

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
 Data File : BF112979.D
 Acq On : 12 Mar 2019 2:59
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 3/12/2019 4:29:08 PM

Quant Time: Mar 12 07:18:49 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	6.73	152	121517	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	492390	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	263743	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	567134	20.00	ng	0.00
76) Chrysene-d12	13.87	240	384201	20.00	ng	0.00
87) Perylene-d12	15.27	264	353319	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	600929	76.93	ng	-0.01
7) Phenol-d6	6.37	99	753936	76.83	ng	0.00
23) Nitrobenzene-d5	7.29	82	670007	81.42	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	280408	100.15	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	1286351	79.42	ng	0.00
79) Terphenyl-d14	12.82	244	1566665	79.05	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.35	88	140704	35.932	ng	98
3) Pyridine	3.06	79	352073	33.607	ng	99
4) n-Nitrosodimethylamine	2.99	42	152420m	33.259	ng	
6) Aniline	6.39	93	490761	36.333	ng	100
8) 2-Chlorophenol	6.52	128	342902	39.432	ng	97
9) Benzaldehyde	6.27	77	222012	31.399	ng	100
10) Phenol	6.38	94	431467	37.679	ng	92
11) bis(2-Chloroethyl)ether	6.47	93	315498	35.896	ng	94
12) 1,3-Dichlorobenzene	6.67	146	361019	40.188	ng	96
13) 1,4-Dichlorobenzene	6.75	146	366683	40.703	ng	97
14) 1,2-Dichlorobenzene	6.90	146	346310	41.934	ng	98
15) Benzyl Alcohol	6.87	79	281553	37.627	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.02	45	519904	35.445	ng	98
17) 2-Methylphenol	6.99	107	265170	37.588	ng	97
18) Hexachloroethane	7.25	117	118483	34.334	ng	96
19) n-Nitroso-di-n-propylamine	7.15	70	222209	35.754	ng	99
20) 3+4-Methylphenols	7.15	107	324933	40.797	ng	98
22) Acetophenone	7.14	105	468656	40.705	ng	# 93
24) Nitrobenzene	7.31	77	353033	38.547	ng	99
25) Isophorone	7.56	82	635852	35.868	ng	97
26) 2-Nitrophenol	7.63	139	172371	45.212	ng	94
27) 2,4-Dimethylphenol	7.67	122	266855	37.623	ng	99
28) bis(2-Chloroethoxy)methane	7.77	93	408760	36.880	ng	98
29) 2,4-Dichlorophenol	7.87	162	293824	42.718	ng	98
30) 1,2,4-Trichlorobenzene	7.96	180	312556	42.291	ng	98
31) Naphthalene	8.04	128	985880	40.329	ng	99
32) Benzoic acid	7.79	122	207897	41.820	ng	89
33) 4-Chloroaniline	8.09	127	422519	40.807	ng	99
34) Hexachlorobutadiene	8.16	225	186611	44.772	ng	99
35) Caprolactam	8.45	113	103937	42.904	ng	94
36) 4-Chloro-3-methylphenol	8.57	107	314109	41.265	ng	98
37) 2-Methylnaphthalene	8.73	142	657665	41.659	ng	99
38) 1-Methylnaphthalene	8.83	142	639934	41.846	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	319677	41.565	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.89	237	33802	8.026	nq	100
43) 2,4,6-Trichlorophenol	9.01	196	211060	39.874	nq	98
44) 2,4,5-Trichlorophenol	9.05	196	238202	41.634	nq	95
46) 1,1'-Biphenyl	9.19	154	841592	39.121	nq	99
47) 2-Chloronaphthalene	9.22	162	644309	38.357	nq	98
48) 2-Nitroaniline	9.31	65	217631	42.625	nq	89
49) Acenaphthylene	9.63	152	1103587	38.700	nq	99
50) Dimethylphthalate	9.49	163	754900	38.818	nq	99
51) 2,6-Dinitrotoluene	9.55	165	172697	42.645	nq #	77
52) Acenaphthene	9.80	154	627373	37.621	nq	99
53) 3-Nitroaniline	9.72	138	217251	44.026	nq	97
54) 2,4-Dinitrophenol	9.83	184	12950	13.640	nq	97
55) Dibenzofuran	9.97	168	974254	40.197	nq	97
56) 4-Nitrophenol	9.88	139	168323	41.971	nq	93
57) 2,4-Dinitrotoluene	9.95	165	241719	48.019	nq	87
58) Fluorene	10.31	166	732502	42.581	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.09	232	208804	43.805	nq	95
60) Diethylphthalate	10.19	149	785975	38.267	nq	100
61) 4-Chlorophenyl-phenylether	10.30	204	358133	44.746	nq	94
62) 4-Nitroaniline	10.32	138	229026	50.227	nq	98
63) Azobenzene	10.46	77	781324	37.139	nq	94
65) 4,6-Dinitro-2-methylphenol	10.36	198	22219	13.231	nq	92
66) n-Nitrosodiphenylamine	10.42	169	678972	37.330	nq	100
67) 4-Bromophenyl-phenylether	10.79	248	247384	41.413	nq	92
68) Hexachlorobenzene	10.86	284	284355	42.228	nq #	91
69) Atrazine	10.95	200	215337	38.656	nq	98
70) Pentachlorophenol	11.06	266	152665	38.519	nq	99
71) Phenanthrene	11.27	178	1130327	40.058	nq	99
72) Anthracene	11.32	178	1181337	39.995	nq	99
73) Carbazole	11.47	167	1122049	38.941	nq	99
74) Di-n-butylphthalate	11.81	149	1393942	40.347	nq	100
75) Fluoranthene	12.45	202	1161341	38.415	nq	96
77) Benzidine	12.57	184	620345	31.121	nq	99
78) Pyrene	12.67	202	1158535	35.769	nq	99
80) Butylbenzylphthalate	13.30	149	540475	37.804	nq	91
81) Benzo(a)anthracene	13.86	228	901237	33.846	nq	99
82) 3,3'-Dichlorobenzidine	13.82	252	398157	39.536	nq	99
83) Chrysene	13.90	228	954922	36.611	nq	98
84) Bis(2-ethylhexyl)phthalate	13.86	149	686309	40.927	nq	99
85) Di-n-octyl phthalate	14.47	149	1204788	41.840	nq	96
86) Indeno(1,2,3-cd)pyrene	16.64	276	731509	32.897	nq	95
88) Benzo(b)fluoranthene	14.87	252	944531	41.329	nq	98
89) Benzo(k)fluoranthene	14.90	252	799569	38.625	nq	98
90) Benzo(a)pyrene	15.22	252	782388	38.704	nq	98
91) Dibenzo(a,h)anthracene	16.65	278	624854	36.225	nq	97
92) Benzo(g,h,i)perylene	17.04	276	559054	32.277	nq	96

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

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