

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF110619\
 Data File : BF117698.D
 Acq On : 6 Nov 2019 13:54
 Operator : JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Manual Integrations
 APPROVED

Sohil
 11/8/2019 3:30:31 PM

Quant Time: Nov 06 15:14:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 06 14:03:21 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	380439	20.00	ng	0.00
21) Naphthalene-d8	8.23	136	1429495	20.00	ng	0.00
39) Acenaphthene-d10	9.99	164	753160	20.00	ng	0.00
64) Phenanthrene-d10	11.47	188	1347008	20.00	ng	0.00
76) Chrysene-d12	14.13	240	898127	20.00	ng	0.00
87) Perylene-d12	15.64	264	837878	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	2080843	90.27	ng	0.00
7) Phenol-d6	6.56	99	2557813	89.87	ng	0.00
23) Nitrobenzene-d5	7.50	82	2408705	103.81	ng	0.00
42) 2,4,6-Tribromophenol	10.78	330	722962	98.50	ng	0.00
45) 2-Fluorobiphenyl	9.30	172	3918634	99.78	ng	0.00
79) Terphenyl-d14	13.07	244	4231819	97.84	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.76	88	526949	49.073	ng	99
3) Pyridine	3.52	79	1447658	49.945	ng	100
4) n-Nitrosodimethylamine	3.49	42	641357	48.857	ng	98
6) Aniline	6.60	93	1827039	47.306	ng	99
8) 2-Chlorophenol	6.72	128	1167617	49.377	ng	97
9) Benzaldehyde	6.49	77	896319	59.849	ng	98
10) Phenol	6.57	94	1406125	49.471	ng	98
11) bis(2-Chloroethyl)ether	6.67	93	1146395	49.137	ng	97
12) 1,3-Dichlorobenzene	6.88	146	1337487	48.993	ng	98
13) 1,4-Dichlorobenzene	6.96	146	1345671	49.083	ng	98
14) 1,2-Dichlorobenzene	7.11	146	1243052	49.492	ng	99
15) Benzyl Alcohol	7.07	79	1046319	50.101	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.21	45	1818543	48.987	ng	100
17) 2-Methylphenol	7.18	107	997229	48.857	ng	99
18) Hexachloroethane	7.46	117	505826	49.806	ng	97
19) n-Nitroso-di-n-propylamine	7.36	70	834722	48.200	ng	98
20) 3+4-Methylphenols	7.34	107	1171894	48.671	ng	90
22) Acetophenone	7.35	105	1591350	49.427	ng	# 98
24) Nitrobenzene	7.52	77	1254844	52.535	ng	97
25) Isophorone	7.76	82	2258993	50.782	ng	99
26) 2-Nitrophenol	7.84	139	602203	57.301	ng	93
27) 2,4-Dimethylphenol	7.87	122	977839	51.425	ng	100
28) bis(2-Chloroethoxy)methane	7.97	93	1462365	52.063	ng	100
29) 2,4-Dichlorophenol	8.07	162	1022247	53.390	ng	98
30) 1,2,4-Trichlorobenzene	8.16	180	1165074	53.089	ng	99
31) Naphthalene	8.25	128	3190618	51.336	ng	99
32) Benzoic acid	7.99	122	824125m	58.152	ng	
33) 4-Chloroaniline	8.29	127	1497668	51.058	ng	97
34) Hexachlorobutadiene	8.37	225	666368	53.050	ng	99
35) Caprolactam	8.67	113	328589m	50.763	ng	
36) 4-Chloro-3-methylphenol	8.77	107	1010696	51.866	ng	99
37) 2-Methylnaphthalene	8.94	142	2170491	51.445	ng	99
38) 1-Methylnaphthalene	9.04	142	2061253	51.572	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.10	216	1021778	51.201	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.10	237	568432	51.592	ng	98
43) 2,4,6-Trichlorophenol	9.22	196	747923	54.312	ng	100
44) 2,4,5-Trichlorophenol	9.25	196	768154	55.450	ng	96
46) 1,1'-Biphenyl	9.41	154	2560898	49.060	ng	99
47) 2-Chloronaphthalene	9.43	162	2154650	51.694	ng	98
48) 2-Nitroaniline	9.52	65	693920	55.378	ng	95
49) Acenaphthylene	9.85	152	3136578	51.112	ng	98
50) Dimethylphthalate	9.71	163	2478197	51.792	ng	99
51) 2,6-Dinitrotoluene	9.77	165	577687	55.515	ng	87
52) Acenaphthene	10.02	154	2012367	51.972	ng	99
53) 3-Nitroaniline	9.94	138	696644	55.002	ng	91
54) 2,4-Dinitrophenol	10.03	184	205012	60.853	ng	# 66
55) Dibenzofuran	10.19	168	2831203	51.949	ng	99
56) 4-Nitrophenol	10.08	139	517876	55.177	ng	98
57) 2,4-Dinitrotoluene	10.17	165	748407	57.272	ng	90
58) Fluorene	10.54	166	2067658	51.175	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.30	232	640264	57.431	ng	100
60) Diethylphthalate	10.41	149	2478011	52.786	ng	99
61) 4-Chlorophenyl-phenylether	10.53	204	1080976	51.729	ng	97
62) 4-Nitroaniline	10.56	138	686384	55.624	ng	99
63) Azobenzene	10.69	77	2329821	51.676	ng	96
65) 4,6-Dinitro-2-methylphenol	10.58	198	290403	59.869	ng	89
66) n-Nitrosodiphenylamine	10.65	169	1983859	50.770	ng	99
67) 4-Bromophenyl-phenylether	11.02	248	733014	51.441	ng	95
68) Hexachlorobenzene	11.09	284	847086	51.254	ng	96
69) Atrazine	11.17	200	651451	80.673	ng	99
70) Pentachlorophenol	11.27	266	497128	59.367	ng	99
71) Phenanthrene	11.50	178	3170061	50.177	ng	99
72) Anthracene	11.56	178	3120975	49.159	ng	98
73) Carbazole	11.70	167	2880784	49.632	ng	98
74) Di-n-butylphthalate	12.04	149	3527467	50.632	ng	98
75) Fluoranthene	12.70	202	3213873	49.895	ng	98
77) Benzidine	12.81	184	1603509	90.833	ng	98
78) Pyrene	12.93	202	3248553	51.454	ng	98
80) Butylbenzylphthalate	13.54	149	1532869	53.423	ng	93
81) Benzo(a)anthracene	14.12	228	2764356	52.225	ng	100
82) 3,3'-Dichlorobenzidine	14.08	252	1057865	55.432	ng	98
83) Chrysene	14.16	228	2739649	52.611	ng	97
84) Bis(2-ethylhexyl)phthalate	14.10	149	1952999	53.182	ng	99
85) Di-n-octyl phthalate	14.73	149	3069864	53.893	ng	100
86) Indeno(1,2,3-cd)pyrene	17.18	276	2498229	51.154	ng	100
88) Benzo(b)fluoranthene	15.19	252	2713962	54.615	ng	98
89) Benzo(k)fluoranthene	15.23	252	2331685	50.712	ng	98
90) Benzo(a)pyrene	15.57	252	2307238	52.847	ng	99
91) Dibenzo(a,h)anthracene	17.20	278	2059697	52.999	ng	98
92) Benzo(g,h,i)perylene	17.64	276	2027744	53.388	ng	100

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

