

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122624\
 Data File : BF140972.D
 Acq On : 26 Dec 2024 17:15
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: Dec 27 00:19:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 27 00:11:18 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	251125	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	1005088	20.000	ng	0.00
39) Acenaphthene-d10	9.857	164	539677	20.000	ng	0.00
64) Phenanthrene-d10	11.345	188	960019	20.000	ng	0.00
76) Chrysene-d12	13.986	240	769052	20.000	ng	0.00
86) Perylene-d12	15.439	264	725753	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	357171	21.439	ng	0.00
7) Phenol-d6	6.440	99	461726	21.514	ng	-0.01
23) Nitrobenzene-d5	7.387	82	277918	17.691	ng	0.00
42) 2,4,6-Tribromophenol	10.645	330	102029	19.456	ng	0.00
45) 2-Fluorobiphenyl	9.181	172	783402	23.128	ng	0.00
79) Terphenyl-d14	12.927	244	908253	19.616	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.563	88	77981	10.339	ng	99
3) Pyridine	3.305	79	196202	10.644	ng	98
4) n-Nitrosodimethylamine	3.234	42	92623	9.918	ng	# 99
6) Aniline	6.487	93	213768	11.168	ng	98
8) 2-Chlorophenol	6.604	128	186317	10.608	ng	99
9) Benzaldehyde	6.375	77	149809	7.569	ng	98
10) Phenol	6.451	94	236093	10.639	ng	97
11) bis(2-Chloroethyl)ether	6.557	93	185749	10.376	ng	100
12) 1,3-Dichlorobenzene	6.769	146	209098	10.787	ng	99
13) 1,4-Dichlorobenzene	6.845	146	212166	10.857	ng	98
14) 1,2-Dichlorobenzene	6.998	146	199494	10.949	ng	99
15) Benzyl Alcohol	6.957	79	160122	10.383	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.104	45	290964	10.738	ng	99
17) 2-Methylphenol	7.069	107	151661	10.393	ng	98
18) Hexachloroethane	7.340	117	71922	10.361	ng	97
19) n-Nitroso-di-n-propyla...	7.234	70	141506	10.763	ng	98
20) 3+4-Methylphenols	7.222	107	203866	11.040	ng	91
22) Acetophenone	7.228	105	268709	10.898	ng	98
24) Nitrobenzene	7.404	77	160105	9.121	ng	100
25) Isophorone	7.640	82	354894	10.381	ng	97
26) 2-Nitrophenol	7.722	139	39980	10.245	ng	99
27) 2,4-Dimethylphenol	7.751	122	128382	10.276	ng	98
28) bis(2-Chloroethoxy)met...	7.857	93	232407	10.730	ng	99
29) 2,4-Dichlorophenol	7.957	162	145068	10.238	ng	100
30) 1,2,4-Trichlorobenzene	8.051	180	168369	10.498	ng	99
31) Naphthalene	8.128	128	579991	10.951	ng	98
32) Benzoic acid	7.810	122	49591	11.743	ng	99
33) 4-Chloroaniline	8.175	127	199889	10.433	ng	99
34) Hexachlorobutadiene	8.251	225	100026	10.579	ng	100
35) Caprolactam	8.504	113	47868	9.936	ng	98
36) 4-Chloro-3-methylphenol	8.639	107	160989	10.230	ng	100
37) 2-Methylnaphthalene	8.816	142	372101	10.982	ng	99
38) 1-Methylnaphthalene	8.916	142	363059	10.903	ng	100
40) 1,2,4,5-Tetrachloroben...	8.981	216	158150	10.652	ng	99
41) Hexachlorocyclopentadiene	8.975	237	47532	9.661	ng	97
43) 2,4,6-Trichlorophenol	9.087	196	96190	9.734	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.122	196	104859	10.371	ng	99
46) 1,1'-Biphenyl	9.281	154	457591	11.074	ng	99
47) 2-Chloronaphthalene	9.304	162	348834	10.884	ng	98
48) 2-Nitroaniline	9.392	65	49140	9.938	ng	99
49) Acenaphthylene	9.722	152	504423	10.933	ng	99
50) Dimethylphthalate	9.575	163	381476	10.746	ng	99
51) 2,6-Dinitrotoluene	9.639	165	46772	9.624	ng	97
52) Acenaphthene	9.892	154	334623	10.729	ng	100
53) 3-Nitroaniline	9.804	138	53244	9.568	ng	# 93
54) 2,4-Dinitrophenol	9.904	184	10078	11.150	ng	# 1
55) Dibenzofuran	10.063	168	480351	10.959	ng	97
56) 4-Nitrophenol	9.945	139	44755	7.370	ng	94
57) 2,4-Dinitrotoluene	10.034	165	50385	9.654	ng	90
58) Fluorene	10.404	166	386427	11.369	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.175	232	81766	9.765	ng	98
60) Diethylphthalate	10.281	149	374095	10.675	ng	99
61) 4-Chlorophenyl-phenyle...	10.404	204	181693	11.109	ng	97
62) 4-Nitroaniline	10.410	138	54741	9.821	ng	91
63) Azobenzene	10.557	77	395397	10.749	ng	99
65) 4,6-Dinitro-2-methylph...	10.439	198	15689	12.644	ng	91
66) n-Nitrosodiphenylamine	10.516	169	324220	10.725	ng	99
67) 4-Bromophenyl-phenylether	10.892	248	110052	10.604	ng	98
68) Hexachlorobenzene	10.951	284	118361	10.343	ng	97
69) Atrazine	11.039	200	98299	11.639	ng	99
70) Pentachlorophenol	11.139	266	65934	9.254	ng	99
71) Phenanthrene	11.369	178	560414	11.091	ng	98
72) Anthracene	11.416	178	545401	11.016	ng	98
73) Carbazole	11.569	167	524879	11.028	ng	99
74) Di-n-butylphthalate	11.910	149	594457	10.981	ng	99
75) Fluoranthene	12.557	202	611729	11.163	ng	98
77) Benzidine	12.680	184	127969	7.873	ng	99
78) Pyrene	12.780	202	642229	9.549	ng	98
80) Butylbenzylphthalate	13.410	149	219532	8.939	ng	97
81) Benzo(a)anthracene	13.974	228	558462	10.419	ng	99
82) 3,3'-Dichlorobenzidine	13.939	252	174840	10.857	ng	97
83) Chrysene	14.010	228	511519	10.587	ng	99
84) Bis(2-ethylhexyl)phtha...	13.974	149	323647	10.159	ng	100
85) Di-n-octyl phthalate	14.592	149	473516	10.269	ng	100
87) Indeno(1,2,3-cd)pyrene	16.880	276	365994	8.004	ng	100
88) Benzo(b)fluoranthene	15.016	252	567548	11.213	ng	99
89) Benzo(k)fluoranthene	15.045	252	406884	10.256	ng	99
90) Benzo(a)pyrene	15.374	252	401924	10.230	ng	99
91) Dibenzo(a,h)anthracene	16.904	278	300782	8.115	ng	99
92) Benzo(g,h,i)perylene	17.315	276	294254	7.654	ng	100

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

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