

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080923\
 Data File : BF134689.D
 Acq On : 09 Aug 2023 14:28
 Operator : CG\JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :

BNA_F

ClientSampleId :

SSTDCCC040

Manual Integrations

APPROVED

Reviewed By :Yogesh Patel 08/10/2023

Supervised By :mohammad ahmed 08/10/2023

Quant Time: Aug 10 01:39:42 2023

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF080123.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Aug 04 09:29:15 2023

Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.804	152	209965	20.000	ng	0.00
21) Naphthalene-d8	8.081	136	810987	20.000	ng	# 0.00
39) Acenaphthene-d10	9.840	164	419030	20.000	ng	0.00
64) Phenanthrene-d10	11.328	188	762786	20.000	ng	# 0.00
76) Chrysene-d12	13.980	240	450629	20.000	ng	# 0.00
86) Perylene-d12	15.451	264	447570	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1033798	74.845	ng	0.00
7) Phenol-d6	6.446	99	1337293	75.738	ng	0.00
23) Nitrobenzene-d5	7.369	82	1131516	87.192	ng	0.00
42) 2,4,6-Tribromophenol	10.634	330	393361	92.087	ng	0.00
45) 2-Fluorobiphenyl	9.163	172	1965903	78.006	ng	0.00
79) Terphenyl-d14	12.922	244	2148297	84.042	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.546	88	235332	37.030	ng	# 94
3) Pyridine	3.305	79	650253	37.760	ng	90
4) n-Nitrosodimethylamine	3.275	42	292131	35.503	ng	# 60
6) Aniline	6.469	93	854139	36.881	ng	93
8) 2-Chlorophenol	6.593	128	550639	39.519	ng	95
9) Benzaldehyde	6.357	77	358385	43.714	ng	98
10) Phenol	6.457	94	675070m	38.797	ng	
11) bis(2-Chloroethyl)ether	6.546	93	540861	39.129	ng	94
12) 1,3-Dichlorobenzene	6.746	146	562817	39.041	ng	94
13) 1,4-Dichlorobenzene	6.822	146	565719	39.155	ng	98
14) 1,2-Dichlorobenzene	6.975	146	520199	38.642	ng	96
15) Benzyl Alcohol	6.946	79	479324	39.351	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.081	45	835802	38.596	ng	89
17) 2-Methylphenol	7.057	107	487577	39.600	ng	92
18) Hexachloroethane	7.310	117	203497	40.131	ng	96
19) n-Nitroso-di-n-propyla...	7.222	70	382180	36.152	ng	92
20) 3+4-Methylphenols	7.210	107	549052	36.976	ng	# 77
22) Acetophenone	7.216	105	677114	36.166	ng	# 91
24) Nitrobenzene	7.387	77	580742	41.422	ng	97
25) Isophorone	7.628	82	1102902	38.965	ng	98
26) 2-Nitrophenol	7.698	139	289674	48.500	ng	# 90
27) 2,4-Dimethylphenol	7.740	122	486489	38.520	ng	84
28) bis(2-Chloroethoxy)met...	7.840	93	678397	40.066	ng	99
29) 2,4-Dichlorophenol	7.940	162	476183	41.542	ng	97
30) 1,2,4-Trichlorobenzene	8.022	180	492651	40.304	ng	99
31) Naphthalene	8.104	128	1564141	38.376	ng	99
32) Benzoic acid	7.869	122	346965	43.080	ng	92
33) 4-Chloroaniline	8.157	127	700320	38.562	ng	97
34) Hexachlorobutadiene	8.222	225	285906	40.550	ng	94
35) Caprolactam	8.534	113	157355	42.550	ng	99
36) 4-Chloro-3-methylphenol	8.640	107	492140	40.541	ng	91
