

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061918\
 Data File : BF106573.D
 Acq On : 19 Jun 2018 10:50
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 6/20/2018 4:58:00 PM

Quant Time: Jun 19 12:18:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 12:10:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	114442	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	458857	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	223290	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	450299	20.00	ng	0.00
75) Chrysene-d12	14.15	240	426953	20.00	ng	0.00
86) Perylene-d12	15.69	264	396226	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	155093	22.94	ng	0.00
7) Phenol-d6	6.60	99	191694	22.44	ng	0.00
23) Nitrobenzene-d5	7.52	82	182772	23.43	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	55771	25.96	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	371975	28.41	ng	0.00
78) Terphenyl-d14	13.08	244	432818	25.18	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.84	88	36623	10.465	ng	98
3) Pyridine	3.62	79	94460m	9.686	ng	
4) n-Nitrosodimethylamine	3.56	42	39919	8.936	ng	94
6) Aniline	6.62	93	126900	10.173	ng	98
8) 2-Chlorophenol	6.75	128	83214	11.444	ng	98
9) Benzaldehyde	6.51	77	71095	10.972	ng	96
10) Phenol	6.61	94	108740	11.660	ng	82
11) bis(2-Chloroethyl)ether	6.69	93	85619	10.724	ng	97
12) 1,3-Dichlorobenzene	6.90	146	95126	11.115	ng	98
13) 1,4-Dichlorobenzene	6.98	146	98210	11.308	ng	96
14) 1,2-Dichlorobenzene	7.13	146	92363	11.298	ng	96
15) Benzyl Alcohol	7.10	79	77117	10.541	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.23	45	113430	9.911	ng	96
17) 2-Methylphenol	7.21	107	67614	10.422	ng	96
18) Hexachloroethane	7.47	117	33227	10.738	ng	93
19) n-Nitroso-di-n-propylamine	7.36	70	66082	10.311	ng	95
20) 3+4-Methylphenols	7.37	107	90202	10.872	ng	# 54
22) Acetophenone	7.37	105	127395	11.017	ng	# 93
24) Nitrobenzene	7.54	77	92747	10.940	ng	92
25) Isophorone	7.77	82	155017	10.175	ng	94
26) 2-Nitrophenol	7.85	139	40525	12.315	ng	92
27) 2,4-Dimethylphenol	7.90	122	67631	10.486	ng	96
28) bis(2-Chloroethoxy)methane	7.98	93	107840	10.917	ng	96
29) 2,4-Dichlorophenol	8.11	162	69697	11.201	ng	97
30) 1,2,4-Trichlorobenzene	8.18	180	77087	11.049	ng	94
31) Naphthalene	8.27	128	247940	11.953	ng	99
32) Benzoic acid	7.97	122	30479m	7.119	ng	
33) 4-Chloroaniline	8.31	127	98483	10.714	ng	97
34) Hexachlorobutadiene	8.38	225	50640	11.298	ng	99
35) Caprolactam	8.66	113	21677	10.300	ng	98
36) 4-Chloro-3-methylphenol	8.80	107	74504	10.825	ng	97
37) 2-Methylnaphthalene	8.96	142	163550	11.576	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.12	216	85038	11.732	ng	97
40) Hexachlorocyclopentadiene	9.11	237	33500	8.376	ng	97

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061918\
 Data File : BF106573.D
 Acq On : 19 Jun 2018 10:50
 Operator : JU/SJ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Sohil
 6/20/2018 4:58:00 PM

Quant Time: Jun 19 12:18:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 19 12:10:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.24	196	54256	11.745	nq	98
43) 2,4,5-Trichlorophenol	9.28	196	56483	11.342	nq	95
45) 1,1'-Biphenyl	9.42	154	208168	12.036	nq	98
46) 2-Chloronaphthalene	9.45	162	159328	11.512	nq	95
47) 2-Nitroaniline	9.54	65	50699	11.644	nq	88
48) Acenaphthylene	9.87	152	256771	12.121	nq	99
49) Dimethylphthalate	9.71	163	184186	11.546	nq	99
50) 2,6-Dinitrotoluene	9.78	165	37851	11.616	nq	95
51) Acenaphthene	10.04	154	158330	11.573	nq	98
52) 3-Nitroaniline	9.95	138	43216	11.174	nq	96
53) 2,4-Dinitrophenol	10.06	184	10839	8.998	nq #	14
54) Dibenzofuran	10.21	168	221091	12.001	nq	99
55) 4-Nitrophenol	10.13	139	29314	9.888	nq	89
56) 2,4-Dinitrotoluene	10.18	165	49135	12.007	nq	97
57) Fluorene	10.55	166	181278	12.419	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.33	232	43738	11.081	nq #	84
59) Diethylphthalate	10.41	149	183340	11.579	nq	97
60) 4-Chlorophenyl-phenylether	10.54	204	91725	11.948	nq #	88
61) 4-Nitroaniline	10.57	138	43395	11.532	nq	99
62) Azobenzene	10.70	77	185126	11.213	nq	97
64) 4,6-Dinitro-2-methylphenol	10.60	198	23153	10.521	nq	78
65) n-Nitrosodiphenylamine	10.66	169	174478	11.529	nq	94
66) 4-Bromophenyl-phenylether	11.04	248	59760	11.668	nq #	81
67) Hexachlorobenzene	11.10	284	60389	11.437	nq #	86
68) Atrazine	11.18	200	53728	11.520	nq	95
69) Pentachlorophenol	11.30	266	21470	6.605	nq	97
70) Phenanthrene	11.52	178	265051	12.019	nq	99
71) Anthracene	11.57	178	268969	12.174	nq	99
72) Carbazole	11.72	167	241128	11.822	nq	99
73) Di-n-butylphthalate	12.04	149	283974	12.510	nq	99
74) Fluoranthene	12.71	202	293687	12.488	nq	95
76) Benzidine	12.83	184	145664	10.251	nq	99
77) Pyrene	12.94	202	303478	11.354	nq	99
79) Butylbenzylphthalate	13.55	149	120704	12.064	nq	97
80) Benzo(a)anthracene	14.14	228	282624	11.124	nq	99
81) 3,3'-Dichlorobenzidine	14.09	252	106325	12.403	nq	99
82) Chrysene	14.18	228	264374	11.397	nq	98
83) Bis(2-ethylhexyl)phthalate	14.11	149	165491	11.535	nq	98
84) Di-n-octyl phthalate	14.74	149	265408	12.833	nq	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	221350	11.956	nq #	93
87) Benzo(b)fluoranthene	15.23	252	255213	10.655	nq	98
88) Benzo(k)fluoranthene	15.27	252	249261	10.508	nq #	95
89) Benzo(a)pyrene	15.62	252	227905	10.579	nq	98
90) Dibenzo(a,h)anthracene	17.27	278	186708	10.395	nq	96
91) Benzo(g,h,i)perylene	17.73	276	176133	10.091	nq #	90

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

