

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091422\
 Data File : BG054939.D
 Acq On : 14 Sep 2022 10:52
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_G

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

 Reviewed By : Christian
 Giraldo

Quant Time: Sep 15 03:17:53 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082522.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Thu Sep 15 03:14:10 2022

Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
Internal Standards							09/15/2022
1) 1,4-Dichlorobenzene-d4	8.135	152	36423	20.000	ng	0.00	Supervised By : Jagrut Padhyay
21) Naphthalene-d8	10.961	136	140238	20.000	ng	# 0.00	
39) Acenaphthene-d10	14.774	164	94240	20.000	ng	0.00	
64) Phenanthrene-d10	17.524	188	214821	20.000	ng	0.00	
76) Chrysene-d12	21.819	240	212006	20.000	ng	0.00	
86) Perylene-d12	25.185	264	232969	20.000	ng	0.00	09/15/2022
System Monitoring Compounds							
5) 2-Fluorophenol	5.673	112	173969	83.173	ng	0.00	
7) Phenol-d6	7.289	99	236866	82.806	ng	0.00	
23) Nitrobenzene-d5	9.310	82	220012	82.359	ng	0.00	
42) 2,4,6-Tribromophenol	16.261	330	109734	93.185	ng	0.00	
45) 2-Fluorobiphenyl	13.393	172	560455	89.386	ng	0.00	
79) Terphenyl-d14	20.121	244	903749	84.464	ng	0.00	
Target Compounds							Qvalue
2) 1,4-Dioxane	3.511	88	37994	38.040	ng	89	
3) Pyridine	3.922	79	106170	38.111	ng	90	
4) n-Nitrosodimethylamine	3.840	42	39302	32.636	ng	82	
6) Aniline	7.453	93	138708	40.813	ng	94	
8) 2-Chlorophenol	7.694	128	96986	42.440	ng	96	
9) Benzaldehyde	7.265	77	68458	36.515	ng	93	
10) Phenol	7.318	94	123000	41.812	ng	94	
11) bis(2-Chloroethyl)ether	7.541	93	93651	39.584	ng	94	
12) 1,3-Dichlorobenzene	8.017	146	111946	42.331	ng	98	
13) 1,4-Dichlorobenzene	8.170	146	113683	42.618	ng	96	
14) 1,2-Dichlorobenzene	8.493	146	108568	42.923	ng	97	
15) Benzyl Alcohol	8.376	79	81143	39.265	ng	93	
16) 2,2'-oxybis(1-Chloropr...	8.658	45	118257	31.944	ng	93	
17) 2-Methylphenol	8.575	107	86168	42.544	ng	95	
18) Hexachloroethane	9.222	117	39782	41.127	ng	93	
19) n-Nitroso-di-n-propyla...	8.940	70	74966	38.111	ng	# 90	
20) 3+4-Methylphenols	8.910	107	120389	42.865	ng	96	
22) Acetophenone	8.963	105	150426	42.875	ng	# 94	
24) Nitrobenzene	9.351	77	106941	39.717	ng	92	
25) Isophorone	9.874	82	202717	40.698	ng	96	
26) 2-Nitrophenol	10.068	139	58760	49.404	ng	98	
27) 2,4-Dimethylphenol	10.121	122	82461	43.758	ng	95	
28) bis(2-Chloroethoxy)met...	10.350	93	116536	41.058	ng	100	
29) 2,4-Dichlorophenol	10.608	162	99222	46.291	ng	97	
30) 1,2,4-Trichlorobenzene	10.820	180	111661	45.586	ng	98	
31) Naphthalene	11.014	128	327066	43.469	ng	100	
32) Benzoic acid	10.338	122	15072m	17.185	ng		
33) 4-Chloroaniline	11.120	127	131498	42.869	ng	97	
34) Hexachlorobutadiene	11.278	225	75298	47.979	ng	98	
35) Caprolactam	11.901	113	32453	42.916	ng	# 76	
36) 4-Chloro-3-methylphenol	12.236	107	102311	43.650	ng	98	
