

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100224\
 Data File : BF139599.D
 Acq On : 02 Oct 2024 11:39
 Operator : RC/JU
 Sample : PB163811BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163811BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/03/2024
 Supervised By :mohammad ahmed 10/04/2024

Quant Time: Oct 02 12:00:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	92492	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	356514	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	207013	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	398612	20.000	ng	0.00
76) Chrysene-d12	14.033	240	227777	20.000	ng	0.00
86) Perylene-d12	15.498	264	194700	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	815381	152.798	ng	0.01
7) Phenol-d6	6.504	99	1101130	153.441	ng	0.00
23) Nitrobenzene-d5	7.433	82	728642	103.478	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	312088	156.171	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1377531	102.156	ng	0.00
79) Terphenyl-d14	12.980	244	1554708	111.996	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.751	88	105251	45.620	ng	99
3) Pyridine	3.493	79	279862	49.877	ng	99
4) n-Nitrosodimethylamine	3.451	42	191756	52.210	ng	98
6) Aniline	6.528	93	346812	54.462	ng	# 77
8) 2-Chlorophenol	6.651	128	312794	54.710	ng	99
9) Benzaldehyde	6.416	77	116880	30.746	ng	98
10) Phenol	6.516	94	407504	54.004	ng	91
11) bis(2-Chloroethyl)ether	6.604	93	311303	50.669	ng	100
12) 1,3-Dichlorobenzene	6.810	146	321112	49.926	ng	99
13) 1,4-Dichlorobenzene	6.886	146	324644	49.837	ng	99
14) 1,2-Dichlorobenzene	7.033	146	314060	51.442	ng	98
15) Benzyl Alcohol	7.010	79	301094	54.913	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.139	45	568260	53.427	ng	95
17) 2-Methylphenol	7.122	107	258160	54.088	ng	99
18) Hexachloroethane	7.375	117	122013	50.217	ng	97
19) n-Nitroso-di-n-propyla...	7.281	70	241474	52.339	ng	98
20) 3+4-Methylphenols	7.275	107	319396	53.277	ng	94
22) Acetophenone	7.275	105	434744	51.373	ng	99
24) Nitrobenzene	7.451	77	364089	49.933	ng	99
25) Isophorone	7.692	82	655389	52.404	ng	100
26) 2-Nitrophenol	7.769	139	166672	53.172	ng	98
27) 2,4-Dimethylphenol	7.804	122	248516m	64.762	ng	
28) bis(2-Chloroethoxy)met...	7.898	93	389118	52.104	ng	99
29) 2,4-Dichlorophenol	8.010	162	267352	53.414	ng	100
30) 1,2,4-Trichlorobenzene	8.092	180	275182	48.615	ng	100
31) Naphthalene	8.175	128	923100	50.641	ng	100
32) Benzoic acid	7.939	122	197201	51.682	ng	98
33) 4-Chloroaniline	8.222	127	162710	26.371	ng	98
34) Hexachlorobutadiene	8.286	225	178674	48.777	ng	99
35) Caprolactam	8.598	113	87675m	53.406	ng	
36) 4-Chloro-3-methylphenol	8.710	107	302151	53.327	ng	99
37) 2-Methylnaphthalene	8.863	142	610767	51.447	ng	99
38) 1-Methylnaphthalene	8.963	142	564440	48.640	ng	100
40) 1,2,4,5-Tetrachloroben...	9.027	216	285529	49.670	ng	99
41) Hexachlorocyclopentadiene	9.016	237	352043	199.726	ng	99
43) 2,4,6-Trichlorophenol	9.145	196	203284	52.458	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	214767	51.391	ng	98
46) 1,1'-Biphenyl	9.327	154	772476	50.171	ng	100
47) 2-Chloronaphthalene	9.351	162	575365	49.847	ng	99
48) 2-Nitroaniline	9.451	65	221879	53.549	ng	99
49) Acenaphthylene	9.769	152	958517	54.604	ng	100
50) Dimethylphthalate	9.633	163	716976	52.908	ng	100
51) 2,6-Dinitrotoluene	9.692	165	158053	51.631	ng	98
52) Acenaphthene	9.939	154	685524m	56.887	ng	
53) 3-Nitroaniline	9.863	138	110197	35.226	ng	98
54) 2,4-Dinitrophenol	9.974	184	187318	113.768	ng	99
55) Dibenzofuran	10.116	168	870342	51.583	ng	100
56) 4-Nitrophenol	10.033	139	255739	107.203	ng	98
57) 2,4-Dinitrotoluene	10.098	165	214697	53.657	ng	98
58) Fluorene	10.457	166	684464	50.932	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.233	232	183171	54.267	ng	100
60) Diethylphthalate	10.333	149	744276	53.174	ng	100
61) 4-Chlorophenyl-phenyle...	10.445	204	336638	50.071	ng	99
62) 4-Nitroaniline	10.480	138	166971	52.181	ng	99
63) Azobenzene	10.610	77	750762	53.336	ng	100
65) 4,6-Dinitro-2-methylph...	10.510	198	120351	53.459	ng	97
66) n-Nitrosodiphenylamine	10.569	169	615392	52.375	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	207315	50.726	ng	99
68) Hexachlorobenzene	11.004	284	228097	50.296	ng	98
69) Atrazine	11.098	200	202789	64.327	ng	99
70) Pentachlorophenol	11.198	266	262609	97.247	ng	99
71) Phenanthrene	11.421	178	990570	52.704	ng	100
72) Anthracene	11.468	178	1013847	54.526	ng	100
73) Carbazole	11.627	167	924076	51.744	ng	100
74) Di-n-butylphthalate	11.951	149	1167404	53.757	ng	100
75) Fluoranthene	12.604	202	1082500	52.688	ng	100
77) Benzidine	12.727	184	432163	107.564	ng	99
78) Pyrene	12.839	202	1085535	53.285	ng	99
80) Butylbenzylphthalate	13.451	149	421640	57.862	ng	100
81) Benzo(a)anthracene	14.021	228	837383	55.917	ng	100
82) 3,3'-Dichlorobenzidine	13.986	252	172924	37.730	ng	99
83) Chrysene	14.062	228	697272	50.928	ng	100
84) Bis(2-ethylhexyl)phtha...	14.009	149	541756	59.539	ng	99
85) Di-n-octyl phthalate	14.621	149	774758	57.957	ng	99
87) Indeno(1,2,3-cd)pyrene	16.980	276	668005	58.292	ng	99
88) Benzo(b)fluoranthene	15.074	252	622970	49.485	ng	100
89) Benzo(k)fluoranthene	15.104	252	611708	57.282	ng	100
90) Benzo(a)pyrene	15.439	252	576420	58.069	ng	99
91) Dibenzo(a,h)anthracene	16.998	278	558324	58.742	ng	99
92) Benzo(g,h,i)perylene	17.427	276	509666	53.467	ng	100

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

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