

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100220\
 Data File : BF121787.D
 Acq On : 2 Oct 2020 11:27
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 10/5/2020 10:27:14 AM

Quant Time: Oct 02 14:21:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 01 14:29:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	151478	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	559420	20.00	ng	0.00
39) Acenaphthene-d10	9.80	164	284139	20.00	ng	0.00
64) Phenanthrene-d10	11.29	188	519201	20.00	ng	0.00
76) Chrysene-d12	13.93	240	389797	20.00	ng	0.00
86) Perylene-d12	15.36	264	348807	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	809891	79.46	ng	0.00
7) Phenol-d6	6.42	99	1047615	78.34	ng	0.00
23) Nitrobenzene-d5	7.34	82	885381	84.98	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	295009	83.54	ng	-0.01
45) 2-Fluorobiphenyl	9.13	172	1464085	78.04	ng	0.00
79) Terphenyl-d14	12.87	244	1629259	84.68	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.49	88	181061	38.666	ng	99
3) Pyridine	3.21	79	519340	40.118	ng	96
4) n-Nitrosodimethylamine	3.18	42	239432	39.874	ng	100
6) Aniline	6.43	93	654748	37.594	ng	# 89
8) 2-Chlorophenol	6.56	128	415338	40.022	ng	95
9) Benzaldehyde	6.32	77	293666	33.593	ng	99
10) Phenol	6.43	94	532064	37.364	ng	90
11) bis(2-Chloroethyl)ether	6.51	93	427006	39.786	ng	96
12) 1,3-Dichlorobenzene	6.70	146	466008	40.231	ng	99
13) 1,4-Dichlorobenzene	6.78	146	467282	40.177	ng	99
14) 1,2-Dichlorobenzene	6.94	146	423163	39.495	ng	99
15) Benzyl Alcohol	6.92	79	362481	39.881	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.05	45	666618	38.598	ng	95
17) 2-Methylphenol	7.03	107	356819	40.410	ng	96
18) Hexachloroethane	7.27	117	171829	41.261	ng	99
19) n-Nitroso-di-n-propylamine	7.19	70	296616	38.487	ng	99
20) 3+4-Methylphenols	7.19	107	406408	38.307	ng	93
22) Acetophenone	7.18	105	559803	39.441	ng	# 97
24) Nitrobenzene	7.36	77	470397	41.743	ng	97
25) Isophorone	7.60	82	839692	40.035	ng	99
26) 2-Nitrophenol	7.67	139	221759	42.132	ng	96
27) 2,4-Dimethylphenol	7.71	122	369518	39.469	ng	99
28) bis(2-Chloroethoxy)methane	7.80	93	518432	39.928	ng	99
29) 2,4-Dichlorophenol	7.92	162	344661	40.419	ng	97
30) 1,2,4-Trichlorobenzene	7.99	180	364492	40.586	ng	99
31) Naphthalene	8.07	128	1171453	39.669	ng	99
32) Benzoic acid	7.86	122	199217	31.335	ng	99
33) 4-Chloroaniline	8.12	127	519306	40.568	ng	98
34) Hexachlorobutadiene	8.19	225	217196	41.203	ng	99
35) Caprolactam	8.51	113	113378	39.926	ng	97
36) 4-Chloro-3-methylphenol	8.62	107	370871	40.374	ng	97
37) 2-Methylnaphthalene	8.76	142	761512	39.349	ng	98
38) 1-Methylnaphthalene	8.86	142	705865	39.190	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.93	216	364062	41.018	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.92	237	174086	34.346	ng	99
43) 2,4,6-Trichlorophenol	9.05	196	240356	39.730	ng	99
44) 2,4,5-Trichlorophenol	9.09	196	260308	40.702	ng	96
46) 1,1'-Biphenyl	9.23	154	905126	39.791	ng	99
47) 2-Chloronaphthalene	9.25	162	712917	39.771	ng	99
48) 2-Nitroaniline	9.35	65	244659	43.922	ng	98
49) Acenaphthylene	9.67	152	1094745	39.748	ng	99
50) Dimethylphthalate	9.53	163	830867	40.345	ng	99
51) 2,6-Dinitrotoluene	9.60	165	188456	42.969	ng	88
52) Acenaphthene	9.84	154	640347	36.811	ng	100
53) 3-Nitroaniline	9.76	138	213894	42.099	ng	97
54) 2,4-Dinitrophenol	9.87	184	89958	40.588	ng	# 89
55) Dibenzofuran	10.01	168	934486	38.146	ng	99
56) 4-Nitrophenol	9.93	139	139474	34.422	ng	85
57) 2,4-Dinitrotoluene	10.00	165	237728	43.085	ng	96
58) Fluorene	10.35	166	741015	39.554	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.13	232	206590	39.734	ng	97
60) Diethylphthalate	10.23	149	790290	39.360	ng	100
61) 4-Chlorophenyl-phenylether	10.35	204	362993	40.448	ng	99
62) 4-Nitroaniline	10.38	138	209564	41.542	ng	99
63) Azobenzene	10.50	77	841315	39.347	ng	99
65) 4,6-Dinitro-2-methylphenol	10.41	198	138509	44.120	ng	90
66) n-Nitrosodiphenylamine	10.46	169	700579	39.331	ng	99
67) 4-Bromophenyl-phenylether	10.83	248	240956	40.762	ng	98
68) Hexachlorobenzene	10.90	284	272488	40.846	ng	# 91
69) Atrazine	10.99	200	170918	38.622	ng	99
70) Pentachlorophenol	11.10	266	117882m	32.176	ng	
71) Phenanthrene	11.32	178	1109800	39.232	ng	100
72) Anthracene	11.36	178	1150284	39.404	ng	100
73) Carbazole	11.52	167	1028874	39.057	ng	100
74) Di-n-butylphthalate	11.85	149	1215619	39.982	ng	99
75) Fluoranthene	12.50	202	1197734	39.663	ng	100
77) Benzidine	12.62	184	748116	57.892	ng	100
78) Pyrene	12.73	202	1201404	42.184	ng	99
80) Butylbenzylphthalate	13.35	149	508743	44.169	ng	98
81) Benzo(a)anthracene	13.92	228	1021952	41.441	ng	100
82) 3,3'-Dichlorobenzidine	13.88	252	405257	42.094	ng	99
83) Chrysene	13.96	228	1009407	40.422	ng	100
84) Bis(2-ethylhexyl)phthalate	13.90	149	670819	43.596	ng	99
85) Di-n-octyl phthalate	14.52	149	1103125	40.623	ng	100
87) Indeno(1,2,3-cd)pyrene	16.78	276	924040	40.679	ng	100
88) Benzo(b)fluoranthene	14.95	252	948390	40.672	ng	99
89) Benzo(k)fluoranthene	14.98	252	895780	41.950	ng	99
90) Benzo(a)pyrene	15.30	252	813460	41.040	ng	99
91) Dibenzo(a,h)anthracene	16.79	278	757197	40.044	ng	99
92) Benzo(g,h,i)perylene	17.21	276	773787	41.630	ng	98

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

