

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091423\
 Data File : BF135268.D
 Acq On : 14 Sep 2023 22:46
 Operator : CG\JU
 Sample : 04283-02MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-2MS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 09/15/2023
 Supervised By :mohammad ahmed 09/15/2023

Quant Time: Sep 15 01:38:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 00:19:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.763	152	119974	20.000	ng	0.00
21) Naphthalene-d8	8.045	136	457701	20.000	ng	0.00
39) Acenaphthene-d10	9.804	164	213284	20.000	ng	0.00
64) Phenanthrene-d10	11.292	188	340011	20.000	ng	0.00
76) Chrysene-d12	13.951	240	209162	20.000	ng	0.00
86) Perylene-d12	15.415	264	223694	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	722507	97.948	ng	0.02
7) Phenol-d6	6.416	99	913017	97.341	ng	0.01
23) Nitrobenzene-d5	7.334	82	573354	68.057	ng	0.00
42) 2,4,6-Tribromophenol	10.598	330	179968	84.910	ng	0.00
45) 2-Fluorobiphenyl	9.122	172	1004536	71.059	ng	0.00
79) Terphenyl-d14	12.886	244	843557	62.520	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.575	88	141225	43.673	ng	99
3) Pyridine	3.328	79	349614	40.171	ng	97
4) n-Nitrosodimethylamine	3.299	42	228022	49.373	ng	95
6) Aniline	6.434	93	419490	34.106	ng	# 82
8) 2-Chlorophenol	6.551	128	402740	51.145	ng	97
9) Benzaldehyde	6.316	77	149415	27.963	ng	96
10) Phenol	6.428	94	463344	45.050	ng	81
11) bis(2-Chloroethyl)ether	6.504	93	380342	49.299	ng	97
12) 1,3-Dichlorobenzene	6.704	146	406741	50.826	ng	99
13) 1,4-Dichlorobenzene	6.781	146	413731	50.815	ng	98
14) 1,2-Dichlorobenzene	6.934	146	388394	51.368	ng	98
15) Benzyl Alcohol	6.916	79	347576	51.641	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.039	45	709439	48.785	ng	90
17) 2-Methylphenol	7.028	107	309949	44.604	ng	96
18) Hexachloroethane	7.269	117	151371	50.317	ng	93
19) n-Nitroso-di-n-propyla...	7.186	70	264832	45.584	ng	96
20) 3+4-Methylphenols	7.181	107	362517	43.385	ng	92
22) Acetophenone	7.175	105	521018	49.790	ng	# 97
24) Nitrobenzene	7.351	77	414291	49.754	ng	98
25) Isophorone	7.592	82	765582	49.489	ng	99
26) 2-Nitrophenol	7.663	139	197415	49.398	ng	97
27) 2,4-Dimethylphenol	7.710	122	316180	47.068	ng	98
28) bis(2-Chloroethoxy)met...	7.798	93	465416	50.266	ng	100
29) 2,4-Dichlorophenol	7.910	162	305592	49.095	ng	97
30) 1,2,4-Trichlorobenzene	7.986	180	336778	51.764	ng	99
31) Naphthalene	8.069	128	1088312	48.939	ng	100
32) Benzoic acid	7.833	122	157112m	31.060	ng	
33) 4-Chloroaniline	8.122	127	323745	33.181	ng	99
34) Hexachlorobutadiene	8.181	225	201555	51.377	ng	99
35) Caprolactam	8.504	113	92552m	44.303	ng	
36) 4-Chloro-3-methylphenol	8.610	107	317554	45.902	ng	97
37) 2-Methylnaphthalene	8.757	142	690005	46.159	ng	99
38) 1-Methylnaphthalene	8.857	142	652608	46.630	ng	100
40) 1,2,4,5-Tetrachloroben...	8.922	216	330634	53.509	ng	100
41) Hexachlorocyclopentadiene	8.904	237	132403	51.526	ng	99
43) 2,4,6-Trichlorophenol	9.039	196	199621	49.342	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.086	196	203379	43.653	ng	97
46) 1,1'-Biphenyl	9.228	154	849386	51.822	ng	98
47) 2-Chloronaphthalene	9.251	162	633296	53.253	ng	98
48) 2-Nitroaniline	9.351	65	239174	51.768	ng	95
49) Acenaphthylene	9.669	152	1013377	51.274	ng	99
50) Dimethylphthalate	9.533	163	758943	52.631	ng	99
51) 2,6-Dinitrotoluene	9.598	165	164280	52.004	ng	89
52) Acenaphthene	9.839	154	582431	49.885	ng	98
53) 3-Nitroaniline	9.769	138	161471	43.545	ng	96
54) 2,4-Dinitrophenol	9.875	184	34343	23.902	ng	# 88
55) Dibenzofuran	10.010	168	855853	48.036	ng	99
56) 4-Nitrophenol	9.939	139	179419	76.132	ng	97
57) 2,4-Dinitrotoluene	10.004	165	186175	47.076	ng	91
58) Fluorene	10.357	166	633882	46.279	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.133	232	151997	42.411	ng	98
60) Diethylphthalate	10.233	149	735160	50.799	ng	100
61) 4-Chlorophenyl-phenyle...	10.345	204	308320	46.861	ng	96
62) 4-Nitroaniline	10.386	138	146185	41.012	ng	95
63) Azobenzene	10.510	77	646747	45.663	ng	99
65) 4,6-Dinitro-2-methylph...	10.416	198	30912	17.117	ng	84
66) n-Nitrosodiphenylamine	10.469	169	577376	56.090	ng	100
67) 4-Bromophenyl-phenylether	10.839	248	187950	55.579	ng	100
68) Hexachlorobenzene	10.904	284	194024	56.318	ng	98
69) Atrazine	11.004	200	168911	60.984	ng	99
70) Pentachlorophenol	11.098	266	126160	61.206	ng	96
71) Phenanthrene	11.322	178	825607	51.340	ng	99
72) Anthracene	11.369	178	847528	51.273	ng	99
73) Carbazole	11.533	167	744737	49.493	ng	99
74) Di-n-butylphthalate	11.863	149	1024765	57.795	ng	100
75) Fluoranthene	12.510	202	832163	49.663	ng	99
77) Benzidine	12.639	184	226879	52.946	ng	98
78) Pyrene	12.745	202	843249	42.345	ng	99
80) Butylbenzylphthalate	13.368	149	359333	51.250	ng	96
81) Benzo(a)anthracene	13.939	228	673138	49.839	ng	99
82) 3,3'-Dichlorobenzidine	13.904	252	220215	52.200	ng	# 99
83) Chrysene	13.980	228	655262	48.797	ng	99
84) Bis(2-ethylhexyl)phtha...	13.933	149	442443	61.494	ng	99
85) Di-n-octyl phthalate	14.551	149	795178	69.545	ng	100
87) Indeno(1,2,3-cd)pyrene	16.904	276	595045	40.571	ng	99
88) Benzo(b)fluoranthene	14.992	252	653041	53.929	ng	99
89) Benzo(k)fluoranthene	15.021	252	630108	52.676	ng	98
90) Benzo(a)pyrene	15.357	252	553508	47.932	ng	99
91) Dibenzo(a,h)anthracene	16.915	278	494890	41.239	ng	99
92) Benzo(g,h,i)perylene	17.345	276	464353	37.050	ng	100

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

