

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030519\
 Data File : BF112857.D
 Acq On : 5 Mar 2019 17:23
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/6/2019 12:24:23 PM

Quant Time: Mar 06 03:31:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	208562	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	832610	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	408380	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	817014	20.00	ng	0.00
76) Chrysene-d12	13.88	240	543858	20.00	ng	0.00
87) Perylene-d12	15.29	264	468930	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.35	112	989596	73.81	ng	0.00
7) Phenol-d6	6.38	99	1290032	76.60	ng	0.01
23) Nitrobenzene-d5	7.30	82	1152097	82.80	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	401852	92.69	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	2062181	83.20	ng	0.00
79) Terphenyl-d14	12.83	244	2170467	77.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.45	88	233921	34.805	ng	98
3) Pyridine	3.16	79	661552	36.792	ng	99
4) n-Nitrosodimethylamine	3.09	42	281898	35.840	ng	97
6) Aniline	6.40	93	849456	36.641	ng	99
8) 2-Chlorophenol	6.53	128	582447	39.025	ng	97
9) Benzaldehyde	6.29	77	395283	32.572	ng	99
10) Phenol	6.39	94	733641	37.328	ng	84
11) bis(2-Chloroethyl)ether	6.48	93	562332	37.277	ng	95
12) 1,3-Dichlorobenzene	6.68	146	618369	40.107	ng	99
13) 1,4-Dichlorobenzene	6.76	146	619555	40.069	ng	99
14) 1,2-Dichlorobenzene	6.92	146	570923	40.279	ng	98
15) Benzyl Alcohol	6.89	79	508158	39.568	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.02	45	917986	36.465	ng	98
17) 2-Methylphenol	7.00	107	471885	38.973	ng	99
18) Hexachloroethane	7.26	117	232503	39.255	ng	93
19) n-Nitroso-di-n-propylamine	7.16	70	400017	37.501	ng	99
20) 3+4-Methylphenols	7.16	107	532694	38.969	ng	95
22) Acetophenone	7.15	105	796052	40.889	ng	100
24) Nitrobenzene	7.32	77	612194	39.530	ng	99
25) Isophorone	7.57	82	1148615	38.317	ng	98
26) 2-Nitrophenol	7.64	139	306840	47.596	ng	97
27) 2,4-Dimethylphenol	7.68	122	466029	38.857	ng	99
28) bis(2-Chloroethoxy)methane	7.77	93	723136	38.584	ng	99
29) 2,4-Dichlorophenol	7.88	162	494703	42.534	ng	99
30) 1,2,4-Trichlorobenzene	7.97	180	528223	42.267	ng	98
31) Naphthalene	8.05	128	1627363	39.368	ng	99
32) Benzoic acid	7.82	122	387724	46.124	ng	93
33) 4-Chloroaniline	8.09	127	697870	39.859	ng	97
34) Hexachlorobutadiene	8.17	225	317124	44.995	ng	98
35) Caprolactam	8.47	113	169302m	41.329	ng	
36) 4-Chloro-3-methylphenol	8.58	107	526290	40.887	ng	99
37) 2-Methylnaphthalene	8.73	142	1083253	40.579	ng	99
38) 1-Methylnaphthalene	8.83	142	1050017	40.605	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	504683	42.379	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	292848	44.907	nq	99
43) 2,4,6-Trichlorophenol	9.02	196	355848	43.417	nq	99
44) 2,4,5-Trichlorophenol	9.05	196	377638	42.628	nq	98
46) 1,1'-Biphenyl	9.20	154	1343566	40.336	nq	100
47) 2-Chloronaphthalene	9.22	162	1027951	39.522	nq	100
48) 2-Nitroaniline	9.32	65	332632	42.075	nq	99
49) Acenaphthylene	9.64	152	1709582	38.718	nq	100
50) Dimethylphthalate	9.50	163	1229408	40.828	nq	100
51) 2,6-Dinitrotoluene	9.56	165	269371	42.958	nq #	80
52) Acenaphthene	9.81	154	1023021	39.619	nq	99
53) 3-Nitroaniline	9.73	138	308531	40.380	nq	94
54) 2,4-Dinitrophenol	9.83	184	124004	50.900	nq	97
55) Dibenzofuran	9.98	168	1443089	38.453	nq	99
56) 4-Nitrophenol	9.89	139	258938	41.699	nq	86
57) 2,4-Dinitrotoluene	9.96	165	346229	44.420	nq	91
58) Fluorene	10.32	166	1065132	39.988	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.10	232	326441	44.229	nq	96
60) Diethylphthalate	10.20	149	1222664	38.445	nq	99
61) 4-Chlorophenyl-phenylether	10.32	204	520417	41.993	nq	95
62) 4-Nitroaniline	10.34	138	309074	43.776	nq	97
63) Azobenzene	10.47	77	1179104	36.197	nq	97
65) 4,6-Dinitro-2-methylphenol	10.37	198	161048	47.976	nq #	73
66) n-Nitrosodiphenylamine	10.43	169	1005864	38.388	nq	100
67) 4-Bromophenyl-phenylether	10.80	248	364119	42.312	nq	92
68) Hexachlorobenzene	10.87	284	414859	42.766	nq #	89
69) Atrazine	10.96	200	316098	39.389	nq	98
70) Pentachlorophenol	11.06	266	261835	45.859	nq	99
71) Phenanthrene	11.28	178	1607459	39.544	nq	99
72) Anthracene	11.33	178	1636655	38.463	nq	100
73) Carbazole	11.48	167	1566968	37.750	nq	100
74) Di-n-butylphthalate	11.82	149	1904843	38.272	nq	100
75) Fluoranthene	12.46	202	1718682	39.463	nq	98
77) Benzidine	12.58	184	968266	34.316	nq	99
78) Pyrene	12.69	202	1750255	38.175	nq	100
80) Butylbenzylphthalate	13.31	149	782456	38.663	nq	91
81) Benzo(a)anthracene	13.87	228	1313083	34.836	nq	100
82) 3,3'-Dichlorobenzidine	13.83	252	561914	39.417	nq	98
83) Chrysene	13.91	228	1353098	36.648	nq	100
84) Bis(2-ethylhexyl)phthalate	13.87	149	871607	36.718	nq	99
85) Di-n-octyl phthalate	14.47	149	1471897	36.110	nq	98
86) Indeno(1,2,3-cd)pyrene	16.65	276	1117565	35.505	nq	94
88) Benzo(b)fluoranthene	14.89	252	1160418	38.258	nq	98
89) Benzo(k)fluoranthene	14.92	252	1096129	39.896	nq	98
90) Benzo(a)pyrene	15.23	252	1041970	38.838	nq	98
91) Dibenzo(a,h)anthracene	16.67	278	916150	40.018	nq	97
92) Benzo(g,h,i)perylene	17.06	276	926622	40.309	nq	95

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

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