

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012420\
 Data File : BF118686.D
 Acq On : 24 Jan 2020 15:28
 Operator : CG/JU
 Sample : L1241-06MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-12MSD

Manual Integrations
 APPROVED

mohammad
 1/28/2020 8:12:13 AM

Quant Time: Jan 27 01:29:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 20 14:45:03 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	266135	20.00	ng	-0.02
21) Naphthalene-d8	8.15	136	994783	20.00	ng	-0.02
39) Acenaphthene-d10	9.90	164	549465	20.00	ng	-0.03
64) Phenanthrene-d10	11.38	188	1007864	20.00	ng	-0.03
76) Chrysene-d12	14.02	240	600016	20.00	ng	-0.03
87) Perylene-d12	15.47	264	627840	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1863837	130.07	ng	0.00
7) Phenol-d6	6.50	99	2253997	121.02	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1448738	81.46	ng	-0.02
42) 2,4,6-Tribromophenol	10.69	330	775012	121.95	ng	-0.02
45) 2-Fluorobiphenyl	9.22	172	2830612	84.67	ng	-0.02
79) Terphenyl-d14	12.97	244	3344931	121.59	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.68	88	311186	44.836	ng	96
3) Pyridine	3.47	79	803087	41.054	ng	95
4) n-Nitrosodimethylamine	3.37	42	362697	43.247	ng	# 78
6) Aniline	6.53	93	726659	29.630	ng	95
8) 2-Chlorophenol	6.65	128	842771	52.399	ng	96
9) Benzaldehyde	6.41	77	418073	36.910	ng	92
10) Phenol	6.51	94	1033285	48.682	ng	98
11) bis(2-Chloroethyl)ether	6.60	93	776171	49.288	ng	95
12) 1,3-Dichlorobenzene	6.80	146	909281	48.301	ng	98
13) 1,4-Dichlorobenzene	6.88	146	916808	48.515	ng	99
14) 1,2-Dichlorobenzene	7.04	146	864487	48.664	ng	98
15) Benzyl Alcohol	7.00	79	712155	48.217	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.14	45	921567	42.014	ng	99
17) 2-Methylphenol	7.12	107	674542	50.349	ng	97
18) Hexachloroethane	7.38	117	318801	46.524	ng	99
19) n-Nitroso-di-n-propylamine	7.27	70	548337	43.602	ng	99
20) 3+4-Methylphenols	7.27	107	794158	48.551	ng	94
22) Acetophenone	7.27	105	1034463	43.897	ng	# 93
24) Nitrobenzene	7.45	77	842669	46.274	ng	93
25) Isophorone	7.69	82	1517842	46.791	ng	98
26) 2-Nitrophenol	7.76	139	452648	58.288	ng	97
27) 2,4-Dimethylphenol	7.80	122	718611	53.588	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	947389	45.326	ng	99
29) 2,4-Dichlorophenol	8.00	162	720006	48.570	ng	96
30) 1,2,4-Trichlorobenzene	8.09	180	812080	47.246	ng	99
31) Naphthalene	8.17	128	2268206	50.474	ng	99
32) Benzoic acid	7.92	122	512516	50.925	ng	95
33) 4-Chloroaniline	8.21	127	305814	14.724	ng	97
34) Hexachlorobutadiene	8.29	225	472952	45.184	ng	100
35) Caprolactam	8.58	113	242979m	50.086	ng	
36) 4-Chloro-3-methylphenol	8.69	107	720204	49.198	ng	96
37) 2-Methylnaphthalene	8.86	142	1591871	50.186	ng	99
38) 1-Methylnaphthalene	8.96	142	1495268	49.619	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	770446	44.813	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	814410	93.890	ng	99
43) 2,4,6-Trichlorophenol	9.13	196	559297	49.024	ng	99
44) 2,4,5-Trichlorophenol	9.17	196	586662	50.351	ng	98
46) 1,1'-Biphenyl	9.32	154	1913549	50.531	ng	97
47) 2-Chloronaphthalene	9.35	162	1545586	48.423	ng	98
48) 2-Nitroaniline	9.44	65	433349	44.166	ng	86
49) Acenaphthylene	9.76	152	2371121	52.805	ng	97
50) Dimethylphthalate	9.62	163	2125683	59.379	ng	99
51) 2,6-Dinitrotoluene	9.68	165	422095	52.655	ng #	81
52) Acenaphthene	9.93	154	1584589	52.774	ng	99
53) 3-Nitroaniline	9.85	138	236456	25.832	ng	92
54) 2,4-Dinitrophenol	9.95	184	382675	119.575	ng #	37
55) Dibenzofuran	10.11	168	2138447	51.392	ng	95
56) 4-Nitrophenol	10.01	139	711822	103.013	ng	89
57) 2,4-Dinitrotoluene	10.09	165	555490	54.968	ng	87
58) Fluorene	10.45	166	1591198	48.711	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.22	232	515210	50.273	ng	97
60) Diethylphthalate	10.32	149	1709805	49.261	ng	99
61) 4-Chlorophenyl-phenylether	10.44	204	846207	47.630	ng	97
62) 4-Nitroaniline	10.46	138	380049	40.477	ng	93
63) Azobenzene	10.60	77	1569355	46.687	ng	96
65) 4,6-Dinitro-2-methylphenol	10.49	198	278406	62.820	ng	84
66) n-Nitrosodiphenylamine	10.56	169	1502689	49.439	ng	98
67) 4-Bromophenyl-phenylether	10.93	248	580487	46.619	ng #	94
68) Hexachlorobenzene	11.00	284	643788	45.603	ng	97
69) Atrazine	11.08	200	495926	47.074	ng	99
70) Pentachlorophenol	11.19	266	711403	90.638	ng	98
71) Phenanthrene	11.41	178	2412145	49.608	ng	95
72) Anthracene	11.46	178	2480638	51.559	ng	96
73) Carbazole	11.61	167	2152555	48.777	ng	97
74) Di-n-butylphthalate	11.94	149	2496042	50.908	ng	98
75) Fluoranthene	12.59	202	2617212	51.034	ng	98
77) Benzidine	12.71	184	1167901	72.848	ng	99
78) Pyrene	12.82	202	2600035	63.827	ng	95
80) Butylbenzylphthalate	13.44	149	1063208	63.774	ng	95
81) Benzo(a)anthracene	14.01	228	1845962	51.143	ng	96
82) 3,3'-Dichlorobenzidine	13.97	252	442588	34.412	ng	97
83) Chrysene	14.04	228	1796961	51.946	ng	96
84) Bis(2-ethylhexyl)phthalate	14.00	149	1260715	61.897	ng	97
85) Di-n-octyl phthalate	14.61	149	1850595	61.383	ng	96
86) Indeno(1,2,3-cd)pyrene	16.95	276	2252716	74.947	ng	98
88) Benzo(b)fluoranthene	15.05	252	1829989	46.371	ng	98
89) Benzo(k)fluoranthene	15.09	252	1610862	44.058	ng	97
90) Benzo(a)pyrene	15.42	252	1622342	47.655	ng	97
91) Dibenzo(a,h)anthracene	16.96	278	1845296	61.279	ng	98
92) Benzo(g,h,i)perylene	17.39	276	1957191	66.939	ng	97

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

