

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG061423\
 Data File : BG057714.D
 Acq On : 14 Jun 2023 19:24
 Operator : CG/JU
 Sample : 03167-01MSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

TP-4-GAMSD

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 06/15/2023

Supervised By :mohammad ahmed 06/16/2023

Quant Time: Jun 15 01:23:37 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG061323.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Wed Jun 14 03:52:08 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.936	152	32035	20.000	ng	0.00
21) Naphthalene-d8	10.739	136	132454	20.000	ng	0.00
39) Acenaphthene-d10	14.570	164	88598	20.000	ng	0.00
64) Phenanthrene-d10	17.314	188	226486	20.000	ng	0.00
76) Chrysene-d12	21.573	240	202182	20.000	ng	0.00
86) Perylene-d12	24.734	264	255788	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	169913	86.364	ng	0.00
7) Phenol-d6	7.096	99	234612	77.624	ng	0.00
23) Nitrobenzene-d5	9.094	82	127212	63.959	ng	-0.01
42) 2,4,6-Tribromophenol	16.056	330	96959	98.092	ng	0.00
45) 2-Fluorobiphenyl	13.195	172	294656	49.574	ng	0.00
79) Terphenyl-d14	19.928	244	554442	50.701	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.418	88	36595	36.536	ng	91
3) Pyridine	3.806	79	94992	33.502	ng	94
4) n-Nitrosodimethylamine	3.724	42	55013	39.781	ng	90
6) Aniline	7.261	93	131953	33.276	ng	98
8) 2-Chlorophenol	7.502	128	88867	47.112	ng	95
9) Benzaldehyde	7.073	77	56926	30.557	ng	96
10) Phenol	7.120	94	127710	43.242	ng	98
11) bis(2-Chloroethyl)ether	7.361	93	89900	38.837	ng	94
12) 1,3-Dichlorobenzene	7.825	146	92786	42.523	ng	97
13) 1,4-Dichlorobenzene	7.972	146	94108	43.500	ng	98
14) 1,2-Dichlorobenzene	8.289	146	88725	42.209	ng	100
15) Benzyl Alcohol	8.171	79	90843	39.245	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.465	45	180903	35.891	ng	98
17) 2-Methylphenol	8.371	107	84636	38.944	ng	96
18) Hexachloroethane	9.017	117	32615	48.784	ng	97
19) n-Nitroso-di-n-propyla...	8.741	70	75433	34.517	ng	97
20) 3+4-Methylphenols	8.700	107	116138	39.084	ng	97
22) Acetophenone	8.759	105	148241	38.525	ng	# 97
24) Nitrobenzene	9.141	77	102762	43.560	ng	100
25) Isophorone	9.658	82	216881	35.327	ng	98
26) 2-Nitrophenol	9.852	139	36689	57.352	ng	91
27) 2,4-Dimethylphenol	9.905	122	87533	42.699	ng	99
28) bis(2-Chloroethoxy)met...	10.146	93	120419	37.732	ng	97
29) 2,4-Dichlorophenol	10.381	162	85208	48.998	ng	97
30) 1,2,4-Trichlorobenzene	10.604	180	89810	39.735	ng	95
31) Naphthalene	10.792	128	273640	38.440	ng	99
32) Benzoic acid	10.034	122	64978	81.110	ng	92
33) 4-Chloroaniline	10.892	127	91878	27.722	ng	99
34) Hexachlorobutadiene	11.074	225	55496	39.045	ng	97
35) Caprolactam	11.661	113	34415m	47.974	ng	
36) 4-Chloro-3-methylphenol	12.014	107	107360	43.578	ng	100
37) 2-Methylnaphthalene	12.396	142	187623	35.626	ng	99
38) 1-Methylnaphthalene	12.619	142	175569	35.535	ng	97
40) 1,2,4,5-Tetrachloroben...	12.760	216	99762	39.777	ng	99
41) Hexachlorocyclopentadiene	12.748	237	119932	125.376	ng	95
43) 2,4,6-Trichlorophenol	13.001	196	74507	46.047	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.072	196	82698	43.423	ng	96
46) 1,1'-Biphenyl	13.406	154	248857	40.653	ng	98
47) 2-Chloronaphthalene	13.448	162	193308	41.781	ng	98
48) 2-Nitroaniline	13.647	65	69395	53.371	ng	99
49) Acenaphthylene	14.294	152	318288	39.669	ng	99
50) Dimethylphthalate	14.029	163	270172	41.589	ng	99
51) 2,6-Dinitrotoluene	14.141	165	51435	52.472	ng	86
52) Acenaphthene	14.634	154	214199m	44.536	ng	
53) 3-Nitroaniline	14.470	138	46910	43.390	ng	# 97
54) 2,4-Dinitrophenol	14.676	184	48804	104.431	ng	96
55) Dibenzofuran	14.969	168	313497	38.832	ng	99
56) 4-Nitrophenol	14.775	139	104473	96.486	ng	93
57) 2,4-Dinitrotoluene	14.928	165	71899	55.322	ng	96
58) Fluorene	15.616	166	251603	39.123	ng	97
59) 2,3,4,6-Tetrachlorophenol	15.193	232	75478	47.251	ng	100
60) Diethylphthalate	15.392	149	279658	41.704	ng	99
61) 4-Chlorophenyl-phenyle...	15.610	204	131711	38.228	ng	98
62) 4-Nitroaniline	15.633	138	62168	50.203	ng	94
63) Azobenzene	15.904	77	295065	40.923	ng	97
65) 4,6-Dinitro-2-methylph...	15.692	198	36621	62.074	ng	85
66) n-Nitrosodiphenylamine	15.827	169	227191	38.373	ng	98
67) 4-Bromophenyl-phenylether	16.509	248	88179	38.236	ng	94
68) Hexachlorobenzene	16.620	284	96170	39.903	ng	97
69) Atrazine	16.773	200	98939	50.228	ng	96
70) Pentachlorophenol	16.961	266	132307	91.440	ng	96
71) Phenanthrene	17.355	178	434423	39.308	ng	99
72) Anthracene	17.449	178	440629	39.356	ng	100
73) Carbazole	17.713	167	439719	42.408	ng	100
74) Di-n-butylphthalate	18.283	149	532704	46.351	ng	99
75) Fluoranthene	19.370	202	545526	40.974	ng	99
77) Benzidine	19.546	184	314526	65.257	ng	97
78) Pyrene	19.728	202	557286	38.740	ng	100
80) Butylbenzylphthalate	20.622	149	240862	46.175	ng	98
81) Benzo(a)anthracene	21.550	228	553818	39.664	ng	100
82) 3,3'-Dichlorobenzidine	21.473	252	158225	34.790	ng	100
83) Chrysene	21.620	228	512978	38.604	ng	99
84) Bis(2-ethylhexyl)phtha...	21.473	149	346132	48.840	ng	100
85) Di-n-octyl phthalate	22.666	149	632784	51.780	ng	97
87) Indeno(1,2,3-cd)pyrene	28.342	276	679376	41.049	ng	98
88) Benzo(b)fluoranthene	23.730	252	586598	42.273	ng	99
89) Benzo(k)fluoranthene	23.800	252	574000	39.884	ng	98
90) Benzo(a)pyrene	24.587	252	525604	38.524	ng	# 98
91) Dibenzo(a,h)anthracene	28.407	278	558826	40.859	ng	97
92) Benzo(g,h,i)perylene	29.470	276	563294	41.264	ng	98

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

