

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031219\
 Data File : BF113023.D
 Acq On : 13 Mar 2019 1:45
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/13/2019 11:40:54 AM

Quant Time: Mar 13 04:54:13 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	194638	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	711706	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	312588	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	540933	20.00	ng	0.00
76) Chrysene-d12	13.87	240	446912	20.00	ng	0.00
87) Perylene-d12	15.27	264	410053	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.34	112	913838	73.03	ng	0.00
7) Phenol-d6	6.37	99	1135051	72.21	ng	0.00
23) Nitrobenzene-d5	7.30	82	1011474	85.04	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	280183	84.43	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1620120	86.19	ng	0.00
79) Terphenyl-d14	12.82	244	1542648	66.91	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.38	88	222431	35.463	ng	97
3) Pyridine	3.09	79	603022	35.936	ng	95
4) n-Nitrosodimethylamine	3.02	42	238319	32.467	ng	92
6) Aniline	6.39	93	722911	33.413	ng	# 70
8) 2-Chlorophenol	6.52	128	522659	37.524	ng	93
9) Benzaldehyde	6.27	77	348544	30.776	ng	98
10) Phenol	6.39	94	630545	34.378	ng	81
11) bis(2-Chloroethyl)ether	6.47	93	502527	35.696	ng	97
12) 1,3-Dichlorobenzene	6.67	146	550369	38.250	ng	97
13) 1,4-Dichlorobenzene	6.75	146	562279	38.967	ng	98
14) 1,2-Dichlorobenzene	6.90	146	510344	38.581	ng	98
15) Benzyl Alcohol	6.87	79	435011	36.295	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.02	45	810658	34.505	ng	97
17) 2-Methylphenol	7.00	107	413877	36.627	ng	97
18) Hexachloroethane	7.25	117	195430	35.356	ng	98
19) n-Nitroso-di-n-propylamine	7.15	70	334813	33.634	ng	99
20) 3+4-Methylphenols	7.15	107	466311	36.553	ng	# 80
22) Acetophenone	7.15	105	689078	41.407	ng	# 95
24) Nitrobenzene	7.32	77	535553	40.456	ng	98
25) Isophorone	7.56	82	962321	37.556	ng	98
26) 2-Nitrophenol	7.63	139	268473	48.719	ng	97
27) 2,4-Dimethylphenol	7.67	122	396773	38.702	ng	98
28) bis(2-Chloroethoxy)methane	7.77	93	613589	38.301	ng	99
29) 2,4-Dichlorophenol	7.87	162	416387	41.882	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	444465	41.607	ng	99
31) Naphthalene	8.04	128	1361561	38.533	ng	100
32) Benzoic acid	7.79	122	280914m	39.095	ng	
33) 4-Chloroaniline	8.09	127	583489	38.988	ng	99
34) Hexachlorobutadiene	8.16	225	260191	43.188	ng	99
35) Caprolactam	8.46	113	131123	37.447	ng	98
36) 4-Chloro-3-methylphenol	8.57	107	397743	36.150	ng	94
37) 2-Methylnaphthalene	8.73	142	856150	37.520	ng	99
38) 1-Methylnaphthalene	8.83	142	817737	36.995	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	411755	45.171	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	67620	13.547	nq	96
43) 2,4,6-Trichlorophenol	9.01	196	259937	41.434	nq	99
44) 2,4,5-Trichlorophenol	9.05	196	278469	41.067	nq	95
46) 1,1'-Biphenyl	9.19	154	1072442	42.062	nq	99
47) 2-Chloronaphthalene	9.22	162	799294	40.148	nq	98
48) 2-Nitroaniline	9.30	65	256279	42.351	nq	98
49) Acenaphthylene	9.63	152	1280548	37.888	nq	99
50) Dimethylphthalate	9.49	163	951793	41.295	nq	99
51) 2,6-Dinitrotoluene	9.55	165	206360	42.995	nq #	78
52) Acenaphthene	9.80	154	732089	37.040	nq	99
53) 3-Nitroaniline	9.72	138	234000	40.011	nq	96
54) 2,4-Dinitrophenol	9.82	184	28286	19.699	nq	92
55) Dibenzofuran	9.97	168	1094383	38.097	nq	97
56) 4-Nitrophenol	9.88	139	163004	34.294	nq	93
57) 2,4-Dinitrotoluene	9.95	165	258138	43.267	nq #	85
58) Fluorene	10.31	166	775192	38.021	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.09	232	218383	38.656	nq #	95
60) Diethylphthalate	10.19	149	916114	37.634	nq	99
61) 4-Chlorophenyl-phenylether	10.30	204	397695	41.924	nq	93
62) 4-Nitroaniline	10.32	138	219141	40.549	nq	96
63) Azobenzene	10.46	77	833281	33.420	nq	95
65) 4,6-Dinitro-2-methylphenol	10.36	198	46943	23.704	nq	85
66) n-Nitrosodiphenylamine	10.42	169	723761	41.720	nq	99
67) 4-Bromophenyl-phenylether	10.79	248	256634	45.042	nq	92
68) Hexachlorobenzene	10.86	284	276546	43.058	nq #	89
69) Atrazine	10.95	200	221308	41.652	nq	99
70) Pentachlorophenol	11.05	266	159154	42.101	nq	99
71) Phenanthrene	11.27	178	1069822	39.750	nq	99
72) Anthracene	11.32	178	1113887	39.538	nq	99
73) Carbazole	11.47	167	1054646	38.375	nq	99
74) Di-n-butylphthalate	11.81	149	1332424	40.434	nq	99
75) Fluoranthene	12.45	202	1171324	40.622	nq	96
77) Benzidine	12.57	184	606617	26.162	nq	99
78) Pyrene	12.67	202	1205472	31.996	nq	99
80) Butylbenzylphthalate	13.30	149	546114	32.839	nq	90
81) Benzo(a)anthracene	13.86	228	1039255	33.552	nq	99
82) 3,3'-Dichlorobenzidine	13.82	252	464478	39.650	nq	99
83) Chrysene	13.90	228	1080711	35.620	nq	98
84) Bis(2-ethylhexyl)phthalate	13.86	149	688685	35.306	nq	99
85) Di-n-octyl phthalate	14.46	149	1353888	40.420	nq	96
86) Indeno(1,2,3-cd)pyrene	16.63	276	845182	32.676	nq	95
88) Benzo(b)fluoranthene	14.87	252	950306	35.829	nq	98
89) Benzo(k)fluoranthene	14.90	252	959708	39.946	nq	98
90) Benzo(a)pyrene	15.22	252	885405	37.741	nq	97
91) Dibenzo(a,h)anthracene	16.65	278	713941	35.663	nq	97
92) Benzo(g,h,i)perylene	17.04	276	673641	33.511	nq	96

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

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