

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110922\
 Data File : BF131085.D
 Acq On : 09 Nov 2022 14:16
 Operator : CG\JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/10/2022
 Supervised By : Jagrut Upadhyay 11/10/2022

Quant Time: Nov 09 15:37:12 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 09 15:21:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	92598	20.000	ng	0.00
21) Naphthalene-d8	8.180	136	324475	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	185248	20.000	ng	0.00
64) Phenanthrene-d10	11.433	188	339200	20.000	ng	0.00
76) Chrysene-d12	14.080	240	224687	20.000	ng	0.00
86) Perylene-d12	15.568	264	158635	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	512465	89.240	ng	0.00
7) Phenol-d6	6.528	99	658125	93.896	ng	0.00
23) Nitrobenzene-d5	7.469	82	669519	95.644	ng	0.00
42) 2,4,6-Tribromophenol	10.733	330	186422	91.404	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	1197791	88.119	ng	0.00
79) Terphenyl-d14	13.021	244	1238870	88.400	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.675	88	118930	43.563	ng	98
3) Pyridine	3.445	79	334742	44.109	ng	99
4) n-Nitrosodimethylamine	3.422	42	192610	45.135	ng	97
6) Aniline	6.563	93	413435	47.429	ng	100
8) 2-Chlorophenol	6.681	128	290559	45.484	ng	95
9) Benzaldehyde	6.451	77	203320	50.218	ng	94
10) Phenol	6.539	94	389788	47.457	ng	100
11) bis(2-Chloroethyl)ether	6.633	93	276995	46.492	ng	98
12) 1,3-Dichlorobenzene	6.833	146	326577	46.453	ng	98
13) 1,4-Dichlorobenzene	6.916	146	330458	45.137	ng	97
14) 1,2-Dichlorobenzene	7.069	146	301684	42.157	ng	98
15) Benzyl Alcohol	7.039	79	260783	42.596	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.169	45	472067	45.562	ng	99
17) 2-Methylphenol	7.151	107	245956	46.019	ng	97
18) Hexachloroethane	7.404	117	126592	45.641	ng	98
19) n-Nitroso-di-n-propyla...	7.316	70	211861	42.414	ng	95
20) 3+4-Methylphenols	7.304	107	297549	43.049	ng	90
22) Acetophenone	7.310	105	386827	44.779	ng	# 93
24) Nitrobenzene	7.486	77	337498	47.009	ng	98
25) Isophorone	7.722	82	537087	44.763	ng	99
26) 2-Nitrophenol	7.798	139	152863	49.994	ng	97
27) 2,4-Dimethylphenol	7.833	122	228196m	48.302	ng	
28) bis(2-Chloroethoxy)met...	7.928	93	304909	46.014	ng	98
29) 2,4-Dichlorophenol	8.039	162	246048	48.356	ng	97
30) 1,2,4-Trichlorobenzene	8.122	180	260298	46.916	ng	98
31) Naphthalene	8.204	128	839703	47.352	ng	99
32) Benzoic acid	7.969	122	161605	50.214	ng	100
33) 4-Chloroaniline	8.257	127	335010	48.310	ng	97
34) Hexachlorobutadiene	8.316	225	176042	48.604	ng	99
35) Caprolactam	8.639	113	74510	47.728	ng	# 88
36) 4-Chloro-3-methylphenol	8.733	107	264830	48.223	ng	99
37) 2-Methylnaphthalene	8.898	142	557246	49.392	ng	97
38) 1-Methylnaphthalene	8.998	142	555912	51.603	ng	99
40) 1,2,4,5-Tetrachloroben...	9.057	216	289642	48.506	ng	99
41) Hexachlorocyclopentadiene	9.039	237	124360	45.183	ng	99
43) 2,4,6-Trichlorophenol	9.175	196	192846	48.426	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.216	196	216720	48.532	ng	94
46) 1,1'-Biphenyl	9.363	154	677790	46.974	ng	99
47) 2-Chloronaphthalene	9.386	162	553073	47.280	ng	100
48) 2-Nitroaniline	9.486	65	202270	47.047	ng	89
49) Acenaphthylene	9.804	152	842370	46.860	ng	100
50) Dimethylphthalate	9.663	163	622890	43.509	ng	100
51) 2,6-Dinitrotoluene	9.727	165	141448	46.444	ng	89
52) Acenaphthene	9.974	154	502426	45.644	ng	98
53) 3-Nitroaniline	9.898	138	163465	49.730	ng	96
54) 2,4-Dinitrophenol	10.004	184	77109	50.281	ng	# 88
55) Dibenzofuran	10.151	168	746965	44.630	ng	95
56) 4-Nitrophenol	10.057	139	115657	49.852	ng	91
57) 2,4-Dinitrotoluene	10.133	165	187603	47.414	ng	90
58) Fluorene	10.492	166	598820	44.989	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	172136	47.318	ng	99
60) Diethylphthalate	10.363	149	594522	39.847	ng	97
61) 4-Chlorophenyl-phenyle...	10.480	204	279307	42.366	ng	98
62) 4-Nitroaniline	10.516	138	166492	49.412	ng	95
63) Azobenzene	10.645	77	626915	43.229	ng	95
65) 4,6-Dinitro-2-methylph...	10.545	198	112638	52.331	ng	93
66) n-Nitrosodiphenylamine	10.604	169	529889	44.209	ng	98
67) 4-Bromophenyl-phenylether	10.974	248	180764	42.414	ng	# 90
68) Hexachlorobenzene	11.039	284	200370	43.343	ng	# 89
69) Atrazine	11.127	200	164745	48.319	ng	98
70) Pentachlorophenol	11.233	266	98767	49.242	ng	97
71) Phenanthrene	11.457	178	878176	45.963	ng	99
72) Anthracene	11.510	178	855517	45.619	ng	99
73) Carbazole	11.668	167	780351	48.635	ng	99
74) Di-n-butylphthalate	11.986	149	881886	41.035	ng	100
75) Fluoranthene	12.651	202	941000	46.400	ng	99
77) Benzidine	12.768	184	330141	54.472	ng	98
78) Pyrene	12.880	202	955940	46.716	ng	99
80) Butylbenzylphthalate	13.492	149	359470	44.590	ng	97
81) Benzo(a)anthracene	14.068	228	772274	47.337	ng	99
82) 3,3'-Dichlorobenzidine	14.033	252	211415	47.688	ng	97
83) Chrysene	14.109	228	730759	47.167	ng	99
84) Bis(2-ethylhexyl)phtha...	14.045	149	386512	41.849	ng	99
85) Di-n-octyl phthalate	14.662	149	493430	37.491	ng	100
87) Indeno(1,2,3-cd)pyrene	17.092	276	609531	52.009	ng	99
88) Benzo(b)fluoranthene	15.133	252	521286	45.219	ng	99
89) Benzo(k)fluoranthene	15.162	252	522180	46.492	ng	99
90) Benzo(a)pyrene	15.509	252	432092	47.145	ng	99
91) Dibenzo(a,h)anthracene	17.109	278	492935	51.236	ng	99
92) Benzo(g,h,i)perylene	17.556	276	521947	53.544	ng	99

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

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