

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122222\
 Data File : BF131779.D
 Acq On : 22 Dec 2022 16:18
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 12/23/2022
 Supervised By : Jagrut Upadhyay 12/24/2022

Quant Time: Dec 23 03:30:25 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122022.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Dec 23 03:03:57 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	87262	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	310688	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	183875	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	352821	20.000	ng	0.00
76) Chrysene-d12	14.051	240	199611	20.000	ng	# 0.00
86) Perylene-d12	15.533	264	169532	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	423513	81.725	ng	0.00
7) Phenol-d6	6.492	99	560736	80.392	ng	0.00
23) Nitrobenzene-d5	7.428	82	618668	80.839	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	165218	82.093	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1073183	81.759	ng	0.00
79) Terphenyl-d14	12.992	244	1176188	91.835	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.569	88	102700	42.409	ng	100
3) Pyridine	3.322	79	286243	41.453	ng	100
4) n-Nitrosodimethylamine	3.299	42	129363	41.408	ng	100
6) Aniline	6.522	93	345752	40.152	ng	99
8) 2-Chlorophenol	6.640	128	225383	40.493	ng	99
9) Benzaldehyde	6.404	77	170762	42.156	ng	99
10) Phenol	6.504	94	334342	40.148	ng	99
11) bis(2-Chloroethyl)ether	6.592	93	243016	40.779	ng	99
12) 1,3-Dichlorobenzene	6.792	146	252842	40.186	ng	98
13) 1,4-Dichlorobenzene	6.875	146	259785	40.393	ng	100
14) 1,2-Dichlorobenzene	7.028	146	237776	39.487	ng	99
15) Benzyl Alcohol	7.004	79	248095	39.823	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.134	45	317350	41.763	ng	99
17) 2-Methylphenol	7.116	107	208237	39.630	ng	97
18) Hexachloroethane	7.369	117	104312	40.916	ng	99
19) n-Nitroso-di-n-propyla...	7.275	70	203772	40.507	ng	98
20) 3+4-Methylphenols	7.269	107	269887	39.908	ng	98
22) Acetophenone	7.275	105	362062	39.933	ng	98
24) Nitrobenzene	7.445	77	317074	40.753	ng	94
25) Isophorone	7.686	82	527074	40.554	ng	99
26) 2-Nitrophenol	7.763	139	120867	39.750	ng	96
27) 2,4-Dimethylphenol	7.798	122	173102	39.971	ng	99
28) bis(2-Chloroethoxy)met...	7.898	93	282441	40.923	ng	100
29) 2,4-Dichlorophenol	8.004	162	214271	40.984	ng	99
30) 1,2,4-Trichlorobenzene	8.086	180	245689	39.974	ng	99
31) Naphthalene	8.169	128	676862	39.891	ng	100
32) Benzoic acid	7.934	122	136662m	39.422	ng	
33) 4-Chloroaniline	8.222	127	257893	38.973	ng	99
34) Hexachlorobutadiene	8.281	225	169249	40.929	ng	99
35) Caprolactam	8.610	113	60799	38.567	ng	95
36) 4-Chloro-3-methylphenol	8.704	107	241047	40.361	ng	100
37) 2-Methylnaphthalene	8.863	142	461234	39.851	ng	99
38) 1-Methylnaphthalene	8.963	142	450916	40.225	ng	100
40) 1,2,4,5-Tetrachloroben...	9.028	216	259013	41.173	ng	100
41) Hexachlorocyclopentadiene	9.010	237	117115	41.091	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	169392	41.220	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.181	196	185004	41.585	ng	98
46) 1,1'-Biphenyl	9.328	154	568217	41.083	ng	99
47) 2-Chloronaphthalene	9.357	162	464209	40.988	ng	99
48) 2-Nitroaniline	9.457	65	169380	41.494	ng	95
49) Acenaphthylene	9.775	152	689908	40.614	ng	100
50) Dimethylphthalate	9.633	163	572956	41.412	ng	100
51) 2,6-Dinitrotoluene	9.698	165	123403	41.480	ng	92
52) Acenaphthene	9.945	154	420306	40.391	ng	99
53) 3-Nitroaniline	9.875	138	122854	40.359	ng	90
54) 2,4-Dinitrophenol	9.975	184	69021	38.720	ng	93
55) Dibenzofuran	10.116	168	651425	40.824	ng	99
56) 4-Nitrophenol	10.033	139	86524	39.713	ng	94
57) 2,4-Dinitrotoluene	10.104	165	163993	41.022	ng	97
58) Fluorene	10.463	166	522926	40.533	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.233	232	152192m	40.978	ng	
60) Diethylphthalate	10.333	149	549377	41.209	ng	99
61) 4-Chlorophenyl-phenyle...	10.451	204	281196	41.009	ng	99
62) 4-Nitroaniline	10.492	138	119918	39.143	ng	95
63) Azobenzene	10.616	77	610095	41.673	ng	99
65) 4,6-Dinitro-2-methylph...	10.510	198	100121	41.333	ng	98
66) n-Nitrosodiphenylamine	10.575	169	440237	40.236	ng	99
67) 4-Bromophenyl-phenylether	10.945	248	170490	40.914	ng	98
68) Hexachlorobenzene	11.004	284	185388	40.451	ng	100
69) Atrazine	11.104	200	146778	40.244	ng	98
70) Pentachlorophenol	11.204	266	98636	40.550	ng	97
71) Phenanthrene	11.427	178	760493	39.529	ng	99
72) Anthracene	11.480	178	743041	39.445	ng	99
73) Carbazole	11.639	167	632529	38.811	ng	100
74) Di-n-butylphthalate	11.963	149	833281	41.062	ng	100
75) Fluoranthene	12.616	202	825398	38.522	ng	98
77) Benzidine	12.739	184	159835	43.668	ng	97
78) Pyrene	12.851	202	823997	45.403	ng	100
80) Butylbenzylphthalate	13.468	149	291291	45.508	ng	95
81) Benzo(a)anthracene	14.039	228	599122	41.675	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	169939	42.340	ng	99
83) Chrysene	14.074	228	571377	42.110	ng	100
84) Bis(2-ethylhexyl)phtha...	14.021	149	348354	45.640	ng	99
85) Di-n-octyl phthalate	14.639	149	487146	45.722	ng	100
87) Indeno(1,2,3-cd)pyrene	17.039	276	527996	46.368	ng	99
88) Benzo(b)fluoranthene	15.098	252	472894	39.427	ng	98
89) Benzo(k)fluoranthene	15.127	252	439253	37.531	ng	97
90) Benzo(a)pyrene	15.474	252	375664	39.541	ng	99
91) Dibenzo(a,h)anthracene	17.056	278	436532	46.388	ng	99
92) Benzo(g,h,i)perylene	17.498	276	440030	47.309	ng	98

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

