

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073024\
 Data File : BF138682.D
 Acq On : 30 Jul 2024 13:56
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/31/2024
 Supervised By :mohammad ahmed 07/31/2024

Quant Time: Jul 30 17:42:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 17:38:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	73349	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	297904	20.000	ng	0.00
39) Acenaphthene-d10	9.880	164	162245	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	269225	20.000	ng	0.00
76) Chrysene-d12	14.004	240	146258	20.000	ng	0.00
86) Perylene-d12	15.468	264	137026	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.469	112	100297	21.108	ng	0.00
7) Phenol-d6	6.475	99	133232	20.884	ng	-0.01
23) Nitrobenzene-d5	7.404	82	123997	20.350	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	26595	20.011	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	231000	21.392	ng	0.00
79) Terphenyl-d14	12.945	244	185079	21.187	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.593	88	20339	9.777	ng	95
3) Pyridine	3.357	79	50057	9.933	ng	98
4) n-Nitrosodimethylamine	3.287	42	29962	9.983	ng	# 98
6) Aniline	6.504	93	59629	10.481	ng	94
8) 2-Chlorophenol	6.628	128	51997	10.401	ng	97
9) Benzaldehyde	6.398	77	41047	10.733	ng	99
10) Phenol	6.492	94	71371	10.626	ng	92
11) bis(2-Chloroethyl)ether	6.575	93	54121	10.471	ng	99
12) 1,3-Dichlorobenzene	6.787	146	59006	10.544	ng	98
13) 1,4-Dichlorobenzene	6.863	146	59226	10.487	ng	97
14) 1,2-Dichlorobenzene	7.016	146	55425	10.501	ng	98
15) Benzyl Alcohol	6.987	79	47206	10.267	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.116	45	94469	10.620	ng	97
17) 2-Methylphenol	7.098	107	42323	10.252	ng	97
18) Hexachloroethane	7.351	117	22530	10.598	ng	97
19) n-Nitroso-di-n-propyla...	7.245	70	39862	10.345	ng	99
20) 3+4-Methylphenols	7.251	107	56989	10.760	ng	97
22) Acetophenone	7.251	105	75892	10.404	ng	99
24) Nitrobenzene	7.422	77	63020	10.164	ng	98
25) Isophorone	7.663	82	105784	10.167	ng	99
26) 2-Nitrophenol	7.739	139	26149	9.803	ng	96
27) 2,4-Dimethylphenol	7.781	122	32386	10.147	ng	98
28) bis(2-Chloroethoxy)met...	7.875	93	64569	10.191	ng	99
29) 2,4-Dichlorophenol	7.987	162	41138	10.031	ng	99
30) 1,2,4-Trichlorobenzene	8.063	180	48272	10.199	ng	98
31) Naphthalene	8.145	128	164167	10.469	ng	99
32) Benzoic acid	7.875	122	20318m	8.102	ng	
33) 4-Chloroaniline	8.198	127	54003	10.260	ng	98
34) Hexachlorobutadiene	8.263	225	29309	10.224	ng	98
35) Caprolactam	8.539	113	12474	10.193	ng	93
36) 4-Chloro-3-methylphenol	8.681	107	47584	10.152	ng	96
37) 2-Methylnaphthalene	8.839	142	104097	10.511	ng	99
38) 1-Methylnaphthalene	8.934	142	102100	10.521	ng	100
40) 1,2,4,5-Tetrachloroben...	9.004	216	45439	10.082	ng	98
41) Hexachlorocyclopentadiene	8.986	237	5161	10.915	ng	96
43) 2,4,6-Trichlorophenol	9.116	196	27439	9.985	ng	98

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44) 2,4,5-Trichlorophenol	9.163	196	30782	10.247	ng	97
46) 1,1'-Biphenyl	9.298	154	132668	10.441	ng	99
47) 2-Chloronaphthalene	9.328	162	97850	10.354	ng	97
48) 2-Nitroaniline	9.422	65	32731	10.216	ng	97
49) Acenaphthylene	9.739	152	141438	10.552	ng	99
50) Dimethylphthalate	9.598	163	105187	10.139	ng	100
51) 2,6-Dinitrotoluene	9.663	165	23351	9.974	ng	95
52) Acenaphthene	9.910	154	95143	10.560	ng	98
53) 3-Nitroaniline	9.833	138	24353	10.062	ng	99
54) 2,4-Dinitrophenol	9.951	184	7908	7.337	ng	88
55) Dibenzofuran	10.080	168	136302	10.717	ng	99
56) 4-Nitrophenol	10.004	139	12643	8.687	ng	95
57) 2,4-Dinitrotoluene	10.069	165	30462	10.198	ng	99
58) Fluorene	10.428	166	107585	10.622	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.204	232	22672	9.872	ng	# 94
60) Diethylphthalate	10.298	149	102108	10.381	ng	99
61) 4-Chlorophenyl-phenyle...	10.416	204	53870	10.814	ng	99
62) 4-Nitroaniline	10.439	138	23653	10.284	ng	98
63) Azobenzene	10.575	77	112880	10.347	ng	99
65) 4,6-Dinitro-2-methylph...	10.475	198	14344	8.733	ng	96
66) n-Nitrosodiphenylamine	10.533	169	86813	10.316	ng	100
67) 4-Bromophenyl-phenylether	10.910	248	29510	10.124	ng	98
68) Hexachlorobenzene	10.975	284	31049	10.317	ng	96
69) Atrazine	11.063	200	23189	10.680	ng	96
70) Pentachlorophenol	11.175	266	10392	7.661	ng	93
71) Phenanthrene	11.392	178	149253	10.766	ng	99
72) Anthracene	11.439	178	145043	10.620	ng	99
73) Carbazole	11.598	167	126321	10.721	ng	100
74) Di-n-butylphthalate	11.922	149	133246	10.060	ng	99
75) Fluoranthene	12.580	202	140690	10.871	ng	99
77) Benzidine	12.704	184	35263	10.080	ng	100
78) Pyrene	12.810	202	143221	10.400	ng	99
80) Butylbenzylphthalate	13.421	149	41771	9.472	ng	99
81) Benzo(a)anthracene	13.998	228	105185	10.444	ng	97
82) 3,3'-Dichlorobenzidine	13.957	252	26584	10.314	ng	98
83) Chrysene	14.033	228	89780	9.881	ng	99
84) Bis(2-ethylhexyl)phtha...	13.980	149	61314	9.495	ng	98
85) Di-n-octyl phthalate	14.592	149	111460	9.329	ng	100
87) Indeno(1,2,3-cd)pyrene	16.933	276	100310	10.215	ng	99
88) Benzo(b)fluoranthene	15.039	252	93263	10.980	ng	99
89) Benzo(k)fluoranthene	15.068	252	71464	9.717	ng	99
90) Benzo(a)pyrene	15.404	252	71724	10.038	ng	98
91) Dibenzo(a,h)anthracene	16.945	278	84356	10.465	ng	97
92) Benzo(g,h,i)perylene	17.374	276	84564	10.110	ng	98

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

