

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051424\
 Data File : BF137917.D
 Acq On : 14 May 2024 14:42
 Operator : CG\JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 05/15/2024
 Supervised By :mohammad ahmed 05/17/2024

Quant Time: May 14 17:20:37 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF051424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 14 15:07:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.045	152	128965	20.000	ng	0.00
21) Naphthalene-d8	8.328	136	476897	20.000	ng	0.00
39) Acenaphthene-d10	10.092	164	261375	20.000	ng	0.00
64) Phenanthrene-d10	11.580	188	432281	20.000	ng	0.00
76) Chrysene-d12	14.233	240	224761	20.000	ng	0.00
86) Perylene-d12	15.792	264	241364	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.669	112	1131180	148.992	ng	0.00
7) Phenol-d6	6.669	99	1382032	147.773	ng	0.01
23) Nitrobenzene-d5	7.616	82	1266206	156.253	ng	0.00
42) 2,4,6-Tribromophenol	10.886	330	419860	159.313	ng	0.00
45) 2-Fluorobiphenyl	9.404	172	2444085	144.040	ng	0.00
79) Terphenyl-d14	13.168	244	2373313	155.623	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.940	88	260462	79.191	ng	99
3) Pyridine	3.716	79	613015	83.066	ng	99
4) n-Nitrosodimethylamine	3.698	42	260271	79.945	ng	95
6) Aniline	6.710	93	675353m	79.656	ng	
8) 2-Chlorophenol	6.834	128	615345	74.893	ng	100
9) Benzaldehyde	6.598	77	331861	65.586	ng	98
10) Phenol	6.681	94	703271	74.499	ng	100
11) bis(2-Chloroethyl)ether	6.781	93	548175	75.630	ng	99
12) 1,3-Dichlorobenzene	6.986	146	706784	74.349	ng	98
13) 1,4-Dichlorobenzene	7.063	146	707819	74.179	ng	98
14) 1,2-Dichlorobenzene	7.222	146	655216	72.959	ng	97
15) Benzyl Alcohol	7.186	79	505998	83.788	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.316	45	669392	75.131	ng	98
17) 2-Methylphenol	7.286	107	504295	76.963	ng	99
18) Hexachloroethane	7.557	117	245687	76.089	ng	98
19) n-Nitroso-di-n-propyla...	7.463	70	392428	74.686	ng	97
20) 3+4-Methylphenols	7.439	107	628192	73.181	ng	95
22) Acetophenone	7.457	105	806451	75.487	ng	99
24) Nitrobenzene	7.633	77	641127	77.989	ng	99
25) Isophorone	7.869	82	1113382	79.164	ng	97
26) 2-Nitrophenol	7.945	139	344500	83.401	ng	96
27) 2,4-Dimethylphenol	7.969	122	422167	79.383	ng	97
28) bis(2-Chloroethoxy)met...	8.069	93	658298	77.229	ng	99
29) 2,4-Dichlorophenol	8.181	162	527887	78.936	ng	100
30) 1,2,4-Trichlorobenzene	8.269	180	573284	76.778	ng	99
31) Naphthalene	8.351	128	1805040	75.277	ng	99
32) Benzoic acid	8.104	122	407527	90.443	ng	99
33) 4-Chloroaniline	8.398	127	627601	79.841	ng	99
34) Hexachlorobutadiene	8.463	225	363833	76.983	ng	99
35) Caprolactam	8.786	113	160757	82.953	ng	99
36) 4-Chloro-3-methylphenol	8.869	107	543237	79.852	ng	96
37) 2-Methylnaphthalene	9.045	142	1159537	75.230	ng	99
38) 1-Methylnaphthalene	9.145	142	1128440	74.470	ng	100
40) 1,2,4,5-Tetrachloroben...	9.204	216	548608	77.205	ng	100
41) Hexachlorocyclopentadiene	9.192	237	239016	88.524	ng	99
43) 2,4,6-Trichlorophenol	9.316	196	369744	79.360	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.357	196	402024	78.862	ng	99
46) 1,1'-Biphenyl	9.510	154	1381260	75.345	ng	99
47) 2-Chloronaphthalene	9.533	162	1106915	75.639	ng	96
48) 2-Nitroaniline	9.627	65	309864	82.687	ng	99
49) Acenaphthylene	9.957	152	1592518	75.682	ng	99
50) Dimethylphthalate	9.810	163	1264950	77.177	ng	100
51) 2,6-Dinitrotoluene	9.869	165	301486	80.144	ng	96
52) Acenaphthene	10.127	154	1074941	77.292	ng	99
53) 3-Nitroaniline	10.045	138	299451	81.426	ng	94
54) 2,4-Dinitrophenol	10.145	184	166997	81.983	ng	# 79
55) Dibenzofuran	10.298	168	1511758	74.690	ng	99
56) 4-Nitrophenol	10.186	139	233018	88.220	ng	97
57) 2,4-Dinitrotoluene	10.274	165	399620	82.173	ng	91
58) Fluorene	10.645	166	1184944	74.016	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.410	232	315911	78.884	ng	100
60) Diethylphthalate	10.504	149	1174205	76.708	ng	99
61) 4-Chlorophenyl-phenyle...	10.627	204	591437	74.913	ng	99
62) 4-Nitroaniline	10.663	138	286852	82.238	ng	97
63) Azobenzene	10.792	77	1068065	76.709	ng	98
65) 4,6-Dinitro-2-methylph...	10.680	198	220895	88.499	ng	88
66) n-Nitrosodiphenylamine	10.751	169	1010704	76.241	ng	99
67) 4-Bromophenyl-phenylether	11.121	248	373544	78.823	ng	97
68) Hexachlorobenzene	11.186	284	403640	77.053	ng	# 91
69) Atrazine	11.269	200	247511	68.374	ng	99
70) Pentachlorophenol	11.374	266	263701	85.938	ng	98
71) Phenanthrene	11.610	178	1575891	73.977	ng	99
72) Anthracene	11.663	178	1576272	74.578	ng	99
73) Carbazole	11.816	167	1451561	75.181	ng	100
74) Di-n-butylphthalate	12.127	149	1645880	76.813	ng	99
75) Fluoranthene	12.804	202	1577028	72.935	ng	99
77) Benzidine	12.916	184	279254	95.069	ng	97
78) Pyrene	13.033	202	1583958	80.574	ng	100
80) Butylbenzylphthalate	13.633	149	507366	85.566	ng	96
81) Benzo(a)anthracene	14.221	228	1141571	80.280	ng	99
82) 3,3'-Dichlorobenzidine	14.180	252	339297	86.100	ng	98
83) Chrysene	14.257	228	1070713	78.732	ng	99
84) Bis(2-ethylhexyl)phtha...	14.186	149	625172	83.873	ng	98
85) Di-n-octyl phthalate	14.815	149	920479	91.411	ng	100
87) Indeno(1,2,3-cd)pyrene	17.427	276	1366262	88.106	ng	99
88) Benzo(b)fluoranthene	15.321	252	1099300	78.661	ng	100
89) Benzo(k)fluoranthene	15.357	252	1019327	74.245	ng	100
90) Benzo(a)pyrene	15.727	252	970441	81.322	ng	98
91) Dibenzo(a,h)anthracene	17.451	278	1103777	86.152	ng	98
92) Benzo(g,h,i)perylene	17.933	276	1181732	87.802	ng	99

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

