

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101323\
 Data File : BF135753.D
 Acq On : 13 Oct 2023 14:55
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 10/16/2023
 Supervised By :mohammad ahmed 10/17/2023

Quant Time: Oct 13 17:06:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 13 16:43:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.740	152	120605	20.000	ng	0.00
21) Naphthalene-d8	8.022	136	469032	20.000	ng	0.00
39) Acenaphthene-d10	9.781	164	241232	20.000	ng	0.00
64) Phenanthrene-d10	11.275	188	457818	20.000	ng	0.00
76) Chrysene-d12	13.945	240	281967	20.000	ng	0.00
86) Perylene-d12	15.415	264	280273	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	874230	116.541	ng	0.00
7) Phenol-d6	6.404	99	1043696	111.509	ng	0.00
23) Nitrobenzene-d5	7.322	82	1015484	119.077	ng	0.00
42) 2,4,6-Tribromophenol	10.580	330	260065	112.348	ng	0.00
45) 2-Fluorobiphenyl	9.104	172	1593586	103.286	ng	0.00
79) Terphenyl-d14	12.874	244	1727360	113.450	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.434	88	214628	59.317	ng	99
3) Pyridine	3.181	79	550699	61.771	ng	96
4) n-Nitrosodimethylamine	3.169	42	304100	63.696	ng	97
6) Aniline	6.416	93	660424	56.069	ng	# 59
8) 2-Chlorophenol	6.534	128	464860	57.466	ng	97
9) Benzaldehyde	6.304	77	266647	50.053	ng	99
10) Phenol	6.416	94	552638	56.349	ng	98
11) bis(2-Chloroethyl)ether	6.487	93	466093	58.982	ng	98
12) 1,3-Dichlorobenzene	6.681	146	464946	56.694	ng	99
13) 1,4-Dichlorobenzene	6.757	146	466981	56.210	ng	99
14) 1,2-Dichlorobenzene	6.916	146	409378	54.727	ng	98
15) Benzyl Alcohol	6.904	79	348330	55.971	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.022	45	794407	55.585	ng	96
17) 2-Methylphenol	7.016	107	394904	56.392	ng	97
18) Hexachloroethane	7.245	117	174974	58.088	ng	94
19) n-Nitroso-di-n-propyla...	7.169	70	309303	53.446	ng	98
20) 3+4-Methylphenols	7.175	107	474912	54.399	ng	92
22) Acetophenone	7.163	105	602998	56.057	ng	95
24) Nitrobenzene	7.340	77	508071	59.362	ng	98
25) Isophorone	7.575	82	962426	59.820	ng	98
26) 2-Nitrophenol	7.645	139	255001	61.051	ng	95
27) 2,4-Dimethylphenol	7.692	122	401504	58.391	ng	98
28) bis(2-Chloroethoxy)met...	7.781	93	557134	58.148	ng	99
29) 2,4-Dichlorophenol	7.892	162	375429	58.262	ng	97
30) 1,2,4-Trichlorobenzene	7.969	180	369666	56.090	ng	99
31) Naphthalene	8.045	128	1272098	55.605	ng	99
32) Benzoic acid	7.869	122	319394m	68.796	ng	
33) 4-Chloroaniline	8.104	127	554276	55.464	ng	99
34) Hexachlorobutadiene	8.157	225	220111	56.424	ng	98
35) Caprolactam	8.510	113	127693m	57.903	ng	
36) 4-Chloro-3-methylphenol	8.604	107	415302	58.117	ng	98
37) 2-Methylnaphthalene	8.739	142	833610	55.306	ng	99
38) 1-Methylnaphthalene	8.839	142	775960	55.816	ng	99
40) 1,2,4,5-Tetrachloroben...	8.904	216	386503	55.699	ng	100
41) Hexachlorocyclopentadiene	8.881	237	68917	61.218	ng	98
43) 2,4,6-Trichlorophenol	9.028	196	256806	57.430	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.075	196	299335	57.554	ng	99
46) 1,1'-Biphenyl	9.204	154	1008747	54.590	ng	97
47) 2-Chloronaphthalene	9.234	162	754484	56.653	ng	98
48) 2-Nitroaniline	9.339	65	314217	59.774	ng	96
49) Acenaphthylene	9.645	152	1203992	54.630	ng	99
50) Dimethylphthalate	9.516	163	938157	57.610	ng	99
51) 2,6-Dinitrotoluene	9.586	165	203629	58.147	ng	97
52) Acenaphthene	9.816	154	728238	55.129	ng	99
53) 3-Nitroaniline	9.763	138	251083	58.488	ng	97
54) 2,4-Dinitrophenol	9.881	184	104456	71.456	ng	97
55) Dibenzofuran	9.992	168	1069770	53.671	ng	99
56) 4-Nitrophenol	9.945	139	114675	59.791	ng	97
57) 2,4-Dinitrotoluene	9.998	165	253625	56.927	ng	# 86
58) Fluorene	10.339	166	846694	54.261	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.116	232	229860	58.481	ng	98
60) Diethylphthalate	10.216	149	925268	55.655	ng	99
61) 4-Chlorophenyl-phenyle...	10.328	204	407986	54.098	ng	93
62) 4-Nitroaniline	10.386	138	236143	59.786	ng	99
63) Azobenzene	10.486	77	923997	57.206	ng	99
65) 4,6-Dinitro-2-methylph...	10.416	198	158989	69.005	ng	95
66) n-Nitrosodiphenylamine	10.451	169	786269	56.347	ng	99
67) 4-Bromophenyl-phenylether	10.816	248	265472	57.765	ng	97
68) Hexachlorobenzene	10.886	284	266911	57.045	ng	99
69) Atrazine	10.986	200	210614	55.490	ng	98
70) Pentachlorophenol	11.092	266	127607	70.133	ng	97
71) Phenanthrene	11.304	178	1181076	53.853	ng	99
72) Anthracene	11.357	178	1228883	54.073	ng	99
73) Carbazole	11.522	167	1142055	53.882	ng	99
74) Di-n-butylphthalate	11.845	149	1393785	53.559	ng	100
75) Fluoranthene	12.504	202	1300357	53.422	ng	97
77) Benzidine	12.627	184	286475	46.560	ng	100
78) Pyrene	12.733	202	1331533	59.426	ng	99
80) Butylbenzylphthalate	13.351	149	582078	59.117	ng	99
81) Benzo(a)anthracene	13.933	228	1068580	57.696	ng	99
82) 3,3'-Dichlorobenzidine	13.898	252	343854	55.980	ng	# 97
83) Chrysene	13.974	228	1023244	57.041	ng	100
84) Bis(2-ethylhexyl)phtha...	13.916	149	776480	58.185	ng	# 97
85) Di-n-octyl phthalate	14.533	149	1199384	56.678	ng	100
87) Indeno(1,2,3-cd)pyrene	16.927	276	963965	63.122	ng	100
88) Benzo(b)fluoranthene	14.992	252	905124	57.703	ng	98
89) Benzo(k)fluoranthene	15.015	252	829756	54.486	ng	99
90) Benzo(a)pyrene	15.357	252	849547	57.596	ng	99
91) Dibenzo(a,h)anthracene	16.927	278	785651	62.756	ng	99
92) Benzo(g,h,i)perylene	17.374	276	829344	64.491	ng	99

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

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