

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101520\
 Data File : BP003623.D
 Acq On : 15 Oct 2020 14:14
 Operator : CG/JU
 Sample : PB132384BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB132384BS

Quant Time: Oct 15 14:55:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 17:03:23 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.540	152	129589	20.00	ng	0.00
21) Naphthalene-d8	11.393	136	450691	20.00	ng	# 0.00
39) Acenaphthene-d10	15.145	164	279607	20.00	ng	0.00
64) Phenanthrene-d10	17.863	188	563729	20.00	ng	0.00
76) Chrysene-d12	21.874	240	501064	20.00	ng	# 0.00
86) Perylene-d12	24.451	264	472213	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.034	112	850674	112.85	ng	0.00
7) Phenol-d6	7.687	99	1104966	100.63	ng	0.00
23) Nitrobenzene-d5	9.710	82	694732	71.47	ng	0.00
42) 2,4,6-Tribromophenol	16.616	330	401618	107.73	ng	0.00
45) 2-Fluorobiphenyl	13.775	172	1410409	66.18	ng	0.00
79) Terphenyl-d14	20.327	244	1883685	68.02	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.699	88	128359	39.75	ng	# 86
3) Pyridine	4.140	79	356391	37.88	ng	# 90
4) n-Nitrosodimethylamine	4.034	42	168952	41.78	ng	# 62
6) Aniline	7.834	93	380352	29.60	ng	# 86
8) 2-Chlorophenol	8.104	128	325701	42.50	ng	# 81
9) Benzaldehyde	7.634	77	190878	32.38	ng	# 79
10) Phenol	7.716	94	438307	41.34	ng	77
11) bis(2-Chloroethyl)ether	7.922	93	340693	40.02	ng	91
12) 1,3-Dichlorobenzene	8.440	146	398872	40.03	ng	# 92
13) 1,4-Dichlorobenzene	8.581	146	400599	40.07	ng	# 93
14) 1,2-Dichlorobenzene	8.916	146	379554	39.59	ng	94
15) Benzyl Alcohol	8.769	79	350960	39.79	ng	# 82
16) 2,2'-oxybis(1-Chloropr...	9.069	45	322904	37.05	ng	99
17) 2-Methylphenol	8.987	107	286238	39.84	ng	# 87
18) Hexachloroethane	9.669	117	156114	40.17	ng	96
19) n-Nitroso-di-n-propyla...	9.346	70	273350	37.35	ng	# 85
20) 3+4-Methylphenols	9.316	107	379205	39.65	ng	91
22) Acetophenone	9.363	105	534296	40.15	ng	# 88
24) Nitrobenzene	9.751	77	417633	43.64	ng	# 75
25) Isophorone	10.281	82	720351	39.71	ng	# 89
26) 2-Nitrophenol	10.481	139	156117	53.94	ng	# 64
27) 2,4-Dimethylphenol	10.534	122	289334	48.40	ng	# 81
28) bis(2-Chloroethoxy)met...	10.751	93	423889	38.01	ng	98
29) 2,4-Dichlorophenol	11.034	162	315456	42.28	ng	95
30) 1,2,4-Trichlorobenzene	11.257	180	391346	39.54	ng	98
31) Naphthalene	11.445	128	974306	40.32	ng	99
32) Benzoic acid	10.675	122	139418	43.79	ng	# 61
33) 4-Chloroaniline	11.522	127	184732	17.71	ng	# 85
34) Hexachlorobutadiene	11.751	225	284632	38.80	ng	98
35) Caprolactam	12.245	113	83568	35.34	ng	# 54
36) 4-Chloro-3-methylphenol	12.622	107	335748	39.42	ng	# 81
37) 2-Methylnaphthalene	13.010	142	702210	39.06	ng	# 95
38) 1-Methylnaphthalene	13.228	142	646462	38.25	ng	# 99
40) 1,2,4,5-Tetrachloroben...	13.381	216	460918	43.48	ng	# 98
41) Hexachlorocyclopentadiene	13.375	237	696720	117.93	ng	99
43) 2,4,6-Trichlorophenol	13.598	196	254724	43.33	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.675	196	272188	44.09	ng	#	94
46) 1,1'-Biphenyl	13.987	154	886718	41.04	ng		98
47) 2-Chloronaphthalene	14.034	162	708473	39.39	ng		98
48) 2-Nitroaniline	14.204	65	197085	39.35	ng	#	57
49) Acenaphthylene	14.869	152	1034677	39.49	ng		100
50) Dimethylphthalate	14.569	163	842765	36.71	ng		99
51) 2,6-Dinitrotoluene	14.681	165	182760	44.20	ng	#	63
52) Acenaphthene	15.210	154	731658	42.87	ng		87
53) 3-Nitroaniline	15.010	138	97086	23.29	ng	#	57
54) 2,4-Dinitrophenol	15.216	184	160065	77.14	ng	#	1
55) Dibenzofuran	15.534	168	1026877	38.71	ng		97
56) 4-Nitrophenol	15.316	139	254675	85.08	ng	#	40
57) 2,4-Dinitrotoluene	15.463	165	236656	39.93	ng	#	64
58) Fluorene	16.181	166	815137	37.98	ng		99
59) 2,3,4,6-Tetrachlorophenol	15.763	232	240136	42.29	ng		93
60) Diethylphthalate	15.910	149	821334	36.40	ng		99
61) 4-Chlorophenyl-phenyle...	16.157	204	484416	38.09	ng		95
62) 4-Nitroaniline	16.163	138	156522	33.74	ng	#	52
63) Azobenzene	16.445	77	812210	37.41	ng		85
65) 4,6-Dinitro-2-methylph...	16.233	198	119719	48.98	ng		89
66) n-Nitrosodiphenylamine	16.363	169	703937	41.81	ng		99
67) 4-Bromophenyl-phenylether	17.039	248	297681	40.23	ng		96
68) Hexachlorobenzene	17.204	284	334378	39.57	ng		97
69) Atrazine	17.280	200	226668	34.54	ng		98
70) Pentachlorophenol	17.528	266	369725	96.17	ng		99
71) Phenanthrene	17.904	178	1236841	39.49	ng		99
72) Anthracene	17.992	178	1254590	41.07	ng		99
73) Carbazole	18.233	167	1050616	38.09	ng		99
74) Di-n-butylphthalate	18.739	149	1278255	38.04	ng	#	98
75) Fluoranthene	19.816	202	1439546	37.13	ng		98
77) Benzidine	19.951	184	534815	45.53	ng		99
78) Pyrene	20.163	202	1439744	43.23	ng		99
80) Butylbenzylphthalate	20.963	149	501406	43.10	ng	#	79
81) Benzo(a)anthracene	21.857	228	1329944	39.47	ng		99
82) 3,3'-Dichlorobenzidine	21.757	252	293808	24.67	ng		97
83) Chrysene	21.916	228	1262087	39.60	ng		98
84) Bis(2-ethylhexyl)phtha...	21.716	149	734761	41.62	ng		100
85) Di-n-octyl phthalate	22.680	149	1186763	41.64	ng	#	92
87) Indeno(1,2,3-cd)pyrene	27.151	276	1364991	38.09	ng	#	91
88) Benzo(b)fluoranthene	23.663	252	1295429	39.48	ng	#	97
89) Benzo(k)fluoranthene	23.715	252	1266924	41.12	ng	#	96
90) Benzo(a)pyrene	24.339	252	1127713	38.16	ng	#	96
91) Dibenzo(a,h)anthracene	27.151	278	1164166	38.95	ng	#	94
92) Benzo(g,h,i)perylene	27.986	276	1112889	39.45	ng	#	91

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

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