

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011720\
 Data File : BF118558.D
 Acq On : 17 Jan 2020 13:26
 Operator : JU
 Sample : SSTDICC100
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC100

Manual Integrations
 APPROVED

Quant Time: Jan 17 15:51:18 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 17 12:09:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	448710	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	1522725	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	838450	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	1448520	20.00	ng	0.00
76) Chrysene-d12	14.06	240	862912	20.00	ng	0.00
87) Perylene-d12	15.53	264	895072	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	3990711	152.25	ng	0.01
7) Phenol-d6	6.53	99	5240050	166.16	ng	0.02
23) Nitrobenzene-d5	7.47	82	4920418	197.88	ng	0.01
42) 2,4,6-Tribromophenol	10.73	330	1688469	166.03	ng	0.01
45) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
79) Terphenyl-d14	13.00	244	6990323	134.25	ng	0.00

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Target Compounds

						Qvalue
2) 1,4-Dioxane	2.68	88	1051033	98.573	ng	99
3) Pyridine	3.45	79	2906101	103.276	ng	99
4) n-Nitrosodimethylamine	3.41	42	1340858	137.317	ng	98
6) Aniline	6.56	93	3605105m	90.605	ng	
8) 2-Chlorophenol	6.69	128	2329001	79.267	ng	98
9) Benzaldehyde	6.44	77	1032740	70.317	ng	99
10) Phenol	6.55	94	2980988	84.067	ng	84
11) bis(2-Chloroethyl)ether	6.64	93	2376541	89.144	ng	98
12) 1,3-Dichlorobenzene	6.84	146	2703819	79.398	ng	98
13) 1,4-Dichlorobenzene	6.92	146	2679432	78.523	ng	98
14) 1,2-Dichlorobenzene	7.07	146	2372009	75.032	ng	97
15) Benzyl Alcohol	7.04	79	2120497	94.341	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	3348653	121.305	ng	97
17) 2-Methylphenol	7.15	107	2048719	89.361	ng	100
18) Hexachloroethane	7.42	117	1018021	84.078	ng	94
19) n-Nitroso-di-n-propylamine	7.33	70	1908074	104.592	ng	98
20) 3+4-Methylphenols	7.31	107	2150917	72.065	ng	# 89
22) Acetophenone	7.32	105	3067208	86.826	ng	98
24) Nitrobenzene	7.49	77	2535939	102.602	ng	96
25) Isophorone	7.73	82	4848240	108.199	ng	99
26) 2-Nitrophenol	7.79	139	1218915	84.051	ng	98
27) 2,4-Dimethylphenol	7.83	122	1989635	102.779	ng	99
28) bis(2-Chloroethoxy)methane	7.93	93	3000054	97.123	ng	98
29) 2,4-Dichlorophenol	8.04	162	2172766	96.486	ng	99
30) 1,2,4-Trichlorobenzene	8.12	180	2388405	93.182	ng	99
31) Naphthalene	8.20	128	5699513	74.611	ng	93
32) Benzoic acid	8.00	122	1620791	115.951	ng	94
33) 4-Chloroaniline	8.25	127	3030119	93.221	ng	99
34) Hexachlorobutadiene	8.32	225	1473614	97.859	ng	98
35) Caprolactam	8.66	113	746060m	107.505	ng	
36) 4-Chloro-3-methylphenol	8.73	107	2091779	90.002	ng	96
37) 2-Methylnaphthalene	8.89	142	4114675	80.575	ng	98
38) 1-Methylnaphthalene	8.99	142	3921717	80.991	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	2264125	88.062	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)	
41) Hexachlorocyclopentadiene	9.05	237	1314117	338.191	nq	97	
43) 2,4,6-Trichlorophenol	9.17	196	1645181	98.814	nq	99	
44) 2,4,5-Trichlorophenol	9.21	196	1615578	97.166	nq	99	
46) 1,1'-Biphenyl	9.36	154	4677086	70.400	nq	94	
47) 2-Chloronaphthalene	9.39	162	4218428	79.084	nq	92	
48) 2-Nitroaniline	9.47	65	1430924	99.886	nq	95	mohammad
49) Acenaphthylene	9.80	152	5709784	73.462	nq	94	
50) Dimethylphthalate	9.67	163	4982404	80.511	nq	97	
51) 2,6-Dinitrotoluene	9.72	165	1194275	87.768	nq	99	
52) Acenaphthene	9.97	154	3912536	81.817	nq	98	
53) 3-Nitroaniline	9.89	138	1296992	82.407	nq	95	1/21/2020 7:58:28 AM
54) 2,4-Dinitrophenol	9.99	184	500335	98.452	nq	1	#
55) Dibenzofuran	10.15	168	5298763	76.364	nq	89	
56) 4-Nitrophenol	10.04	139	996108	138.079	nq	95	
57) 2,4-Dinitrotoluene	10.13	165	1515165	87.113	nq	91	
58) Fluorene	10.49	166	3893064	69.069	nq	99	
59) 2,3,4,6-Tetrachlorophenol	10.26	232	1345707	98.971	nq	99	
60) Diethylphthalate	10.36	149	4605797	74.693	nq	93	
61) 4-Chlorophenyl-phenylether	10.47	204	2172366	76.606	nq	96	
62) 4-Nitroaniline	10.52	138	1271122	81.787	nq	96	
63) Azobenzene	10.64	77	4370590	85.573	nq	93	
65) 4,6-Dinitro-2-methylphenol	10.53	198	681188	89.123	nq	83	
66) n-Nitrosodiphenylamine	10.60	169	3996629	84.037	nq	96	
67) 4-Bromophenyl-phenylether	10.97	248	1728835	99.008	nq	97	
68) Hexachlorobenzene	11.04	284	1892587	91.413	nq	95	
70) Pentachlorophenol	11.22	266	1008156	176.440	nq	99	
71) Phenanthrene	11.45	178	5708451	73.208	nq	95	
72) Anthracene	11.50	178	5682793	73.122	nq	95	
73) Carbazole	11.65	167	5220778	75.194	nq	96	
77) Benzidine	12.75	184	1510338	78.417	nq	95	#
78) Pyrene	12.86	202	5764553	79.059	nq	94	
80) Butylbenzylphthalate	13.47	149	2593612	79.463	nq	95	
81) Benzo(a)anthracene	14.05	228	4581704	76.600	nq	96	
82) 3,3'-Dichlorobenzidine	14.00	252	1535774	75.622	nq	98	
83) Chrysene	14.09	228	4536148	79.124	nq	96	
84) Bis(2-ethylhexyl)phthalate	14.04	149	2961926	67.923	nq	96	#
85) Di-n-octyl phthalate	14.65	149	4669899	71.573	nq	99	
86) Indeno(1,2,3-cd)pyrene	17.03	276	6258834	116.178	nq	99	
88) Benzo(b)fluoranthene	15.10	252	5127155	89.380	nq	98	
89) Benzo(k)fluoranthene	15.13	252	4392161	84.527	nq	96	
90) Benzo(a)pyrene	15.47	252	4571952	89.439	nq	96	
91) Dibenzo(a,h)anthracene	17.05	278	4993919	107.591	nq	98	
92) Benzo(g,h,i)perylene	17.48	276	5059311	112.043	nq	98	

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

