

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042519\
 Data File : BF113975.D
 Acq On : 25 Apr 2019 14:13
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/29/2019 5:49:42 AM

Quant Time: Apr 25 16:48:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	122030	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	483472	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	259948	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	520184	20.00	ng	0.00
76) Chrysene-d12	14.07	240	435130	20.00	ng	0.00
87) Perylene-d12	15.58	264	443323	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.52	112	573855	80.43	ng	0.00
7) Phenol-d6	6.53	99	723847	80.86	ng	0.00
23) Nitrobenzene-d5	7.46	82	668253	85.34	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	260717	82.33	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1341701	80.73	ng	0.00
79) Terphenyl-d14	13.00	244	1718721	80.98	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.67	88	127645	37.953	ng	97
3) Pyridine	3.44	79	359015	39.108	ng	97
4) n-Nitrosodimethylamine	3.39	42	117999	37.549	ng	94
6) Aniline	6.56	93	489422	39.817	ng	98
8) 2-Chlorophenol	6.69	128	332255	40.789	ng	98
9) Benzaldehyde	6.45	77	200687	39.350	ng	95
10) Phenol	6.54	94	428729	40.054	ng	98
11) bis(2-Chloroethyl)ether	6.63	93	320053	39.735	ng	95
12) 1,3-Dichlorobenzene	6.84	146	375513	40.703	ng	99
13) 1,4-Dichlorobenzene	6.92	146	379192	40.864	ng	99
14) 1,2-Dichlorobenzene	7.07	146	354037	41.515	ng	98
15) Benzyl Alcohol	7.03	79	255824	42.434	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.17	45	359697	38.255	ng	92
17) 2-Methylphenol	7.15	107	284567	39.854	ng	99
18) Hexachloroethane	7.42	117	132910	40.860	ng	100
19) n-Nitroso-di-n-propylamine	7.31	70	227297	39.210	ng	99
20) 3+4-Methylphenols	7.30	107	339794	42.120	ng	90
22) Acetophenone	7.30	105	436822	40.601	ng	# 94
24) Nitrobenzene	7.48	77	356937	41.254	ng	94
25) Isophorone	7.72	82	624901	39.003	ng	99
26) 2-Nitrophenol	7.79	139	180750	45.793	ng	95
27) 2,4-Dimethylphenol	7.83	122	243139	37.808	ng	97
28) bis(2-Chloroethoxy)methane	7.92	93	406507	39.816	ng	98
29) 2,4-Dichlorophenol	8.04	162	303172	41.225	ng	97
30) 1,2,4-Trichlorobenzene	8.12	180	332521	40.552	ng	99
31) Naphthalene	8.20	128	948970	41.318	ng	99
32) Benzoic acid	7.95	122	186663m	42.981	ng	
33) 4-Chloroaniline	8.25	127	394325	40.575	ng	99
34) Hexachlorobutadiene	8.32	225	207095	40.514	ng	99
35) Caprolactam	8.62	113	93631	40.024	ng	# 85
36) 4-Chloro-3-methylphenol	8.73	107	296131	40.645	ng	97
37) 2-Methylnaphthalene	8.89	142	632475	40.837	ng	98
38) 1-Methylnaphthalene	8.99	142	615948	41.205	ng	97
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	338067	40.037	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	168425	35.769	nq	99
43) 2,4,6-Trichlorophenol	9.17	196	215156	40.127	nq	97
44) 2,4,5-Trichlorophenol	9.21	196	238794	39.705	nq	98
46) 1,1'-Biphenyl	9.36	154	803536	40.464	nq	99
47) 2-Chloronaphthalene	9.39	162	646811	40.053	nq	99
48) 2-Nitroaniline	9.47	65	182486	43.106	nq	86
49) Acenaphthylene	9.80	152	1035827	40.972	nq	99
50) Dimethylphthalate	9.66	163	759099	40.378	nq	99
51) 2,6-Dinitrotoluene	9.72	165	176865	43.851	nq	99
52) Acenaphthene	9.97	154	613998	41.095	nq	97
53) 3-Nitroaniline	9.88	138	183326	43.255	nq	97
54) 2,4-Dinitrophenol	9.99	184	73660	46.141	nq	# 1
55) Dibenzofuran	10.15	168	919603	41.090	nq	99
56) 4-Nitrophenol	10.04	139	143307	42.705	nq	99
57) 2,4-Dinitrotoluene	10.12	165	238048	46.760	nq	96
58) Fluorene	10.49	166	699788	41.531	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.26	232	209622	39.662	nq	100
60) Diethylphthalate	10.35	149	724118	40.561	nq	100
61) 4-Chlorophenyl-phenylether	10.47	204	367166	41.146	nq	100
62) 4-Nitroaniline	10.49	138	185282	44.798	nq	99
63) Azobenzene	10.63	77	697691	40.323	nq	98
65) 4,6-Dinitro-2-methylphenol	10.53	198	103746	44.196	nq	94
66) n-Nitrosodiphenylamine	10.59	169	620412	39.749	nq	99
67) 4-Bromophenyl-phenylether	10.96	248	246152	39.171	nq	97
68) Hexachlorobenzene	11.04	284	270408	39.158	nq	97
69) Atrazine	11.12	200	204791	40.334	nq	100
70) Pentachlorophenol	11.23	266	154876	39.612	nq	97
71) Phenanthrene	11.44	178	1087288	41.591	nq	100
72) Anthracene	11.50	178	1090300	41.300	nq	99
73) Carbazole	11.64	167	1005876	42.709	nq	100
74) Di-n-butylphthalate	11.98	149	1174515	42.394	nq	99
75) Fluoranthene	12.63	202	1190662	41.746	nq	99
77) Benzidine	12.74	184	525806	40.556	nq	99
78) Pyrene	12.86	202	1213794	39.889	nq	99
80) Butylbenzylphthalate	13.47	149	512981	42.847	nq	99
81) Benzo(a)anthracene	14.06	228	1142240	44.554	nq	99
82) 3,3'-Dichlorobenzidine	14.02	252	444515	47.305	nq	99
83) Chrysene	14.10	228	1107073	44.339	nq	98
84) Bis(2-ethylhexyl)phthalate	14.04	149	704539	46.252	nq	98
85) Di-n-octyl phthalate	14.69	149	1187760	49.817	nq	98
86) Indeno(1,2,3-cd)pyrene	17.08	276	971627	41.083	nq	98
88) Benzo(b)fluoranthene	15.14	252	1099821	39.211	nq	99
89) Benzo(k)fluoranthene	15.17	252	1011707	41.933	nq	99
90) Benzo(a)pyrene	15.52	252	979497	40.374	nq	100
91) Dibenzo(a,h)anthracene	17.10	278	809357	35.538	nq	100
92) Benzo(g,h,i)perylene	17.53	276	749822	32.625	nq	98

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

