

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042024\
 Data File : BG060975.D
 Acq On : 20 Apr 2024 19:41
 Operator : MA/JU
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Quant Time: Apr 22 04:15:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG042024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 22 04:10:09 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.943	152	20618	20.000	ng	0.00
21) Naphthalene-d8	10.740	136	82022	20.000	ng	0.00
39) Acenaphthene-d10	14.570	164	72860	20.000	ng	0.00
64) Phenanthrene-d10	17.320	188	190054	20.000	ng	0.00
76) Chrysene-d12	21.580	240	189201	20.000	ng	0.00
86) Perylene-d12	24.700	264	221856	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.528	112	79000	77.504	ng	0.00
7) Phenol-d6	7.109	99	120973	79.989	ng	0.00
23) Nitrobenzene-d5	9.095	82	138496	75.864	ng	0.00
42) 2,4,6-Tribromophenol	16.063	330	122075	76.299	ng	0.00
45) 2-Fluorobiphenyl	13.196	172	405065	76.944	ng	0.00
79) Terphenyl-d14	19.941	244	925313	75.914	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.448	88	13923	39.468	ng	100
3) Pyridine	3.836	79	42344	38.832	ng	100
4) n-Nitrosodimethylamine	3.754	42	38880	39.644	ng	100
6) Aniline	7.267	93	71868	39.539	ng	100
8) 2-Chlorophenol	7.508	128	48574	38.246	ng	100
9) Benzaldehyde	7.085	77	29092	37.397	ng	100
10) Phenol	7.132	94	59648	39.697	ng	100
11) bis(2-Chloroethyl)ether	7.361	93	41519	38.621	ng	100
12) 1,3-Dichlorobenzene	7.831	146	54235	38.508	ng	100
13) 1,4-Dichlorobenzene	7.978	146	56247	38.771	ng	100
14) 1,2-Dichlorobenzene	8.295	146	55340	39.032	ng	100
15) Benzyl Alcohol	8.178	79	55538	38.894	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.472	45	91243	38.981	ng	100
17) 2-Methylphenol	8.384	107	48422	39.911	ng	100
18) Hexachloroethane	9.024	117	20850	39.886	ng	100
19) n-Nitroso-di-n-propyla...	8.742	70	43282	38.402	ng	100
20) 3+4-Methylphenols	8.707	107	65500	38.603	ng	100
22) Acetophenone	8.760	105	85413	37.293	ng	100
24) Nitrobenzene	9.136	77	66768	38.061	ng	100
25) Isophorone	9.664	82	115516	37.810	ng	100
26) 2-Nitrophenol	9.847	139	31552	38.270	ng	100
27) 2,4-Dimethylphenol	9.911	122	48515	36.712	ng	100
28) bis(2-Chloroethoxy)met...	10.146	93	58976	36.695	ng	100
29) 2,4-Dichlorophenol	10.387	162	57443	37.663	ng	100
30) 1,2,4-Trichlorobenzene	10.599	180	63327	37.631	ng	100
31) Naphthalene	10.787	128	167849	37.379	ng	100
32) Benzoic acid	10.046	122	39833	39.587	ng	100
33) 4-Chloroaniline	10.887	127	76925	37.711	ng	100
34) Hexachlorobutadiene	11.075	225	55464	37.477	ng	100
35) Caprolactam	11.650	113	19492	40.124	ng	100
36) 4-Chloro-3-methylphenol	12.020	107	67185	37.769	ng	100
37) 2-Methylnaphthalene	12.397	142	135019	38.242	ng	100
38) 1-Methylnaphthalene	12.614	142	127344	38.651	ng	100
40) 1,2,4,5-Tetrachloroben...	12.761	216	101120	39.054	ng	100
41) Hexachlorocyclopentadiene	12.749	237	60088	39.440	ng	100
43) 2,4,6-Trichlorophenol	13.002	196	65079	38.236	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.072	196	76369	37.931	ng	100
46) 1,1'-Biphenyl	13.407	154	189175	38.546	ng	100
47) 2-Chloronaphthalene	13.448	162	143782	38.210	ng	100
48) 2-Nitroaniline	13.642	65	57311	37.737	ng	100
49) Acenaphthylene	14.294	152	246099	38.833	ng	100
50) Dimethylphthalate	14.030	163	229143	38.139	ng	100
51) 2,6-Dinitrotoluene	14.142	165	48474	38.667	ng	100
52) Acenaphthene	14.635	154	152021	39.478	ng	100
53) 3-Nitroaniline	14.471	138	45874	37.111	ng	100
54) 2,4-Dinitrophenol	14.676	184	27723	37.148	ng	100
55) Dibenzofuran	14.970	168	254416	38.260	ng	100
56) 4-Nitrophenol	14.782	139	35145	39.685	ng	100
57) 2,4-Dinitrotoluene	14.929	165	70073	38.810	ng	100
58) Fluorene	15.616	166	211141	37.619	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.193	232	77625	38.125	ng	100
60) Diethylphthalate	15.393	149	223961	37.851	ng	100
61) 4-Chlorophenyl-phenyle...	15.616	204	130949	37.830	ng	100
62) 4-Nitroaniline	15.634	138	50669	38.014	ng	100
63) Azobenzene	15.910	77	180829	37.574	ng	100
65) 4,6-Dinitro-2-methylph...	15.693	198	46387	39.528	ng	100
66) n-Nitrosodiphenylamine	15.828	169	199844	38.167	ng	100
67) 4-Bromophenyl-phenylether	16.509	248	94890	37.831	ng	100
68) Hexachlorobenzene	16.627	284	107014	38.626	ng	100
69) Atrazine	16.780	200	77115	39.350	ng	100
70) Pentachlorophenol	16.968	266	72774	40.058	ng	100
71) Phenanthrene	17.361	178	366609	37.791	ng	100
72) Anthracene	17.449	178	382135	38.247	ng	100
73) Carbazole	17.720	167	315624	38.038	ng	100
74) Di-n-butylphthalate	18.290	149	382433	37.962	ng	100
75) Fluoranthene	19.371	202	487219	37.844	ng	100
77) Benzidine	19.553	184	150880	43.039	ng	100
78) Pyrene	19.735	202	489843	38.024	ng	100
80) Butylbenzylphthalate	20.634	149	163883	38.167	ng	100
81) Benzo(a)anthracene	21.562	228	505091	37.971	ng	100
82) 3,3'-Dichlorobenzidine	21.480	252	183446	38.160	ng	100
83) Chrysene	21.621	228	478813	37.807	ng	100
84) Bis(2-ethylhexyl)phtha...	21.486	149	235210	37.677	ng	100
85) Di-n-octyl phthalate	22.673	149	393022	37.920	ng	100
87) Indeno(1,2,3-cd)pyrene	28.243	276	595413	38.780	ng	100
88) Benzo(b)fluoranthene	23.719	252	484591	38.544	ng	100
89) Benzo(k)fluoranthene	23.783	252	485296	38.203	ng	100
90) Benzo(a)pyrene	24.559	252	467229	38.394	ng	100
91) Dibenzo(a,h)anthracene	28.313	278	497648	39.578	ng	100
92) Benzo(g,h,i)perylene	29.353	276	478816	38.405	ng	100

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

