

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010923\
 Data File : BF131962.D
 Acq On : 09 Jan 2023 14:33
 Operator : CG\JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

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

Quant Time: Jan 10 00:02:18 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF010923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 09 23:53:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.792	152	96506	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	320418	20.000	ng	0.00
39) Acenaphthene-d10	9.845	164	186179	20.000	ng	0.00
64) Phenanthrene-d10	11.333	188	344230	20.000	ng	0.00
76) Chrysene-d12	13.986	240	185475	20.000	ng	# 0.00
86) Perylene-d12	15.445	264	150238	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.398	112	858953	146.661	ng	0.00
7) Phenol-d6	6.445	99	1103047	142.048	ng	0.01
23) Nitrobenzene-d5	7.381	82	1176494	152.390	ng	0.01
42) 2,4,6-Tribromophenol	10.645	330	339644	158.664	ng	0.00
45) 2-Fluorobiphenyl	9.169	172	2093815	151.820	ng	0.00
79) Terphenyl-d14	12.933	244	2216353	171.270	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.469	88	189673	74.151	ng	100
3) Pyridine	3.210	79	544180	75.802	ng	99
4) n-Nitrosodimethylamine	3.199	42	256121	75.912	ng	90
6) Aniline	6.469	93	729491	71.399	ng	97
8) 2-Chlorophenol	6.587	128	429577	73.106	ng	93
10) Phenol	6.457	94	627626	72.453	ng	97
11) bis(2-Chloroethyl)ether	6.539	93	471652	72.227	ng	98
12) 1,3-Dichlorobenzene	6.734	146	477584	72.970	ng	99
13) 1,4-Dichlorobenzene	6.810	146	480606	72.574	ng	99
14) 1,2-Dichlorobenzene	6.969	146	452191	71.897	ng	99
15) Benzyl Alcohol	6.951	79	474181	70.123	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.075	45	562144	68.214	ng	88
17) 2-Methylphenol	7.063	107	418098	71.898	ng	91
18) Hexachloroethane	7.304	117	202658	74.000	ng	93
19) n-Nitroso-di-n-propyla...	7.234	70	399667	71.149	ng	98
20) 3+4-Methylphenols	7.228	107	524691	70.281	ng	# 77
22) Acetophenone	7.222	105	713215	76.450	ng	# 98
24) Nitrobenzene	7.398	77	599003	76.213	ng	96
25) Isophorone	7.639	82	1120238	77.672	ng	98
26) 2-Nitrophenol	7.704	139	242863	78.532	ng	97
27) 2,4-Dimethylphenol	7.751	122	399267	77.348	ng	100
28) bis(2-Chloroethoxy)met...	7.839	93	607131	76.060	ng	99
29) 2,4-Dichlorophenol	7.951	162	408846	77.326	ng	96
30) 1,2,4-Trichlorobenzene	8.028	180	474625	78.381	ng	98
31) Naphthalene	8.110	128	1274168	75.246	ng	99
32) Benzoic acid	7.933	122	258783	87.854	ng	93
33) 4-Chloroaniline	8.163	127	509179	72.868	ng	99
34) Hexachlorobutadiene	8.216	225	330229	78.658	ng	100
35) Caprolactam	8.581	113	129910m	83.585	ng	
36) 4-Chloro-3-methylphenol	8.657	107	471989	76.059	ng	98
37) 2-Methylnaphthalene	8.804	142	938599	77.177	ng	99
38) 1-Methylnaphthalene	8.898	142	855411	75.919	ng	97
40) 1,2,4,5-Tetrachloroben...	8.969	216	518767	79.074	ng	99
41) Hexachlorocyclopentadiene	8.945	237	263280	100.479	ng	97
43) 2,4,6-Trichlorophenol	9.080	196	311775	78.952	ng	99
44) 2,4,5-Trichlorophenol	9.133	196	365476	79.497	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.269	154	1112894	77.446	ng	99
47) 2-Chloronaphthalene	9.298	162	858572	76.980	ng	97
48) 2-Nitroaniline	9.398	65	313090	77.287	ng	98
49) Acenaphthylene	9.710	152	1303451	75.312	ng	100
50) Dimethylphthalate	9.586	163	1106873	76.271	ng	99
51) 2,6-Dinitrotoluene	9.651	165	233143	76.383	ng	97
52) Acenaphthene	9.886	154	798310	76.841	ng	100
53) 3-Nitroaniline	9.822	138	219179	74.015	ng	94
54) 2,4-Dinitrophenol	9.927	184	134420	87.710	ng	99
55) Dibenzofuran	10.057	168	1212998	74.071	ng	100
56) 4-Nitrophenol	9.992	139	137105	82.198	ng	88
57) 2,4-Dinitrotoluene	10.057	165	299767	73.457	ng	96
58) Fluorene	10.398	166	984477	75.255	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.175	232	274394m	80.976	ng	
60) Diethylphthalate	10.280	149	1054147	73.870	ng	100
61) 4-Chlorophenyl-phenyle...	10.392	204	571417	76.302	ng	99
62) 4-Nitroaniline	10.451	138	195501	70.859	ng	94
63) Azobenzene	10.551	77	1113215	73.538	ng	99
65) 4,6-Dinitro-2-methylph...	10.469	198	187256	84.354	ng	87
66) n-Nitrosodiphenylamine	10.516	169	851143	78.812	ng	99
67) 4-Bromophenyl-phenylether	10.880	248	339372	78.809	ng	97
68) Hexachlorobenzene	10.945	284	341005	78.281	ng	97
69) Atrazine	11.051	200	296878	75.593	ng	99
70) Pentachlorophenol	11.145	266	184890	90.286	ng	96
71) Phenanthrene	11.369	178	1372559	76.488	ng	99
72) Anthracene	11.421	178	1412560	76.464	ng	99
73) Carbazole	11.574	167	1147497	75.029	ng	98
74) Di-n-butylphthalate	11.904	149	1622034	76.167	ng	100
75) Fluoranthene	12.557	202	1460397	72.363	ng	99
78) Pyrene	12.786	202	1472553	87.049	ng	100
80) Butylbenzylphthalate	13.404	149	578143	84.670	ng	97
81) Benzo(a)anthracene	13.980	228	1067080	79.079	ng	99
82) 3,3'-Dichlorobenzidine	13.945	252	296089	70.306	ng	99
83) Chrysene	14.015	228	1011100	80.177	ng	99
84) Bis(2-ethylhexyl)phtha...	13.963	149	771285	83.376	ng	# 97
85) Di-n-octyl phthalate	14.574	149	1051539	83.068	ng	100
87) Indeno(1,2,3-cd)pyrene	16.921	276	848203	92.055	ng	100
88) Benzo(b)fluoranthene	15.027	252	774610	75.172	ng	99
89) Benzo(k)fluoranthene	15.057	252	778419	75.692	ng	99
90) Benzo(a)pyrene	15.392	252	727464	78.355	ng	99
91) Dibenzo(a,h)anthracene	16.939	278	713484	92.373	ng	98
92) Benzo(g,h,i)perylene	17.368	276	692994	91.104	ng	98

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

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