

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122023\
 Data File : BF136643.D
 Acq On : 20 Dec 2023 10:27
 Operator : CG\JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 12/21/2023
 Supervised By :mohammad ahmed 12/21/2023

Quant Time: Dec 20 17:10:37 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 20 17:02:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.792	152	69657	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	288713	20.000	ng	0.00
39) Acenaphthene-d10	9.827	164	154919	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	315486	20.000	ng	0.00
76) Chrysene-d12	13.968	240	207092	20.000	ng	0.00
86) Perylene-d12	15.445	264	160972	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	95009	21.282	ng	0.00
7) Phenol-d6	6.428	99	120190	20.777	ng	-0.01
23) Nitrobenzene-d5	7.351	82	112134	19.874	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	38567	21.138	ng	0.00
45) 2-Fluorobiphenyl	9.145	172	247780	21.512	ng	0.00
79) Terphenyl-d14	12.910	244	299741	20.227	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.540	88	19462	10.463	ng	96
3) Pyridine	3.310	79	47312	10.425	ng	98
4) n-Nitrosodimethylamine	3.246	42	23401	9.655	ng	# 97
6) Aniline	6.457	93	60918	10.533	ng	97
8) 2-Chlorophenol	6.581	128	50546	10.346	ng	92
9) Benzaldehyde	6.345	77	36985	11.240	ng	99
10) Phenol	6.445	94	64735	10.483	ng	87
11) bis(2-Chloroethyl)ether	6.528	93	48857	10.403	ng	99
12) 1,3-Dichlorobenzene	6.734	146	55138	10.346	ng	96
13) 1,4-Dichlorobenzene	6.810	146	56788	10.428	ng	96
14) 1,2-Dichlorobenzene	6.957	146	53403	10.539	ng	96
15) Benzyl Alcohol	6.934	79	40248	10.313	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.063	45	79144	10.308	ng	98
17) 2-Methylphenol	7.045	107	41733	10.311	ng	97
18) Hexachloroethane	7.298	117	20226	10.227	ng	98
19) n-Nitroso-di-n-propyla...	7.192	70	36594	10.436	ng	97
20) 3+4-Methylphenols	7.198	107	54198	10.719	ng	98
22) Acetophenone	7.198	105	72848	10.067	ng	# 93
24) Nitrobenzene	7.369	77	56147	9.934	ng	97
25) Isophorone	7.604	82	102784	10.012	ng	99
26) 2-Nitrophenol	7.686	139	25687	9.499	ng	97
27) 2,4-Dimethylphenol	7.728	122	32538	9.884	ng	91
28) bis(2-Chloroethoxy)met...	7.816	93	62523	10.019	ng	99
29) 2,4-Dichlorophenol	7.928	162	42456	10.017	ng	98
30) 1,2,4-Trichlorobenzene	8.010	180	47347	10.026	ng	99
31) Naphthalene	8.086	128	160315	10.107	ng	99
32) Benzoic acid	7.810	122	26030	8.233	ng	96
33) 4-Chloroaniline	8.139	127	59145	10.100	ng	98
34) Hexachlorobutadiene	8.204	225	28933	10.084	ng	98
35) Caprolactam	8.480	113	14361	10.268	ng	96
36) 4-Chloro-3-methylphenol	8.622	107	47900	10.039	ng	98
37) 2-Methylnaphthalene	8.780	142	106159	10.400	ng	100
38) 1-Methylnaphthalene	8.880	142	103304	10.315	ng	99
40) 1,2,4,5-Tetrachloroben...	8.945	216	49072	10.228	ng	98
41) Hexachlorocyclopentadiene	8.933	237	8610	10.182	ng	94
43) 2,4,6-Trichlorophenol	9.063	196	30750	9.994	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.104	196	35414	10.328	ng	99
46) 1,1'-Biphenyl	9.245	154	136598	10.507	ng	100
47) 2-Chloronaphthalene	9.269	162	101624	10.283	ng	99
48) 2-Nitroaniline	9.369	65	31728	10.252	ng	92
49) Acenaphthylene	9.686	152	158935	10.521	ng	99
50) Dimethylphthalate	9.545	163	117479	10.382	ng	99
51) 2,6-Dinitrotoluene	9.610	165	25852	10.067	ng	93
52) Acenaphthene	9.857	154	99726	10.650	ng	97
53) 3-Nitroaniline	9.780	138	28199	10.366	ng	91
54) 2,4-Dinitrophenol	9.886	184	9207	9.902	ng	# 86
55) Dibenzofuran	10.027	168	151972	10.707	ng	99
56) 4-Nitrophenol	9.951	139	18101	9.857	ng	97
57) 2,4-Dinitrotoluene	10.016	165	35563	10.667	ng	# 85
58) Fluorene	10.374	166	123850	10.997	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.151	232	26667m	10.033	ng	
60) Diethylphthalate	10.245	149	125203	10.664	ng	99
61) 4-Chlorophenyl-phenyle...	10.369	204	59324	10.837	ng	97
62) 4-Nitroaniline	10.386	138	29147	10.846	ng	95
63) Azobenzene	10.527	77	115625	10.544	ng	98
65) 4,6-Dinitro-2-methylph...	10.422	198	15638	8.649	ng	97
66) n-Nitrosodiphenylamine	10.486	169	105105	10.256	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	35661	10.178	ng	98
68) Hexachlorobenzene	10.922	284	40193	10.294	ng	98
69) Atrazine	11.016	200	29722	11.339	ng	99
70) Pentachlorophenol	11.121	266	17544	9.640	ng	97
71) Phenanthrene	11.339	178	171865	10.617	ng	100
72) Anthracene	11.392	178	171611	10.468	ng	99
73) Carbazole	11.551	167	165178	10.687	ng	100
74) Di-n-butylphthalate	11.880	149	202133	10.717	ng	100
75) Fluoranthene	12.533	202	191976	11.087	ng	99
77) Benzidine	12.657	184	61659	10.098	ng	98
78) Pyrene	12.763	202	195313	9.607	ng	99
80) Butylbenzylphthalate	13.386	149	70222	9.493	ng	94
81) Benzo(a)anthracene	13.957	228	146955	10.219	ng	100
82) 3,3'-Dichlorobenzidine	13.921	252	41088	10.061	ng	# 98
83) Chrysene	13.992	228	143268	10.188	ng	99
84) Bis(2-ethylhexyl)phtha...	13.951	149	89950	9.665	ng	100
85) Di-n-octyl phthalate	14.568	149	131613	9.142	ng	99
87) Indeno(1,2,3-cd)pyrene	16.939	276	105637	9.089	ng	98
88) Benzo(b)fluoranthene	15.009	252	107910	10.037	ng	99
89) Benzo(k)fluoranthene	15.039	252	110311	10.862	ng	100
90) Benzo(a)pyrene	15.380	252	91108	10.039	ng	99
91) Dibenzo(a,h)anthracene	16.956	278	87889	9.218	ng	99
92) Benzo(g,h,i)perylene	17.392	276	87519	8.962	ng	99

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

