

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031819\
 Data File : BF113109.D
 Acq On : 19 Mar 2019 00:17
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
Client Sampled :
 SSTDCCC040

Manual Integrations
APPROVED
 Sohil
 3/20/2019 6:46:24 PM

Quant Time: Mar 20 05:44:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	144479	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	580271	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	237392	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	424952	20.00	ng	0.00
76) Chrysene-d12	13.87	240	369417	20.00	ng	0.00
87) Perylene-d12	15.28	264	281895	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.35	112	625756	67.37	ng	0.00
7) Phenol-d6	6.38	99	755271	64.73	ng	0.00
23) Nitrobenzene-d5	7.30	82	735644	75.86	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	211560	83.95	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1291331	92.18	ng	0.00
79) Terphenyl-d14	12.82	244	1401495	73.54	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.37	88	176675	37.947	ng	97
3) Pyridine	3.09	79	422365	33.909	ng	98
4) n-Nitrosodimethylamine	3.02	42	182333	33.463	ng	99
6) Aniline	6.39	93	474117	29.522	ng	# 38
8) 2-Chlorophenol	6.52	128	353466	34.187	ng	94
9) Benzaldehyde	6.27	77	226552	26.949	ng	96
10) Phenol	6.39	94	429155	31.521	ng	# 67
11) bis(2-Chloroethyl)ether	6.47	93	339899	32.526	ng	95
12) 1,3-Dichlorobenzene	6.67	146	401444	37.586	ng	97
13) 1,4-Dichlorobenzene	6.75	146	433881	40.507	ng	97
14) 1,2-Dichlorobenzene	6.90	146	398782	40.613	ng	98
15) Benzyl Alcohol	6.87	79	308740	34.703	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.01	45	597206	34.244	ng	81
17) 2-Methylphenol	7.00	107	300352	35.809	ng	98
18) Hexachloroethane	7.25	117	122280	29.803	ng	91
19) n-Nitroso-di-n-propylamine	7.15	70	258066	34.924	ng	97
20) 3+4-Methylphenols	7.15	107	366076	38.658	ng	# 70
22) Acetophenone	7.14	105	538770	39.708	ng	# 89
24) Nitrobenzene	7.32	77	389221	36.062	ng	95
25) Isophorone	7.55	82	691676	33.108	ng	99
26) 2-Nitrophenol	7.63	139	103963	23.139	ng	93
27) 2,4-Dimethylphenol	7.67	122	315155	37.704	ng	98
28) bis(2-Chloroethoxy)methane	7.76	93	503188	38.524	ng	100
29) 2,4-Dichlorophenol	7.88	162	316694	39.070	ng	96
30) 1,2,4-Trichlorobenzene	7.96	180	364015	41.794	ng	99
31) Naphthalene	8.03	128	1101035	38.218	ng	99
32) Benzoic acid	7.76	122	106024m	18.097	ng	
33) 4-Chloroaniline	8.09	127	454562	37.253	ng	100
34) Hexachlorobutadiene	8.16	225	225776	45.965	ng	99
35) Caprolactam	8.45	113	88092	30.856	ng	96
36) 4-Chloro-3-methylphenol	8.57	107	308634	34.405	ng	99
37) 2-Methylnaphthalene	8.73	142	690023	37.089	ng	99
38) 1-Methylnaphthalene	8.83	142	654780	36.332	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.89	216	327694	47.337	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.88	237	5313	1.402	ng	97
43) 2,4,6-Trichlorophenol	9.01	196	182901	38.390	ng	100
44) 2,4,5-Trichlorophenol	9.06	196	194602	37.789	ng	94
46) 1,1'-Biphenyl	9.19	154	803052	41.474	ng	99
47) 2-Chloronaphthalene	9.22	162	582529	38.529	ng	96
48) 2-Nitroaniline	9.31	65	185506	40.366	ng	92
49) Acenaphthylene	9.63	152	962376	37.494	ng	98
50) Dimethylphthalate	9.49	163	711711	40.660	ng	99
51) 2,6-Dinitrotoluene	9.55	165	129538	35.538	ng #	74
52) Acenaphthene	9.80	154	544561	36.280	ng	99
53) 3-Nitroaniline	9.72	138	187903	42.306	ng	97
55) Dibenzofuran	9.97	168	827164	37.916	ng	97
56) 4-Nitrophenol	9.89	139	100505	27.843	ng	94
57) 2,4-Dinitrotoluene	9.95	165	157315	34.720	ng	88
58) Fluorene	10.31	166	602003	38.880	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.09	232	159490	37.174	ng	93
60) Diethylphthalate	10.18	149	720342	38.965	ng	99
61) 4-Chlorophenyl-phenylether	10.30	204	319574	44.360	ng	89
62) 4-Nitroaniline	10.32	138	188231	45.863	ng	97
63) Azobenzene	10.46	77	613241	32.385	ng	93
65) 4,6-Dinitro-2-methylphenol	10.55	198	299	4.768	ng #	1
66) n-Nitrosodiphenylamine	10.42	169	545211	40.005	ng	99
67) 4-Bromophenyl-phenylether	10.79	248	195850	43.755	ng #	90
68) Hexachlorobenzene	10.86	284	216589	42.926	ng #	90
69) Atrazine	10.94	200	150431	36.040	ng	98
70) Pentachlorophenol	11.06	266	86639	29.174	ng	97
71) Phenanthrene	11.27	178	846747	40.049	ng	99
72) Anthracene	11.32	178	872214	39.409	ng	98
73) Carbazole	11.47	167	838282	38.827	ng	98
74) Di-n-butylphthalate	11.80	149	1009411	38.992	ng	100
75) Fluoranthene	12.45	202	1014600	44.790	ng	96
77) Benzidine	12.57	184	568954	29.686	ng	99
78) Pyrene	12.68	202	1045934	33.585	ng	99
80) Butylbenzylphthalate	13.30	149	478528	34.811	ng	89
81) Benzo(a)anthracene	13.86	228	832770	32.526	ng	99
82) 3,3'-Dichlorobenzidine	13.83	252	386025	39.865	ng #	98
83) Chrysene	13.90	228	863328	34.424	ng	99
84) Bis(2-ethylhexyl)phthalate	13.86	149	626605	38.862	ng	99
85) Di-n-octyl phthalate	14.47	149	1175337	42.451	ng	97
86) Indeno(1,2,3-cd)pyrene	16.66	276	658928	30.819	ng	95
88) Benzo(b)fluoranthene	14.88	252	711596	39.026	ng #	96
89) Benzo(k)fluoranthene	14.91	252	573464m	34.721	ng	
90) Benzo(a)pyrene	15.22	252	597722	37.061	ng	98
91) Dibenzo(a,h)anthracene	16.66	278	550616	40.009	ng	97
92) Benzo(g,h,i)perylene	17.06	276	530307	38.374	ng	95

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

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