

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091018\
 Data File : BF108887.D
 Acq On : 10 Sep 2018 12:37
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 9/11/2018 9:25:47 AM

Quant Time: Sep 11 02:15:00 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 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	199029	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	790496	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	374705	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	740081	20.00	ng	0.00
75) Chrysene-d12	13.87	240	503208	20.00	ng	0.00
86) Perylene-d12	15.27	264	447747	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.34	112	952497	79.28	ng	0.00
7) Phenol-d6	6.37	99	1222914	79.11	ng	0.00
23) Nitrobenzene-d5	7.30	82	1075083	85.11	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	304309	85.07	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1900638	86.53	ng	0.00
78) Terphenyl-d14	12.82	244	1644373	76.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.41	88	236380	40.941	ng	93
3) Pyridine	3.12	79	649392	41.070	ng	90
4) n-Nitrosodimethylamine	3.05	42	236295	38.908	ng	# 92
6) Aniline	6.40	93	810538	38.986	ng	98
8) 2-Chlorophenol	6.52	128	526860	39.620	ng	99
9) Benzaldehyde	6.28	77	423659	38.633	ng	98
10) Phenol	6.38	94	691418	39.053	ng	91
11) bis(2-Chloroethyl)ether	6.47	93	545656	38.473	ng	98
12) 1,3-Dichlorobenzene	6.68	146	604831	39.858	ng	98
13) 1,4-Dichlorobenzene	6.75	146	607887	40.058	ng	99
14) 1,2-Dichlorobenzene	6.91	146	568695	40.657	ng	100
15) Benzyl Alcohol	6.88	79	474761	40.021	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.02	45	808892	38.154	ng	97
17) 2-Methylphenol	7.00	107	430700	38.326	ng	98
18) Hexachloroethane	7.25	117	218847	39.995	ng	98
19) n-Nitroso-di-n-propylamine	7.15	70	396058	36.861	ng	99
20) 3+4-Methylphenols	7.15	107	511592	38.862	ng	# 59
22) Acetophenone	7.15	105	682715	38.055	ng	# 91
24) Nitrobenzene	7.32	77	564926	41.682	ng	99
25) Isophorone	7.56	82	934912	37.106	ng	98
26) 2-Nitrophenol	7.64	139	216755	40.594	ng	94
27) 2,4-Dimethylphenol	7.68	122	423252	39.034	ng	96
28) bis(2-Chloroethoxy)methane	7.77	93	649939	38.582	ng	99
29) 2,4-Dichlorophenol	7.88	162	445052	39.806	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	477666	39.422	ng	98
31) Naphthalene	8.04	128	1437490	40.419	ng	99
32) Benzoic acid	7.79	122	235616m	35.493	ng	
33) 4-Chloroaniline	8.09	127	652683	39.959	ng	97
34) Hexachlorobutadiene	8.17	225	291829	40.316	ng	99
35) Caprolactam	8.46	113	146869	38.766	ng	# 79
36) 4-Chloro-3-methylphenol	8.57	107	462613	38.841	ng	97
37) 2-Methylnaphthalene	8.73	142	931355	39.640	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	465019	41.516	ng	98
40) Hexachlorocyclopentadiene	8.89	237	223617	39.684	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.01	196	301687	39.543	nq	99
43) 2,4,5-Trichlorophenol	9.04	196	327119	41.364	nq	98
45) 1,1'-Biphenyl	9.20	154	1178393	40.849	nq	99
46) 2-Chloronaphthalene	9.22	162	953170	40.856	nq	98
47) 2-Nitroaniline	9.31	65	281471	44.163	nq	96
48) Acenaphthylene	9.63	152	1422788	40.914	nq	98
49) Dimethylphthalate	9.50	163	1049641	40.050	nq	99
50) 2,6-Dinitrotoluene	9.55	165	224541	44.963	nq	96
51) Acenaphthene	9.80	154	874221	40.624	nq	99
52) 3-Nitroaniline	9.72	138	276193	46.461	nq	98
53) 2,4-Dinitrophenol	9.82	184	63157	39.725	nq #	21
54) Dibenzofuran	9.97	168	1271275	40.527	nq	97
55) 4-Nitrophenol	9.87	139	188978	42.523	nq	94
56) 2,4-Dinitrotoluene	9.95	165	291852	45.698	nq #	94
57) Fluorene	10.31	166	873293	43.382	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.09	232	270404	42.400	nq #	86
59) Diethylphthalate	10.20	149	1046683	38.700	nq	96
60) 4-Chlorophenyl-phenylether	10.31	204	467500	43.484	nq	96
61) 4-Nitroaniline	10.32	138	260794	48.427	nq	98
62) Azobenzene	10.47	77	1096608	39.280	nq	98
64) 4,6-Dinitro-2-methylphenol	10.36	198	108408	38.787	nq	86
65) n-Nitrosodiphenylamine	10.42	169	929156	37.761	nq	96
66) 4-Bromophenyl-phenylether	10.80	248	323263	38.673	nq #	89
67) Hexachlorobenzene	10.87	284	342924	38.431	nq #	89
68) Atrazine	10.95	200	305485	38.781	nq	99
69) Pentachlorophenol	11.05	266	212236	38.904	nq	99
70) Phenanthrene	11.27	178	1379007	39.019	nq	98
71) Anthracene	11.32	178	1413040	41.759	nq	100
72) Carbazole	11.47	167	1418178	39.796	nq	99
73) Di-n-butylphthalate	11.81	149	1598992	40.405	nq	99
74) Fluoranthene	12.45	202	1452506	41.509	nq	97
76) Benzidine	12.57	184	839417	35.196	nq	99
77) Pyrene	12.68	202	1516868	39.174	nq	99
79) Butylbenzylphthalate	13.30	149	659874	37.843	nq	95
80) Benzo(a)anthracene	13.86	228	1181330	36.626	nq	98
81) 3,3'-Dichlorobenzidine	13.82	252	475926	37.873	nq	98
82) Chrysene	13.90	228	1193402	38.220	nq	100
83) Bis(2-ethylhexyl)phthalate	13.87	149	780274	35.607	nq	99
84) Di-n-octyl phthalate	14.48	149	1310531	35.872	nq	100
85) Indeno(1,2,3-cd)pyrene	16.63	276	942431	34.105	nq	95
87) Benzo(b)fluoranthene	14.88	252	965745	39.667	nq	98
88) Benzo(k)fluoranthene	14.91	252	965098	43.045	nq	97
89) Benzo(a)pyrene	15.22	252	866537	39.445	nq	99
90) Dibenzo(a,h)anthracene	16.65	278	781518	40.373	nq	97
91) Benzo(g,h,i)perylene	17.04	276	798192	41.183	nq	93

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

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