

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010819\
 Data File : BF111932.D
 Acq On : 9 Jan 2019 3:35
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 1/9/2019 10:36:58 AM

Quant Time: Jan 09 04:07:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	112806	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	357191	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	139886	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	282614	20.00	ng	0.00
76) Chrysene-d12	14.06	240	184594	20.00	ng	0.00
87) Perylene-d12	15.54	264	180101	20.00	ng	0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.52	112	543046	82.36	ng	0.00
7) Phenol-d6	6.53	99	578265	68.04	ng	0.00
23) Nitrobenzene-d5	7.45	82	578412	86.55	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	142539	78.34	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	788059	82.15	ng	0.00
79) Terphenyl-d14	13.00	244	968918	99.64	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.64	88	139358	48.336	ng	100
3) Pyridine	3.41	79	360620	48.258	ng	99
4) n-Nitrosodimethylamine	3.35	42	169145	45.493	ng	99
6) Aniline	6.55	93	418637	40.740	ng	# 69
8) 2-Chlorophenol	6.68	128	277089	37.845	ng	96
9) Benzaldehyde	6.43	77	169144	33.634	ng	96
10) Phenol	6.55	94	363383	39.341	ng	97
11) bis(2-Chloroethyl)ether	6.62	93	273467	38.270	ng	97
12) 1,3-Dichlorobenzene	6.83	146	272663	32.124	ng	99
13) 1,4-Dichlorobenzene	6.90	146	335226	38.920	ng	98
14) 1,2-Dichlorobenzene	7.06	146	330432	40.133	ng	98
15) Benzyl Alcohol	7.03	79	225966	33.605	ng	100
16) 2,2'-oxybis(1-Chloropropan	7.16	45	487288	39.996	ng	91
17) 2-Methylphenol	7.15	107	227849	38.720	ng	98
18) Hexachloroethane	7.40	117	58227	19.362	ng	90
19) n-Nitroso-di-n-propylamine	7.30	70	208117	37.690	ng	99
20) 3+4-Methylphenols	7.30	107	288408	39.381	ng	94
22) Acetophenone	7.29	105	402209	44.478	ng	# 97
24) Nitrobenzene	7.47	77	288167	42.623	ng	98
25) Isophorone	7.70	82	455142	42.056	ng	99
26) 2-Nitrophenol	7.79	139	90687	27.243	ng	91
27) 2,4-Dimethylphenol	7.83	122	206756	42.957	ng	98
28) bis(2-Chloroethoxy)methane	7.91	93	311674	43.145	ng	99
29) 2,4-Dichlorophenol	8.03	162	222410	40.185	ng	97
30) 1,2,4-Trichlorobenzene	8.12	180	251380	41.171	ng	98
31) Naphthalene	8.19	128	633378	37.354	ng	99
32) Benzoic acid	7.92	122	72736m	23.287	ng	
33) 4-Chloroaniline	8.24	127	274088	37.394	ng	100
34) Hexachlorobutadiene	8.32	225	126548	33.293	ng	99
35) Caprolactam	8.60	113	55047m	30.442	ng	
36) 4-Chloro-3-methylphenol	8.73	107	166359	30.231	ng	100
37) 2-Methylnaphthalene	8.89	142	343079	32.653	ng	99
38) 1-Methylnaphthalene	8.99	142	330488	32.795	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.05	216	186618	41.846	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.04	237	273	0.106	ng	# 27
43) 2,4,6-Trichlorophenol	9.16	196	118182	39.071	ng	99
44) 2,4,5-Trichlorophenol	9.21	196	122490	38.651	ng	97
46) 1,1'-Biphenyl	9.35	154	470338	40.196	ng	99
47) 2-Chloronaphthalene	9.37	162	364295	39.668	ng	98
48) 2-Nitroaniline	9.46	65	126656	45.110	ng	98
49) Acenaphthylene	9.79	152	525530	38.707	ng	98
50) Dimethylphthalate	9.64	163	406913	38.723	ng	99
51) 2,6-Dinitrotoluene	9.70	165	59609	25.367	ng	96
52) Acenaphthene	9.96	154	312199	35.668	ng	99
53) 3-Nitroaniline	9.87	138	117656	46.754	ng	99
54) 2,4-Dinitrophenol	9.99	184	125	0.126	ng	# 1
55) Dibenzofuran	10.13	168	489986	38.492	ng	100
56) 4-Nitrophenol	10.05	139	60331	33.930	ng	94
57) 2,4-Dinitrotoluene	10.11	165	68086	22.423	ng	# 97
58) Fluorene	10.47	166	366264	39.543	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.26	232	105441	39.488	ng	99
60) Diethylphthalate	10.34	149	398428	39.179	ng	99
61) 4-Chlorophenyl-phenylether	10.46	204	201166	38.730	ng	97
62) 4-Nitroaniline	10.49	138	133055	55.744	ng	99
63) Azobenzene	10.63	77	362408	37.305	ng	95
65) 4,6-Dinitro-2-methylphenol	10.52	198	801	0.459	ng	# 64
66) n-Nitrosodiphenylamine	10.58	169	348209	36.716	ng	99
67) 4-Bromophenyl-phenylether	10.96	248	128755	35.174	ng	95
68) Hexachlorobenzene	11.03	284	145030	35.750	ng	100
69) Atrazine	11.11	200	102478	30.755	ng	98
70) Pentachlorophenol	11.23	266	72253	31.299	ng	99
71) Phenanthrene	11.44	178	590987	39.112	ng	99
72) Anthracene	11.49	178	599316	39.072	ng	99
73) Carbazole	11.65	167	590422	39.572	ng	98
74) Di-n-butylphthalate	11.97	149	663732	38.298	ng	100
75) Fluoranthene	12.63	202	588086	37.912	ng	99
77) Benzdine	12.74	184	356319	64.600	ng	97
78) Pyrene	12.86	202	589560	46.418	ng	98
80) Butylbenzylphthalate	13.47	149	271642	50.588	ng	99
81) Benzo(a)anthracene	14.05	228	414751	39.040	ng	98
82) 3,3'-Dichlorobenzidine	14.00	252	186427	47.533	ng	99
83) Chrysene	14.09	228	386214	38.128	ng	97
84) Bis(2-ethylhexyl)phthalate	14.03	149	366814	52.016	ng	# 99
85) Di-n-octyl phthalate	14.64	149	500446	49.589	ng	99
86) Indeno(1,2,3-cd)pyrene	17.05	276	352588	47.746	ng	99
88) Benzo(b)fluoranthene	15.11	252	412310	38.271	ng	100
89) Benzo(k)fluoranthene	15.14	252	412237	40.707	ng	99
90) Benzo(a)pyrene	15.48	252	373878	39.468	ng	99
91) Dibenzo(a,h)anthracene	17.06	278	308072	40.435	ng	98
92) Benzo(g,h,i)perylene	17.50	276	298562	40.065	ng	98

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

