

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061522\
 Data File : BP010733.D
 Acq On : 15 Jun 2022 17:15
 Operator : CG/JU
 Sample : N3139-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COMP-2MS

Quant Time: Jun 16 00:05:31 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP060722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 16 00:00:39 2022
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 06/16/2022
 Supervised By :Yogesh Patel 06/16/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	64095	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	292586	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	232625	20.000	ng	0.00
64) Phenanthrene-d10	17.222	188	586638	20.000	ng	0.00
76) Chrysene-d12	21.316	240	708170	20.000	ng	# 0.00
86) Perylene-d12	23.710	264	790344	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.417	112	491728	132.408	ng	0.00
7) Phenol-d6	7.011	99	711481	140.086	ng	0.00
23) Nitrobenzene-d5	8.993	82	484951	78.054	ng	0.00
42) 2,4,6-Tribromophenol	15.969	330	516199	162.951	ng	0.00
45) 2-Fluorobiphenyl	13.093	172	1223946	75.689	ng	0.00
79) Terphenyl-d14	19.786	244	2790870	73.692	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.317	88	52416	35.114	ng	99
3) Pyridine	3.723	79	150189	35.584	ng	98
4) n-Nitrosodimethylamine	3.628	42	121152	51.440	ng	# 95
6) Aniline	7.158	93	286965	48.329	ng	97
8) 2-Chlorophenol	7.399	128	225567	55.939	ng	99
9) Benzaldehyde	6.975	77	123146	40.126	ng	99
10) Phenol	7.034	94	299445	57.147	ng	94
11) bis(2-Chloroethyl)ether	7.258	93	195730	48.611	ng	96
12) 1,3-Dichlorobenzene	7.716	146	221475	47.981	ng	99
13) 1,4-Dichlorobenzene	7.863	146	228206	48.313	ng	98
14) 1,2-Dichlorobenzene	8.181	146	222776	48.993	ng	98
15) Benzyl Alcohol	8.075	79	237685m	58.995	ng	
16) 2,2'-oxybis(1-Chloropr...	8.358	45	342005	50.529	ng	99
17) 2-Methylphenol	8.281	107	204497	57.797	ng	97
18) Hexachloroethane	8.905	117	86614	47.254	ng	89
19) n-Nitroso-di-n-propyla...	8.640	70	194296	54.245	ng	97
20) 3+4-Methylphenols	8.611	107	289687	59.648	ng	99
22) Acetophenone	8.658	105	364223	46.709	ng	# 98
24) Nitrobenzene	9.034	77	291955	46.543	ng	99
25) Isophorone	9.563	82	542840	51.054	ng	97
26) 2-Nitrophenol	9.746	139	130962	50.800	ng	97
27) 2,4-Dimethylphenol	9.810	122	241475	63.895	ng	99
28) bis(2-Chloroethoxy)met...	10.040	93	301866	52.713	ng	98
29) 2,4-Dichlorophenol	10.287	162	259043	58.197	ng	96
30) 1,2,4-Trichlorobenzene	10.493	180	249266	48.451	ng	98
31) Naphthalene	10.675	128	717855	45.978	ng	99
32) Benzoic acid	9.981	122	185766	58.200	ng	97
33) 4-Chloroaniline	10.799	127	160290	26.137	ng	98
34) Hexachlorobutadiene	10.963	225	160590	46.005	ng	97
35) Caprolactam	11.604	113	96908m	63.917	ng	
36) 4-Chloro-3-methylphenol	11.934	107	322157	60.629	ng	97
37) 2-Methylnaphthalene	12.287	142	557718	51.231	ng	98
38) 1-Methylnaphthalene	12.510	142	542944	50.687	ng	97
40) 1,2,4,5-Tetrachloroben...	12.663	216	335704	46.910	ng	99
41) Hexachlorocyclopentadiene	12.640	237	378947	105.515	ng	99
43) 2,4,6-Trichlorophenol	12.904	196	246355	53.058	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.981	196	276135	53.468	ng	98
46) 1,1'-Biphenyl	13.304	154	803537	46.837	ng	98
47) 2-Chloronaphthalene	13.346	162	616498	45.925	ng	100
48) 2-Nitroaniline	13.551	65	252557	50.765	ng	97
49) Acenaphthylene	14.193	152	1067251	50.148	ng	100
50) Dimethylphthalate	13.934	163	893330	50.515	ng	99
51) 2,6-Dinitrotoluene	14.051	165	199360	52.890	ng	99
52) Acenaphthene	14.534	154	607527	47.077	ng	98
53) 3-Nitroaniline	14.381	138	150919	39.122	ng	98
54) 2,4-Dinitrophenol	14.598	184	287054	115.227	ng	97
55) Dibenzofuran	14.869	168	1044994	50.369	ng	99
56) 4-Nitrophenol	14.710	139	361717	125.293	ng	93
57) 2,4-Dinitrotoluene	14.845	165	287604	54.825	ng	94
58) Fluorene	15.522	166	881777	51.007	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.104	232	260596	55.412	ng	96
60) Diethylphthalate	15.298	149	955657	52.781	ng	100
61) 4-Chlorophenyl-phenyle...	15.516	204	470175	53.582	ng	94
62) 4-Nitroaniline	15.551	138	212100	53.582	ng	94
63) Azobenzene	15.810	77	889131	48.958	ng	96
65) 4,6-Dinitro-2-methylph...	15.616	198	184278	50.059	ng	93
66) n-Nitrosodiphenylamine	15.734	169	787859	45.987	ng	99
67) 4-Bromophenyl-phenylether	16.410	248	314235	49.208	ng	97
68) Hexachlorobenzene	16.534	284	372764	50.478	ng	# 68
69) Atrazine	16.692	200	331721	54.444	ng	99
70) Pentachlorophenol	16.881	266	508682	126.404	ng	96
71) Phenanthrene	17.269	178	1599957	49.603	ng	100
72) Anthracene	17.357	178	1547782	49.626	ng	99
73) Carbazole	17.634	167	1424801	51.184	ng	99
74) Di-n-butylphthalate	18.181	149	1801975	49.356	ng	99
75) Fluoranthene	19.239	202	2152762	56.395	ng	98
77) Benzidine	19.422	184	1086696	78.526	ng	99
78) Pyrene	19.592	202	2180607	45.641	ng	99
80) Butylbenzylphthalate	20.463	149	906228	44.248	ng	93
81) Benzo(a)anthracene	21.298	228	2349887	49.910	ng	99
82) 3,3'-Dichlorobenzidine	21.233	252	682083	49.637	ng	# 98
83) Chrysene	21.351	228	2161331	48.074	ng	99
84) Bis(2-ethylhexyl)phtha...	21.222	149	1393385	47.905	ng	98
85) Di-n-octyl phthalate	22.139	149	2356571	48.214	ng	99
87) Indeno(1,2,3-cd)pyrene	26.221	276	2879104	50.064	ng	# 95
88) Benzo(b)fluoranthene	22.986	252	2618024	53.085	ng	# 98
89) Benzo(k)fluoranthene	23.033	252	2347219	46.697	ng	# 98
90) Benzo(a)pyrene	23.610	252	2414612	58.484	ng	# 97
91) Dibenzo(a,h)anthracene	26.227	278	2522638	52.444	ng	# 97
92) Benzo(g,h,i)perylene	26.980	276	2474137	51.893	ng	# 95

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

