

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051221\
 Data File : BF123922.D
 Acq On : 12 May 2021 13:58
 Operator : JU/CG
 Sample : PB136281BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB136281BSD

Manual Integrations
 APPROVED

Sohil
 5/13/2021 2:51:24 PM

Quant Time: May 12 14:24:10 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 11 18:02:17 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.757	152	174865	20.00	ng	0.00
21) Naphthalene-d8	8.045	136	741889	20.00	ng	# 0.00
39) Acenaphthene-d10	9.798	164	349017	20.00	ng	0.00
64) Phenanthrene-d10	11.286	188	589618	20.00	ng	# 0.00
76) Chrysene-d12	13.927	240	388671	20.00	ng	# 0.00
86) Perylene-d12	15.363	264	276631	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	1096387	75.07	ng	0.00
7) Phenol-d6	6.404	99	1420524	73.39	ng	0.00
23) Nitrobenzene-d5	7.328	82	1138095	68.34	ng	0.00
42) 2,4,6-Tribromophenol	10.592	330	187388	77.94	ng	0.00
45) 2-Fluorobiphenyl	9.122	172	1429154	64.50	ng	0.00
79) Terphenyl-d14	12.874	244	1289451	59.98	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.552	88	310687	37.34	ng	# 74
3) Pyridine	3.263	79	804194	41.77	ng	# 68
4) n-Nitrosodimethylamine	3.228	42	348390	44.15	ng	# 62
6) Aniline	6.422	93	694985	31.89	ng	# 47
8) 2-Chlorophenol	6.551	128	606660	41.79	ng	88
9) Benzaldehyde	6.304	77	94630	10.65	ng	93
10) Phenol	6.422	94	805229	39.81	ng	87
11) bis(2-Chloroethyl)ether	6.498	93	661861	43.44	ng	84
12) 1,3-Dichlorobenzene	6.698	146	602246	42.19	ng	96
13) 1,4-Dichlorobenzene	6.775	146	607828	42.31	ng	99
14) 1,2-Dichlorobenzene	6.928	146	574285	42.84	ng	98
15) Benzyl Alcohol	6.910	79	529325	43.85	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.040	45	1050436	41.44	ng	88
17) 2-Methylphenol	7.028	107	513802	41.62	ng	98
18) Hexachloroethane	7.269	117	251839	42.07	ng	92
19) n-Nitroso-di-n-propyla...	7.181	70	466819	41.64	ng	88
20) 3+4-Methylphenols	7.181	107	615773	40.34	ng	# 84
22) Acetophenone	7.169	105	866289	41.98	ng	# 89
24) Nitrobenzene	7.345	77	717736	41.32	ng	97
25) Isophorone	7.587	82	1320763	41.43	ng	99
26) 2-Nitrophenol	7.663	139	319669	42.37	ng	# 80
27) 2,4-Dimethylphenol	7.704	122	533780	43.07	ng	93
28) bis(2-Chloroethoxy)met...	7.798	93	797112	44.03	ng	99
29) 2,4-Dichlorophenol	7.910	162	440008	39.43	ng	98
30) 1,2,4-Trichlorobenzene	7.987	180	451879	41.04	ng	98
31) Naphthalene	8.069	128	1763088	41.66	ng	99
32) Benzoic acid	7.851	122	308851	51.24	ng	95
33) 4-Chloroaniline	8.116	127	404117	24.33	ng	# 93
34) Hexachlorobutadiene	8.187	225	227783	40.73	ng	100
35) Caprolactam	8.492	113	173397m	44.78	ng	
36) 4-Chloro-3-methylphenol	8.610	107	564447	42.71	ng	96
37) 2-Methylnaphthalene	8.757	142	1085260	43.09	ng	98
38) 1-Methylnaphthalene	8.857	142	1006647	41.95	ng	100
40) 1,2,4,5-Tetrachloroben...	8.928	216	362835	41.67	ng	97
41) Hexachlorocyclopentadiene	8.916	237	318025	154.41	ng	95
43) 2,4,6-Trichlorophenol	9.039	196	238943	39.83	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.092	196	259294	40.66	ng	#	91
46) 1,1'-Biphenyl	9.222	154	1227556	40.63	ng		98
47) 2-Chloronaphthalene	9.245	162	899260	41.03	ng		97
48) 2-Nitroaniline	9.351	65	354064	43.10	ng	#	78
49) Acenaphthylene	9.663	152	1378075	41.75	ng		99
50) Dimethylphthalate	9.534	163	1025730	42.39	ng		99
51) 2,6-Dinitrotoluene	9.592	165	231672	41.83	ng	#	81
52) Acenaphthene	9.834	154	878451	40.21	ng		97
53) 3-Nitroaniline	9.757	138	215955	30.40	ng		97
54) 2,4-Dinitrophenol	9.869	184	218144	98.52	ng		96
55) Dibenzofuran	10.010	168	1184581	38.77	ng		99
56) 4-Nitrophenol	9.939	139	361835	87.65	ng		85
57) 2,4-Dinitrotoluene	9.998	165	294080	40.46	ng		93
58) Fluorene	10.351	166	909463	40.11	ng		96
59) 2,3,4,6-Tetrachlorophenol	10.128	232	193392	40.70	ng		90
60) Diethylphthalate	10.228	149	997987	41.73	ng		97
61) 4-Chlorophenyl-phenyle...	10.345	204	400779	41.22	ng		97
62) 4-Nitroaniline	10.381	138	283143	41.81	ng		88
63) Azobenzene	10.504	77	1270599	41.70	ng		92
65) 4,6-Dinitro-2-methylph...	10.404	198	158958	40.60	ng		89
66) n-Nitrosodiphenylamine	10.463	169	811450	38.46	ng		99
67) 4-Bromophenyl-phenylether	10.833	248	232042	39.09	ng		93
68) Hexachlorobenzene	10.898	284	224474	39.96	ng	#	86
69) Atrazine	10.992	200	220739	35.10	ng		96
70) Pentachlorophenol	11.098	266	216276	91.53	ng		96
71) Phenanthrene	11.310	178	1410915	39.39	ng		100
72) Anthracene	11.363	178	1459317	40.77	ng		100
73) Carbazole	11.522	167	1309452	39.49	ng		98
74) Di-n-butylphthalate	11.851	149	1627959	41.06	ng		100
75) Fluoranthene	12.498	202	1459226	45.26	ng		94
77) Benzidine	12.622	184	495856	41.38	ng		98
78) Pyrene	12.727	202	1460670	40.18	ng		97
80) Butylbenzylphthalate	13.351	149	719179	42.93	ng		93
81) Benzo(a)anthracene	13.916	228	1003300	38.67	ng		99
82) 3,3'-Dichlorobenzidine	13.880	252	273693	33.95	ng	#	96
83) Chrysene	13.957	228	1070818	39.70	ng		98
84) Bis(2-ethylhexyl)phtha...	13.910	149	874436	39.76	ng	#	99
85) Di-n-octyl phthalate	14.521	149	1592449	42.25	ng		97
87) Indeno(1,2,3-cd)pyrene	16.786	276	807119	44.26	ng	#	79
88) Benzo(b)fluoranthene	14.951	252	847437	41.32	ng	#	93
89) Benzo(k)fluoranthene	14.980	252	861219	46.05	ng	#	95
90) Benzo(a)pyrene	15.304	252	708507	39.78	ng	#	92
91) Dibenzo(a,h)anthracene	16.798	278	660166	42.85	ng	#	90
92) Benzo(g,h,i)perylene	17.215	276	670618	42.40	ng	#	81

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

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