

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031022\
 Data File : BF127289.D
 Acq On : 10 Mar 2022 15:32
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/11/2022
 Supervised By : Jagrut Upadhyay 03/11/2022

Quant Time: Mar 10 15:54:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 10 14:28:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	82412	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	287227	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	162536	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	292128	20.000	ng	# 0.00
76) Chrysene-d12	14.057	240	207243	20.000	ng	# 0.00
86) Perylene-d12	15.539	264	156014	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	546496	98.213	ng	0.00
7) Phenol-d6	6.510	99	755663	101.128	ng	0.00
23) Nitrobenzene-d5	7.445	82	782142	108.922	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	203508	112.562	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	1242885	102.383	ng	0.00
79) Terphenyl-d14	12.992	244	1373235	109.333	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.646	88	154875	55.418	ng	# 82
3) Pyridine	3.410	79	414592	56.599	ng	# 76
4) n-Nitrosodimethylamine	3.399	42	234701	57.988	ng	# 63
6) Aniline	6.545	93	457287	51.611	ng	93
8) 2-Chlorophenol	6.663	128	289223	51.580	ng	93
9) Benzaldehyde	6.428	77	194392	45.155	ng	93
10) Phenol	6.528	94	413903	50.289	ng	# 66
11) bis(2-Chloroethyl)ether	6.616	93	311471	51.736	ng	93
12) 1,3-Dichlorobenzene	6.810	146	322415	51.302	ng	95
13) 1,4-Dichlorobenzene	6.892	146	321795	51.103	ng	99
14) 1,2-Dichlorobenzene	7.045	146	296518	49.409	ng	98
15) Benzyl Alcohol	7.022	79	299582	53.915	ng	90
16) 2,2'-oxybis(1-Chloropr...	7.145	45	576204	50.321	ng	100
17) 2-Methylphenol	7.128	107	245676	51.771	ng	# 89
18) Hexachloroethane	7.381	117	130041	54.087	ng	# 78
19) n-Nitroso-di-n-propyla...	7.292	70	231661	49.798	ng	# 86
20) 3+4-Methylphenols	7.287	107	284917	46.876	ng	# 83
22) Acetophenone	7.287	105	392868	48.780	ng	# 81
24) Nitrobenzene	7.463	77	369448	54.517	ng	92
25) Isophorone	7.698	82	617939	54.064	ng	# 96
26) 2-Nitrophenol	7.775	139	154645	55.736	ng	# 73
27) 2,4-Dimethylphenol	7.810	122	226437	53.986	ng	93
28) bis(2-Chloroethoxy)met...	7.904	93	386488	53.056	ng	97
29) 2,4-Dichlorophenol	8.016	162	261159	54.046	ng	98
30) 1,2,4-Trichlorobenzene	8.098	180	293125	53.251	ng	98
31) Naphthalene	8.181	128	842243	52.772	ng	99
32) Benzoic acid	7.963	122	173289	79.262	ng	88
33) 4-Chloroaniline	8.234	127	320994	51.682	ng	94
34) Hexachlorobutadiene	8.287	225	196680	53.441	ng	98
35) Caprolactam	8.628	113	79388m	56.153	ng	
36) 4-Chloro-3-methylphenol	8.716	107	275412	52.833	ng	97
37) 2-Methylnaphthalene	8.869	142	540447	52.166	ng	97
38) 1-Methylnaphthalene	8.969	142	527466	53.053	ng	100
40) 1,2,4,5-Tetrachloroben...	9.034	216	285264	53.218	ng	# 95
41) Hexachlorocyclopentadiene	9.016	237	127202	87.609	ng	95
43) 2,4,6-Trichlorophenol	9.151	196	190795	57.579	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	199345	56.446	ng	92
46) 1,1'-Biphenyl	9.334	154	685609	53.396	ng	98
47) 2-Chloronaphthalene	9.363	162	537340	54.159	ng	100
48) 2-Nitroaniline	9.463	65	213125	56.638	ng	88
49) Acenaphthylene	9.781	152	799344	52.080	ng	98
50) Dimethylphthalate	9.639	163	643799	53.534	ng	99
51) 2,6-Dinitrotoluene	9.710	165	134598	55.651	ng	95
52) Acenaphthene	9.951	154	474661	51.888	ng	96
53) 3-Nitroaniline	9.881	138	148526	56.199	ng	96
54) 2,4-Dinitrophenol	9.986	184	79327	81.608	ng	95
55) Dibenzofuran	10.122	168	758284	53.135	ng	98
56) 4-Nitrophenol	10.045	139	95487	61.906	ng	# 69
57) 2,4-Dinitrotoluene	10.116	165	180748	56.568	ng	# 90
58) Fluorene	10.469	166	562985	50.598	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.239	232	171993	59.634	ng	# 80
60) Diethylphthalate	10.339	149	595714	53.867	ng	97
61) 4-Chlorophenyl-phenyle...	10.457	204	314573	52.829	ng	# 87
62) 4-Nitroaniline	10.504	138	142036	53.917	ng	# 81
63) Azobenzene	10.616	77	664116	51.895	ng	87
65) 4,6-Dinitro-2-methylph...	10.528	198	111204	62.634	ng	90
66) n-Nitrosodiphenylamine	10.580	169	483357	50.598	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	191755	51.994	ng	99
68) Hexachlorobenzene	11.010	284	212907	53.040	ng	# 91
69) Atrazine	11.104	200	163289	49.212	ng	93
70) Pentachlorophenol	11.210	266	100236	73.190	ng	98
71) Phenanthrene	11.433	178	847864	52.467	ng	99
72) Anthracene	11.486	178	821444	51.183	ng	100
73) Carbazole	11.645	167	773336	51.555	ng	98
74) Di-n-butylphthalate	11.957	149	941867	53.771	ng	99
75) Fluoranthene	12.622	202	926857	51.359	ng	93
77) Benzidine	12.745	184	291552	55.408	ng	96
78) Pyrene	12.857	202	920076	54.572	ng	98
80) Butylbenzylphthalate	13.463	149	359431	56.500	ng	# 96
81) Benzo(a)anthracene	14.045	228	785878	54.524	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	230994	50.205	ng	# 94
83) Chrysene	14.080	228	724234	53.975	ng	98
84) Bis(2-ethylhexyl)phtha...	14.016	149	478285	55.630	ng	# 98
85) Di-n-octyl phthalate	14.627	149	705566	54.748	ng	97
87) Indeno(1,2,3-cd)pyrene	17.051	276	589278	61.716	ng	# 83
88) Benzo(b)fluoranthene	15.104	252	555433	53.057	ng	# 91
89) Benzo(k)fluoranthene	15.133	252	539294	53.599	ng	# 93
90) Benzo(a)pyrene	15.480	252	460767	56.028	ng	# 91
91) Dibenzo(a,h)anthracene	17.068	278	479070	59.854	ng	# 90
92) Benzo(g,h,i)perylene	17.515	276	496488	63.809	ng	# 84

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

