

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010220\
 Data File : BF118431.D
 Acq On : 2 Jan 2020 12:58
 Operator : JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 1/3/2020 11:04:01 AM

Quant Time: Jan 02 14:27:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	87872	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	361660	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	176921	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	323005	20.00	ng	0.00
76) Chrysene-d12	14.05	240	256340	20.00	ng	0.00
87) Perylene-d12	15.54	264	223531	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	335744	65.22	ng	0.00
7) Phenol-d6	6.49	99	497235	77.86	ng	0.00
23) Nitrobenzene-d5	7.42	82	492434	77.88	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	177125	81.07	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	1024465	88.39	ng	0.00
79) Terphenyl-d14	12.98	244	1212768	78.49	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.59	88	67210	32.843	ng	96
3) Pyridine	3.35	79	169629m	31.112	ng	
4) n-Nitrosodimethylamine	3.32	42	88946	39.458	ng	# 70
6) Aniline	6.52	93	310033	38.233	ng	94
8) 2-Chlorophenol	6.64	128	231586	39.722	ng	98
9) Benzaldehyde	6.40	77	129066	34.473	ng	97
10) Phenol	6.50	94	280370	39.062	ng	82
11) bis(2-Chloroethyl)ether	6.59	93	198158	38.604	ng	100
12) 1,3-Dichlorobenzene	6.79	146	267770	39.879	ng	99
13) 1,4-Dichlorobenzene	6.87	146	275972	40.464	ng	98
14) 1,2-Dichlorobenzene	7.02	146	258802	40.193	ng	99
15) Benzyl Alcohol	7.00	79	204564	42.072	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.12	45	217857	39.773	ng	68
17) 2-Methylphenol	7.11	107	185004	39.563	ng	94
18) Hexachloroethane	7.36	117	102938	41.124	ng	94
19) n-Nitroso-di-n-propylamine	7.27	70	160831	40.882	ng	96
20) 3+4-Methylphenols	7.27	107	245745	41.061	ng	99
22) Acetophenone	7.26	105	346384	39.078	ng	# 92
24) Nitrobenzene	7.44	77	241709	38.159	ng	98
25) Isophorone	7.68	82	418975	38.019	ng	99
26) 2-Nitrophenol	7.76	139	131373	39.157	ng	95
27) 2,4-Dimethylphenol	7.79	122	173572	37.918	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	278076	38.583	ng	98
29) 2,4-Dichlorophenol	8.00	162	204228	38.762	ng	97
30) 1,2,4-Trichlorobenzene	8.08	180	240754	39.226	ng	97
31) Naphthalene	8.16	128	754305	41.035	ng	99
32) Benzoic acid	7.95	122	121620	33.419	ng	98
33) 4-Chloroaniline	8.21	127	305094	38.716	ng	98
34) Hexachlorobutadiene	8.27	225	141810	38.815	ng	98
35) Caprolactam	8.59	113	59542m	36.198	ng	
36) 4-Chloro-3-methylphenol	8.70	107	215658	38.124	ng	99
37) 2-Methylnaphthalene	8.85	142	476080	37.828	ng	99
38) 1-Methylnaphthalene	8.95	142	448766	37.848	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	219474	41.288	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.00	237	65982	35.186	nq	99
43) 2,4,6-Trichlorophenol	9.14	196	145362	40.932	nq	97
44) 2,4,5-Trichlorophenol	9.19	196	146017	39.969	nq	97
46) 1,1'-Biphenyl	9.32	154	623869	44.583	nq	97
47) 2-Chloronaphthalene	9.35	162	500940	44.479	nq	99
48) 2-Nitroaniline	9.45	65	142158	44.688	nq	98
49) Acenaphthylene	9.76	152	706378	42.290	nq	99
50) Dimethylphthalate	9.62	163	560253	42.187	nq	99
51) 2,6-Dinitrotoluene	9.69	165	120481	42.080	nq	92
52) Acenaphthene	9.93	154	410067	40.305	nq	99
53) 3-Nitroaniline	9.86	138	132513	39.433	nq	87
54) 2,4-Dinitrophenol	9.98	184	60599	44.441	nq	98
55) Dibenzofuran	10.11	168	618932	41.171	nq	98
56) 4-Nitrophenol	10.04	139	101820	49.080	nq	98
57) 2,4-Dinitrotoluene	10.10	165	155166	40.558	nq	# 77
58) Fluorene	10.45	166	493791	40.689	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.23	232	124798	40.205	nq	98
60) Diethylphthalate	10.32	149	536836	40.860	nq	99
61) 4-Chlorophenyl-phenylether	10.44	204	243193	40.143	nq	100
62) 4-Nitroaniline	10.48	138	135989	38.954	nq	89
63) Azobenzene	10.60	77	476917	42.379	nq	97
65) 4,6-Dinitro-2-methylphenol	10.52	198	69630	39.960	nq	88
66) n-Nitrosodiphenylamine	10.56	169	431367	41.669	nq	99
67) 4-Bromophenyl-phenylether	10.93	248	148039	39.163	nq	98
68) Hexachlorobenzene	11.00	284	171860	38.625	nq	94
69) Atrazine	11.09	200	123141	38.541	nq	98
70) Pentachlorophenol	11.20	266	79252	41.127	nq	99
71) Phenanthrene	11.42	178	742470	42.165	nq	99
72) Anthracene	11.47	178	759034	43.065	nq	99
73) Carbazole	11.63	167	641841	41.040	nq	99
74) Di-n-butylphthalate	11.95	149	888179	44.789	nq	99
75) Fluoranthene	12.62	202	772893	43.492	nq	99
77) Benzidine	12.73	184	249929	35.583	nq	100
78) Pyrene	12.85	202	789860	37.733	nq	99
80) Butylbenzylphthalate	13.45	149	358532	38.040	nq	95
81) Benzo(a)anthracene	14.04	228	711565	39.657	nq	96
82) 3,3'-Dichlorobenzidine	14.00	252	220857	36.303	nq	98
83) Chrysene	14.07	228	680091	39.584	nq	99
84) Bis(2-ethylhexyl)phthalate	14.01	149	480735	38.789	nq	99
85) Di-n-octyl phthalate	14.63	149	717133	39.337	nq	99
86) Indeno(1,2,3-cd)pyrene	17.06	276	490607	33.513	nq	100
88) Benzo(b)fluoranthene	15.10	252	587727	39.610	nq	100
89) Benzo(k)fluoranthene	15.13	252	543917	38.177	nq	98
90) Benzo(a)pyrene	15.47	252	508387	38.894	nq	99
91) Dibenzo(a,h)anthracene	17.06	278	410702	36.208	nq	99
92) Benzo(g,h,i)perylene	17.50	276	369589	33.731	nq	98

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

