

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF103020\
 Data File : BF122180.D
 Acq On : 30 Oct 2020 14:33
 Operator : JU/CG
 Sample : PB132732BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB132732BS

Manual Integrations
 APPROVED

mohammad
 11/2/2020 12:35:56 PM

Quant Time: Oct 30 15:10:57 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF101220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 10:16:54 2020
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.710	152	115392	20.00	ng	0.00
21) Naphthalene-d8	7.992	136	474071	20.00	ng	0.00
39) Acenaphthene-d10	9.751	164	242518	20.00	ng	0.00
64) Phenanthrene-d10	11.239	188	449974	20.00	ng	0.00
76) Chrysene-d12	13.886	240	314698	20.00	ng	0.00
86) Perylene-d12	15.315	264	263616	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.346	112	944063	120.75	ng	0.01
7) Phenol-d6	6.375	99	1130091	112.29	ng	0.00
23) Nitrobenzene-d5	7.287	82	776752	84.03	ng	0.00
42) 2,4,6-Tribromophenol	10.545	330	370222	123.41	ng	0.00
45) 2-Fluorobiphenyl	9.075	172	1292099	78.39	ng	0.00
79) Terphenyl-d14	12.822	244	1362838	82.53	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.481	88	119632	34.79	ng	98
3) Pyridine	3.216	79	326559	36.12	ng	99
4) n-Nitrosodimethylamine	3.169	42	200552	45.89	ng	93
6) Aniline	6.381	93	403931	32.59	ng	# 22
8) 2-Chlorophenol	6.504	128	368744	46.66	ng	97
9) Benzaldehyde	6.269	77	119403	17.66	ng	98
10) Phenol	6.387	94	445121	42.86	ng	99
11) bis(2-Chloroethyl)ether	6.451	93	364404	45.38	ng	100
12) 1,3-Dichlorobenzene	6.651	146	383073	43.10	ng	98
13) 1,4-Dichlorobenzene	6.728	146	392029	43.93	ng	99
14) 1,2-Dichlorobenzene	6.881	146	359137	44.02	ng	97
15) Benzyl Alcohol	6.869	79	326171	50.11	ng	95
16) 2,2'-oxybis(1-Chloropr...	6.987	45	524207	42.42	ng	# 11
17) 2-Methylphenol	6.987	107	296751	43.92	ng	# 86
18) Hexachloroethane	7.210	117	144653	44.87	ng	98
19) n-Nitroso-di-n-propyla...	7.140	70	271090	47.31	ng	97
20) 3+4-Methylphenols	7.145	107	390772	47.96	ng	# 68
22) Acetophenone	7.134	105	508299	41.67	ng	# 89
24) Nitrobenzene	7.304	77	409801	41.47	ng	100
25) Isophorone	7.545	82	733950	41.99	ng	99
26) 2-Nitrophenol	7.622	139	194275	42.70	ng	95
27) 2,4-Dimethylphenol	7.663	122	320498	41.32	ng	96
28) bis(2-Chloroethoxy)met...	7.751	93	455621	41.66	ng	100
29) 2,4-Dichlorophenol	7.869	162	310532	42.75	ng	98
30) 1,2,4-Trichlorobenzene	7.939	180	316186	40.94	ng	99
31) Naphthalene	8.016	128	1032678	40.66	ng	99
32) Benzoic acid	7.851	122	165189	38.98	ng	94
33) 4-Chloroaniline	8.081	127	290301	26.84	ng	98
34) Hexachlorobutadiene	8.128	225	195854	42.77	ng	99
35) Caprolactam	8.469	113	94639m	41.03	ng	
36) 4-Chloro-3-methylphenol	8.581	107	340085	43.59	ng	99
37) 2-Methylnaphthalene	8.710	142	680623	41.38	ng	98
38) 1-Methylnaphthalene	8.804	142	624460	41.00	ng	100
40) 1,2,4,5-Tetrachloroben...	8.875	216	326454	41.71	ng	100
41) Hexachlorocyclopentadiene	8.857	237	134115	62.32	ng	98
43) 2,4,6-Trichlorophenol	8.998	196	198734	41.15	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.051	196	198570	38.07	ng	98
46) 1,1'-Biphenyl	9.175	154	825712	41.87	ng	97
47) 2-Chloronaphthalene	9.198	162	636415	41.50	ng	99
48) 2-Nitroaniline	9.310	65	222057	43.90	ng	95
49) Acenaphthylene	9.610	152	941920	39.99	ng	99
50) Dimethylphthalate	9.481	163	746567	42.10	ng	100
51) 2,6-Dinitrotoluene	9.551	165	165963	42.66	ng	93
52) Acenaphthene	9.781	154	581677	41.52	ng	99
53) 3-Nitroaniline	9.728	138	122563	27.70	ng	96
54) 2,4-Dinitrophenol	9.851	184	136173	69.84	ng	91
55) Dibenzofuran	9.957	168	848375	41.41	ng	97
56) 4-Nitrophenol	9.916	139	116332	48.34	ng	# 78
57) 2,4-Dinitrotoluene	9.963	165	212633	44.40	ng	# 87
58) Fluorene	10.298	166	687065	41.62	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.086	232	170462	42.24	ng	99
60) Diethylphthalate	10.181	149	743806	43.50	ng	99
61) 4-Chlorophenyl-phenyle...	10.286	204	342155	43.22	ng	95
62) 4-Nitroaniline	10.351	138	160213	36.25	ng	93
63) Azobenzene	10.451	77	773258	42.76	ng	98
65) 4,6-Dinitro-2-methylph...	10.381	198	118074	39.98	ng	93
66) n-Nitrosodiphenylamine	10.416	169	638842	41.21	ng	99
67) 4-Bromophenyl-phenylether	10.775	248	222835	42.86	ng	96
68) Hexachlorobenzene	10.851	284	254255	43.21	ng	95
69) Atrazine	10.945	200	174286	45.99	ng	98
70) Pentachlorophenol	11.057	266	154908	68.25	ng	98
71) Phenanthrene	11.263	178	1000038	40.53	ng	99
72) Anthracene	11.316	178	1038403	40.58	ng	100
73) Carbazole	11.475	167	922031	40.52	ng	99
74) Di-n-butylphthalate	11.792	149	1124665	42.96	ng	99
75) Fluoranthene	12.451	202	1061449	40.12	ng	98
77) Benzidine	12.580	184	458702	47.28	ng	99
78) Pyrene	12.680	202	1074999	44.83	ng	99
80) Butylbenzylphthalate	13.292	149	452048	47.53	ng	99
81) Benzo(a)anthracene	13.874	228	880519	42.70	ng	100
82) 3,3'-Dichlorobenzidine	13.833	252	201661	26.36	ng	98
83) Chrysene	13.910	228	848846	41.93	ng	99
84) Bis(2-ethylhexyl)phtha...	13.845	149	636002	49.26	ng	99
85) Di-n-octyl phthalate	14.457	149	984911	47.16	ng	99
87) Indeno(1,2,3-cd)pyrene	16.739	276	740278	43.25	ng	98
88) Benzo(b)fluoranthene	14.910	252	751164	42.15	ng	99
89) Benzo(k)fluoranthene	14.939	252	733304	43.32	ng	98
90) Benzo(a)pyrene	15.263	252	647009	42.04	ng	99
91) Dibenzo(a,h)anthracene	16.733	278	614847	43.63	ng	98
92) Benzo(g,h,i)perylene	17.151	276	622419	44.13	ng	97

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

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Abundance

TIC: BF122180.D\data.ms