

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100322\
 Data File : BF130495.D
 Acq On : 03 Oct 2022 13:21
 Operator : CG\JU
 Sample : PB148034BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB148034BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/04/2022
 Supervised By :mohammad ahmed 10/06/2022

Quant Time: Oct 04 03:13:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.640	152	88024	20.000	ng	0.00
21) Naphthalene-d8	7.922	136	355837	20.000	ng	0.00
39) Acenaphthene-d10	9.675	164	186066	20.000	ng	0.00
64) Phenanthrene-d10	11.151	188	301167	20.000	ng	0.00
76) Chrysene-d12	13.780	240	141761	20.000	ng	-0.01
86) Perylene-d12	15.169	264	132297	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.252	112	732457	144.327	ng	0.00
7) Phenol-d6	6.293	99	909046	143.157	ng	0.00
23) Nitrobenzene-d5	7.210	82	608394	87.349	ng	0.00
42) 2,4,6-Tribromophenol	10.463	330	235239	131.944	ng	0.00
45) 2-Fluorobiphenyl	8.998	172	1063648	83.243	ng	-0.01
79) Terphenyl-d14	12.739	244	904662	105.711	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.316	88	149456	64.123	ng	94
3) Pyridine	2.993	79	284531	46.798	ng	99
4) n-Nitrosodimethylamine	2.958	42	191552	57.926	ng	99
6) Aniline	6.304	93	338633	43.281	ng	# 34
8) 2-Chlorophenol	6.422	128	271280	46.925	ng	98
9) Benzaldehyde	6.187	77	128212	30.143	ng	99
10) Phenol	6.304	94	348070	46.774	ng	# 59
11) bis(2-Chloroethyl)ether	6.381	93	256211	47.734	ng	99
12) 1,3-Dichlorobenzene	6.575	146	286350	45.016	ng	97
13) 1,4-Dichlorobenzene	6.657	146	292144	45.036	ng	99
14) 1,2-Dichlorobenzene	6.804	146	275959	45.540	ng	98
15) Benzyl Alcohol	6.793	79	224286	44.548	ng	99
16) 2,2'-oxybis(1-Chloropr...	6.922	45	380228	48.569	ng	92
17) 2-Methylphenol	6.910	107	214565	45.657	ng	# 93
18) Hexachloroethane	7.145	117	114295	44.699	ng	96
19) n-Nitroso-di-n-propyla...	7.063	70	208843	48.740	ng	99
20) 3+4-Methylphenols	7.063	107	280717	45.982	ng	# 76
22) Acetophenone	7.057	105	407533	46.844	ng	99
24) Nitrobenzene	7.228	77	318763	45.072	ng	97
25) Isophorone	7.469	82	547998	48.815	ng	98
26) 2-Nitrophenol	7.545	139	125914	43.431	ng	95
27) 2,4-Dimethylphenol	7.587	122	223145	49.988	ng	98
28) bis(2-Chloroethoxy)met...	7.681	93	318906	50.450	ng	98
29) 2,4-Dichlorophenol	7.787	162	211158	43.166	ng	98
30) 1,2,4-Trichlorobenzene	7.863	180	216792	40.991	ng	94
31) Naphthalene	7.945	128	781731	42.642	ng	99
32) Benzoic acid	7.745	122	110812	32.100	ng	98
33) 4-Chloroaniline	7.998	127	262602	37.663	ng	96
34) Hexachlorobutadiene	8.057	225	139421	41.248	ng	99
35) Caprolactam	8.381	113	66821m	47.648	ng	
36) 4-Chloro-3-methylphenol	8.492	107	253037	46.667	ng	94
37) 2-Methylnaphthalene	8.634	142	509315	43.574	ng	99
38) 1-Methylnaphthalene	8.734	142	488566	43.511	ng	98
40) 1,2,4,5-Tetrachloroben...	8.798	216	224580	44.227	ng	98
41) Hexachlorocyclopentadiene	8.787	237	220673	81.171	ng	99
43) 2,4,6-Trichlorophenol	8.916	196	140711	39.706	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.963	196	150033	39.232	ng	# 91
46) 1,1'-Biphenyl	9.098	154	619518	44.582	ng	99
47) 2-Chloronaphthalene	9.122	162	476216	42.364	ng	99
48) 2-Nitroaniline	9.228	65	167416	44.653	ng	97
49) Acenaphthylene	9.534	152	741442	43.991	ng	99
50) Dimethylphthalate	9.410	163	559191	43.508	ng	100
51) 2,6-Dinitrotoluene	9.469	165	121901	45.350	ng	95
52) Acenaphthene	9.704	154	441894	39.904	ng	99
53) 3-Nitroaniline	9.639	138	104625	34.931	ng	99
54) 2,4-Dinitrophenol	9.751	184	95776	66.803	ng	98
55) Dibenzofuran	9.881	168	664662	43.152	ng	99
56) 4-Nitrophenol	9.816	139	148192	67.930	ng	99
57) 2,4-Dinitrotoluene	9.875	165	160925	46.845	ng	# 88
58) Fluorene	10.222	166	545213	43.282	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.998	232	116387	38.883	ng	99
60) Diethylphthalate	10.104	149	568681	44.122	ng	99
61) 4-Chlorophenyl-phenyle...	10.216	204	249239	45.103	ng	99
62) 4-Nitroaniline	10.257	138	120593m	40.079	ng	
63) Azobenzene	10.375	77	589452	44.414	ng	98
65) 4,6-Dinitro-2-methylph...	10.281	198	65076	38.817	ng	95
66) n-Nitrosodiphenylamine	10.339	169	448028	44.518	ng	99
67) 4-Bromophenyl-phenylether	10.704	248	151153	45.496	ng	99
68) Hexachlorobenzene	10.763	284	169552	45.020	ng	93
69) Atrazine	10.869	200	135318	46.009	ng	99
70) Pentachlorophenol	10.963	266	137947	68.249	ng	99
71) Phenanthrene	11.181	178	723628	42.796	ng	100
72) Anthracene	11.228	178	735530	45.077	ng	99
73) Carbazole	11.392	167	636322	43.210	ng	98
74) Di-n-butylphthalate	11.722	149	778559	43.843	ng	100
75) Fluoranthene	12.363	202	648737	39.705	ng	97
77) Benzidine	12.492	184	199619	48.508	ng	99
78) Pyrene	12.586	202	637699	51.422	ng	99
80) Butylbenzylphthalate	13.216	149	208026	45.280	ng	98
81) Benzo(a)anthracene	13.774	228	407890	43.252	ng	99
82) 3,3'-Dichlorobenzidine	13.745	252	122047	47.542	ng	98
83) Chrysene	13.810	228	383916	42.875	ng	99
84) Bis(2-ethylhexyl)phtha...	13.774	149	250840	47.667	ng	100
85) Di-n-octyl phthalate	14.386	149	365128	45.364	ng	100
87) Indeno(1,2,3-cd)pyrene	16.498	276	465230	49.589	ng	99
88) Benzo(b)fluoranthene	14.786	252	389870	45.150	ng	98
89) Benzo(k)fluoranthene	14.810	252	368882	43.428	ng	100
90) Benzo(a)pyrene	15.116	252	336965	49.257	ng	98
91) Dibenzo(a,h)anthracene	16.510	278	409767	53.242	ng	99
92) Benzo(g,h,i)perylene	16.892	276	382622	49.352	ng	99

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

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