

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072423\
 Data File : BF134429.D
 Acq On : 24 Jul 2023 14:53
 Operator : CG\JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/25/2023
 Supervised By :mohammad ahmed 07/25/2023

Quant Time: Jul 25 01:32:26 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072423.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 25 01:30:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	50610	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	188067	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	92918	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	181268	20.000	ng	0.00
76) Chrysene-d12	14.027	240	156648	20.000	ng	0.00
86) Perylene-d12	15.533	264	147998	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	62321	20.345	ng	0.00
7) Phenol-d6	6.469	99	79436	21.166	ng	-0.01
23) Nitrobenzene-d5	7.392	82	74456	20.030	ng	-0.01
42) 2,4,6-Tribromophenol	10.663	330	25469	19.292	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	148116	20.927	ng	0.00
79) Terphenyl-d14	12.957	244	175291	19.534	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.593	88	13081	10.068	ng	95
3) Pyridine	3.404	79	28555m	9.807	ng	
4) n-Nitrosodimethylamine	3.334	42	16180m	9.387	ng	
6) Aniline	6.498	93	47148	10.659	ng	91
8) 2-Chlorophenol	6.622	128	32417	10.446	ng	98
9) Benzaldehyde	6.392	77	23560	11.722	ng	92
10) Phenol	6.487	94	39697	10.166	ng	91
11) bis(2-Chloroethyl)ether	6.569	93	28084	10.302	ng	91
12) 1,3-Dichlorobenzene	6.775	146	34234	10.259	ng	99
13) 1,4-Dichlorobenzene	6.851	146	35417	10.313	ng	99
14) 1,2-Dichlorobenzene	7.004	146	33882	10.503	ng	98
15) Benzyl Alcohol	6.975	79	28384m	9.892	ng	
16) 2,2'-oxybis(1-Chloropr...	7.110	45	42826	10.668	ng	99
17) 2-Methylphenol	7.092	107	28141	10.168	ng	97
18) Hexachloroethane	7.339	117	14139	10.717	ng	94
19) n-Nitroso-di-n-propyla...	7.239	70	24433	10.458	ng	97
20) 3+4-Methylphenols	7.245	107	38425	10.844	ng	92
22) Acetophenone	7.239	105	49129	10.504	ng	91
24) Nitrobenzene	7.416	77	36702	9.928	ng	96
25) Isophorone	7.651	82	63024	10.173	ng	97
26) 2-Nitrophenol	7.734	139	16963	9.977	ng	98
27) 2,4-Dimethylphenol	7.769	122	27579	10.278	ng	94
28) bis(2-Chloroethoxy)met...	7.863	93	34833	10.102	ng	100
29) 2,4-Dichlorophenol	7.981	162	27560	10.275	ng	94
30) 1,2,4-Trichlorobenzene	8.057	180	29839	10.313	ng	96
31) Naphthalene	8.133	128	97997	10.470	ng	99
32) Benzoic acid	7.857	122	15774m	8.903	ng	
33) 4-Chloroaniline	8.186	127	39253	10.146	ng	94
34) Hexachlorobutadiene	8.245	225	20206	10.425	ng	97
35) Caprolactam	8.533	113	8605	10.512	ng	# 90
36) 4-Chloro-3-methylphenol	8.675	107	31575	10.202	ng	98
37) 2-Methylnaphthalene	8.828	142	66817	10.274	ng	100
38) 1-Methylnaphthalene	8.928	142	60604	10.045	ng	96
40) 1,2,4,5-Tetrachloroben...	8.992	216	30709	10.429	ng	100
41) Hexachlorocyclopentadiene	8.975	237	2796	10.522	ng	93
43) 2,4,6-Trichlorophenol	9.110	196	19015	10.066	ng	98

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44) 2,4,5-Trichlorophenol	9.157	196	20999	9.541	ng	90
46) 1,1'-Biphenyl	9.292	154	77991	10.184	ng	97
47) 2-Chloronaphthalene	9.316	162	58001	10.273	ng	98
48) 2-Nitroaniline	9.422	65	19410	9.546	ng	94
49) Acenaphthylene	9.733	152	98204	10.296	ng	96
50) Dimethylphthalate	9.586	163	69138	10.037	ng	98
51) 2,6-Dinitrotoluene	9.657	165	14700	9.726	ng	100
52) Acenaphthene	9.904	154	55738	9.892	ng	97
53) 3-Nitroaniline	9.833	138	17407	10.045	ng	# 97
54) 2,4-Dinitrophenol	9.957	184	4115	11.333	ng	# 82
55) Dibenzofuran	10.080	168	86959	10.028	ng	98
56) 4-Nitrophenol	10.016	139	7121	7.706	ng	# 84
57) 2,4-Dinitrotoluene	10.069	165	19626	9.668	ng	95
58) Fluorene	10.422	166	71096	10.162	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.204	232	16213m	9.552	ng	
60) Diethylphthalate	10.292	149	71766	10.063	ng	98
61) 4-Chlorophenyl-phenyle...	10.416	204	33175	9.927	ng	96
62) 4-Nitroaniline	10.445	138	17192	9.674	ng	86
63) Azobenzene	10.575	77	64868	9.882	ng	98
65) 4,6-Dinitro-2-methylph...	10.480	198	8586	8.257	ng	94
66) n-Nitrosodiphenylamine	10.533	169	59573	10.330	ng	96
67) 4-Bromophenyl-phenylether	10.904	248	20065	10.103	ng	# 89
68) Hexachlorobenzene	10.975	284	23486	10.501	ng	93
69) Atrazine	11.063	200	17231	10.716	ng	97
70) Pentachlorophenol	11.180	266	8149m	8.682	ng	
71) Phenanthrene	11.392	178	100027	10.561	ng	99
72) Anthracene	11.445	178	102739	10.510	ng	97
73) Carbazole	11.604	167	95125	10.334	ng	98
74) Di-n-butylphthalate	11.927	149	108468	10.265	ng	99
75) Fluoranthene	12.586	202	114301	10.355	ng	99
77) Benzidine	12.716	184	38741	11.583	ng	95
78) Pyrene	12.821	202	118659	9.527	ng	99
80) Butylbenzylphthalate	13.439	149	47165	9.453	ng	95
81) Benzo(a)anthracene	14.015	228	110076	10.072	ng	97
82) 3,3'-Dichlorobenzidine	13.980	252	35952	10.386	ng	97
83) Chrysene	14.057	228	104527	9.999	ng	97
84) Bis(2-ethylhexyl)phtha...	13.998	149	57616	9.505	ng	# 99
85) Di-n-octyl phthalate	14.621	149	84313	10.009	ng	# 98
87) Indeno(1,2,3-cd)pyrene	17.080	276	82469	8.529	ng	99
88) Benzo(b)fluoranthene	15.086	252	99156	10.709	ng	100
89) Benzo(k)fluoranthene	15.115	252	91763	10.075	ng	98
90) Benzo(a)pyrene	15.468	252	86692	10.130	ng	100
91) Dibenzo(a,h)anthracene	17.086	278	64285	8.233	ng	95
92) Benzo(g,h,i)perylene	17.545	276	65626	8.185	ng	98

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

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