

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG022123\
 Data File : BG056681.D
 Acq On : 21 Feb 2023 11:43
 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: Feb 21 14:03:37 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG021523.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Feb 21 14:02:56 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.429	152	14758	20.000	ng	0.00
21) Naphthalene-d8	11.273	136	58858	20.000	ng	0.00
39) Acenaphthene-d10	15.033	164	45046	20.000	ng	0.00
64) Phenanthrene-d10	17.771	188	112872	20.000	ng	0.00
76) Chrysene-d12	22.090	240	103162	20.000	ng	0.00
86) Perylene-d12	25.621	264	114711	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.932	112	76192	84.002	ng	0.00
7) Phenol-d6	7.554	99	112353	83.857	ng	0.00
23) Nitrobenzene-d5	9.604	82	119285	103.767	ng	0.00
42) 2,4,6-Tribromophenol	16.514	330	57907	93.893	ng	0.00
45) 2-Fluorobiphenyl	13.664	172	289479	86.052	ng	0.00
79) Terphenyl-d14	20.339	244	545287	86.481	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.752	88	15843	39.841	ng	97
3) Pyridine	4.169	79	52673	41.876	ng	90
4) n-Nitrosodimethylamine	4.075	42	28801	49.123	ng	# 79
6) Aniline	7.742	93	77887	41.829	ng	95
8) 2-Chlorophenol	7.982	128	36657	40.847	ng	92
9) Benzaldehyde	7.554	77	28710	40.519	ng	91
10) Phenol	7.577	94	58329	42.559	ng	96
11) bis(2-Chloroethyl)ether	7.830	93	43286	42.093	ng	92
12) 1,3-Dichlorobenzene	8.317	146	43421	40.610	ng	95
13) 1,4-Dichlorobenzene	8.464	146	43952	40.817	ng	95
14) 1,2-Dichlorobenzene	8.793	146	41697	40.286	ng	94
15) Benzyl Alcohol	8.658	79	52033	43.372	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.952	45	62965	45.607	ng	97
17) 2-Methylphenol	8.852	107	41086	42.658	ng	94
18) Hexachloroethane	9.534	117	15382	45.095	ng	94
19) n-Nitroso-di-n-propyla...	9.234	70	46792	46.650	ng	92
20) 3+4-Methylphenols	9.181	107	56953	43.046	ng	95
22) Acetophenone	9.257	105	76599	42.593	ng	# 96
24) Nitrobenzene	9.651	77	63165	47.221	ng	96
25) Isophorone	10.168	82	122788	43.889	ng	98
26) 2-Nitrophenol	10.362	139	23070	48.740	ng	92
27) 2,4-Dimethylphenol	10.403	122	44758	41.380	ng	97
28) bis(2-Chloroethoxy)met...	10.644	93	60790	42.636	ng	97
29) 2,4-Dichlorophenol	10.897	162	45668	42.600	ng	97
30) 1,2,4-Trichlorobenzene	11.126	180	53236	41.688	ng	95
31) Naphthalene	11.320	128	137825	42.251	ng	100
32) Benzoic acid	10.515	122	32558m	46.702	ng	
33) 4-Chloroaniline	11.414	127	60666	40.706	ng	98
34) Hexachlorobutadiene	11.596	225	38820	42.538	ng	94
35) Caprolactam	12.172	113	15176	44.015	ng	# 68
36) 4-Chloro-3-methylphenol	12.495	107	50827	42.994	ng	97
37) 2-Methylnaphthalene	12.894	142	108923	42.290	ng	99
38) 1-Methylnaphthalene	13.112	142	102333	42.579	ng	98
40) 1,2,4,5-Tetrachloroben...	13.247	216	73124	43.667	ng	100
41) Hexachlorocyclopentadiene	13.229	237	43971	48.413	ng	95
43) 2,4,6-Trichlorophenol	13.476	196	47322	46.577	ng	98

02/22/2023

Supervised By :Jagrut

padhyay

02/22/2023

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.547	196	54944	45.121	ng	98	02/22/2023
46) 1,1'-Biphenyl	13.876	154	147517	43.231	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.928	162	115850	42.903	ng	98	
48) 2-Nitroaniline	14.111	65	40842	51.493	ng	85	
49) Acenaphthylene	14.763	152	186121	42.420	ng	98	
50) Dimethylphthalate	14.475	163	162668	43.683	ng	99	
51) 2,6-Dinitrotoluene	14.592	165	31612	43.838	ng	# 81	02/22/2023
52) Acenaphthene	15.098	154	120422	44.495	ng	98	
53) 3-Nitroaniline	14.927	138	34679	47.976	ng	# 96	
54) 2,4-Dinitrophenol	15.121	184	20544	47.754	ng	# 92	
55) Dibenzofuran	15.427	168	196073	42.883	ng	99	
56) 4-Nitrophenol	15.198	139	26552	47.944	ng	86	
57) 2,4-Dinitrotoluene	15.374	165	45755	50.381	ng	# 76	
58) Fluorene	16.073	166	156113	43.183	ng	93	
59) 2,3,4,6-Tetrachlorophenol	15.644	232	50765m	47.567	ng		
60) Diethylphthalate	15.820	149	163675	43.574	ng	100	
61) 4-Chlorophenyl-phenyle...	16.055	204	95723	43.342	ng	98	
62) 4-Nitroaniline	16.079	138	36410	48.332	ng	89	
63) Azobenzene	16.349	77	171032	45.495	ng	97	
65) 4,6-Dinitro-2-methylph...	16.132	198	29101	49.927	ng	93	
66) n-Nitrosodiphenylamine	16.267	169	142616	43.433	ng	98	
67) 4-Bromophenyl-phenylether	16.948	248	63754	43.994	ng	93	
68) Hexachlorobenzene	17.078	284	64182	43.471	ng	98	
69) Atrazine	17.201	200	55762	44.225	ng	96	
70) Pentachlorophenol	17.407	266	47374	48.451	ng	95	
71) Phenanthrene	17.812	178	266828	43.452	ng	99	
72) Anthracene	17.906	178	274325	43.658	ng	99	
73) Carbazole	18.165	167	227009	41.910	ng	97	
74) Di-n-butylphthalate	18.693	149	258609	42.663	ng	99	
75) Fluoranthene	19.798	202	338618	41.791	ng	98	
77) Benzidine	19.963	184	117053	47.440	ng	100	
78) Pyrene	20.156	202	338184	43.407	ng	99	
80) Butylbenzylphthalate	21.020	149	109351	46.740	ng	96	
81) Benzo(a)anthracene	22.066	228	330270	43.049	ng	97	
82) 3,3'-Dichlorobenzidine	21.960	252	112056	45.019	ng	97	
83) Chrysene	22.137	228	307065	42.814	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.937	149	160787	45.863	ng	99	
85) Di-n-octyl phthalate	23.259	149	267470	45.903	ng	97	
87) Indeno(1,2,3-cd)pyrene	29.710	276	389168	43.319	ng	# 99	
88) Benzo(b)fluoranthene	24.493	252	321309	42.875	ng	99	
89) Benzo(k)fluoranthene	24.569	252	324165m	43.987	ng		
90) Benzo(a)pyrene	25.462	252	314315	43.633	ng	98	
91) Dibenzo(a,h)anthracene	29.786	278	323431	43.469	ng	# 97	
92) Benzo(g,h,i)perylene	30.997	276	314163	43.211	ng	94	

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

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