

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110424\
 Data File : BF140218.D
 Acq On : 04 Nov 2024 14:46
 Operator : RC/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICCC040

Quant Time: Nov 04 16:54:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 16:40:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	218244	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	806521	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	468102	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	833437	20.000	ng	0.00
76) Chrysene-d12	14.057	240	519714	20.000	ng	0.00
86) Perylene-d12	15.545	264	459164	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	998052	77.490	ng	0.00
7) Phenol-d6	6.504	99	1271918	76.962	ng	0.00
23) Nitrobenzene-d5	7.439	82	1244930	78.089	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	412406	79.905	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	2294804	76.627	ng	0.00
79) Terphenyl-d14	12.992	244	2571293	78.937	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.669	88	222058	39.059	ng	100
3) Pyridine	3.434	79	552579	39.569	ng	100
4) n-Nitrosodimethylamine	3.393	42	310902	39.743	ng	100
6) Aniline	6.545	93	563966	39.096	ng	100
8) 2-Chlorophenol	6.663	128	524393	38.626	ng	100
9) Benzaldehyde	6.428	77	399291	37.185	ng	100
10) Phenol	6.516	94	682268	38.815	ng	100
11) bis(2-Chloroethyl)ether	6.616	93	521054	38.558	ng	100
12) 1,3-Dichlorobenzene	6.822	146	622778	38.853	ng	100
13) 1,4-Dichlorobenzene	6.898	146	621933	38.409	ng	100
14) 1,2-Dichlorobenzene	7.051	146	580529	38.363	ng	100
15) Benzyl Alcohol	7.016	79	513086	39.636	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.151	45	752863	38.709	ng	100
17) 2-Methylphenol	7.128	107	429040	39.045	ng	100
18) Hexachloroethane	7.392	117	236329	39.203	ng	100
19) n-Nitroso-di-n-propyla...	7.292	70	410465	38.038	ng	100
20) 3+4-Methylphenols	7.281	107	545639	38.591	ng	100
22) Acetophenone	7.286	105	761763	38.537	ng	100
24) Nitrobenzene	7.463	77	647639	38.913	ng	100
25) Isophorone	7.698	82	1056162	38.998	ng	100
26) 2-Nitrophenol	7.775	139	264370	40.474	ng	100
27) 2,4-Dimethylphenol	7.804	122	353217	39.484	ng	100
28) bis(2-Chloroethoxy)met...	7.904	93	636323	38.996	ng	100
29) 2,4-Dichlorophenol	8.016	162	446674	39.483	ng	100
30) 1,2,4-Trichlorobenzene	8.098	180	538036	38.735	ng	100
31) Naphthalene	8.181	128	1573759	38.706	ng	100
32) Benzoic acid	7.922	122	292435	41.112	ng	100
33) 4-Chloroaniline	8.228	127	511942	37.990	ng	100
34) Hexachlorobutadiene	8.298	225	377324	39.144	ng	100
35) Caprolactam	8.604	113	134745	40.138	ng	100
36) 4-Chloro-3-methylphenol	8.704	107	480969	39.039	ng	100
37) 2-Methylnaphthalene	8.869	142	1044965	38.512	ng	100
38) 1-Methylnaphthalene	8.969	142	1017797	38.322	ng	100
40) 1,2,4,5-Tetrachloroben...	9.033	216	556861	38.543	ng	100
41) Hexachlorocyclopentadiene	9.022	237	173564	43.803	ng	100
43) 2,4,6-Trichlorophenol	9.145	196	358015	40.581	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	356101	38.633	ng	100
46) 1,1'-Biphenyl	9.333	154	1308608	38.514	ng	100
47) 2-Chloronaphthalene	9.363	162	992821	38.511	ng	100
48) 2-Nitroaniline	9.457	65	325782	40.655	ng	100
49) Acenaphthylene	9.780	152	1400520	38.599	ng	100
50) Dimethylphthalate	9.639	163	1110672	38.691	ng	100
51) 2,6-Dinitrotoluene	9.698	165	268465	39.246	ng	100
52) Acenaphthene	9.951	154	1006352	38.911	ng	100
53) 3-Nitroaniline	9.869	138	250085	40.066	ng	100
54) 2,4-Dinitrophenol	9.975	184	121375	38.827	ng	100
55) Dibenzofuran	10.122	168	1413860	38.702	ng	100
56) 4-Nitrophenol	10.022	139	192314	41.638	ng	100
57) 2,4-Dinitrotoluene	10.104	165	360025	40.205	ng	100
58) Fluorene	10.469	166	1119281	38.313	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.239	232	305647	40.021	ng	100
60) Diethylphthalate	10.339	149	1136253	38.856	ng	100
61) 4-Chlorophenyl-phenyle...	10.457	204	571115	37.992	ng	100
62) 4-Nitroaniline	10.486	138	253973	40.023	ng	100
63) Azobenzene	10.616	77	1140730	38.469	ng	100
65) 4,6-Dinitro-2-methylph...	10.516	198	184843	42.358	ng	100
66) n-Nitrosodiphenylamine	10.574	169	942552	39.262	ng	100
67) 4-Bromophenyl-phenylether	10.951	248	367721	39.178	ng	100
68) Hexachlorobenzene	11.016	284	417351	39.402	ng	100
69) Atrazine	11.104	200	276820	41.217	ng	100
70) Pentachlorophenol	11.210	266	242711	43.026	ng	100
71) Phenanthrene	11.433	178	1568592	38.717	ng	100
72) Anthracene	11.480	178	1543534	38.876	ng	100
73) Carbazole	11.639	167	1414089	39.269	ng	100
74) Di-n-butylphthalate	11.969	149	1670622	39.557	ng	100
75) Fluoranthene	12.621	202	1694379	39.045	ng	100
77) Benzidine	12.739	184	312126	34.623	ng	100
78) Pyrene	12.851	202	1685783	39.497	ng	100
80) Butylbenzylphthalate	13.468	149	608675	40.662	ng	100
81) Benzo(a)anthracene	14.045	228	1337026	39.223	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	406483	39.524	ng	100
83) Chrysene	14.080	228	1221565	38.958	ng	100
84) Bis(2-ethylhexyl)phtha...	14.033	149	827560	39.751	ng	100
85) Di-n-octyl phthalate	14.657	149	1230269	39.849	ng	100
87) Indeno(1,2,3-cd)pyrene	17.056	276	1164443	40.291	ng	100
88) Benzo(b)fluoranthene	15.109	252	1176440	40.620	ng	100
89) Benzo(k)fluoranthene	15.139	252	952334	36.899	ng	100
90) Benzo(a)pyrene	15.486	252	919956	39.410	ng	100
91) Dibenzo(a,h)anthracene	17.074	278	958091	40.705	ng	100
92) Benzo(g,h,i)perylene	17.515	276	971858	40.439	ng	100

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

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