

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
 Data File : BF131963.D
 Acq On : 09 Jan 2023 15:05
 Operator : CG\JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ICVBF010923

Manual Integrations
 APPROVED

Reviewed By :Christian
 Giraldo
 01/10/2023
 Supervised By :Jagrut
 Upadhyay
 01/10/2023

Quant Time: Jan 10 00:35:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 10 00:32:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.787	152	92081	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	326733	20.000	ng	0.00
39) Acenaphthene-d10	9.839	164	194879	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	387917	20.000	ng	0.00
76) Chrysene-d12	13.986	240	242222	20.000	ng	0.00
86) Perylene-d12	15.445	264	177654	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.392	112	440018	78.741	ng	0.00
7) Phenol-d6	6.428	99	584991	78.954	ng	0.00
23) Nitrobenzene-d5	7.369	82	635392	80.711	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	188774	84.248	ng	0.00
45) 2-Fluorobiphenyl	9.163	172	1170800	81.103	ng	0.00
79) Terphenyl-d14	12.927	244	1397242	82.677	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.463	88	97582	39.982	ng	99
3) Pyridine	3.193	79	280543	40.956	ng	99
4) n-Nitrosodimethylamine	3.169	42	129466	40.216	ng	92
6) Aniline	6.457	93	386557	39.653	ng	99
8) 2-Chlorophenol	6.575	128	226694	40.433	ng	98
9) Benzaldehyde	6.339	77	189169	38.371	ng	98
10) Phenol	6.445	94	329615	39.879	ng	99
11) bis(2-Chloroethyl)ether	6.534	93	249037	39.969	ng	98
12) 1,3-Dichlorobenzene	6.728	146	250018	40.036	ng	99
13) 1,4-Dichlorobenzene	6.804	146	252501	39.961	ng	99
14) 1,2-Dichlorobenzene	6.957	146	236215	39.362	ng	97
15) Benzyl Alcohol	6.939	79	258742	40.102	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.069	45	315209	40.087	ng	100
17) 2-Methylphenol	7.051	107	221978	40.007	ng	97
18) Hexachloroethane	7.298	117	103430	39.582	ng	96
19) n-Nitroso-di-n-propyla...	7.216	70	213527	39.839	ng	98
20) 3+4-Methylphenols	7.210	107	281223	39.479	ng	97
22) Acetophenone	7.210	105	383291	40.291	ng	100
24) Nitrobenzene	7.386	77	325051	40.558	ng	99
25) Isophorone	7.622	82	600579	40.837	ng	98
26) 2-Nitrophenol	7.698	139	130107	41.258	ng	98
27) 2,4-Dimethylphenol	7.739	122	214480	40.747	ng	100
28) bis(2-Chloroethoxy)met...	7.833	93	326990	40.173	ng	97
29) 2,4-Dichlorophenol	7.945	162	219059	40.630	ng	99
30) 1,2,4-Trichlorobenzene	8.022	180	249424	40.394	ng	100
31) Naphthalene	8.104	128	693689	40.174	ng	99
32) Benzoic acid	7.892	122	126709	41.522	ng	96
33) 4-Chloroaniline	8.157	127	291150	40.861	ng	97
34) Hexachlorobutadiene	8.216	225	170451	39.815	ng	98
35) Caprolactam	8.551	113	63554m	40.136	ng	
36) 4-Chloro-3-methylphenol	8.645	107	258785	40.896	ng	99
37) 2-Methylnaphthalene	8.798	142	509689	41.100	ng	100
38) 1-Methylnaphthalene	8.892	142	464913	40.464	ng	98
40) 1,2,4,5-Tetrachloroben...	8.963	216	281940	41.056	ng	99
41) Hexachlorocyclopentadiene	8.939	237	121984	40.042	ng	100
43) 2,4,6-Trichlorophenol	9.075	196	170433	41.232	ng	99

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44) 2,4,5-Trichlorophenol	9.122	196	199062	41.366	ng	98
46) 1,1'-Biphenyl	9.263	154	614955	40.884	ng	99
47) 2-Chloronaphthalene	9.286	162	485223	41.563	ng	100
48) 2-Nitroaniline	9.392	65	174860	41.238	ng	96
49) Acenaphthylene	9.704	152	739226	40.805	ng	100
50) Dimethylphthalate	9.575	163	629528	41.442	ng	100
51) 2,6-Dinitrotoluene	9.639	165	131188	41.061	ng	98
52) Acenaphthene	9.875	154	447226	41.126	ng	97
53) 3-Nitroaniline	9.810	138	125713	40.557	ng	96
54) 2,4-Dinitrophenol	9.916	184	69684	43.440	ng	94
55) Dibenzofuran	10.051	168	696806	40.650	ng	99
56) 4-Nitrophenol	9.975	139	75868	43.454	ng	98
57) 2,4-Dinitrotoluene	10.045	165	177643	41.587	ng	99
58) Fluorene	10.392	166	556164	40.616	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.169	232	153448	42.795	ng	94
60) Diethylphthalate	10.269	149	613892	41.099	ng	100
61) 4-Chlorophenyl-phenyle...	10.386	204	320235	40.852	ng	99
62) 4-Nitroaniline	10.427	138	118941m	42.674	ng	
63) Azobenzene	10.545	77	648622	40.935	ng	99
65) 4,6-Dinitro-2-methylph...	10.457	198	107461	42.957	ng	100
66) n-Nitrosodiphenylamine	10.510	169	485377	39.882	ng	100
67) 4-Bromophenyl-phenylether	10.874	248	193944	39.966	ng	99
68) Hexachlorobenzene	10.939	284	194061	39.532	ng	97
69) Atrazine	11.039	200	178649	40.366	ng	99
70) Pentachlorophenol	11.139	266	96708	39.288	ng	99
71) Phenanthrene	11.357	178	799420	39.532	ng	99
72) Anthracene	11.410	178	835684	40.142	ng	100
73) Carbazole	11.569	167	690003	40.035	ng	100
74) Di-n-butylphthalate	11.898	149	979936	40.833	ng	99
75) Fluoranthene	12.551	202	901823	39.653	ng	99
77) Benzidine	12.674	184	300209	37.306	ng	99
78) Pyrene	12.780	202	918464	41.574	ng	100
80) Butylbenzylphthalate	13.398	149	375141	42.069	ng	98
81) Benzo(a)anthracene	13.974	228	710326	40.308	ng	99
82) 3,3'-Dichlorobenzidine	13.939	252	213042	38.735	ng	98
83) Chrysene	14.015	228	664428	40.344	ng	98
84) Bis(2-ethylhexyl)phtha...	13.957	149	498521	41.265	ng	99
85) Di-n-octyl phthalate	14.574	149	687228	41.570	ng	100
87) Indeno(1,2,3-cd)pyrene	16.909	276	455295	41.787	ng	100
88) Benzo(b)fluoranthene	15.021	252	489784	40.196	ng	99
89) Benzo(k)fluoranthene	15.051	252	503911	41.438	ng	99
90) Benzo(a)pyrene	15.386	252	446520	40.672	ng	98
91) Dibenzo(a,h)anthracene	16.921	278	383786	42.020	ng	99
92) Benzo(g,h,i)perylene	17.351	276	368001	38.390	ng	98

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