37) 2-Methylnaphthalene	8.798	142	1070265	39.615	ng	99
38) 1-Methylnaphthalene	8.898	142	995160	39.613	ng	97
40) 1,2,4,5-Tetrachloroben...	8.963	216	496436	40.741	ng	99
41) Hexachlorocyclopentadiene	8.945	237	223681	37.749	ng	97
43) 2,4,6-Trichlorophenol	9.075	196	337031	42.263	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.116	196	384396	42.175	ng	90
46) 1,1'-Biphenyl	9.263	154	1268424	38.927	ng	98
47) 2-Chloronaphthalene	9.287	162	960355	39.333	ng	98
48) 2-Nitroaniline	9.387	65	325801	43.144	ng	86
49) Acenaphthylene	9.704	152	1483800	38.661	ng	99
50) Dimethylphthalate	9.569	163	1180591	41.337	ng	98
51) 2,6-Dinitrotoluene	9.628	165	262662	45.048	ng	87
52) Acenaphthene	9.875	154	984172m	39.570	ng	
53) 3-Nitroaniline	9.798	138	290426	45.624	ng	96
54) 2,4-Dinitrophenol	9.904	184	111102	55.698	ng	87
55) Dibenzofuran	10.045	168	1376110	38.867	ng	95
56) 4-Nitrophenol	9.957	139	219929	42.242	ng	# 73
57) 2,4-Dinitrotoluene	10.034	165	324761	46.071	ng	# 89
58) Fluorene	10.392	166	1004298	38.943	ng	96
59) 2,3,4,6-Tetrachlorophenol	10.163	232	307964m	42.709	ng	
60) Diethylphthalate	10.269	149	1099604	40.620	ng	95
61) 4-Chlorophenyl-phenyle...	10.381	204	522115	41.189	ng	95
62) 4-Nitroaniline	10.416	138	286414	45.904	ng	# 85
63) Azobenzene	10.545	77	1108912	38.126	ng	98
65) 4,6-Dinitro-2-methylph...	10.445	198	171217	54.170	ng	# 79
66) n-Nitrosodiphenylamine	10.504	169	963255	39.675	ng	99
67) 4-Bromophenyl-phenylether	10.875	248	355416	42.564	ng	95
68) Hexachlorobenzene	10.939	284	374407	42.462	ng	93
69) Atrazine	11.034	200	266921	40.789	ng	98
70) Pentachlorophenol	11.134	266	236009	43.229	ng	97
71) Phenanthrene	11.357	178	1579427	38.983	ng	100
72) Anthracene	11.410	178	1610245	38.909	ng	98
73) Carbazole	11.563	167	1405277	38.030	ng	99
74) Di-n-butylphthalate	11.898	149	1642219	41.832	ng	98
75) Fluoranthene	12.545	202	1614843	37.732	ng	96
77) Benzidine	12.669	184	361266	36.931	ng	99
78) Pyrene	12.775	202	1632439	41.864	ng	97
80) Butylbenzylphthalate	13.404	149	601906	45.649	ng	# 98
81) Benzo(a)anthracene	13.969	228	1184301	36.754	ng	99
82) 3,3'-Dichlorobenzidine	13.933	252	400145	42.842	ng	# 97
83) Chrysene	14.010	228	1185889	37.776	ng	100
84) Bis(2-ethylhexyl)phtha...	13.963	149	649671	41.869	ng	100
85) Di-n-octyl phthalate	14.580	149	1128181	47.812	ng	99
87) Indeno(1,2,3-cd)pyrene	16.951	276	1263108	47.986	ng	# 91
88) Benzo(b)fluoranthene	15.022	252	1029746	37.724	ng	# 97
89) Benzo(k)fluoranthene	15.051	252	1055302	38.809	ng	# 98
90) Benzo(a)pyrene	15.392	252	1020840	40.110	ng	# 96
91) Dibenzo(a,h)anthracene	16.974	278	1033949	48.604	ng	# 96
92) Benzo(g,h,i)perylene	17.410	276	1058890	48.797	ng	# 92

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

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