

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012023\
 Data File : BM038497.D
 Acq On : 20 Jan 2023 13:26
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/23/2023
 Supervised By : Jagrut Upadhyay 01/23/2023

Quant Time: Jan 21 00:05:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 21 00:02:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.969	152	66063	20.000	ng	0.00
21) Naphthalene-d8	10.781	136	220423	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	136626	20.000	ng	0.00
64) Phenanthrene-d10	17.345	188	310068	20.000	ng	0.00
76) Chrysene-d12	21.521	240	291763	20.000	ng	0.00
86) Perylene-d12	23.933	264	385704	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	78693	19.692	ng	0.00
7) Phenol-d6	7.128	99	98212	19.262	ng	0.00
23) Nitrobenzene-d5	9.140	82	97721	19.252	ng	0.00
42) 2,4,6-Tribromophenol	16.092	330	34630	17.601	ng	0.00
45) 2-Fluorobiphenyl	13.228	172	210295	18.960	ng	0.00
79) Terphenyl-d14	19.957	244	309575	20.685	ng	0.00
Target Compounds						
2) 1,4-Dioxane	3.375	88	20316m	10.868	ng	Qvalue
3) Pyridine	3.799	79	49119m	10.393	ng	
4) n-Nitrosodimethylamine	3.705	42	25283m	10.996	ng	96
6) Aniline	7.299	93	61998	9.796	ng	
8) 2-Chlorophenol	7.528	128	39919	9.778	ng	98
9) Benzaldehyde	7.110	77	34538m	11.163	ng	92
10) Phenol	7.157	94	49690	9.753	ng	
11) bis(2-Chloroethyl)ether	7.393	93	38648	9.614	ng	94
12) 1,3-Dichlorobenzene	7.857	146	44872	9.668	ng	97
13) 1,4-Dichlorobenzene	8.005	146	45036	9.604	ng	97
14) 1,2-Dichlorobenzene	8.322	146	43493	9.575	ng	96
15) Benzyl Alcohol	8.216	79	36764	9.644	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.493	45	42863	10.104	ng	99
17) 2-Methylphenol	8.416	107	33893	9.399	ng	95
18) Hexachloroethane	9.046	117	16557	9.497	ng	91
19) n-Nitroso-di-n-propyla...	8.781	70	32787	10.013	ng	99
20) 3+4-Methylphenols	8.746	107	43707	9.043	ng	97
22) Acetophenone	8.799	105	59360	9.776	ng	96
24) Nitrobenzene	9.187	77	51053	9.940	ng	99
25) Isophorone	9.710	82	77546	9.257	ng	99
26) 2-Nitrophenol	9.898	139	18238	8.946	ng	92
27) 2,4-Dimethylphenol	9.951	122	31315	9.230	ng	91
28) bis(2-Chloroethoxy)met...	10.193	93	47584	9.710	ng	96
29) 2,4-Dichlorophenol	10.428	162	36646	9.207	ng	96
30) 1,2,4-Trichlorobenzene	10.640	180	44056	9.678	ng	99
31) Naphthalene	10.828	128	111155	9.282	ng	99
32) Benzoic acid	10.040	122	16577	6.856	ng	92
33) 4-Chloroaniline	10.951	127	45757	9.155	ng	98
34) Hexachlorobutadiene	11.104	225	28826	9.627	ng	93
35) Caprolactam	11.728	113	7635	7.912	ng	88
36) 4-Chloro-3-methylphenol	12.063	107	32092	8.866	ng	90
37) 2-Methylnaphthalene	12.434	142	77126	8.962	ng	97
38) 1-Methylnaphthalene	12.651	142	72459	8.928	ng	98
40) 1,2,4,5-Tetrachloroben...	12.798	216	50623	9.850	ng	99
41) Hexachlorocyclopentadiene	12.775	237	24909	8.818	ng	96
43) 2,4,6-Trichlorophenol	13.045	196	30679	9.325	ng	95

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.116	196	35941	9.327	ng	98
46) 1,1'-Biphenyl	13.439	154	102196	9.340	ng	100
47) 2-Chloronaphthalene	13.486	162	79947	9.360	ng	98
48) 2-Nitroaniline	13.698	65	21918	8.738	ng	96
49) Acenaphthylene	14.328	152	121659	9.104	ng	99
50) Dimethylphthalate	14.063	163	96257	9.152	ng	100
51) 2,6-Dinitrotoluene	14.186	165	19762	8.929	ng	90
52) Acenaphthene	14.669	154	75324	9.242	ng	99
53) 3-Nitroaniline	14.522	138	18274	8.308	ng	89
54) 2,4-Dinitrophenol	14.733	184	11255	7.352	ng	98
55) Dibenzofuran	14.998	168	128434	9.426	ng	98
56) 4-Nitrophenol	14.828	139	13639	7.530	ng	96
57) 2,4-Dinitrotoluene	14.975	165	25330	8.442	ng	# 91
58) Fluorene	15.645	166	100404	9.087	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.228	232	29173	9.262	ng	99
60) Diethylphthalate	15.416	149	93013	9.031	ng	99
61) 4-Chlorophenyl-phenyle...	15.639	204	56246	9.696	ng	96
62) 4-Nitroaniline	15.680	138	17570	7.775	ng	94
63) Azobenzene	15.933	77	101009	9.257	ng	99
65) 4,6-Dinitro-2-methylph...	15.733	198	17161	8.527	ng	93
66) n-Nitrosodiphenylamine	15.857	169	85854	9.223	ng	99
67) 4-Bromophenyl-phenylether	16.533	248	31927	9.283	ng	100
68) Hexachlorobenzene	16.645	284	34867	9.395	ng	98
69) Atrazine	16.804	200	28815	9.279	ng	94
70) Pentachlorophenol	16.992	266	22873	8.664	ng	96
71) Phenanthrene	17.386	178	161367	9.319	ng	98
72) Anthracene	17.480	178	161271	9.153	ng	99
73) Carbazole	17.751	167	139181	8.970	ng	99
74) Di-n-butylphthalate	18.304	149	146897	8.617	ng	99
75) Fluoranthene	19.398	202	191188	9.187	ng	98
77) Benzidine	19.586	184	69702	10.543	ng	99
78) Pyrene	19.757	202	203390	10.012	ng	99
80) Butylbenzylphthalate	20.645	149	60482	8.609	ng	94
81) Benzo(a)anthracene	21.504	228	196773	9.518	ng	99
82) 3,3'-Dichlorobenzidine	21.439	252	62881	9.469	ng	98
83) Chrysene	21.557	228	198352	9.811	ng	99
84) Bis(2-ethylhexyl)phtha...	21.415	149	85187	8.367	ng	99
85) Di-n-octyl phthalate	22.339	149	164944	9.111	ng	97
87) Indeno(1,2,3-cd)pyrene	26.450	276	264703	9.274	ng	99
88) Benzo(b)fluoranthene	23.192	252	210396	9.043	ng	99
89) Benzo(k)fluoranthene	23.245	252	223196	9.271	ng	99
90) Benzo(a)pyrene	23.827	252	209663	9.089	ng	99
91) Dibenzo(a,h)anthracene	26.462	278	220939	9.300	ng	98
92) Benzo(g,h,i)perylene	27.227	276	228741	9.400	ng	97

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

