

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042222\
 Data File : BG053267.D
 Acq On : 22 Apr 2022 18:57
 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: Apr 24 01:15:51 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG040822.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Sun Apr 24 01:13:30 2022

Response via : Initial Calibration

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

1) 1,4-Dichlorobenzene-d4	8.114	152	28751	20.000	ng	0.00
21) Naphthalene-d8	10.922	136	119039	20.000	ng	0.00
39) Acenaphthene-d10	14.734	164	89313	20.000	ng	0.00
64) Phenanthrene-d10	17.477	188	211290	20.000	ng	0.00
76) Chrysene-d12	21.766	240	211911	20.000	ng	0.00
86) Perylene-d12	25.055	264	221064	20.000	ng	0.00

04/25/2022

 Supervised By :Jagrut
 padhyay

System Monitoring Compounds

5) 2-Fluorophenol	5.676	112	149185	88.947	ng	0.00
7) Phenol-d6	7.274	99	224912	86.957	ng	0.00
23) Nitrobenzene-d5	9.271	82	235311	80.263	ng	0.00
42) 2,4,6-Tribromophenol	16.220	330	116375	81.900	ng	0.00
45) 2-Fluorobiphenyl	13.360	172	530251	81.866	ng	0.00
79) Terphenyl-d14	20.080	244	1007449	80.292	ng	0.00

04/25/2022

Target Compounds

Compound	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.555	88	35027	43.730	ng	95
3) Pyridine	3.960	79	92534	43.802	ng	# 86
4) n-Nitrosodimethylamine	3.866	42	44766	42.792	ng	89
6) Aniline	7.432	93	129344	43.150	ng	98
8) 2-Chlorophenol	7.679	128	79281	43.784	ng	97
9) Benzaldehyde	7.244	77	63628m	41.155	ng	
10) Phenol	7.297	94	111237	43.206	ng	90
11) bis(2-Chloroethyl)ether	7.520	93	79475	42.823	ng	93
12) 1,3-Dichlorobenzene	8.002	146	92051	43.748	ng	96
13) 1,4-Dichlorobenzene	8.143	146	92803	43.263	ng	99
14) 1,2-Dichlorobenzene	8.466	146	85661	41.411	ng	96
15) Benzyl Alcohol	8.343	79	100468	43.701	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.631	45	129333	41.741	ng	98
17) 2-Methylphenol	8.554	107	72965	41.881	ng	94
18) Hexachloroethane	9.195	117	35003	43.514	ng	96
19) n-Nitroso-di-n-propyla...	8.913	70	80094	42.298	ng	# 95
20) 3+4-Methylphenols	8.877	107	106711	43.388	ng	96
22) Acetophenone	8.930	105	135583	41.055	ng	95
24) Nitrobenzene	9.312	77	114910	40.296	ng	# 88
25) Isophorone	9.835	82	203880	40.914	ng	97
26) 2-Nitrophenol	10.029	139	48324	40.526	ng	# 86
27) 2,4-Dimethylphenol	10.087	122	70953	41.547	ng	91
28) bis(2-Chloroethoxy)met...	10.317	93	117225	41.873	ng	98
29) 2,4-Dichlorophenol	10.569	162	88311	41.134	ng	98
30) 1,2,4-Trichlorobenzene	10.787	180	99633	41.003	ng	93
31) Naphthalene	10.975	128	256826	40.439	ng	99
32) Benzoic acid	10.240	122	42791	39.219	ng	# 83
33) 4-Chloroaniline	11.074	127	110230	40.530	ng	97
34) Hexachlorobutadiene	11.256	225	74765	41.422	ng	98
35) Caprolactam	11.844	113	30393	40.162	ng	# 87
36) 4-Chloro-3-methylphenol	12.196	107	101036	40.423	ng	96
37) 2-Methylnaphthalene	12.566	142	189562	40.337	ng	96
38) 1-Methylnaphthalene	12.790	142	185314m	40.398	ng	
40) 1,2,4,5-Tetrachloroben...	12.937	216	123983	40.873	ng	100
41) Hexachlorocyclopentadiene	12.919	237	57994	37.036	ng	93
43) 2,4,6-Trichlorophenol	13.177	196	85596	40.927	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.254	196	87943	39.460	ng	96	04/25/2022
46) 1,1'-Biphenyl	13.571	154	260223	40.459	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.612	162	205540	40.055	ng	97	
48) 2-Nitroaniline	13.812	65	81028	41.609	ng	91	
49) Acenaphthylene	14.458	152	330092	41.094	ng	98	
50) Dimethylphthalate	14.182	163	289077	40.467	ng	98	
51) 2,6-Dinitrotoluene	14.299	165	60076	40.312	ng	85	04/25/2022
52) Acenaphthene	14.799	154	216965m	39.829	ng		
53) 3-Nitroaniline	14.634	138	64467	40.910	ng	94	
54) 2,4-Dinitrophenol	14.846	184	32980	34.780	ng	#	86
55) Dibenzofuran	15.128	168	343721	40.297	ng		98
56) 4-Nitrophenol	14.946	139	49962	41.571	ng	#	79
57) 2,4-Dinitrotoluene	15.092	165	88200	41.570	ng		88
58) Fluorene	15.780	166	275342	39.967	ng		98
59) 2,3,4,6-Tetrachlorophenol	15.357	232	88673	40.970	ng		99
60) Diethylphthalate	15.539	149	296649	39.634	ng		99
61) 4-Chlorophenyl-phenyle...	15.768	204	158058	40.473	ng		97
62) 4-Nitroaniline	15.797	138	67759	40.749	ng	#	79
63) Azobenzene	16.062	77	299501	39.689	ng		98
65) 4,6-Dinitro-2-methylph...	15.856	198	58600	39.753	ng		93
66) n-Nitrosodiphenylamine	15.980	169	249220	41.017	ng		97
67) 4-Bromophenyl-phenylether	16.661	248	110426	40.787	ng		96
68) Hexachlorobenzene	16.790	284	123328	42.063	ng		96
69) Atrazine	16.919	200	92832	39.113	ng		99
70) Pentachlorophenol	17.137	266	68621	42.824	ng		94
71) Phenanthrene	17.524	178	472957	40.980	ng		98
72) Anthracene	17.613	178	464063	40.588	ng		99
73) Carbazole	17.877	167	456795	41.100	ng		99
74) Di-n-butylphthalate	18.429	149	518672	41.148	ng		98
75) Fluoranthene	19.528	202	610599	41.274	ng		99
77) Benzidine	19.698	184	239256	39.572	ng		99
78) Pyrene	19.892	202	614700	40.729	ng		97
80) Butylbenzylphthalate	20.761	149	233927	41.575	ng		96
81) Benzo(a)anthracene	21.742	228	597269	41.222	ng	100	
82) 3,3'-Dichlorobenzidine	21.654	252	218317	40.971	ng		99
83) Chrysene	21.813	228	556452	41.382	ng		99
84) Bis(2-ethylhexyl)phtha...	21.637	149	317257	41.714	ng	100	
85) Di-n-octyl phthalate	22.870	149	530763	41.727	ng		97
87) Indeno(1,2,3-cd)pyrene	28.844	276	670257	40.810	ng	#	96
88) Benzo(b)fluoranthene	24.010	252	573816	41.140	ng		99
89) Benzo(k)fluoranthene	24.080	252	563682	41.120	ng		99
90) Benzo(a)pyrene	24.903	252	488360	41.100	ng		97
91) Dibenzo(a,h)anthracene	28.903	278	550077	40.693	ng		98
92) Benzo(g,h,i)perylene	30.013	276	543759	40.812	ng		99

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

