

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092920\
 Data File : BF121727.D
 Acq On : 29 Sep 2020 13:00
 Operator : JU/CG
 Sample : L4158-02MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1-IMS

Manual Integrations
 APPROVED

mohammad
 10/1/2020 11:45:05 AM

Quant Time: Sep 29 13:42:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 29 11:12:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	131958	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	509004	20.00	ng	0.00
39) Acenaphthene-d10	9.82	164	257088	20.00	ng	0.00
64) Phenanthrene-d10	11.30	188	469729	20.00	ng	0.00
76) Chrysene-d12	13.94	240	372764	20.00	ng	0.00
86) Perylene-d12	15.38	264	327011	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	574936	64.75	ng	0.00
7) Phenol-d6	6.42	99	740086	63.53	ng	0.00
23) Nitrobenzene-d5	7.35	82	499435	52.68	ng	-0.01
42) 2,4,6-Tribromophenol	10.61	330	199926	62.57	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	876301	51.62	ng	0.00
79) Terphenyl-d14	12.89	244	831270	45.18	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.57	88	150991	37.014	ng	99
3) Pyridine	3.30	79	425054	37.691	ng	99
4) n-Nitrosodimethylamine	3.25	42	205458	39.278	ng	100
6) Aniline	6.45	93	259996	17.137	ng	100
8) 2-Chlorophenol	6.57	128	377811	41.791	ng	99
9) Benzaldehyde	6.33	77	118593	15.573	ng	94
10) Phenol	6.43	94	482909	38.929	ng	82
11) bis(2-Chloroethyl)ether	6.52	93	367478	39.305	ng	99
12) 1,3-Dichlorobenzene	6.72	146	403106	39.949	ng	99
13) 1,4-Dichlorobenzene	6.80	146	406195	40.091	ng	99
14) 1,2-Dichlorobenzene	6.95	146	384951	41.243	ng	99
15) Benzyl Alcohol	6.93	79	333487	42.118	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.06	45	581645	38.660	ng	100
17) 2-Methylphenol	7.05	107	312550	40.632	ng	98
18) Hexachloroethane	7.29	117	147839	40.751	ng	96
19) n-Nitroso-di-n-propylamine	7.20	70	260190	38.755	ng	97
20) 3+4-Methylphenols	7.20	107	374594	40.532	ng	94
22) Acetophenone	7.19	105	495522	38.370	ng	98
24) Nitrobenzene	7.37	77	412169	40.199	ng	99
25) Isophorone	7.61	82	730692	38.289	ng	99
26) 2-Nitrophenol	7.69	139	193983	40.604	ng	92
27) 2,4-Dimethylphenol	7.73	122	329127	38.637	ng	99
28) bis(2-Chloroethoxy)methane	7.82	93	459648	38.907	ng	100
29) 2,4-Dichlorophenol	7.93	162	303805	39.157	ng	99
30) 1,2,4-Trichlorobenzene	8.01	180	327128	40.033	ng	99
31) Naphthalene	8.09	128	1054539	39.247	ng	99
32) Benzoic acid	7.86	122	17250m	9.258	ng	
33) 4-Chloroaniline	8.14	127	171261	14.704	ng	100
34) Hexachlorobutadiene	8.20	225	192680	40.173	ng	99
35) Caprolactam	8.51	113	89994m	34.831	ng	
36) 4-Chloro-3-methylphenol	8.63	107	329317	39.401	ng	99
37) 2-Methylnaphthalene	8.78	142	694372	39.434	ng	100
38) 1-Methylnaphthalene	8.88	142	640654	39.093	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.95	216	323114	40.235	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.93	237	290601	63.366	nq	98
43) 2,4,6-Trichlorophenol	9.06	196	211845	38.702	nq	98
44) 2,4,5-Trichlorophenol	9.10	196	222995	38.536	nq	95
46) 1,1'-Biphenyl	9.25	154	815790	39.637	nq	99
47) 2-Chloronaphthalene	9.27	162	642232	39.598	nq	99
48) 2-Nitroaniline	9.37	65	217001	43.056	nq	93
49) Acenaphthylene	9.69	152	1004411	40.306	nq	99
50) Dimethylphthalate	9.55	163	844667	45.331	nq	98
51) 2,6-Dinitrotoluene	9.61	165	168422	42.442	nq	94
52) Acenaphthene	9.86	154	585832	37.220	nq	99
53) 3-Nitroaniline	9.77	138	108431	23.587	nq	93
54) 2,4-Dinitrophenol	9.89	184	52326	29.324	nq	95
55) Dibenzofuran	10.03	168	883154	39.844	nq	99
56) 4-Nitrophenol	9.95	139	246323	67.189	nq	97
57) 2,4-Dinitrotoluene	10.02	165	217109	43.488	nq	99
58) Fluorene	10.37	166	683248	40.308	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.15	232	166669	35.428	nq	99
60) Diethylphthalate	10.25	149	719516	39.605	nq	99
61) 4-Chlorophenyl-phenylether	10.36	204	323216	39.805	nq	97
62) 4-Nitroaniline	10.39	138	177079	38.796	nq	98
63) Azobenzene	10.52	77	751438	38.841	nq	98
65) 4,6-Dinitro-2-methylphenol	10.42	198	69250	27.447	nq	99
66) n-Nitrosodiphenylamine	10.48	169	632303	39.236	nq	99
67) 4-Bromophenyl-phenylether	10.85	248	214164	40.045	nq	95
68) Hexachlorobenzene	10.92	284	242479	40.176	nq	94
69) Atrazine	11.01	200	152809	38.167	nq	99
70) Pentachlorophenol	11.12	266	158756	47.896	nq	98
71) Phenanthrene	11.33	178	1014514	39.641	nq	99
72) Anthracene	11.38	178	1042356	39.467	nq	99
73) Carbazole	11.54	167	949168	39.826	nq	99
74) Di-n-butylphthalate	11.87	149	1074336	39.056	nq	99
75) Fluoranthene	12.52	202	1079219	39.503	nq	99
77) Benzidine	12.64	184	420892	34.059	nq	100
78) Pyrene	12.74	202	1101765	40.454	nq	99
80) Butylbenzylphthalate	13.36	149	443512	40.265	nq	93
81) Benzo(a)anthracene	13.93	228	941077	39.905	nq	100
82) 3,3'-Dichlorobenzidine	13.89	252	101611	11.037	nq	98
83) Chrysene	13.97	228	953096	39.911	nq	98
84) Bis(2-ethylhexyl)phthalate	13.92	149	570935	38.800	nq	# 96
85) Di-n-octyl phthalate	14.53	149	966415	37.214	nq	100
87) Indeno(1,2,3-cd)pyrene	16.81	276	903780	42.439	nq	100
88) Benzo(b)fluoranthene	14.97	252	871042	39.844	nq	98
89) Benzo(k)fluoranthene	15.00	252	866937	43.306	nq	98
90) Benzo(a)pyrene	15.33	252	777629	41.847	nq	99
91) Dibenzo(a,h)anthracene	16.82	278	736870	41.567	nq	100
92) Benzo(g,h,i)perylene	17.24	276	767131	44.023	nq	99

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

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