

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060723\
 Data File : BF133622.D
 Acq On : 07 Jun 2023 13:43
 Operator : CG\JU
 Sample : PB153153BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

PB153153BS

Manual Integrations

APPROVED

Reviewed By :Yogesh

Patel

06/08/2023

Supervised By :mohammad

ahmed

06/08/2023

Quant Time: Jun 07 14:34:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 07 11:27:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	148995	20.000	ng	0.00
21) Naphthalene-d8	8.122	136	572792	20.000	ng	0.00
39) Acenaphthene-d10	9.880	164	275820	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	527994	20.000	ng	0.00
76) Chrysene-d12	14.015	240	289832	20.000	ng	# 0.00
86) Perylene-d12	15.474	264	306879	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	997119	111.658	ng	0.02
7) Phenol-d6	6.475	99	1222429	111.201	ng	0.00
23) Nitrobenzene-d5	7.404	82	807743	94.156	ng	0.00
42) 2,4,6-Tribromophenol	10.674	330	370209	134.650	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	1402006	78.454	ng	0.00
79) Terphenyl-d14	12.957	244	1504478	92.293	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.657	88	134249	33.287	ng	99
3) Pyridine	3.404	79	363990	32.188	ng	96
4) n-Nitrosodimethylamine	3.363	42	238336	38.646	ng	93
6) Aniline	6.498	93	551312	36.936	ng	100
8) 2-Chlorophenol	6.622	128	406007	46.142	ng	95
9) Benzaldehyde	6.387	77	240642	45.047	ng	98
10) Phenol	6.487	94	477442m	43.294	ng	
11) bis(2-Chloroethyl)ether	6.575	93	375707	41.206	ng	98
12) 1,3-Dichlorobenzene	6.781	146	415975	42.582	ng	97
13) 1,4-Dichlorobenzene	6.857	146	424055	43.299	ng	98
14) 1,2-Dichlorobenzene	7.004	146	404540	44.761	ng	99
15) Benzyl Alcohol	6.981	79	364424	43.665	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.116	45	671026	38.803	ng	97
17) 2-Methylphenol	7.092	107	333380	40.008	ng	98
18) Hexachloroethane	7.351	117	152654	45.077	ng	# 89
19) n-Nitroso-di-n-propyla...	7.257	70	276274	38.843	ng	98
20) 3+4-Methylphenols	7.245	107	392602	41.675	ng	91
22) Acetophenone	7.251	105	530187	40.194	ng	# 99
24) Nitrobenzene	7.428	77	431737	47.232	ng	94
25) Isophorone	7.663	82	787766	39.401	ng	99
26) 2-Nitrophenol	7.739	139	218129	51.674	ng	97
27) 2,4-Dimethylphenol	7.775	122	354714	41.697	ng	98
28) bis(2-Chloroethoxy)met...	7.875	93	456665	39.390	ng	99
29) 2,4-Dichlorophenol	7.981	162	329529	43.056	ng	96
30) 1,2,4-Trichlorobenzene	8.063	180	347269	41.949	ng	98
31) Naphthalene	8.145	128	1110811	40.001	ng	99
32) Benzoic acid	7.904	122	254793	45.857	ng	96
33) 4-Chloroaniline	8.192	127	295962	24.867	ng	96
34) Hexachlorobutadiene	8.257	225	211347	41.970	ng	98
35) Caprolactam	8.575	113	118826m	47.340	ng	
36) 4-Chloro-3-methylphenol	8.675	107	370731	43.242	ng	100
37) 2-Methylnaphthalene	8.833	142	733651	38.694	ng	99
38) 1-Methylnaphthalene	8.933	142	673473	38.105	ng	100
40) 1,2,4,5-Tetrachloroben...	9.004	216	338650	41.425	ng	99
41) Hexachlorocyclopentadiene	8.986	237	368078	87.861	ng	100
43) 2,4,6-Trichlorophenol	9.110	196	238170	45.530	ng	98

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44) 2,4,5-Trichlorophenol	9.151	196	254056	41.843	ng	98
46) 1,1'-Biphenyl	9.304	154	900477	40.183	ng	99
47) 2-Chloronaphthalene	9.328	162	667411	41.095	ng	98
48) 2-Nitroaniline	9.422	65	238551	45.937	ng	99
49) Acenaphthylene	9.745	152	1113331	40.170	ng	99
50) Dimethylphthalate	9.604	163	798163	41.173	ng	99
51) 2,6-Dinitrotoluene	9.669	165	174948	47.506	ng	# 83
52) Acenaphthene	9.916	154	752824m	44.841	ng	
53) 3-Nitroaniline	9.833	138	172232	37.984	ng	# 95
54) 2,4-Dinitrophenol	9.945	184	179658	90.810	ng	92
55) Dibenzofuran	10.086	168	982213	39.324	ng	98
56) 4-Nitrophenol	9.998	139	311094	93.797	ng	95
57) 2,4-Dinitrotoluene	10.075	165	230296	50.110	ng	# 81
58) Fluorene	10.427	166	742885	40.595	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.204	232	212395	46.448	ng	99
60) Diethylphthalate	10.304	149	807993	40.150	ng	100
61) 4-Chlorophenyl-phenyle...	10.422	204	353801	40.265	ng	97
62) 4-Nitroaniline	10.457	138	206847	44.646	ng	96
63) Azobenzene	10.580	77	779070	38.674	ng	97
65) 4,6-Dinitro-2-methylph...	10.480	198	115899	52.647	ng	# 77
66) n-Nitrosodiphenylamine	10.539	169	682658	39.825	ng	99
67) 4-Bromophenyl-phenylether	10.910	248	228958	40.832	ng	97
68) Hexachlorobenzene	10.974	284	255591	43.588	ng	98
69) Atrazine	11.069	200	227388	48.347	ng	98
70) Pentachlorophenol	11.174	266	278466	79.422	ng	98
71) Phenanthrene	11.392	178	1068299	39.948	ng	99
72) Anthracene	11.445	178	1097667	39.882	ng	100
73) Carbazole	11.598	167	1057779	40.639	ng	100
74) Di-n-butylphthalate	11.927	149	1223344	39.747	ng	99
75) Fluoranthene	12.580	202	1136703	38.481	ng	99
77) Benzidine	12.704	184	598563	70.611	ng	98
78) Pyrene	12.810	202	1132327	45.911	ng	99
80) Butylbenzylphthalate	13.433	149	475385	48.716	ng	97
81) Benzo(a)anthracene	14.004	228	809643	41.934	ng	99
82) 3,3'-Dichlorobenzidine	13.962	252	257261	42.623	ng	98
83) Chrysene	14.039	228	782082	40.125	ng	99
84) Bis(2-ethylhexyl)phtha...	13.992	149	593908	48.691	ng	98
85) Di-n-octyl phthalate	14.610	149	920372	45.296	ng	97
87) Indeno(1,2,3-cd)pyrene	16.939	276	835409	42.737	ng	98
88) Benzo(b)fluoranthene	15.051	252	803156	40.192	ng	99
89) Benzo(k)fluoranthene	15.080	252	753612	38.677	ng	99
90) Benzo(a)pyrene	15.415	252	697813	38.376	ng	99
91) Dibenzo(a,h)anthracene	16.962	278	695299	43.871	ng	97
92) Benzo(g,h,i)perylene	17.386	276	648264	40.565	ng	99

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

Manual Integrations

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[illegible]