

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032720\
 Data File : BF119605.D
 Acq On : 28 Mar 2020 9:08
 Operator : CG/JU
 Sample : L2053-04MSD
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-13-SAMSD

Manual Integrations
 APPROVED

mohammad
 3/31/2020 8:14:03 AM

Quant Time: Mar 29 23:34:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 27 07:52:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	280180	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1029993	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	526606	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	1010319	20.00	ng	0.00
76) Chrysene-d12	14.01	240	638461	20.00	ng	0.00
87) Perylene-d12	15.47	264	533386	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	2392214	135.92	ng	0.02
7) Phenol-d6	6.46	99	2712623	120.65	ng	0.00
23) Nitrobenzene-d5	7.40	82	2053276	108.34	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	869585	160.06	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	3461015	97.37	ng	0.00
79) Terphenyl-d14	12.96	244	3832439	117.60	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.80	88	374402	43.071	ng	# 83
3) Pyridine	3.48	79	730470	30.503	ng	93
4) n-Nitrosodimethylamine	3.42	42	539220	46.173	ng	94
6) Aniline	6.49	93	562237	18.119	ng	99
8) 2-Chlorophenol	6.62	128	1051937	56.907	ng	99
9) Benzaldehyde	6.39	77	432615	30.332	ng	97
10) Phenol	6.48	94	1128748	47.051	ng	92
11) bis(2-Chloroethyl)ether	6.57	93	1033454	53.862	ng	97
12) 1,3-Dichlorobenzene	6.77	146	982422	49.105	ng	96
13) 1,4-Dichlorobenzene	6.85	146	1005885	50.265	ng	98
14) 1,2-Dichlorobenzene	7.01	146	948952	50.733	ng	97
15) Benzyl Alcohol	6.97	79	892233	52.757	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.11	45	1877346	48.508	ng	99
17) 2-Methylphenol	7.09	107	837627	49.456	ng	98
18) Hexachloroethane	7.35	117	376816	51.382	ng	96
19) n-Nitroso-di-n-propylamine	7.26	70	737182	54.398	ng	97
20) 3+4-Methylphenols	7.24	107	983997	47.406	ng	# 87
22) Acetophenone	7.25	105	1375530	55.890	ng	# 90
24) Nitrobenzene	7.42	77	1124924	55.542	ng	95
25) Isophorone	7.66	82	2088085	54.314	ng	99
26) 2-Nitrophenol	7.73	139	574870	72.087	ng	92
27) 2,4-Dimethylphenol	7.77	122	904642	54.861	ng	99
28) bis(2-Chloroethoxy)methane	7.87	93	1300939	57.226	ng	99
29) 2,4-Dichlorophenol	7.97	162	829183	54.727	ng	99
30) 1,2,4-Trichlorobenzene	8.06	180	864911	51.618	ng	100
31) Naphthalene	8.14	128	2757053	52.015	ng	100
32) Benzoic acid	7.89	122	652418	61.508	ng	95
33) 4-Chloroaniline	8.19	127	331611	13.653	ng	95
34) Hexachlorobutadiene	8.26	225	480763	51.281	ng	99
35) Caprolactam	8.56	113	201960m	40.729	ng	
36) 4-Chloro-3-methylphenol	8.67	107	915954	55.312	ng	97
37) 2-Methylnaphthalene	8.83	142	1924826	55.333	ng	99
38) 1-Methylnaphthalene	8.93	142	1837184	55.797	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	805247	52.169	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	843360	96.108	nq	99
43) 2,4,6-Trichlorophenol	9.11	196	621355	56.253	nq	99
44) 2,4,5-Trichlorophenol	9.15	196	657338	53.123	nq	98
46) 1,1'-Biphenyl	9.30	154	2258686	52.663	nq	99
47) 2-Chloronaphthalene	9.33	162	1795610	53.448	nq	98
48) 2-Nitroaniline	9.42	65	688226	59.350	nq	94
49) Acenaphthylene	9.75	152	2811776	53.296	nq	100
50) Dimethylphthalate	9.61	163	2234546	59.739	nq	100
51) 2,6-Dinitrotoluene	9.66	165	492301	61.403	nq	89
52) Acenaphthene	9.92	154	1906940	57.980	nq	99
53) 3-Nitroaniline	9.83	138	259090	25.597	nq	98
54) 2,4-Dinitrophenol	9.94	184	403503	113.008	nq	97
55) Dibenzofuran	10.09	168	2522467	53.493	nq	98
56) 4-Nitrophenol	9.99	139	749052	93.808	nq	98
57) 2,4-Dinitrotoluene	10.07	165	649860	64.343	nq	98
58) Fluorene	10.43	166	1935912	55.516	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.20	232	533702	57.702	nq	99
60) Diethylphthalate	10.31	149	2116379	55.747	nq	99
61) 4-Chlorophenyl-phenylether	10.42	204	908145	53.256	nq	99
62) 4-Nitroaniline	10.45	138	507858	52.322	nq	99
63) Azobenzene	10.58	77	2078661	54.372	nq	99
65) 4,6-Dinitro-2-methylphenol	10.47	198	295544	69.761	nq	84
66) n-Nitrosodiphenylamine	10.55	169	1790711	53.990	nq	97
67) 4-Bromophenyl-phenylether	10.91	248	603853	52.481	nq	98
68) Hexachlorobenzene	10.98	284	658111	55.884	nq	# 91
69) Atrazine	11.07	200	543078	59.726	nq	99
70) Pentachlorophenol	11.17	266	714172	103.426	nq	97
71) Phenanthrene	11.39	178	2785543	53.172	nq	99
72) Anthracene	11.45	178	2848397	53.497	nq	99
73) Carbazole	11.60	167	2696303	51.162	nq	99
74) Di-n-butylphthalate	11.93	149	3148067	55.841	nq	99
75) Fluoranthene	12.58	202	2991015	52.755	nq	99
77) Benzidine	12.70	184	913685	33.815	nq	98
78) Pyrene	12.81	202	2960637	56.503	nq	100
80) Butylbenzylphthalate	13.43	149	1325831	62.561	nq	98
81) Benzo(a)anthracene	14.00	228	2120840	51.082	nq	98
82) 3,3'-Dichlorobenzidine	13.96	252	408145	24.932	nq	98
83) Chrysene	14.04	228	2230487	52.055	nq	100
84) Bis(2-ethylhexyl)phthalate	13.99	149	1647802	65.225	nq	98
85) Di-n-octyl phthalate	14.60	149	2608807	58.716	nq	99
86) Indeno(1,2,3-cd)pyrene	16.94	276	2017001	49.982	nq	100
88) Benzo(b)fluoranthene	15.04	252	2057480	57.746	nq	98
89) Benzo(k)fluoranthene	15.08	252	1762629	51.953	nq	99
90) Benzo(a)pyrene	15.41	252	1725220	53.280	nq	100
91) Dibenzo(a,h)anthracene	16.96	278	1636772	57.792	nq	99
92) Benzo(g,h,i)perylene	17.38	276	1705754	59.950	nq	97

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

