

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012023\
 Data File : BM038503.D
 Acq On : 20 Jan 2023 17:14
 Operator : CG/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM012023

Quant Time: Jan 20 17:50:23 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 20 17:18:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	74634	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	254900	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	164427	20.000	ng	0.00
64) Phenanthrene-d10	17.345	188	370656	20.000	ng	0.00
76) Chrysene-d12	21.521	240	382768	20.000	ng	0.00
86) Perylene-d12	23.933	264	443043	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	355393	78.719	ng	0.00
7) Phenol-d6	7.134	99	454517	78.906	ng	0.00
23) Nitrobenzene-d5	9.145	82	464412	79.118	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	211416	89.288	ng	0.00
45) 2-Fluorobiphenyl	13.233	172	1106560	82.899	ng	0.00
79) Terphenyl-d14	19.962	244	1626537	82.843	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.375	88	76525	36.235	ng	98
3) Pyridine	3.787	79	184622	34.577	ng	99
4) n-Nitrosodimethylamine	3.705	42	83554	32.167	ng	97
6) Aniline	7.298	93	278987	39.019	ng	99
8) 2-Chlorophenol	7.528	128	186244	40.382	ng	98
9) Benzaldehyde	7.110	77	131611	37.653	ng	96
10) Phenol	7.157	94	224413	38.988	ng	98
11) bis(2-Chloroethyl)ether	7.393	93	173951	38.303	ng	100
12) 1,3-Dichlorobenzene	7.857	146	211448	40.328	ng	98
13) 1,4-Dichlorobenzene	8.004	146	215716	40.718	ng	99
14) 1,2-Dichlorobenzene	8.322	146	209802	40.885	ng	99
15) Benzyl Alcohol	8.216	79	170122	39.500	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.504	45	177444	37.024	ng	98
17) 2-Methylphenol	8.416	107	163954	40.245	ng	98
18) Hexachloroethane	9.045	117	80127	40.683	ng	95
19) n-Nitroso-di-n-propyla...	8.781	70	152737	38.826	ng	97
20) 3+4-Methylphenols	8.745	107	218775	40.068	ng	96
22) Acetophenone	8.804	105	280151	39.896	ng	99
24) Nitrobenzene	9.187	77	232339	39.119	ng	96
25) Isophorone	9.710	82	387936	40.047	ng	99
26) 2-Nitrophenol	9.898	139	101776	43.171	ng	93
27) 2,4-Dimethylphenol	9.951	122	163963	41.790	ng	99
28) bis(2-Chloroethoxy)met...	10.192	93	229080	40.425	ng	100
29) 2,4-Dichlorophenol	10.428	162	193858	42.116	ng	99
30) 1,2,4-Trichlorobenzene	10.645	180	221674	42.112	ng	98
31) Naphthalene	10.834	128	566520	40.911	ng	99
32) Benzoic acid	10.092	122	117931	39.505	ng	96
33) 4-Chloroaniline	10.951	127	245490	42.473	ng	96
34) Hexachlorobutadiene	11.104	225	145212	41.938	ng	97
35) Caprolactam	11.745	113	47457	42.526	ng	92
36) 4-Chloro-3-methylphenol	12.069	107	173000	41.331	ng	99
37) 2-Methylnaphthalene	12.433	142	418001	42.003	ng	98
38) 1-Methylnaphthalene	12.657	142	392092	41.776	ng	100
40) 1,2,4,5-Tetrachloroben...	12.804	216	260955	42.190	ng	100
41) Hexachlorocyclopentadiene	12.775	237	152449	44.845	ng	97
43) 2,4,6-Trichlorophenol	13.045	196	167912	42.408	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.116	196	196913	42.462	ng	99
46) 1,1'-Biphenyl	13.439	154	547468	41.574	ng	99
47) 2-Chloronaphthalene	13.486	162	426199	41.461	ng	98
48) 2-Nitroaniline	13.698	65	123810	41.012	ng	93
49) Acenaphthylene	14.327	152	669666	41.638	ng	99
50) Dimethylphthalate	14.069	163	529463	41.831	ng	100
51) 2,6-Dinitrotoluene	14.192	165	113392	42.569	ng	98
52) Acenaphthene	14.669	154	402916	41.078	ng	99
53) 3-Nitroaniline	14.522	138	114782	43.362	ng	95
54) 2,4-Dinitrophenol	14.733	184	77644	39.819	ng	97
55) Dibenzofuran	15.004	168	675847	41.213	ng	99
56) 4-Nitrophenol	14.822	139	89046	40.849	ng	96
57) 2,4-Dinitrotoluene	14.974	165	155865	43.166	ng	100
58) Fluorene	15.651	166	551180	41.449	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.227	232	160567	42.359	ng	100
60) Diethylphthalate	15.421	149	512318	41.332	ng	99
61) 4-Chlorophenyl-phenyle...	15.639	204	287238	41.142	ng	98
62) 4-Nitroaniline	15.680	138	115504	38.543	ng	97
63) Azobenzene	15.933	77	511426	38.943	ng	98
65) 4,6-Dinitro-2-methylph...	15.739	198	103945	43.203	ng	94
66) n-Nitrosodiphenylamine	15.857	169	465830	41.864	ng	98
67) 4-Bromophenyl-phenylether	16.533	248	177554	43.186	ng	100
68) Hexachlorobenzene	16.651	284	194674	43.880	ng	97
69) Atrazine	16.810	200	160226	43.160	ng	97
70) Pentachlorophenol	16.998	266	143424	45.444	ng	98
71) Phenanthrene	17.386	178	847692	40.953	ng	99
72) Anthracene	17.480	178	871133	41.360	ng	100
73) Carbazole	17.757	167	774204	41.740	ng	99
74) Di-n-butylphthalate	18.304	149	853345	41.876	ng	100
75) Fluoranthene	19.398	202	1041291	41.857	ng	99
77) Benzidine	19.586	184	385024	44.392	ng	99
78) Pyrene	19.762	202	1102836	41.379	ng	99
80) Butylbenzylphthalate	20.651	149	391435	42.472	ng	98
81) Benzo(a)anthracene	21.503	228	1127059	41.555	ng	99
82) 3,3'-Dichlorobenzidine	21.439	252	377558	43.338	ng	99
83) Chrysene	21.562	228	1093128	41.213	ng	99
84) Bis(2-ethylhexyl)phtha...	21.415	149	572795	42.885	ng	99
85) Di-n-octyl phthalate	22.345	149	985932	41.513	ng	99
87) Indeno(1,2,3-cd)pyrene	26.462	276	1353633	41.289	ng	98
88) Benzo(b)fluoranthene	23.197	252	1113492	41.663	ng	100
89) Benzo(k)fluoranthene	23.244	252	1159685	41.935	ng	99
90) Benzo(a)pyrene	23.833	252	1104239	41.674	ng	99
91) Dibenzo(a,h)anthracene	26.474	278	1127812	41.329	ng	97
92) Benzo(g,h,i)perylene	27.238	276	1167537	41.769	ng	98

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