37) 2-Methylnaphthalene	12.612	142	233569	44.291	ng	99	
38) 1-Methylnaphthalene	12.829	142	225530	43.957	ng	100	
40) 1,2,4,5-Tetrachloroben...	12.970	216	134172	47.230	ng	98	
41) Hexachlorocyclopentadiene	12.947	237	62868	41.729	ng	100	
43) 2,4,6-Trichlorophenol	13.211	196	78382	41.569	ng	97	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.294	196	88163	41.634	ng	96	09/15/2022
46) 1,1'-Biphenyl	13.605	154	303225	43.328	ng	99	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.658	162	230177	43.308	ng	98	
48) 2-Nitroaniline	13.858	65	66296	39.801	ng	89	
49) Acenaphthylene	14.498	152	380906	43.430	ng	100	
50) Dimethylphthalate	14.222	163	314546	43.324	ng	99	
51) 2,6-Dinitrotoluene	14.345	165	66903	45.052	ng	# 83	09/15/2022
52) Acenaphthene	14.839	154	241649m	42.893	ng		
53) 3-Nitroaniline	14.686	138	72037	43.333	ng	88	
54) 2,4-Dinitrophenol	14.898	184	27187	36.422	ng	94	
55) Dibenzofuran	15.174	168	364427	44.021	ng	97	
56) 4-Nitrophenol	14.997	139	46042	35.789	ng	87	
57) 2,4-Dinitrotoluene	15.138	165	94996	46.248	ng	# 80	
58) Fluorene	15.820	166	303751	43.555	ng	99	
59) 2,3,4,6-Tetrachlorophenol	15.397	232	72465	37.962	ng	97	
60) Diethylphthalate	15.573	149	300544	41.455	ng	96	
61) 4-Chlorophenyl-phenyle...	15.802	204	154108	44.794	ng	97	
62) 4-Nitroaniline	15.843	138	74793	44.066	ng	93	
63) Azobenzene	16.096	77	255677	37.108	ng	93	
65) 4,6-Dinitro-2-methylph...	15.902	198	48932	44.748	ng	92	
66) n-Nitrosodiphenylamine	16.020	169	265476	42.853	ng	99	
67) 4-Bromophenyl-phenylether	16.701	248	110266	46.751	ng	96	
68) Hexachlorobenzene	16.825	284	125982	47.515	ng	# 92	
69) Atrazine	16.960	200	95767	43.818	ng	99	
70) Pentachlorophenol	17.171	266	51943m	36.056	ng		
71) Phenanthrene	17.565	178	491402	42.941	ng	99	
72) Anthracene	17.659	178	483868	43.490	ng	99	
73) Carbazole	17.929	167	435690	42.503	ng	99	
74) Di-n-butylphthalate	18.464	149	531099	40.944	ng	99	
75) Fluoranthene	19.569	202	605498	43.495	ng	98	
77) Benzidine	19.745	184	190459	36.270	ng	99	
78) Pyrene	19.933	202	614784	41.587	ng	99	
80) Butylbenzylphthalate	20.802	149	243147	39.432	ng	91	
81) Benzo(a)anthracene	21.795	228	613766	42.698	ng	99	
82) 3,3'-Dichlorobenzidine	21.701	252	226653	44.972	ng	99	
83) Chrysene	21.866	228	582955	42.415	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.666	149	352277	39.653	ng	98	
85) Di-n-octyl phthalate	22.923	149	596356	39.583	ng	# 91	
87) Indeno(1,2,3-cd)pyrene	29.087	276	761901	43.079	ng	# 95	
88) Benzo(b)fluoranthene	24.110	252	616567	42.189	ng	99	
89) Benzo(k)fluoranthene	24.181	252	624374	42.452	ng	98	
90) Benzo(a)pyrene	25.033	252	530784	42.511	ng	98	
91) Dibenzo(a,h)anthracene	29.145	278	627521	43.551	ng	99	
92) Benzo(g,h,i)perylene	30.309	276	624523	42.904	ng	# 96	

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

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