

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061322\
 Data File : BM035459.D
 Acq On : 13 Jun 2022 16:11
 Operator : CG/JU
 Sample : N3331-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 B-220(26-28)MS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 06/14/2022
 Supervised By : Jagrut Upadhyay 06/14/2022

Quant Time: Jun 14 04:29:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM061022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 10 16:05:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	226483	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	934708	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	543563	20.000	ng	0.00
64) Phenanthrene-d10	17.209	188	1095319	20.000	ng	# 0.00
76) Chrysene-d12	21.397	240	1132797	20.000	ng	0.00
86) Perylene-d12	23.738	264	1230415	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	2002709	149.952	ng	0.00
7) Phenol-d6	7.004	99	2687282	145.197	ng	0.00
23) Nitrobenzene-d5	8.981	82	1578020	91.745	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	869983	150.331	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	3453770	90.773	ng	0.00
79) Terphenyl-d14	19.833	244	5005356	84.366	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.299	88	243400	47.781	ng	95
3) Pyridine	3.704	79	666344	46.771	ng	95
4) n-Nitrosodimethylamine	3.616	42	388312	59.051	ng	84
6) Aniline	7.145	93	1149837	57.477	ng	100
8) 2-Chlorophenol	7.387	128	872703	56.923	ng	98
9) Benzaldehyde	6.963	77	431574	46.450	ng	97
10) Phenol	7.034	94	1068061	59.405	ng	93
11) bis(2-Chloroethyl)ether	7.245	93	777834	53.974	ng	100
12) 1,3-Dichlorobenzene	7.704	146	873530	53.411	ng	98
13) 1,4-Dichlorobenzene	7.851	146	890821	53.205	ng	100
14) 1,2-Dichlorobenzene	8.169	146	853567	52.232	ng	99
15) Benzyl Alcohol	8.069	79	707753m	57.083	ng	
16) 2,2'-oxybis(1-Chloropr...	8.345	45	1236291	49.394	ng	97
17) 2-Methylphenol	8.281	107	713247	57.809	ng	95
18) Hexachloroethane	8.892	117	319181	53.154	ng	95
19) n-Nitroso-di-n-propyla...	8.633	70	610021	52.437	ng	96
20) 3+4-Methylphenols	8.610	107	972319	56.219	ng	96
22) Acetophenone	8.645	105	1227573	54.224	ng	# 99
24) Nitrobenzene	9.022	77	893943	56.763	ng	99
25) Isophorone	9.557	82	1627201	58.172	ng	97
26) 2-Nitrophenol	9.733	139	463535	57.512	ng	92
27) 2,4-Dimethylphenol	9.804	122	800748	67.799	ng	100
28) bis(2-Chloroethoxy)met...	10.033	93	1026232	52.728	ng	99
29) 2,4-Dichlorophenol	10.280	162	764168	58.061	ng	98
30) 1,2,4-Trichlorobenzene	10.480	180	760752	51.993	ng	99
31) Naphthalene	10.663	128	2514653	53.711	ng	100
32) Benzoic acid	9.998	122	632451	56.457	ng	97
33) 4-Chloroaniline	10.786	127	892686	46.863	ng	98
34) Hexachlorobutadiene	10.951	225	414464	50.334	ng	98
35) Caprolactam	11.598	113	268179m	60.114	ng	
36) 4-Chloro-3-methylphenol	11.927	107	844931	56.626	ng	99
37) 2-Methylnaphthalene	12.280	142	1722484	53.424	ng	98
38) 1-Methylnaphthalene	12.498	142	1658921	52.643	ng	100
40) 1,2,4,5-Tetrachloroben...	12.657	216	799315	52.176	ng	99
41) Hexachlorocyclopentadiene	12.633	237	705686	113.694	ng	99
43) 2,4,6-Trichlorophenol	12.904	196	581067	55.296	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.980	196	640322	56.484	ng	97
46) 1,1'-Biphenyl	13.298	154	2231618	52.715	ng	98
47) 2-Chloronaphthalene	13.339	162	1699087	53.367	ng	98
48) 2-Nitroaniline	13.551	65	572145	59.365	ng	98
49) Acenaphthylene	14.180	152	2892513	58.000	ng	99
50) Dimethylphthalate	13.927	163	2127086	53.178	ng	99
51) 2,6-Dinitrotoluene	14.045	165	489553	58.466	ng	94
52) Acenaphthene	14.527	154	1628626	56.064	ng	99
53) 3-Nitroaniline	14.380	138	438952	48.848	ng	97
54) 2,4-Dinitrophenol	14.598	184	555357	103.174	ng	90
55) Dibenzofuran	14.863	168	2554469	54.886	ng	97
56) 4-Nitrophenol	14.710	139	908794	124.741	ng	89
57) 2,4-Dinitrotoluene	14.839	165	667586	62.783	ng	98
58) Fluorene	15.510	166	2066856	56.846	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.098	232	519746	57.651	ng	98
60) Diethylphthalate	15.292	149	2178790	55.659	ng	99
61) 4-Chlorophenyl-phenyle...	15.504	204	954698	54.263	ng	99
62) 4-Nitroaniline	15.545	138	552143	63.488	ng	89
63) Azobenzene	15.798	77	2035357	54.895	ng	98
65) 4,6-Dinitro-2-methylph...	15.610	198	359560	52.138	ng	94
66) n-Nitrosodiphenylamine	15.721	169	1819269	55.776	ng	99
67) 4-Bromophenyl-phenylether	16.398	248	595946	52.731	ng	98
68) Hexachlorobenzene	16.521	284	667140	53.993	ng	94
69) Atrazine	16.680	200	624281	63.048	ng	99
70) Pentachlorophenol	16.868	266	863685	132.750	ng	99
71) Phenanthrene	17.251	178	3183791	54.769	ng	99
72) Anthracene	17.345	178	3257133	57.425	ng	100
73) Carbazole	17.615	167	3129798	56.909	ng	99
74) Di-n-butylphthalate	18.180	149	3835268	58.305	ng	99
75) Fluoranthene	19.268	202	3732193	61.827	ng	97
77) Benzidine	19.456	184	2459663	118.117	ng	99
78) Pyrene	19.633	202	3865870	48.621	ng	100
80) Butylbenzylphthalate	20.533	149	1825982	51.844	ng	97
81) Benzo(a)anthracene	21.380	228	3887922	55.092	ng	99
82) 3,3'-Dichlorobenzidine	21.315	252	1225917	75.456	ng	99
83) Chrysene	21.433	228	3628405	53.229	ng	100
84) Bis(2-ethylhexyl)phtha...	21.309	149	2706949	53.588	ng	100
85) Di-n-octyl phthalate	22.215	149	4834485	58.907	ng	99
87) Indeno(1,2,3-cd)pyrene	26.174	276	4476622	52.666	ng #	94
88) Benzo(b)fluoranthene	23.027	252	4032636	55.642	ng #	96
89) Benzo(k)fluoranthene	23.074	252	3998362	54.495	ng #	97
90) Benzo(a)pyrene	23.638	252	3930463	64.280	ng #	97
91) Dibenzo(a,h)anthracene	26.179	278	3893463	55.369	ng	97
92) Benzo(g,h,i)perylene	26.909	276	3930313	56.191	ng #	95

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

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