

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032321\
 Data File : BF123408.D
 Acq On : 23 Mar 2021 17:40
 Operator : JU/CG
 Sample : PB135297BSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB135297BSD

Manual Integrations
 APPROVED

mohammad
 3/24/2021 1:33:10 PM

Quant Time: Mar 23 18:53:30 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 23 15:58:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	220736	20.00	ng	0.00
21) Naphthalene-d8	8.198	136	871561	20.00	ng	# 0.00
39) Acenaphthene-d10	9.963	164	502516	20.00	ng	0.00
64) Phenanthrene-d10	11.457	188	942212	20.00	ng	0.00
76) Chrysene-d12	14.110	240	692321	20.00	ng	0.00
86) Perylene-d12	15.609	264	561751	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.534	112	1160339	86.00	ng	0.01
7) Phenol-d6	6.534	99	1514836	84.34	ng	0.00
23) Nitrobenzene-d5	7.475	82	1207247	74.76	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	494884	88.09	ng	0.00
45) 2-Fluorobiphenyl	9.280	172	2234237	63.52	ng	0.00
79) Terphenyl-d14	13.045	244	2544939	62.30	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.810	88	267368	41.20	ng	# 67
3) Pyridine	3.557	79	721170	42.26	ng	# 60
4) n-Nitrosodimethylamine	3.510	42	457408	44.80	ng	# 65
6) Aniline	6.575	93	787279	35.82	ng	93
8) 2-Chlorophenol	6.698	128	669327	44.99	ng	87
9) Benzaldehyde	6.463	77	208447	20.52	ng	99
10) Phenol	6.551	94	848070	46.00	ng	95
11) bis(2-Chloroethyl)ether	6.651	93	666755	45.36	ng	84
12) 1,3-Dichlorobenzene	6.857	146	705981m	43.97	ng	
13) 1,4-Dichlorobenzene	6.934	146	707703	44.08	ng	98
14) 1,2-Dichlorobenzene	7.086	146	682966	45.79	ng	98
15) Benzyl Alcohol	7.051	79	637798	45.35	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.186	45	1506562	45.66	ng	85
17) 2-Methylphenol	7.163	107	583705	47.72	ng	99
18) Hexachloroethane	7.434	117	276615	45.98	ng	98
19) n-Nitroso-di-n-propyla...	7.328	70	525040	45.19	ng	# 72
20) 3+4-Methylphenols	7.310	107	702857	46.13	ng	# 82
22) Acetophenone	7.322	105	929751	42.49	ng	# 89
24) Nitrobenzene	7.498	77	774241	43.37	ng	97
25) Isophorone	7.739	82	1444517	41.61	ng	98
26) 2-Nitrophenol	7.810	139	337020	52.81	ng	# 86
27) 2,4-Dimethylphenol	7.845	122	666478	46.96	ng	97
28) bis(2-Chloroethoxy)met...	7.945	93	885188	47.31	ng	100
29) 2,4-Dichlorophenol	8.051	162	611161	43.88	ng	98
30) 1,2,4-Trichlorobenzene	8.139	180	634650	41.77	ng	99
31) Naphthalene	8.222	128	1929460	41.18	ng	99
32) Benzoic acid	7.969	122	469318	56.05	ng	96
33) 4-Chloroaniline	8.263	127	465589	24.22	ng	# 94
34) Hexachlorobutadiene	8.345	225	372593	40.20	ng	97
35) Caprolactam	8.639	113	187167m	44.68	ng	
36) 4-Chloro-3-methylphenol	8.745	107	670244	45.85	ng	94
37) 2-Methylnaphthalene	8.916	142	1365531	42.51	ng	100
38) 1-Methylnaphthalene	9.016	142	1287800	42.70	ng	98
40) 1,2,4,5-Tetrachloroben...	9.080	216	591964	39.59	ng	# 98
41) Hexachlorocyclopentadiene	9.075	237	787027	94.27	ng	96
43) 2,4,6-Trichlorophenol	9.192	196	453193	43.28	ng	97

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44) 2,4,5-Trichlorophenol	9.227	196	470631	42.78	ng	97
46) 1,1'-Biphenyl	9.380	154	1623975	41.94	ng	99
47) 2-Chloronaphthalene	9.404	162	1262766	39.97	ng	99
48) 2-Nitroaniline	9.498	65	479847	50.05	ng	# 74
49) Acenaphthylene	9.827	152	2018746	40.85	ng	99
50) Dimethylphthalate	9.686	163	1574965	44.13	ng	99
51) 2,6-Dinitrotoluene	9.745	165	351067	47.09	ng	# 82
52) Acenaphthene	9.998	154	1422631	46.12	ng	99
53) 3-Nitroaniline	9.910	138	269361	34.21	ng	86
54) 2,4-Dinitrophenol	10.016	184	347660	96.04	ng	# 81
55) Dibenzofuran	10.174	168	1795522	40.49	ng	100
56) 4-Nitrophenol	10.069	139	612791	96.50	ng	# 62
57) 2,4-Dinitrotoluene	10.151	165	464531	49.92	ng	# 89
58) Fluorene	10.516	166	1376054	41.10	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.286	232	407824	45.41	ng	93
60) Diethylphthalate	10.392	149	1556746	44.68	ng	99
61) 4-Chlorophenyl-phenyle...	10.504	204	699217	42.73	ng	95
62) 4-Nitroaniline	10.533	138	366588	49.13	ng	94
63) Azobenzene	10.669	77	1589308	41.45	ng	91
65) 4,6-Dinitro-2-methylph...	10.557	198	239124	47.27	ng	# 68
66) n-Nitrosodiphenylamine	10.627	169	1269781	40.22	ng	99
67) 4-Bromophenyl-phenylether	10.998	248	475667	42.62	ng	98
68) Hexachlorobenzene	11.069	284	512340	41.66	ng	94
69) Atrazine	11.157	200	401936	40.06	ng	98
70) Pentachlorophenol	11.257	266	630526	85.92	ng	99
71) Phenanthrene	11.480	178	2101678	40.35	ng	99
72) Anthracene	11.533	178	2150456	41.05	ng	99
73) Carbazole	11.686	167	1939017	39.17	ng	99
74) Di-n-butylphthalate	12.021	149	2483395	44.07	ng	99
75) Fluoranthene	12.674	202	2301398	40.81	ng	98
77) Benzidine	12.786	184	912508	44.25	ng	99
78) Pyrene	12.904	202	2349256	39.34	ng	100
80) Butylbenzylphthalate	13.521	149	1029581	46.48	ng	90
81) Benzo(a)anthracene	14.098	228	1810661	39.96	ng	99
82) 3,3'-Dichlorobenzidine	14.057	252	459403	31.17	ng	100
83) Chrysene	14.133	228	1873993	40.06	ng	99
84) Bis(2-ethylhexyl)phtha...	14.086	149	1363738	46.77	ng	100
85) Di-n-octyl phthalate	14.709	149	2406833	52.22	ng	# 87
87) Indeno(1,2,3-cd)pyrene	17.133	276	1408268	40.34	ng	99
88) Benzo(b)fluoranthene	15.168	252	1633369	42.15	ng	99
89) Benzo(k)fluoranthene	15.204	252	1618315m	44.17	ng	
90) Benzo(a)pyrene	15.545	252	1412515	42.22	ng	99
91) Dibenzo(a,h)anthracene	17.156	278	1182919	39.92	ng	99
92) Benzo(g,h,i)perylene	17.598	276	1139154	38.40	ng	99

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

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