

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092022\
 Data File : BP011759.D
 Acq On : 20 Sep 2022 12:51
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_P

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Christian Giraldo 09/21/2022
 Supervised By :Jagrut Upadhyay 09/21/2022

Quant Time: Sep 21 06:44:13 2022

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Sep 20 04:42:52 2022

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	490516	20.000	ng	0.00
21) Naphthalene-d8	10.746	136	2027995	20.000	ng	0.00
39) Acenaphthene-d10	14.575	164	1162504	20.000	ng	0.00
64) Phenanthrene-d10	17.328	188	2265078	20.000	ng	# 0.00
76) Chrysene-d12	21.410	240	2031436	20.000	ng	# 0.00
86) Perylene-d12	23.863	264	1719345	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	2301039	85.387	ng	0.00
7) Phenol-d6	7.093	99	3149522	87.676	ng	0.00
23) Nitrobenzene-d5	9.105	82	3164557	92.545	ng	0.00
42) 2,4,6-Tribromophenol	16.069	330	707846	77.930	ng	0.00
45) 2-Fluorobiphenyl	13.204	172	6134809	82.031	ng	0.00
79) Terphenyl-d14	19.881	244	7843614	83.963	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.376	88	499257	42.068	ng	# 85
3) Pyridine	3.781	79	1547873	45.241	ng	# 80
4) n-Nitrosodimethylamine	3.687	42	600615	45.791	ng	# 48
6) Aniline	7.264	93	1965362	44.414	ng	91
8) 2-Chlorophenol	7.499	128	1350387	42.260	ng	94
9) Benzaldehyde	7.069	77	970006	42.186	ng	93
10) Phenol	7.122	94	1766669	43.725	ng	83
11) bis(2-Chloroethyl)ether	7.358	93	1402623	43.578	ng	95
12) 1,3-Dichlorobenzene	7.822	146	1441696	40.684	ng	94
13) 1,4-Dichlorobenzene	7.969	146	1471365	40.828	ng	97
14) 1,2-Dichlorobenzene	8.287	146	1413626	41.148	ng	97
15) Benzyl Alcohol	8.175	79	1152439	44.939	ng	91
16) 2,2'-oxybis(1-Chloropr...	8.469	45	2127809	49.494	ng	96
17) 2-Methylphenol	8.375	107	1161767	44.286	ng	# 92
18) Hexachloroethane	9.022	117	567119	45.497	ng	91
19) n-Nitroso-di-n-propyla...	8.752	70	1096003	47.122	ng	88
20) 3+4-Methylphenols	8.710	107	1597796	44.349	ng	93
22) Acetophenone	8.763	105	2048347	43.086	ng	# 90
24) Nitrobenzene	9.146	77	1614147	45.369	ng	91
25) Isophorone	9.675	82	2852102	47.150	ng	# 96
26) 2-Nitrophenol	9.857	139	660622	40.980	ng	# 85
27) 2,4-Dimethylphenol	9.916	122	1079017	42.666	ng	90
28) bis(2-Chloroethoxy)met...	10.157	93	1687451	43.625	ng	99
29) 2,4-Dichlorophenol	10.393	162	1171054	41.801	ng	97
30) 1,2,4-Trichlorobenzene	10.610	180	1224647	39.485	ng	98
31) Naphthalene	10.799	128	4360583	41.434	ng	100
32) Benzoic acid	10.063	122	847024	40.991	ng	89
33) 4-Chloroaniline	10.910	127	1788620	43.318	ng	93
34) Hexachlorobutadiene	11.081	225	643319	37.653	ng	98
35) Caprolactam	11.716	113	424152	46.725	ng	# 86
36) 4-Chloro-3-methylphenol	12.028	107	1327080	44.047	ng	94
37) 2-Methylnaphthalene	12.404	142	2944130	41.802	ng	97
38) 1-Methylnaphthalene	12.622	142	2860610	41.543	ng	100
40) 1,2,4,5-Tetrachloroben...	12.769	216	1179856	38.455	ng	# 97
41) Hexachlorocyclopentadiene	12.751	237	614432	40.121	ng	99
43) 2,4,6-Trichlorophenol	13.010	196	858732	41.146	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.081	196	959993	41.140	ng	97
46) 1,1'-Biphenyl	13.410	154	3505042	41.446	ng	98
47) 2-Chloronaphthalene	13.451	162	2747558	41.293	ng	99
48) 2-Nitroaniline	13.657	65	984101	52.739	ng	90
49) Acenaphthylene	14.298	152	4476187	43.476	ng	99
50) Dimethylphthalate	14.040	163	3433424	42.336	ng	100
51) 2,6-Dinitrotoluene	14.151	165	732766	44.557	ng	95
52) Acenaphthene	14.640	154	2884304m	42.116	ng	
53) 3-Nitroaniline	14.487	138	891488	43.035	ng	92
54) 2,4-Dinitrophenol	14.687	184	329037	38.366	ng	93
55) Dibenzofuran	14.975	168	4084880	41.420	ng	100
56) 4-Nitrophenol	14.781	139	712584	44.594	ng	# 79
57) 2,4-Dinitrotoluene	14.940	165	995597	42.071	ng	92
58) Fluorene	15.628	166	3451177	43.144	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.198	232	768997m	40.114	ng	
60) Diethylphthalate	15.404	149	3590288	44.467	ng	98
61) 4-Chlorophenyl-phenyle...	15.622	204	1506200	40.734	ng	99
62) 4-Nitroaniline	15.651	138	931121	43.919	ng	# 82
63) Azobenzene	15.910	77	3802143	49.002	ng	92
65) 4,6-Dinitro-2-methylph...	15.704	198	465497	39.324	ng	91
66) n-Nitrosodiphenylamine	15.834	169	2855803	42.374	ng	97
67) 4-Bromophenyl-phenylether	16.516	248	837487	39.497	ng	94
68) Hexachlorobenzene	16.634	284	872302	38.198	ng	# 92
69) Atrazine	16.792	200	880523	42.802	ng	95
70) Pentachlorophenol	16.975	266	568111	39.215	ng	99
71) Phenanthrene	17.375	178	5178935	42.020	ng	97
72) Anthracene	17.463	178	5062147	43.427	ng	98
73) Carbazole	17.734	167	4838457	43.998	ng	98
74) Di-n-butylphthalate	18.281	149	6207017	46.828	ng	99
75) Fluoranthene	19.333	202	5728533	43.096	ng	92
77) Benzidine	19.516	184	1660006	34.361	ng	98
78) Pyrene	19.686	202	5931319	43.290	ng	97
80) Butylbenzylphthalate	20.557	149	2831523	49.316	ng	96
81) Benzo(a)anthracene	21.398	228	5547581	42.369	ng	96
82) 3,3'-Dichlorobenzidine	21.327	252	1663406	41.828	ng	# 98
83) Chrysene	21.451	228	5231583	41.818	ng	96
84) Bis(2-ethylhexyl)phtha...	21.316	149	4104615	49.695	ng	# 99
85) Di-n-octyl phthalate	22.257	149	6564509	46.315	ng	95
87) Indeno(1,2,3-cd)pyrene	26.433	276	4674475	37.737	ng	# 87
88) Benzo(b)fluoranthene	23.116	252	4789332	45.039	ng	# 96
89) Benzo(k)fluoranthene	23.163	252	4809385	44.881	ng	# 95
90) Benzo(a)pyrene	23.751	252	3867756	44.029	ng	# 94
91) Dibenzo(a,h)anthracene	26.451	278	3921879	38.130	ng	# 93
92) Benzo(g,h,i)perylene	27.215	276	3599364	35.895	ng	# 87

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

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