

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103124\
 Data File : BF140178.D
 Acq On : 31 Oct 2024 10:15
 Operator : RC/JU
 Sample : PB164540BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

PB164540BS

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 11/04/2024

Supervised By :mohammad ahmed 11/04/2024

Quant Time: Oct 31 12:10:15 2024

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Oct 18 15:07:50 2024

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	149764	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	588319	20.000	ng	0.00
39) Acenaphthene-d10	9.934	164	343514	20.000	ng	0.00
64) Phenanthrene-d10	11.428	188	638189	20.000	ng	0.01
76) Chrysene-d12	14.092	240	408760	20.000	ng	0.04
86) Perylene-d12	15.615	264	343987	20.000	ng	0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	909411	95.061	ng	0.02
7) Phenol-d6	6.516	99	1172309	94.615	ng	0.00
23) Nitrobenzene-d5	7.451	82	798703	75.243	ng	0.00
42) 2,4,6-Tribromophenol	10.728	330	387013	120.448	ng	0.01
45) 2-Fluorobiphenyl	9.245	172	1477734	71.096	ng	0.00
79) Terphenyl-d14	13.022	244	1759884	70.156	ng	0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.769	88	113749	25.502	ng	98
3) Pyridine	3.516	79	286326	25.898	ng	99
4) n-Nitrosodimethylamine	3.475	42	195354	32.887	ng	92
6) Aniline	6.551	93	385036	33.344	ng	99
8) 2-Chlorophenol	6.675	128	355629	36.363	ng	98
9) Benzaldehyde	6.440	77	62537	8.972	ng	98
10) Phenol	6.534	94	440337m	34.472	ng	
11) bis(2-Chloroethyl)ether	6.622	93	342938	34.734	ng	99
12) 1,3-Dichlorobenzene	6.828	146	378473	33.712	ng	99
13) 1,4-Dichlorobenzene	6.904	146	381982	34.057	ng	99
14) 1,2-Dichlorobenzene	7.057	146	367763	35.132	ng	99
15) Benzyl Alcohol	7.028	79	341566	37.581	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.157	45	565941	33.617	ng	96
17) 2-Methylphenol	7.134	107	298820	35.832	ng	97
18) Hexachloroethane	7.398	117	141896	35.477	ng	97
19) n-Nitroso-di-n-propyla...	7.298	70	265745	35.363	ng	99
20) 3+4-Methylphenols	7.287	107	369008	34.594	ng	# 85
22) Acetophenone	7.298	105	498653	34.605	ng	99
24) Nitrobenzene	7.469	77	407175	34.986	ng	99
25) Isophorone	7.710	82	732113	36.338	ng	99
26) 2-Nitrophenol	7.781	139	190854	43.469	ng	97
27) 2,4-Dimethylphenol	7.816	122	301422	41.172	ng	96
28) bis(2-Chloroethoxy)met...	7.916	93	426467	34.938	ng	99
29) 2,4-Dichlorophenol	8.022	162	308490	36.994	ng	100
30) 1,2,4-Trichlorobenzene	8.110	180	314619	34.098	ng	99
31) Naphthalene	8.192	128	1061127	34.972	ng	100
32) Benzoic acid	7.940	122	229900	36.004	ng	99
33) 4-Chloroaniline	8.234	127	209860	20.284	ng	100
34) Hexachlorobutadiene	8.304	225	208389	35.945	ng	99
35) Caprolactam	8.616	113	96843m	36.524	ng	
36) 4-Chloro-3-methylphenol	8.722	107	350066	37.913	ng	99
37) 2-Methylnaphthalene	8.881	142	675332	36.342	ng	99
38) 1-Methylnaphthalene	8.981	142	641523	35.187	ng	100
40) 1,2,4,5-Tetrachloroben...	9.051	216	327707	35.361	ng	98
41) Hexachlorocyclopentadiene	9.034	237	419426	130.069	ng	100
43) 2,4,6-Trichlorophenol	9.163	196	250539	38.185	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	240421	35.516	ng	98
46) 1,1'-Biphenyl	9.351	154	835336	34.931	ng	99
47) 2-Chloronaphthalene	9.375	162	665137	34.534	ng	99
48) 2-Nitroaniline	9.469	65	243003	41.252	ng	100
49) Acenaphthylene	9.792	152	1043569	37.478	ng	99
50) Dimethylphthalate	9.657	163	806826	37.708	ng	100
51) 2,6-Dinitrotoluene	9.716	165	179607	38.768	ng	99
52) Acenaphthene	9.969	154	755838	41.962	ng	98
53) 3-Nitroaniline	9.886	138	134444	28.141	ng	97
54) 2,4-Dinitrophenol	9.998	184	219749	111.377	ng #	1
55) Dibenzofuran	10.139	168	926029	35.832	ng	98
56) 4-Nitrophenol	10.051	139	297191	80.854	ng	92
57) 2,4-Dinitrotoluene	10.128	165	244383	42.896	ng	95
58) Fluorene	10.486	166	727555	36.962	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.257	232	205849	38.790	ng	99
60) Diethylphthalate	10.363	149	795973	37.888	ng	100
61) 4-Chlorophenyl-phenyle...	10.475	204	364593	36.677	ng	98
62) 4-Nitroaniline	10.510	138	193417	42.865	ng	95
63) Azobenzene	10.639	77	809448	36.343	ng	98
65) 4,6-Dinitro-2-methylph...	10.539	198	146695	56.353	ng	95
66) n-Nitrosodiphenylamine	10.598	169	678341	35.514	ng	99
67) 4-Bromophenyl-phenylether	10.969	248	236140	35.917	ng	98
68) Hexachlorobenzene	11.033	284	265113	35.854	ng	99
69) Atrazine	11.128	200	233794	45.185	ng	98
70) Pentachlorophenol	11.233	266	331054	73.771	ng	99
71) Phenanthrene	11.457	178	1087142	36.063	ng	100
72) Anthracene	11.504	178	1108125	37.675	ng	100
73) Carbazole	11.657	167	993450	36.318	ng	99
74) Di-n-butylphthalate	11.986	149	1253855	39.675	ng	100
75) Fluoranthene	12.645	202	1159597	38.051	ng	99
77) Benzidine	12.769	184	439202	65.344	ng	99
78) Pyrene	12.874	202	1177466	32.908	ng	99
80) Butylbenzylphthalate	13.498	149	491798	45.847	ng	99
81) Benzo(a)anthracene	14.080	228	955949	35.926	ng	100
82) 3,3'-Dichlorobenzidine	14.039	252	209783	27.040	ng	100
83) Chrysene	14.121	228	917850	37.621	ng	99
84) Bis(2-ethylhexyl)phtha...	14.063	149	652028	54.177	ng	98
85) Di-n-octyl phthalate	14.710	149	1017141	46.465	ng	98
87) Indeno(1,2,3-cd)pyrene	17.157	276	730882	33.027	ng	97
88) Benzo(b)fluoranthene	15.174	252	752340	35.904	ng	99
89) Benzo(k)fluoranthene	15.204	252	718371	39.752	ng	99
90) Benzo(a)pyrene	15.551	252	667766	38.729	ng	99
91) Dibenzo(a,h)anthracene	17.180	278	599184	32.457	ng	98
92) Benzo(g,h,i)perylene	17.627	276	549722	29.811	ng	97

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

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