

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF083018\
 Data File : BF108638.D
 Acq On : 30 Aug 2018 13:27
 Operator : JU/SJ
 Sample : SSTDICC080
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC080

Manual Integrations
 APPROVED

Sohil
 8/31/2018 12:18:03 PM

Quant Time: Aug 30 15:03:41 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 12:33:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.01	152	126089	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	522661	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	270608	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	558243	20.00	ng	0.00
75) Chrysene-d12	14.21	240	347795	20.00	ng	0.00
86) Perylene-d12	15.76	264	271119	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.65	112	1151643	165.36	ng	0.00
7) Phenol-d6	6.67	99	1546618	167.11	ng	0.02
23) Nitrobenzene-d5	7.58	82	1549923	165.48	ng	0.01
41) 2,4,6-Tribromophenol	10.85	330	537394	199.09	ng	0.00
44) 2-Fluorobiphenyl	9.37	172	2764386	164.01	ng	0.00
78) Terphenyl-d14	13.14	244	2788515	203.86	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.91	88	258391	72.204	ng	90
3) Pyridine	3.70	79	618331	67.075	ng	86
4) n-Nitrosodimethylamine	3.68	42	300276	57.888	ng	# 81
6) Aniline	6.70	93	878681	69.799	ng	# 82
8) 2-Chlorophenol	6.81	128	668872	85.273	ng	96
10) Phenol	6.68	94	922377	96.351	ng	# 67
11) bis(2-Chloroethyl)ether	6.75	93	767766	99.202	ng	94
12) 1,3-Dichlorobenzene	6.95	146	786429	86.710	ng	99
13) 1,4-Dichlorobenzene	7.02	146	844092	92.631	ng	98
14) 1,2-Dichlorobenzene	7.18	146	737234	86.978	ng	99
15) Benzyl Alcohol	7.16	79	589912	75.715	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.27	45	1162180	72.892	ng	53
17) 2-Methylphenol	7.27	107	549217	81.648	ng	# 73
18) Hexachloroethane	7.51	117	287718	88.688	ng	98
19) n-Nitroso-di-n-propylamine	7.42	70	527921	84.353	ng	95
20) 3+4-Methylphenols	7.42	107	651520	75.583	ng	# 61
22) Acetophenone	7.42	105	977151	79.956	ng	# 83
24) Nitrobenzene	7.60	77	809018	85.416	ng	93
25) Isophorone	7.84	82	1346573	75.426	ng	97
26) 2-Nitrophenol	7.91	139	373647	104.462	ng	# 84
27) 2,4-Dimethylphenol	7.94	122	520821	65.730	ng	95
28) bis(2-Chloroethoxy)methane	8.03	93	820489	78.726	ng	99
29) 2,4-Dichlorophenol	8.15	162	633470	86.874	ng	99
30) 1,2,4-Trichlorobenzene	8.22	180	687422	85.506	ng	97
31) Naphthalene	8.31	128	1835680	77.228	ng	98
32) Benzoic acid	8.13	122	340201m	80.965	ng	
33) 4-Chloroaniline	8.37	127	815650	78.741	ng	97
34) Hexachlorobutadiene	8.41	225	447754	87.977	ng	99
35) Caprolactam	8.78	113	143759	62.912	ng	# 73
36) 4-Chloro-3-methylphenol	8.86	107	669837	85.153	ng	96
37) 2-Methylnaphthalene	9.00	142	1248922	74.265	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.17	216	721447	86.816	ng	97
40) Hexachlorocyclopentadiene	9.14	237	326801	60.884	ng	99
42) 2,4,6-Trichlorophenol	9.29	196	455303	88.291	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.34	196	482921	85.070	nq	95
45) 1,1'-Biphenyl	9.47	154	1701650	85.288	nq	98
46) 2-Chloronaphthalene	9.50	162	1353003	85.800	nq	99
47) 2-Nitroaniline	9.61	65	470340	92.181	nq	85
48) Acenaphthylene	9.92	152	1914037	76.776	nq	98
49) Dimethylphthalate	9.78	163	1610617	85.444	nq	# 99
50) 2,6-Dinitrotoluene	9.85	165	352932	89.490	nq	# 87
51) Acenaphthene	10.09	154	1186542	77.707	nq	98
52) 3-Nitroaniline	10.03	138	383766	88.937	nq	94
53) 2,4-Dinitrophenol	10.13	184	142679	115.112	nq	# 26
54) Dibenzofuran	10.27	168	1767729	79.908	nq	96
56) 2,4-Dinitrotoluene	10.26	165	445962	87.849	nq	# 80
57) Fluorene	10.61	166	1398223	79.842	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.38	232	416674	87.818	nq	# 86
59) Diethylphthalate	10.47	149	1549757	84.903	nq	95
60) 4-Chlorophenyl-phenylether	10.59	204	797884	87.438	nq	95
61) 4-Nitroaniline	10.67	138	350694	82.204	nq	88
62) Azobenzene	10.75	77	1521883	80.734	nq	92
64) 4,6-Dinitro-2-methylphenol	10.68	198	208499	108.361	nq	88
65) n-Nitrosodiphenylamine	10.72	169	1378335	83.553	nq	97
66) 4-Bromophenyl-phenylether	11.08	248	521237	85.919	nq	# 88
67) Hexachlorobenzene	11.15	284	564324	88.410	nq	# 88
68) Atrazine	11.24	200	279395	50.496	nq	97
69) Pentachlorophenol	11.35	266	319058	104.431	nq	96
70) Phenanthrene	11.58	178	2025676	76.933	nq	99
71) Anthracene	11.63	178	2089606	77.431	nq	97
72) Carbazole	11.79	167	2010366	83.846	nq	98
73) Di-n-butylphthalate	12.09	149	2276920	83.839	nq	# 98
74) Fluoranthene	12.77	202	2142703	74.564	nq	95
77) Pyrene	13.01	202	2155457	97.891	nq	98
79) Butylbenzylphthalate	13.59	149	887487	101.504	nq	94
80) Benzo(a)anthracene	14.19	228	1603440	80.333	nq	99
81) 3,3'-Dichlorobenzidine	14.15	252	490905	68.628	nq	# 98
82) Chrysene	14.24	228	1500372	81.992	nq	99
83) Bis(2-ethylhexyl)phthalate	14.14	149	1111463	93.872	nq	99
84) Di-n-octyl phthalate	14.77	149	1538291	85.189	nq	95
85) Indeno(1,2,3-cd)pyrene	17.40	276	1209108	76.259	nq	# 89
87) Benzo(b)fluoranthene	15.29	252	1193316	73.659	nq	# 95
88) Benzo(k)fluoranthene	15.33	252	1246335	83.691	nq	# 95
89) Benzo(a)pyrene	15.69	252	1144171	78.870	nq	97
90) Dibenzo(a,h)anthracene	17.41	278	1010018	84.146	nq	# 93
91) Benzo(g,h,i)perylene	17.89	276	1002092	84.784	nq	# 88

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

