

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062622\
 Data File : BF129015.D
 Acq On : 26 Jun 2022 20:25
 Operator : CG\JU
 Sample : PB145630BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

PB145630BS

Manual Integrations

APPROVED

Reviewed By : Christian Giraldo 06/27/2022
 Supervised By : Jagrut Upadhyay 06/27/2022

Quant Time: Jun 27 06:32:15 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062622.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Mon Jun 27 05:55:00 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.704	152	93311	20.000	ng	0.00
21) Naphthalene-d8	7.986	136	359895	20.000	ng	0.00
39) Acenaphthene-d10	9.745	164	201357	20.000	ng	0.00
64) Phenanthrene-d10	11.233	188	376196	20.000	ng	0.00
76) Chrysene-d12	13.880	240	274892	20.000	ng	# 0.00
86) Perylene-d12	15.304	264	205558	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.345	112	715954	119.116	ng	0.01
7) Phenol-d6	6.369	99	839219	119.444	ng	0.00
23) Nitrobenzene-d5	7.281	82	660609	88.368	ng	0.00
42) 2,4,6-Tribromophenol	10.539	330	194954	93.257	ng	0.00
45) 2-Fluorobiphenyl	9.069	172	1231492	88.617	ng	0.00
79) Terphenyl-d14	12.821	244	1409688	95.062	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.493	88	91330	55.979	ng	97
3) Pyridine	3.193	79	321875	51.107	ng	94
4) n-Nitrosodimethylamine	3.169	42	237625	63.122	ng	# 84
6) Aniline	6.369	93	387661	51.033	ng	# 62
8) 2-Chlorophenol	6.498	128	303406	51.427	ng	97
9) Benzaldehyde	6.257	77	208698	54.496	ng	95
10) Phenol	6.381	94	473773	63.847	ng	82
11) bis(2-Chloroethyl)ether	6.445	93	292424	53.052	ng	98
12) 1,3-Dichlorobenzene	6.645	146	317352	46.854	ng	99
13) 1,4-Dichlorobenzene	6.722	146	317272	46.510	ng	98
14) 1,2-Dichlorobenzene	6.875	146	300646	47.920	ng	99
15) Benzyl Alcohol	6.863	79	269345	48.875	ng	95
16) 2,2'-oxybis(1-Chloropr...	6.981	45	405120	44.475	ng	# 6
17) 2-Methylphenol	7.139	107	308442	48.744	ng	# 80
18) Hexachloroethane	7.210	117	119133	48.166	ng	97
19) n-Nitroso-di-n-propyla...	7.134	70	219115	50.529	ng	96
20) 3+4-Methylphenols	7.139	107	308442	48.744	ng	# 82
22) Acetophenone	7.122	105	417041	43.907	ng	95
24) Nitrobenzene	7.298	77	330595	47.783	ng	97
25) Isophorone	7.539	82	576871	50.916	ng	97
26) 2-Nitrophenol	7.610	139	154126	47.454	ng	89
27) 2,4-Dimethylphenol	7.663	122	256143	53.478	ng	95
28) bis(2-Chloroethoxy)met...	7.745	93	333467	45.567	ng	98
29) 2,4-Dichlorophenol	7.863	162	247281	45.951	ng	99
30) 1,2,4-Trichlorobenzene	7.933	180	270381	45.484	ng	100
31) Naphthalene	8.010	128	843933	46.689	ng	99
32) Benzoic acid	7.833	122	87941	41.427	ng	95
33) 4-Chloroaniline	8.069	127	217268	31.193	ng	97
34) Hexachlorobutadiene	8.122	225	181472	47.457	ng	98
35) Caprolactam	8.457	113	73921m	50.821	ng	
36) 4-Chloro-3-methylphenol	8.575	107	278495	52.564	ng	96
37) 2-Methylnaphthalene	8.704	142	536434	53.185	ng	99
38) 1-Methylnaphthalene	8.804	142	514914	53.129	ng	100
40) 1,2,4,5-Tetrachloroben...	8.869	216	284779	44.253	ng	98
41) Hexachlorocyclopentadiene	8.851	237	158512	76.433	ng	98
43) 2,4,6-Trichlorophenol	8.992	196	166573	45.149	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.045	196	171660	44.687	ng	# 95
46) 1,1'-Biphenyl	9.169	154	674589	46.126	ng	99
47) 2-Chloronaphthalene	9.192	162	530503	46.611	ng	98
48) 2-Nitroaniline	9.304	65	194071	47.923	ng	98
49) Acenaphthylene	9.610	152	844012	49.372	ng	100
50) Dimethylphthalate	9.480	163	634217	47.409	ng	100
51) 2,6-Dinitrotoluene	9.545	165	136365	48.666	ng	# 80
52) Acenaphthene	9.780	154	486121	47.065	ng	99
53) 3-Nitroaniline	9.716	138	104172	33.769	ng	97
54) 2,4-Dinitrophenol	9.839	184	112325	82.112	ng	87
55) Dibenzofuran	9.951	168	723257	45.206	ng	93
56) 4-Nitrophenol	9.916	139	112430	101.606	ng	80
57) 2,4-Dinitrotoluene	9.957	165	179109	48.620	ng	# 85
58) Fluorene	10.298	166	589004	46.769	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.080	232	135085	48.260	ng	96
60) Diethylphthalate	10.180	149	618016	47.463	ng	98
61) 4-Chlorophenyl-phenyle...	10.286	204	294032	45.780	ng	91
62) 4-Nitroaniline	10.339	138	130582	44.841	ng	88
63) Azobenzene	10.451	77	613727	46.026	ng	98
65) 4,6-Dinitro-2-methylph...	10.369	198	91340	42.368	ng	98
66) n-Nitrosodiphenylamine	10.416	169	543771	47.025	ng	99
67) 4-Bromophenyl-phenylether	10.774	248	187795	45.490	ng	93
68) Hexachlorobenzene	10.845	284	217185	47.757	ng	93
69) Atrazine	10.945	200	184093	53.084	ng	97
70) Pentachlorophenol	11.051	266	94998	88.749	ng	98
71) Phenanthrene	11.257	178	874549	46.908	ng	98
72) Anthracene	11.310	178	891949	47.960	ng	99
73) Carbazole	11.474	167	793655	45.031	ng	99
74) Di-n-butylphthalate	11.798	149	946620	47.306	ng	99
75) Fluoranthene	12.451	202	943565	49.246	ng	95
77) Benzidine	12.574	184	357377	Below Cal		98
78) Pyrene	12.680	202	964891	49.076	ng	98
80) Butylbenzylphthalate	13.298	149	387400	49.153	ng	92
81) Benzo(a)anthracene	13.868	228	811191	47.542	ng	100
82) 3,3'-Dichlorobenzidine	13.833	252	184386	47.681	ng	# 97
83) Chrysene	13.910	228	768101	47.065	ng	99
84) Bis(2-ethylhexyl)phtha...	13.851	149	505096	49.373	ng	99
85) Di-n-octyl phthalate	14.468	149	779796	49.508	ng	100
87) Indeno(1,2,3-cd)pyrene	16.715	276	575269	47.629	ng	# 93
88) Benzo(b)fluoranthene	14.904	252	683591m	54.723	ng	
89) Benzo(k)fluoranthene	14.927	252	596757	48.493	ng	# 98
90) Benzo(a)pyrene	15.251	252	597374	58.939	ng	# 97
91) Dibenzo(a,h)anthracene	16.721	278	500073	50.245	ng	# 97
92) Benzo(g,h,i)perylene	17.139	276	501020	50.703	ng	# 93

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

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