

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG042322\
 Data File : BG053286.D
 Acq On : 23 Apr 2022 10:33
 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 02:47:56 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	8.107	152	31398	20.000	ng	0.00
21) Naphthalene-d8	10.921	136	123117	20.000	ng	0.00
39) Acenaphthene-d10	14.733	164	91638	20.000	ng	0.00
64) Phenanthrene-d10	17.477	188	217099	20.000	ng	0.00
76) Chrysene-d12	21.759	240	221260	20.000	ng	0.00
86) Perylene-d12	25.055	264	230486	20.000	ng	0.00

04/25/2022

Supervised By : Jagrut
Padhyay

System Monitoring Compounds

5) 2-Fluorophenol	5.675	112	157529	86.003	ng	0.00
7) Phenol-d6	7.267	99	236158	83.607	ng	0.00
23) Nitrobenzene-d5	9.270	82	246670	81.351	ng	0.00
42) 2,4,6-Tribromophenol	16.220	330	118978	81.608	ng	0.00
45) 2-Fluorobiphenyl	13.359	172	555713	83.620	ng	0.00
79) Terphenyl-d14	20.079	244	1011665	77.222	ng	0.00

04/25/2022

Target Compounds

Compound	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.554	88	35525	40.613	ng	95
3) Pyridine	3.960	79	100702	43.650	ng	# 87
4) n-Nitrosodimethylamine	3.872	42	49499	43.327	ng	86
6) Aniline	7.431	93	136609	41.731	ng	94
8) 2-Chlorophenol	7.672	128	84184	42.572	ng	92
9) Benzaldehyde	7.244	77	66739m	39.528	ng	
10) Phenol	7.296	94	119436	42.480	ng	88
11) bis(2-Chloroethyl)ether	7.520	93	83088	40.996	ng	91
12) 1,3-Dichlorobenzene	8.001	146	93136	40.532	ng	95
13) 1,4-Dichlorobenzene	8.142	146	97234	41.507	ng	98
14) 1,2-Dichlorobenzene	8.465	146	91985	40.719	ng	99
15) Benzyl Alcohol	8.342	79	101178	40.299	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.630	45	139047	41.093	ng	98
17) 2-Methylphenol	8.548	107	80071	42.085	ng	96
18) Hexachloroethane	9.194	117	36156	41.158	ng	94
19) n-Nitroso-di-n-propyla...	8.912	70	84857	41.035	ng	# 96
20) 3+4-Methylphenols	8.877	107	112173	41.764	ng	95
22) Acetophenone	8.929	105	141752	41.501	ng	# 94
24) Nitrobenzene	9.311	77	122724	41.611	ng	93
25) Isophorone	9.834	82	209890	40.725	ng	100
26) 2-Nitrophenol	10.028	139	51898	42.081	ng	93
27) 2,4-Dimethylphenol	10.087	122	75517	42.755	ng	94
28) bis(2-Chloroethoxy)met...	10.310	93	119477	41.264	ng	98
29) 2,4-Dichlorophenol	10.568	162	93419	42.072	ng	98
30) 1,2,4-Trichlorobenzene	10.780	180	103682	41.256	ng	96
31) Naphthalene	10.968	128	276714	42.127	ng	99
32) Benzoic acid	10.239	122	45599m	40.298	ng	
33) 4-Chloroaniline	11.074	127	115331	41.001	ng	98
34) Hexachlorobutadiene	11.256	225	75182	40.273	ng	98
35) Caprolactam	11.849	113	31823	40.659	ng	# 90
36) 4-Chloro-3-methylphenol	12.190	107	107247	41.487	ng	90
37) 2-Methylnaphthalene	12.566	142	201560	41.470	ng	94
38) 1-Methylnaphthalene	12.783	142	193333	40.751	ng	96
40) 1,2,4,5-Tetrachloroben...	12.936	216	129303	41.545	ng	99
41) Hexachlorocyclopentadiene	12.918	237	75134	46.764	ng	98
43) 2,4,6-Trichlorophenol	13.171	196	88459	41.223	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	13.247	196	92431	40.422	ng	96	04/25/2022
46) 1,1'-Biphenyl	13.564	154	276768	41.940	ng	98	Supervised By :Jagrut Upadhyay
47) 2-Chloronaphthalene	13.611	162	216047	41.035	ng	96	
48) 2-Nitroaniline	13.811	65	82355	41.218	ng	92	
49) Acenaphthylene	14.457	152	347128	42.118	ng	99	
50) Dimethylphthalate	14.181	163	296042	40.390	ng	98	
51) 2,6-Dinitrotoluene	14.299	165	61041	39.920	ng	# 85	04/25/2022
52) Acenaphthene	14.798	154	228710m	40.920	ng		
53) 3-Nitroaniline	14.634	138	67055	41.472	ng	97	
54) 2,4-Dinitrophenol	14.839	184	36116	37.121	ng	# 87	
55) Dibenzofuran	15.127	168	360759	41.221	ng	97	
56) 4-Nitrophenol	14.945	139	48611	39.421	ng	# 78	
57) 2,4-Dinitrotoluene	15.086	165	90608	41.622	ng	94	
58) Fluorene	15.779	166	289192	40.913	ng	98	
59) 2,3,4,6-Tetrachlorophenol	15.356	232	91101	41.024	ng	99	
60) Diethylphthalate	15.538	149	303029	39.459	ng	99	
61) 4-Chlorophenyl-phenyle...	15.761	204	164150	40.966	ng	97	
62) 4-Nitroaniline	15.797	138	69663	40.831	ng	# 82	
63) Azobenzene	16.055	77	315428	40.739	ng	97	
65) 4,6-Dinitro-2-methylph...	15.855	198	64319	42.465	ng	94	
66) n-Nitrosodiphenylamine	15.979	169	257734	41.283	ng	99	
67) 4-Bromophenyl-phenylether	16.660	248	114461	41.146	ng	96	
68) Hexachlorobenzene	16.789	284	122029	40.507	ng	97	
69) Atrazine	16.925	200	92889	38.090	ng	97	
70) Pentachlorophenol	17.130	266	68687	41.718	ng	92	
71) Phenanthrene	17.518	178	483401	40.765	ng	98	
72) Anthracene	17.612	178	480985	40.942	ng	98	
73) Carbazole	17.876	167	459476	40.235	ng	99	
74) Di-n-butylphthalate	18.422	149	521104	40.235	ng	98	
75) Fluoranthene	19.527	202	619063	40.727	ng	100	
77) Benzidine	19.697	184	215444	34.128	ng	99	
78) Pyrene	19.885	202	613460	38.930	ng	98	
80) Butylbenzylphthalate	20.761	149	235057	40.010	ng	97	
81) Benzo(a)anthracene	21.742	228	602314	39.814	ng	99	
82) 3,3'-Dichlorobenzidine	21.648	252	217812	39.149	ng	99	
83) Chrysene	21.806	228	566356	40.339	ng	99	
84) Bis(2-ethylhexyl)phtha...	21.630	149	325633	41.006	ng	99	
85) Di-n-octyl phthalate	22.864	149	541317	40.758	ng	96	
87) Indeno(1,2,3-cd)pyrene	28.838	276	698921	40.816	ng	99	
88) Benzo(b)fluoranthene	24.009	252	580846	39.942	ng	100	
89) Benzo(k)fluoranthene	24.080	252	589410	41.239	ng	99	
90) Benzo(a)pyrene	24.902	252	503875	40.672	ng	98	
91) Dibenzo(a,h)anthracene	28.897	278	571987	40.584	ng	99	
92) Benzo(g,h,i)perylene	30.013	276	561840	40.445	ng	97	

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

