

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090823\
 Data File : BF135150.D
 Acq On : 08 Sep 2023 16:55
 Operator : CG\JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Quant Time: Sep 09 02:26:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 09 02:14:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.763	152	111022	20.000	ng	0.00
21) Naphthalene-d8	8.045	136	453697	20.000	ng	0.00
39) Acenaphthene-d10	9.798	164	240575	20.000	ng	0.00
64) Phenanthrene-d10	11.292	188	445038	20.000	ng	0.00
76) Chrysene-d12	13.945	240	268569	20.000	ng	0.00
86) Perylene-d12	15.404	264	195391	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	161052	22.449	ng	0.00
7) Phenol-d6	6.398	99	202509	22.433	ng	-0.01
23) Nitrobenzene-d5	7.322	82	178477	21.413	ng	-0.01
42) 2,4,6-Tribromophenol	10.592	330	58994	21.806	ng	0.00
45) 2-Fluorobiphenyl	9.122	172	351328	22.932	ng	0.00
79) Terphenyl-d14	12.886	244	381238	21.858	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.475	88	32408	10.902	ng	97
3) Pyridine	3.234	79	83588	10.603	ng	96
4) n-Nitrosodimethylamine	3.169	42	40371	10.229	ng	100
6) Aniline	6.428	93	123508	11.138	ng	99
8) 2-Chlorophenol	6.545	128	82269	11.111	ng	99
9) Benzaldehyde	6.316	77	60121	11.877	ng	99
10) Phenol	6.410	94	108086	11.444	ng	85
11) bis(2-Chloroethyl)ether	6.504	93	76747	10.973	ng	94
12) 1,3-Dichlorobenzene	6.704	146	83348	11.015	ng	96
13) 1,4-Dichlorobenzene	6.781	146	82991	10.919	ng	97
14) 1,2-Dichlorobenzene	6.934	146	80271	11.282	ng	95
15) Benzyl Alcohol	6.904	79	72404	11.305	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.039	45	126307	11.119	ng	99
17) 2-Methylphenol	7.016	107	72215	11.060	ng	98
18) Hexachloroethane	7.269	117	29322	10.525	ng	95
19) n-Nitroso-di-n-propyla...	7.169	70	59763	10.977	ng	96
20) 3+4-Methylphenols	7.175	107	91480	11.628	ng	97
22) Acetophenone	7.169	105	115510	11.038	ng	# 97
24) Nitrobenzene	7.345	77	91156	10.768	ng	97
25) Isophorone	7.581	82	164913	10.594	ng	99
26) 2-Nitrophenol	7.663	139	42067	10.181	ng	92
27) 2,4-Dimethylphenol	7.698	122	69917	10.711	ng	99
28) bis(2-Chloroethoxy)met...	7.798	93	97167	10.678	ng	98
29) 2,4-Dichlorophenol	7.904	162	67630	10.622	ng	94
30) 1,2,4-Trichlorobenzene	7.986	180	75494	10.874	ng	97
31) Naphthalene	8.063	128	245924	10.943	ng	99
32) Benzoic acid	7.786	122	47548	9.445	ng	90
33) 4-Chloroaniline	8.116	127	103960	10.620	ng	98
34) Hexachlorobutadiene	8.181	225	43660	10.927	ng	99
35) Caprolactam	8.463	113	20916	10.152	ng	97
36) 4-Chloro-3-methylphenol	8.598	107	76809	10.743	ng	96
37) 2-Methylnaphthalene	8.757	142	171901	11.275	ng	98
38) 1-Methylnaphthalene	8.857	142	157333	11.141	ng	99
40) 1,2,4,5-Tetrachloroben...	8.922	216	74423	10.905	ng	99
41) Hexachlorocyclopentadiene	8.910	237	15288	11.658	ng	99
43) 2,4,6-Trichlorophenol	9.033	196	48810	10.779	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.075	196	55333	10.561	ng	99
46) 1,1'-Biphenyl	9.222	154	208379	11.165	ng	99
47) 2-Chloronaphthalene	9.245	162	152057	10.974	ng	100
48) 2-Nitroaniline	9.345	65	50870	10.392	ng	97
49) Acenaphthylene	9.657	152	242473	11.322	ng	99
50) Dimethylphthalate	9.522	163	185833	10.778	ng	99
51) 2,6-Dinitrotoluene	9.586	165	40589	10.751	ng	96
52) Acenaphthene	9.833	154	148522	11.050	ng	100
53) 3-Nitroaniline	9.757	138	44882	10.552	ng	98
54) 2,4-Dinitrophenol	9.869	184	13246	7.518	ng	94
55) Dibenzofuran	10.004	168	222889	11.413	ng	98
56) 4-Nitrophenol	9.922	139	28485	10.084	ng	95
57) 2,4-Dinitrotoluene	9.992	165	52775	11.156	ng	93
58) Fluorene	10.351	166	174651	11.397	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.127	232	45516	11.233	ng	99
60) Diethylphthalate	10.227	149	177964	10.996	ng	98
61) 4-Chlorophenyl-phenyle...	10.345	204	86740	11.511	ng	93
62) 4-Nitroaniline	10.363	138	43675	10.505	ng	96
63) Azobenzene	10.504	77	174042	10.831	ng	99
65) 4,6-Dinitro-2-methylph...	10.404	198	24348	9.279	ng	85
66) n-Nitrosodiphenylamine	10.463	169	151462	10.890	ng	99
67) 4-Bromophenyl-phenylether	10.833	248	50893	10.764	ng	92
68) Hexachlorobenzene	10.898	284	55603	10.976	ng	97
69) Atrazine	10.992	200	44353	13.282	ng	98
70) Pentachlorophenol	11.092	266	29337	10.671	ng	90
71) Phenanthrene	11.316	178	261567	11.251	ng	98
72) Anthracene	11.369	178	266966	11.249	ng	99
73) Carbazole	11.521	167	227644	11.283	ng	99
74) Di-n-butylphthalate	11.863	149	270022	11.232	ng	99
75) Fluoranthene	12.510	202	262877	11.447	ng	97
77) Benzidine	12.633	184	70731	12.995	ng	99
78) Pyrene	12.739	202	265957	10.216	ng	99
80) Butylbenzylphthalate	13.368	149	91914	10.105	ng	96
81) Benzo(a)anthracene	13.933	228	196281	10.782	ng	99
82) 3,3'-Dichlorobenzidine	13.898	252	58888	10.598	ng	98
83) Chrysene	13.968	228	187617	10.554	ng	99
84) Bis(2-ethylhexyl)phtha...	13.933	149	95358	9.931	ng	# 98
85) Di-n-octyl phthalate	14.545	149	131322	8.909	ng	100
87) Indeno(1,2,3-cd)pyrene	16.868	276	139044	9.795	ng	99
88) Benzo(b)fluoranthene	14.974	252	133416	11.232	ng	97
89) Benzo(k)fluoranthene	15.004	252	129484	11.226	ng	# 96
90) Benzo(a)pyrene	15.339	252	116704	10.506	ng	98
91) Dibenzo(a,h)anthracene	16.886	278	117691	10.157	ng	99
92) Benzo(g,h,i)perylene	17.315	276	118715	9.930	ng	98

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

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