

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061019\
 Data File : BF114925.D
 Acq On : 10 Jun 2019 10:30
 Operator : HP/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

mohammad
 6/12/2019 9:13:10 AM

Quant Time: Jun 10 12:42:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 10 12:30:05 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	274439	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1126670	20.00	ng	0.00
39) Acenaphthene-d10	9.69	164	568824	20.00	ng	0.00
64) Phenanthrene-d10	11.16	188	1123927	20.00	ng	0.00
76) Chrysene-d12	13.79	240	841048	20.00	ng	0.00
87) Perylene-d12	15.16	264	886931	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.26	112	394011	25.09	ng	-0.01
7) Phenol-d6	6.29	99	479543	25.09	ng	-0.02
23) Nitrobenzene-d5	7.22	82	417884	22.83	ng	-0.01
42) 2,4,6-Tribromophenol	10.47	330	146223	23.85	ng	-0.01
45) 2-Fluorobiphenyl	9.02	172	919967	30.30	ng	0.00
79) Terphenyl-d14	12.74	244	973460	22.51	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	2.28	88	84781	11.306	ng	98
3) Pyridine	2.98	79	217031	10.531	ng	99
4) n-Nitrosodimethylamine	2.89	42	77488	10.274	ng	99
6) Aniline	6.32	93	323729	11.938	ng	# 81
8) 2-Chlorophenol	6.44	128	217209	11.970	ng	96
9) Benzaldehyde	6.20	77	158036	11.921	ng	99
10) Phenol	6.30	94	266862	12.211	ng	# 76
11) bis(2-Chloroethyl)ether	6.39	93	197955	11.606	ng	97
12) 1,3-Dichlorobenzene	6.60	146	255879	12.101	ng	99
13) 1,4-Dichlorobenzene	6.67	146	255605	12.013	ng	99
14) 1,2-Dichlorobenzene	6.83	146	245637	12.783	ng	96
15) Benzyl Alcohol	6.80	79	172610	11.944	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.95	45	278006	11.397	ng	98
17) 2-Methylphenol	6.92	107	174256	11.234	ng	96
18) Hexachloroethane	7.17	117	90655	11.320	ng	96
19) n-Nitroso-di-n-propylamine	7.07	70	142700	11.280	ng	95
20) 3+4-Methylphenols	7.07	107	221751	12.961	ng	97
22) Acetophenone	7.06	105	303211	12.224	ng	93
24) Nitrobenzene	7.23	77	215134	11.346	ng	99
25) Isophorone	7.48	82	394663	10.907	ng	99
26) 2-Nitrophenol	7.56	139	110604	10.770	ng	99
27) 2,4-Dimethylphenol	7.60	122	177113	11.347	ng	99
28) bis(2-Chloroethoxy)methane	7.69	93	263374	11.387	ng	98
29) 2,4-Dichlorophenol	7.80	162	185885	11.320	ng	97
30) 1,2,4-Trichlorobenzene	7.89	180	216007	11.464	ng	96
31) Naphthalene	7.96	128	659632	13.005	ng	97
32) Benzoic acid	7.69	122	90675	7.976	ng	99
33) 4-Chloroaniline	8.01	127	287568	11.884	ng	98
34) Hexachlorobutadiene	8.09	225	122411	11.427	ng	100
35) Caprolactam	8.35	113	59265	11.048	ng	98
36) 4-Chloro-3-methylphenol	8.49	107	186708	11.426	ng	97
37) 2-Methylnaphthalene	8.65	142	436094	12.395	ng	97
38) 1-Methylnaphthalene	8.75	142	415623m	12.457	ng	
40) 1,2,4,5-Tetrachlorobenzene	8.82	216	197273	12.034	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.81	237	84638	8.513	nq	98
43) 2,4,6-Trichlorophenol	8.93	196	136083	10.676	nq	99
44) 2,4,5-Trichlorophenol	8.97	196	141456	11.019	nq	98
46) 1,1'-Biphenyl	9.12	154	547811	12.994	nq	95
47) 2-Chloronaphthalene	9.13	162	434904	12.140	nq	96
48) 2-Nitroaniline	9.23	65	117881	11.027	nq	96
49) Acenaphthylene	9.55	152	680227	13.094	nq	97
50) Dimethylphthalate	9.41	163	498310	11.995	nq	99
51) 2,6-Dinitrotoluene	9.47	165	108577	11.113	nq	96
52) Acenaphthene	9.72	154	404178	11.877	nq	98
53) 3-Nitroaniline	9.63	138	130561	11.439	nq	99
54) 2,4-Dinitrophenol	9.75	184	36549	8.267	nq	93
55) Dibenzofuran	9.89	168	604155	12.767	nq	92
56) 4-Nitrophenol	9.80	139	82833	9.904	nq	93
57) 2,4-Dinitrotoluene	9.87	165	142620	11.612	nq	99
58) Fluorene	10.23	166	463395	13.653	nq	97
59) 2,3,4,6-Tetrachlorophenol	10.01	232	116593	11.086	nq	99
60) Diethylphthalate	10.11	149	503938	12.088	nq	97
61) 4-Chlorophenyl-phenylether	10.23	204	228794	12.624	nq	97
62) 4-Nitroaniline	10.24	138	128777	11.633	nq	96
63) Azobenzene	10.39	77	436043	12.288	nq	97
65) 4,6-Dinitro-2-methylphenol	10.28	198	54224	9.288	nq	93
66) n-Nitrosodiphenylamine	10.34	169	405752	11.939	nq	100
67) 4-Bromophenyl-phenylether	10.71	248	142052	11.128	nq	99
68) Hexachlorobenzene	10.78	284	161175	11.537	nq	97
69) Atrazine	10.87	200	136171	11.521	nq	100
70) Pentachlorophenol	10.97	266	70482	8.727	nq	97
71) Phenanthrene	11.19	178	716994	13.363	nq	96
72) Anthracene	11.23	178	720940	12.768	nq	96
73) Carbazole	11.39	167	647785	13.382	nq	97
74) Di-n-butylphthalate	11.73	149	815391	12.807	nq	97
75) Fluoranthene	12.36	202	741595	13.280	nq	97
77) Benzidine	12.49	184	428263	10.286	nq	96
78) Pyrene	12.59	202	762162	10.446	nq	96
80) Butylbenzylphthalate	13.22	149	338191	9.179	nq	94
81) Benzo(a)anthracene	13.77	228	670695	10.348	nq	97
82) 3,3'-Dichlorobenzidine	13.74	252	264623	10.069	nq	# 97
83) Chrysene	13.81	228	660881	10.482	nq	96
84) Bis(2-ethylhexyl)phthalate	13.79	149	459510	9.863	nq	99
85) Di-n-octyl phthalate	14.39	149	805182	10.171	nq	98
86) Indeno(1,2,3-cd)pyrene	16.46	276	576157	8.009	nq	99
88) Benzo(b)fluoranthene	14.77	252	677913	12.006	nq	98
89) Benzo(k)fluoranthene	14.80	252	566762	11.820	nq	98
90) Benzo(a)pyrene	15.10	252	573406	11.733	nq	98
91) Dibenzo(a,h)anthracene	16.47	278	470578	10.168	nq	98
92) Benzo(g,h,i)perylene	16.85	276	446870	9.602	nq	98

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

