

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120722\
 Data File : BF131545.D
 Acq On : 07 Dec 2022 20:07
 Operator : CG\JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 12/08/2022
 Supervised By : Jagrut Upadhyay 12/08/2022

Quant Time: Dec 08 03:47:53 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF120722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 08 03:39:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.910	152	69070	20.000	ng	0.00
21) Naphthalene-d8	8.204	136	238137	20.000	ng	0.00
39) Acenaphthene-d10	9.969	164	145702	20.000	ng	0.00
64) Phenanthrene-d10	11.457	188	264455	20.000	ng	# 0.00
76) Chrysene-d12	14.110	240	151407	20.000	ng	0.00
86) Perylene-d12	15.621	264	121615	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	392300	96.992	ng	0.00
7) Phenol-d6	6.540	99	524617	103.868	ng	0.00
23) Nitrobenzene-d5	7.487	82	605449	116.688	ng	0.00
42) 2,4,6-Tribromophenol	10.763	330	169787	102.342	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	1136551	102.332	ng	0.00
79) Terphenyl-d14	13.051	244	1156061	115.308	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.675	88	92333	53.328	ng	98
3) Pyridine	3.446	79	246893	51.083	ng	99
4) n-Nitrosodimethylamine	3.422	42	131426	53.062	ng	99
6) Aniline	6.581	93	312117	50.057	ng	98
8) 2-Chlorophenol	6.698	128	208961	48.379	ng	93
9) Benzaldehyde	6.463	77	149816	46.472	ng	98
10) Phenol	6.551	94	292810	51.248	ng	94
11) bis(2-Chloroethyl)ether	6.651	93	219369	51.983	ng	97
12) 1,3-Dichlorobenzene	6.851	146	241299	47.863	ng	98
13) 1,4-Dichlorobenzene	6.928	146	244258	47.406	ng	98
14) 1,2-Dichlorobenzene	7.087	146	233974	48.738	ng	98
15) Benzyl Alcohol	7.057	79	248950	58.929	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.187	45	275608	66.628	ng	99
17) 2-Methylphenol	7.163	107	193666	53.680	ng	99
18) Hexachloroethane	7.428	117	99488	51.973	ng	99
19) n-Nitroso-di-n-propyla...	7.334	70	200831	58.751	ng	98
20) 3+4-Methylphenols	7.322	107	266473	54.617	ng	94
22) Acetophenone	7.328	105	365275	56.527	ng	# 94
24) Nitrobenzene	7.504	77	303198	57.916	ng	96
25) Isophorone	7.745	82	503703	62.409	ng	99
26) 2-Nitrophenol	7.816	139	109968	57.371	ng	96
27) 2,4-Dimethylphenol	7.851	122	175274	55.694	ng	97
28) bis(2-Chloroethoxy)met...	7.951	93	261897	58.532	ng	99
29) 2,4-Dichlorophenol	8.057	162	198725	54.517	ng	97
30) 1,2,4-Trichlorobenzene	8.140	180	244251	53.625	ng	99
31) Naphthalene	8.228	128	646081	49.075	ng	99
32) Benzoic acid	7.987	122	133715	70.455	ng	95
33) 4-Chloroaniline	8.275	127	245100	49.706	ng	99
34) Hexachlorobutadiene	8.339	225	178317	54.813	ng	99
35) Caprolactam	8.669	113	51945m	50.764	ng	
36) 4-Chloro-3-methylphenol	8.757	107	225679	57.916	ng	96
37) 2-Methylnaphthalene	8.922	142	455672	51.723	ng	98
38) 1-Methylnaphthalene	9.022	142	439626	51.932	ng	97
40) 1,2,4,5-Tetrachloroben...	9.086	216	268095	49.810	ng	99
41) Hexachlorocyclopentadiene	9.069	237	142763	58.882	ng	98
43) 2,4,6-Trichlorophenol	9.198	196	169242	52.007	ng	98

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44) 2,4,5-Trichlorophenol	9.239	196	177354	51.103	ng	97
46) 1,1'-Biphenyl	9.386	154	564118	49.101	ng	97
47) 2-Chloronaphthalene	9.416	162	464638	49.509	ng	97
48) 2-Nitroaniline	9.510	65	164882	60.804	ng	92
49) Acenaphthylene	9.834	152	682776	48.704	ng	99
50) Dimethylphthalate	9.692	163	552178	52.434	ng	100
51) 2,6-Dinitrotoluene	9.757	165	113881	50.610	ng	97
52) Acenaphthene	10.004	154	458841m	50.448	ng	
53) 3-Nitroaniline	9.928	138	107053	45.452	ng	92
54) 2,4-Dinitrophenol	10.028	184	64986	57.559	ng	96
55) Dibenzofuran	10.175	168	635662	47.874	ng	98
56) 4-Nitrophenol	10.081	139	78015	47.466	ng	98
57) 2,4-Dinitrotoluene	10.157	165	154086	52.189	ng	87
58) Fluorene	10.522	166	526379	48.041	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.292	232	153819m	52.250	ng	
60) Diethylphthalate	10.392	149	541112	52.115	ng	100
61) 4-Chlorophenyl-phenyle...	10.510	204	295094	52.064	ng	96
62) 4-Nitroaniline	10.545	138	104618	44.550	ng	97
63) Azobenzene	10.675	77	589693	57.595	ng	97
65) 4,6-Dinitro-2-methylph...	10.569	198	90358	62.270	ng	99
66) n-Nitrosodiphenylamine	10.633	169	422123	50.859	ng	99
67) 4-Bromophenyl-phenylether	11.004	248	175860	53.151	ng	93
68) Hexachlorobenzene	11.063	284	184594	50.341	ng	# 91
69) Atrazine	11.157	200	136994	55.820	ng	99
70) Pentachlorophenol	11.257	266	105335	53.649	ng	98
71) Phenanthrene	11.486	178	732417	49.842	ng	99
72) Anthracene	11.539	178	699273	48.402	ng	98
73) Carbazole	11.692	167	571415	46.983	ng	99
74) Di-n-butylphthalate	12.022	149	792581	58.369	ng	99
75) Fluoranthene	12.680	202	750872	46.428	ng	99
77) Benzidine	12.798	184	91672	38.861	ng	99
78) Pyrene	12.910	202	743480	54.825	ng	100
80) Butylbenzylphthalate	13.527	149	255158	69.754	ng	99
81) Benzo(a)anthracene	14.104	228	539432	52.171	ng	99
82) 3,3'-Dichlorobenzidine	14.063	252	149591	53.424	ng	99
83) Chrysene	14.139	228	501867	50.489	ng	100
84) Bis(2-ethylhexyl)phtha...	14.086	149	321448	78.634	ng	# 96
85) Di-n-octyl phthalate	14.704	149	468828	97.536	ng	100
87) Indeno(1,2,3-cd)pyrene	17.162	276	455603	54.530	ng	99
88) Benzo(b)fluoranthene	15.174	252	424796	53.772	ng	100
89) Benzo(k)fluoranthene	15.210	252	406300	47.379	ng	99
90) Benzo(a)pyrene	15.557	252	335562	51.185	ng	100
91) Dibenzo(a,h)anthracene	17.186	278	382636	57.080	ng	99
92) Benzo(g,h,i)perylene	17.633	276	373038	52.527	ng	96

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

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