

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090518\
 Data File : BF108775.D
 Acq On : 5 Sep 2018 15:41
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Quant Time: Sep 05 16:38:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 16:28:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	290973	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	1179960	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	583037	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	1124262	20.00	ng	0.00
75) Chrysene-d12	13.87	240	804946	20.00	ng	0.00
86) Perylene-d12	15.27	264	781206	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	898852	53.66	ng	0.00
7) Phenol-d6	6.36	99	1157859	48.27	ng	0.00
23) Nitrobenzene-d5	7.30	82	976330	42.00	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	293513	38.13	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1912824	44.11	ng	0.00
78) Terphenyl-d14	12.82	244	1654199	39.93	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.42	88	211396	27.157	ng	93
3) Pyridine	3.12	79	596138	33.147	ng	88
4) n-Nitrosodimethylamine	3.05	42	220808	24.613	ng	91
6) Aniline	6.39	93	788899	30.475	ng	98
8) 2-Chlorophenol	6.51	128	508643	24.837	ng	100
9) Benzaldehyde	6.28	77	416113	28.064	ng	99
10) Phenol	6.37	94	667680	23.867	ng	99
11) bis(2-Chloroethyl)ether	6.47	93	522485	21.210	ng	98
12) 1,3-Dichlorobenzene	6.67	146	565997	23.836	ng	98
13) 1,4-Dichlorobenzene	6.75	146	569479	21.971	ng	99
14) 1,2-Dichlorobenzene	6.91	146	536704	22.546	ng	97
15) Benzyl Alcohol	6.87	79	455191	26.378	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.01	45	792324	20.966	ng	99
17) 2-Methylphenol	6.98	107	413383	25.962	ng	97
18) Hexachloroethane	7.25	117	203996	22.885	ng	94
19) n-Nitroso-di-n-propylamine	7.15	70	392487	24.431	ng	98
20) 3+4-Methylphenols	7.14	107	498091	24.967	ng	# 71
22) Acetophenone	7.14	105	678364	23.029	ng	94
24) Nitrobenzene	7.31	77	527190	22.366	ng	99
25) Isophorone	7.55	82	938744	24.309	ng	99
26) 2-Nitrophenol	7.63	139	185383	17.373	ng	97
27) 2,4-Dimethylphenol	7.67	122	413232	26.628	ng	96
28) bis(2-Chloroethoxy)methane	7.77	93	643001	26.358	ng	99
29) 2,4-Dichlorophenol	7.87	162	430097	23.361	ng	97
30) 1,2,4-Trichlorobenzene	7.96	180	464342	22.554	ng	97
31) Naphthalene	8.04	128	1390818	24.747	ng	99
32) Benzoic acid	7.77	122	228666	22.977	ng	99
33) 4-Chloroaniline	8.08	127	637264	25.686	ng	99
34) Hexachlorobutadiene	8.17	225	272897	20.188	ng	99
35) Caprolactam	8.44	113	144201	29.394	ng	91
36) 4-Chloro-3-methylphenol	8.56	107	462974	23.430	ng	99
37) 2-Methylnaphthalene	8.72	142	920419	24.205	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	451992	21.837	ng	99
40) Hexachlorocyclopentadiene	8.89	237	221463	30.209	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.00	196	311925	25.448	nq	99
43) 2,4,5-Trichlorophenol	9.04	196	317103	22.185	nq	96
45) 1,1'-Biphenyl	9.19	154	1184414	23.335	nq	99
46) 2-Chloronaphthalene	9.21	162	948014	23.858	nq	98
47) 2-Nitroaniline	9.30	65	244189	18.374	nq	96
48) Acenaphthylene	9.62	152	1433027	24.633	nq	99
49) Dimethylphthalate	9.49	163	1065545	23.209	nq	99
50) 2,6-Dinitrotoluene	9.54	165	208641	20.617	nq	88
51) Acenaphthene	9.80	154	872207	25.075	nq	97
52) 3-Nitroaniline	9.71	138	251888	23.008	nq	# 87
53) 2,4-Dinitrophenol	9.81	184	45978	13.361	nq	# 16
54) Dibenzofuran	9.97	168	1287212	23.650	nq	97
55) 4-Nitrophenol	9.86	139	182042	30.237	nq	95
56) 2,4-Dinitrotoluene	9.94	165	245862	18.348	nq	# 89
57) Fluorene	10.31	166	875888	20.765	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.08	232	260809	20.368	nq	91
59) Diethylphthalate	10.19	149	1096124	23.697	nq	97
60) 4-Chlorophenyl-phenylether	10.31	204	469069	19.598	nq	93
61) 4-Nitroaniline	10.32	138	217757	20.555	nq	96
62) Azobenzene	10.47	77	1134820	24.683	nq	94
64) 4,6-Dinitro-2-methylphenol	10.35	198	85284	17.096	nq	92
65) n-Nitrosodiphenylamine	10.42	169	953006	25.514	nq	96
66) 4-Bromophenyl-phenylether	10.79	248	324151	23.459	nq	92
67) Hexachlorobenzene	10.86	284	341896	22.967	nq	# 88
68) Atrazine	10.95	200	299531	28.684	nq	98
69) Pentachlorophenol	11.04	266	218438	27.615	nq	98
70) Phenanthrene	11.27	178	1393231	24.552	nq	99
71) Anthracene	11.31	178	1426053	23.917	nq	99
72) Carbazole	11.47	167	1390336	24.115	nq	99
73) Di-n-butylphthalate	11.81	149	1657307	24.985	nq	99
74) Fluoranthene	12.44	202	1465250	22.939	nq	100
76) Benzidine	12.57	184	940758	41.519	nq	99
77) Pyrene	12.67	202	1494041	24.691	nq	99
79) Butylbenzylphthalate	13.30	149	672122	26.416	nq	98
80) Benzo(a)anthracene	13.85	228	1276708	26.017	nq	99
81) 3,3'-Dichlorobenzidine	13.82	252	515645	28.975	nq	97
82) Chrysene	13.89	228	1223848	25.934	nq	99
83) Bis(2-ethylhexyl)phthalate	13.87	149	855709	26.165	nq	100
84) Di-n-octyl phthalate	14.47	149	1458156	30.325	nq	100
85) Indeno(1,2,3-cd)pyrene	16.61	276	983271	29.966	nq	97
87) Benzo(b)fluoranthene	14.87	252	1125817	23.460	nq	97
88) Benzo(k)fluoranthene	14.89	252	989438	20.813	nq	98
89) Benzo(a)pyrene	15.21	252	976395	22.514	nq	98
90) Dibenzo(a,h)anthracene	16.63	278	828943	23.933	nq	98
91) Benzo(g,h,i)perylene	17.02	276	819861	24.354	nq	94

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

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