

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062419\
 Data File : BF115153.D
 Acq On : 24 Jun 2019 18:03
 Operator : HP/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

mohammad
 6/28/2019 8:25:17 AM

Quant Time: Jun 25 08:07:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	255930	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1050577	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	526208	20.00	ng	0.00
64) Phenanthrene-d10	11.11	188	1029556	20.00	ng	0.00
76) Chrysene-d12	13.74	240	735290	20.00	ng	-0.01
87) Perylene-d12	15.10	264	904547	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	1323110	79.78	ng	0.00
7) Phenol-d6	6.25	99	1624199	80.69	ng	0.00
23) Nitrobenzene-d5	7.18	82	1455286	85.21	ng	0.00
42) 2,4,6-Tribromophenol	10.42	330	478420	86.22	ng	0.00
45) 2-Fluorobiphenyl	8.98	172	2552460	82.39	ng	0.00
79) Terphenyl-d14	12.70	244	2664063	77.03	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.19	88	316334	40.415	ng	99
3) Pyridine	2.84	79	861097	39.033	ng	99
4) n-Nitrosodimethylamine	2.79	42	332806	38.482	ng	98
6) Aniline	6.28	93	1141512	41.261	ng	100
8) 2-Chlorophenol	6.40	128	745390	39.970	ng	98
9) Benzaldehyde	6.15	77	443430	33.765	ng	99
10) Phenol	6.27	94	913455	38.740	ng	95
11) bis(2-Chloroethyl)ether	6.35	93	698252	40.173	ng	98
12) 1,3-Dichlorobenzene	6.55	146	854459	41.285	ng	98
13) 1,4-Dichlorobenzene	6.63	146	862412	41.554	ng	99
14) 1,2-Dichlorobenzene	6.79	146	795476	41.389	ng	98
15) Benzyl Alcohol	6.76	79	591431	40.349	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1021494	40.520	ng	100
17) 2-Methylphenol	6.88	107	611380	39.718	ng	98
18) Hexachloroethane	7.13	117	306868	41.013	ng	98
19) n-Nitroso-di-n-propylamine	7.04	70	496087	38.494	ng	98
20) 3+4-Methylphenols	7.03	107	705566	39.696	ng	90
22) Acetophenone	7.03	105	977547	40.644	ng	# 99
24) Nitrobenzene	7.20	77	770082	40.930	ng	98
25) Isophorone	7.44	82	1363221	38.965	ng	100
26) 2-Nitrophenol	7.51	139	404639	43.549	ng	99
27) 2,4-Dimethylphenol	7.56	122	603756	39.687	ng	98
28) bis(2-Chloroethoxy)methane	7.65	93	896091	40.524	ng	99
29) 2,4-Dichlorophenol	7.75	162	635111	40.454	ng	98
30) 1,2,4-Trichlorobenzene	7.84	180	729100	42.043	ng	99
31) Naphthalene	7.92	128	2078197	40.298	ng	99
32) Benzoic acid	7.68	122	457694	42.148	ng	99
33) 4-Chloroaniline	7.97	127	976963	41.452	ng	99
34) Hexachlorobutadiene	8.05	225	406554	42.780	ng	99
35) Caprolactam	8.35	113	218412	41.379	ng	89
36) 4-Chloro-3-methylphenol	8.45	107	639027	40.192	ng	95
37) 2-Methylnaphthalene	8.61	142	1418077	40.783	ng	100
38) 1-Methylnaphthalene	8.71	142	1337049	40.262	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	623889	41.957	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.77	237	361569	41.007	nq	99
43) 2,4,6-Trichlorophenol	8.88	196	478114	41.648	nq	99
44) 2,4,5-Trichlorophenol	8.92	196	493169	41.716	nq	97
46) 1,1'-Biphenyl	9.07	154	1681888	39.946	nq	99
47) 2-Chloronaphthalene	9.10	162	1404052	41.111	nq	100
48) 2-Nitroaniline	9.18	65	431601	43.346	nq	99
49) Acenaphthylene	9.51	152	2096169	39.393	nq	100
50) Dimethylphthalate	9.38	163	1585277	40.948	nq	100
51) 2,6-Dinitrotoluene	9.43	165	370972	41.505	nq	97
52) Acenaphthene	9.68	154	1353820	40.659	nq	99
53) 3-Nitroaniline	9.60	138	448798	41.147	nq	99
54) 2,4-Dinitrophenol	9.70	184	190810	45.023	nq	90
55) Dibenzofuran	9.85	168	1842301	38.550	nq	100
56) 4-Nitrophenol	9.75	139	356314	40.748	nq	90
57) 2,4-Dinitrotoluene	9.83	165	496829	43.050	nq	99
58) Fluorene	10.19	166	1305886	40.363	nq	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	407799	41.137	nq	99
60) Diethylphthalate	10.08	149	1594053	40.962	nq	99
61) 4-Chlorophenyl-phenylether	10.18	204	647039	42.398	nq	97
62) 4-Nitroaniline	10.20	138	466952	42.675	nq	98
63) Azobenzene	10.34	77	1442144	39.675	nq	98
65) 4,6-Dinitro-2-methylphenol	10.24	198	236735	43.709	nq	94
66) n-Nitrosodiphenylamine	10.30	169	1294189	40.348	nq	100
67) 4-Bromophenyl-phenylether	10.67	248	467556	41.886	nq	93
68) Hexachlorobenzene	10.74	284	515353	42.534	nq	97
69) Atrazine	10.83	200	401715	38.933	nq	99
70) Pentachlorophenol	10.92	266	329976	42.749	nq	98
71) Phenanthrene	11.14	178	2133126	38.481	nq	100
72) Anthracene	11.19	178	2187662	39.096	nq	98
73) Carbazole	11.34	167	2001356	40.054	nq	99
74) Di-n-butylphthalate	11.70	149	2387699	41.353	nq	99
75) Fluoranthene	12.32	202	2233334	40.380	nq	98
77) Benzidine	12.44	184	1168736	35.796	nq	99
78) Pyrene	12.54	202	2286768	40.468	nq	98
80) Butylbenzylphthalate	13.18	149	1087900	43.625	nq	97
81) Benzo(a)anthracene	13.72	228	2170949	41.148	nq	99
82) 3,3'-Dichlorobenzidine	13.69	252	861895	45.238	nq	99
83) Chrysene	13.77	228	2015386	41.702	nq	99
84) Bis(2-ethylhexyl)phthalate	13.75	149	1436367	43.958	nq	# 98
85) Di-n-octyl phthalate	14.35	149	2448344	44.596	nq	99
86) Indeno(1,2,3-cd)pyrene	16.38	276	2502401	43.758	nq	99
88) Benzo(b)fluoranthene	14.73	252	2308018	40.893	nq	99
89) Benzo(k)fluoranthene	14.76	252	1931038m	38.151	nq	
90) Benzo(a)pyrene	15.05	252	2059141	40.478	nq	99
91) Dibenzo(a,h)anthracene	16.40	278	2007853	42.278	nq	98
92) Benzo(g,h,i)perylene	16.76	276	2005735	41.728	nq	99

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

