	14.639	139	203813	104.33	ng	# 57
57) 2,4-Dinitrotoluene	14.781	165	157999	48.17	ng	# 63
58) Fluorene	15.469	166	494045	44.05	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.045	232	145432m	47.58	ng	
60) Diethylphthalate	15.251	149	523663	43.98	ng	99
61) 4-Chlorophenyl-phenyle...	15.469	204	261512	42.00	ng	96
62) 4-Nitroaniline	15.492	138	112952	41.53	ng	# 60
63) Azobenzene	15.763	77	492330	44.02	ng	86
65) 4,6-Dinitro-2-methylph...	15.545	198	98121	54.91	ng	82
66) n-Nitrosodiphenylamine	15.681	169	435770	45.82	ng	96
67) 4-Bromophenyl-phenylether	16.363	248	158842	42.19	ng	# 91
68) Hexachlorobenzene	16.475	284	174716	42.59	ng	# 90
69) Atrazine	16.639	200	139895	44.84	ng	100
70) Pentachlorophenol	16.822	266	243519	106.07	ng	99
71) Phenanthrene	17.216	178	768583	42.92	ng	100
72) Anthracene	17.304	178	791703	45.98	ng	98
73) Carbazole	17.580	167	709893	44.94	ng	99
74) Di-n-butylphthalate	18.139	149	879558	45.38	ng	# 98
75) Fluoranthene	19.186	202	917401	43.00	ng	100
77) Benzidine	19.369	184	310282	53.53	ng	98
78) Pyrene	19.539	202	930580	44.28	ng	99
80) Butylbenzylphthalate	20.416	149	394558	47.77	ng	# 79
81) Benzo(a)anthracene	21.245	228	916967	43.39	ng	98
82) 3,3'-Dichlorobenzidine	21.180	252	214717	32.05	ng	98
83) Chrysene	21.298	228	884619	43.22	ng	99
84) Bis(2-ethylhexyl)phtha...	21.174	149	580341	46.55	ng	100
85) Di-n-octyl phthalate	22.068	149	984897	48.71	ng	# 93
87) Indeno(1,2,3-cd)pyrene	25.951	276	1043195	45.18	ng	96
88) Benzo(b)fluoranthene	22.868	252	933182	44.06	ng	98
89) Benzo(k)fluoranthene	22.915	252	893821	43.96	ng	98
90) Benzo(a)pyrene	23.468	252	831121	43.42	ng	# 98
91) Dibenzo(a,h)anthracene	25.956	278	878109	45.59	ng	# 97
92) Benzo(g,h,i)perylene	26.674	276	842965	45.61	ng	# 95

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

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