

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090922\
 Data File : BG054900.D
 Acq On : 9 Sep 2022 17:00
 Operator : CG/JU
 Sample : N4528-01MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PIT-10MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 09/12/2022
 Supervised By : Jagrut Upadhyay 09/12/2022

Quant Time: Sep 10 01:59:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 10 01:53:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.144	152	33719	20.000	ng	0.00
21) Naphthalene-d8	10.970	136	142424	20.000	ng	0.00
39) Acenaphthene-d10	14.783	164	104987	20.000	ng	0.00
64) Phenanthrene-d10	17.527	188	251278	20.000	ng	0.00
76) Chrysene-d12	21.828	240	241576	20.000	ng	0.00
86) Perylene-d12	25.206	264	257844	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.688	112	260122	134.336	ng	0.00
7) Phenol-d6	7.304	99	335639	126.746	ng	0.00
23) Nitrobenzene-d5	9.319	82	225916	83.271	ng	0.00
42) 2,4,6-Tribromophenol	16.270	330	201245	153.402	ng	0.00
45) 2-Fluorobiphenyl	13.403	172	592522	84.827	ng	0.00
79) Terphenyl-d14	20.130	244	1122202	92.042	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.526	88	32177	34.800	ng	# 73
3) Pyridine	3.943	79	77336	29.987	ng	92
4) n-Nitrosodimethylamine	3.849	42	40369	36.210	ng	93
6) Aniline	7.463	93	100065	31.804	ng	96
8) 2-Chlorophenol	7.709	128	107305	50.721	ng	96
9) Benzaldehyde	7.275	77	47249m	27.223	ng	
10) Phenol	7.333	94	119377	43.835	ng	94
11) bis(2-Chloroethyl)ether	7.557	93	97286	44.418	ng	97
12) 1,3-Dichlorobenzene	8.033	146	104669	42.753	ng	98
13) 1,4-Dichlorobenzene	8.179	146	106740	43.224	ng	98
14) 1,2-Dichlorobenzene	8.503	146	105494	45.052	ng	98
15) Benzyl Alcohol	8.385	79	93553m	48.901	ng	
16) 2,2'-oxybis(1-Chloropr...	8.667	45	129680	37.839	ng	95
17) 2-Methylphenol	8.591	107	94099	50.186	ng	96
18) Hexachloroethane	9.237	117	36833	41.132	ng	93
19) n-Nitroso-di-n-propyla...	8.949	70	83795	46.015	ng	94
20) 3+4-Methylphenols	8.920	107	129243	49.708	ng	96
22) Acetophenone	8.973	105	165797	46.531	ng	# 95
24) Nitrobenzene	9.366	77	116642	42.655	ng	94
25) Isophorone	9.883	82	245052	48.442	ng	98
26) 2-Nitrophenol	10.077	139	62462	51.710	ng	99
27) 2,4-Dimethylphenol	10.130	122	109749	57.345	ng	98
28) bis(2-Chloroethoxy)met...	10.359	93	143016	49.615	ng	99
29) 2,4-Dichlorophenol	10.618	162	114895	52.781	ng	96
30) 1,2,4-Trichlorobenzene	10.829	180	108710	43.700	ng	98
31) Naphthalene	11.023	128	332863	43.561	ng	100
32) Benzoic acid	10.377	122	10902m	14.485	ng	
33) 4-Chloroaniline	11.135	127	36233	11.631	ng	96
34) Hexachlorobutadiene	11.293	225	71053	44.579	ng	95
35) Caprolactam	11.910	113	36213m	47.153	ng	
36) 4-Chloro-3-methylphenol	12.251	107	124893	52.466	ng	99
37) 2-Methylnaphthalene	12.621	142	246382	46.004	ng	100
38) 1-Methylnaphthalene	12.839	142	238625	45.795	ng	99
40) 1,2,4,5-Tetrachloroben...	12.980	216	142909	45.156	ng	99
41) Hexachlorocyclopentadiene	12.956	237	94596	56.362	ng	99
43) 2,4,6-Trichlorophenol	13.221	196	96501	45.939	ng	99

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44) 2,4,5-Trichlorophenol	13.303	196	106885	45.308	ng	98
46) 1,1'-Biphenyl	13.620	154	348891	44.750	ng	99
47) 2-Chloronaphthalene	13.667	162	259012	43.745	ng	98
48) 2-Nitroaniline	13.873	65	80406	43.331	ng	88
49) Acenaphthylene	14.507	152	440622	45.096	ng	99
50) Dimethylphthalate	14.231	163	381689	47.190	ng	99
51) 2,6-Dinitrotoluene	14.355	165	82726	50.004	ng	# 84
52) Acenaphthene	14.848	154	277000m	44.135	ng	
53) 3-Nitroaniline	14.689	138	50035	27.017	ng	92
54) 2,4-Dinitrophenol	14.901	184	27148	33.342	ng	98
55) Dibenzofuran	15.183	168	426137	46.206	ng	98
56) 4-Nitrophenol	15.007	139	110946	77.412	ng	92
57) 2,4-Dinitrotoluene	15.142	165	118102	51.611	ng	# 83
58) Fluorene	15.829	166	360127	46.353	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.406	232	92161	43.338	ng	96
60) Diethylphthalate	15.582	149	383273	47.454	ng	97
61) 4-Chlorophenyl-phenyle...	15.812	204	188206	49.105	ng	97
62) 4-Nitroaniline	15.853	138	81932	43.331	ng	97
63) Azobenzene	16.105	77	319107	41.573	ng	94
65) 4,6-Dinitro-2-methylph...	15.912	198	35785	27.977	ng	93
66) n-Nitrosodiphenylamine	16.029	169	326802	45.098	ng	99
67) 4-Bromophenyl-phenylether	16.711	248	134899	48.897	ng	96
68) Hexachlorobenzene	16.828	284	149845	48.316	ng	96
69) Atrazine	16.975	200	125700	49.170	ng	98
70) Pentachlorophenol	17.181	266	92676	54.998	ng	98
71) Phenanthrene	17.574	178	595314	44.474	ng	100
72) Anthracene	17.662	178	607293	46.664	ng	99
73) Carbazole	17.933	167	561207	46.805	ng	99
74) Di-n-butylphthalate	18.467	149	680986	44.882	ng	99
75) Fluoranthene	19.578	202	740147	45.453	ng	99
77) Benzidine	19.754	184	94357	15.769	ng	99
78) Pyrene	19.942	202	750295	44.541	ng	99
80) Butylbenzylphthalate	20.806	149	301034	42.844	ng	96
81) Benzo(a)anthracene	21.805	228	728298	44.464	ng	99
82) 3,3'-Dichlorobenzidine	21.711	252	151717	26.419	ng	98
83) Chrysene	21.875	228	681820	43.536	ng	100
84) Bis(2-ethylhexyl)phtha...	21.681	149	444514	43.911	ng	98
85) Di-n-octyl phthalate	22.939	149	747534	43.544	ng	# 93
87) Indeno(1,2,3-cd)pyrene	29.108	276	886175	45.272	ng	98
88) Benzo(b)fluoranthene	24.125	252	774189	47.864	ng	98
89) Benzo(k)fluoranthene	24.196	252	754285	46.337	ng	98
90) Benzo(a)pyrene	25.048	252	680008	49.208	ng	98
91) Dibenzo(a,h)anthracene	29.172	278	769499	48.253	ng	99
92) Benzo(g,h,i)perylene	30.336	276	763414	47.386	ng	95

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

