

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020222\
 Data File : BF126779.D
 Acq On : 02 Feb 2022 11:42
 Operator : CG/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/04/2022
 Supervised By : Jagrut Upadhyay 02/04/2022

Quant Time: Feb 02 14:32:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 14:30:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.945	152	126779	20.000	ng	0.00
21) Naphthalene-d8	8.233	136	489644	20.000	ng	0.00
39) Acenaphthene-d10	9.992	164	274349	20.000	ng	0.00
64) Phenanthrene-d10	11.480	188	475701	20.000	ng	0.00
76) Chrysene-d12	14.127	240	361933	20.000	ng	# 0.00
86) Perylene-d12	15.645	264	256395	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.557	112	209887	21.785	ng	0.00
7) Phenol-d6	6.557	99	286042	21.369	ng	-0.01
23) Nitrobenzene-d5	7.504	82	280164	22.396	ng	-0.01
42) 2,4,6-Tribromophenol	10.774	330	58757	22.885	ng	-0.01
45) 2-Fluorobiphenyl	9.304	172	398760	22.260	ng	0.00
79) Terphenyl-d14	13.068	244	460080	21.321	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.787	88	51376	10.820	ng	# 88
3) Pyridine	3.551	79	138128	10.385	ng	# 84
4) n-Nitrosodimethylamine	3.504	42	41004	8.712	ng	# 74
6) Aniline	6.610	93	173027	10.315	ng	97
8) 2-Chlorophenol	6.728	128	93814	10.616	ng	97
9) Benzaldehyde	6.498	77	87004	8.350	ng	86
10) Phenol	6.569	94	153771	10.801	ng	97
11) bis(2-Chloroethyl)ether	6.675	93	121651	10.530	ng	93
12) 1,3-Dichlorobenzene	6.886	146	105982	10.803	ng	# 95
13) 1,4-Dichlorobenzene	6.963	146	109193	11.022	ng	97
14) 1,2-Dichlorobenzene	7.116	146	104228	11.160	ng	97
15) Benzyl Alcohol	7.081	79	108855	9.121	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.216	45	130226	9.880	ng	83
17) 2-Methylphenol	7.186	107	90172	10.379	ng	96
18) Hexachloroethane	7.463	117	41132	10.283	ng	# 92
19) n-Nitroso-di-n-propyla...	7.345	70	94405	9.663	ng	# 85
20) 3+4-Methylphenols	7.334	107	120057	10.658	ng	# 71
22) Acetophenone	7.351	105	157392	10.431	ng	# 88
24) Nitrobenzene	7.528	77	138129	10.696	ng	89
25) Isophorone	7.763	82	234206	9.654	ng	# 94
26) 2-Nitrophenol	7.845	139	43831	13.011	ng	# 82
27) 2,4-Dimethylphenol	7.869	122	80318	10.047	ng	# 83
28) bis(2-Chloroethoxy)met...	7.975	93	152413	10.073	ng	# 97
29) 2,4-Dichlorophenol	8.081	162	78790	10.390	ng	99
30) 1,2,4-Trichlorobenzene	8.169	180	86334	10.616	ng	97
31) Naphthalene	8.251	128	282830	10.692	ng	99
32) Benzoic acid	7.939	122	54006	11.909	ng	# 70
33) 4-Chloroaniline	8.298	127	111900	10.588	ng	# 92
34) Hexachlorobutadiene	8.369	225	53658	10.373	ng	98
35) Caprolactam	8.633	113	26835	9.909	ng	# 82
36) 4-Chloro-3-methylphenol	8.763	107	100742	10.159	ng	86
37) 2-Methylnaphthalene	8.945	142	173289	10.534	ng	97
38) 1-Methylnaphthalene	9.045	142	169032	10.601	ng	98
40) 1,2,4,5-Tetrachloroben...	9.104	216	84376	10.965	ng	96
41) Hexachlorocyclopentadiene	9.092	237	44671	9.535	ng	97
43) 2,4,6-Trichlorophenol	9.216	196	58296	10.627	ng	93

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.251	196	61020m	10.740	ng	
46) 1,1'-Biphenyl	9.404	154	224344	10.854	ng	96
47) 2-Chloronaphthalene	9.433	162	172938	10.650	ng	98
48) 2-Nitroaniline	9.522	65	67064	11.961	ng	# 88
49) Acenaphthylene	9.851	152	270026	10.601	ng	97
50) Dimethylphthalate	9.698	163	205512	10.562	ng	99
51) 2,6-Dinitrotoluene	9.763	165	40566	12.065	ng	# 74
52) Acenaphthene	10.022	154	169585	10.890	ng	94
53) 3-Nitroaniline	9.933	138	46241	12.028	ng	# 75
54) 2,4-Dinitrophenol	10.039	184	15812	12.907	ng	88
55) Dibenzofuran	10.192	168	255375	10.912	ng	98
56) 4-Nitrophenol	10.080	139	32662	10.670	ng	80
57) 2,4-Dinitrotoluene	10.169	165	53572	13.148	ng	95
58) Fluorene	10.539	166	194329	10.998	ng	96
59) 2,3,4,6-Tetrachlorophenol	10.304	232	51274	11.039	ng	# 84
60) Diethylphthalate	10.404	149	202410	10.464	ng	97
61) 4-Chlorophenyl-phenyle...	10.527	204	96912	10.982	ng	91
62) 4-Nitroaniline	10.545	138	40866	10.953	ng	# 75
63) Azobenzene	10.686	77	285604	9.977	ng	89
65) 4,6-Dinitro-2-methylph...	10.574	198	24987	12.389	ng	93
66) n-Nitrosodiphenylamine	10.645	169	169165	10.535	ng	99
67) 4-Bromophenyl-phenylether	11.022	248	57618	10.509	ng	94
68) Hexachlorobenzene	11.080	284	64262	10.986	ng	96
69) Atrazine	11.169	200	53204	10.375	ng	94
70) Pentachlorophenol	11.274	266	34590	10.169	ng	98
71) Phenanthrene	11.504	178	291491	10.725	ng	99
72) Anthracene	11.557	178	290126	10.640	ng	98
73) Carbazole	11.710	167	267679	10.456	ng	99
74) Di-n-butylphthalate	12.033	149	309563	10.028	ng	# 99
75) Fluoranthene	12.698	202	291496	10.859	ng	97
77) Benzidine	12.816	184	81041	6.491	ng	98
78) Pyrene	12.927	202	296732	10.138	ng	99
80) Butylbenzylphthalate	13.545	149	114869	9.033	ng	# 96
81) Benzo(a)anthracene	14.115	228	250005	10.163	ng	99
82) 3,3'-Dichlorobenzidine	14.080	252	69749	8.631	ng	# 97
83) Chrysene	14.157	228	239242	10.107	ng	100
84) Bis(2-ethylhexyl)phtha...	14.098	149	155213	9.013	ng	# 98
85) Di-n-octyl phthalate	14.721	149	203618	7.723	ng	# 93
87) Indeno(1,2,3-cd)pyrene	17.186	276	137913	8.706	ng	# 88
88) Benzo(b)fluoranthene	15.192	252	182124	10.711	ng	# 94
89) Benzo(k)fluoranthene	15.221	252	169591	10.159	ng	# 95
90) Benzo(a)pyrene	15.574	252	134028	9.688	ng	# 94
91) Dibenzo(a,h)anthracene	17.209	278	115705	8.817	ng	# 93
92) Benzo(g,h,i)perylene	17.656	276	111719	8.682	ng	# 86

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

