

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011821\
 Data File : BF122981.D
 Acq On : 18 Jan 2021 10:50
 Operator : JU/CG
 Sample : SP5424
 Misc : 8270 SPIKE
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP5424

Manual Integrations
 APPROVED

mohammad
 1/20/2021 2:15:08 PM

Quant Time: Jan 18 13:28:15 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF011621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 18 13:14:08 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.898	152	237360	20.00	ng	0.00
21) Naphthalene-d8	8.180	136	967610	20.00	ng	# 0.00
39) Acenaphthene-d10	9.939	164	504608	20.00	ng	0.00
64) Phenanthrene-d10	11.427	188	844990	20.00	ng	# 0.00
76) Chrysene-d12	14.074	240	529602	20.00	ng	# 0.00
86) Perylene-d12	15.551	264	482772	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	0.000	112	0	0.00	ng	
7) Phenol-d6	0.000	99	0d	0.00	ng	
23) Nitrobenzene-d5	0.000	82	0d	0.00	ng	
42) 2,4,6-Tribromophenol	0.000	330	0d	0.00	ng	
45) 2-Fluorobiphenyl	0.000	172	0d	0.00	ng	
79) Terphenyl-d14	0.000	244	0	0.00	ng	
Target Compounds						Qvalue
2) 1,4-Dioxane	2.828	88	443728	58.18	ng	# 85
3) Pyridine	3.557	79	789805	38.26	ng	# 68
4) n-Nitrosodimethylamine	3.528	42	501514	54.43	ng	# 68
6) Aniline	6.557	93	869549	34.16	ng	96
8) 2-Chlorophenol	6.681	128	768604	46.35	ng	89
9) Benzaldehyde	6.445	77	551739	54.17	ng	92
10) Phenol	6.528	94	997160	47.05	ng	94
11) bis(2-Chloroethyl)ether	6.628	93	793525	46.24	ng	85
12) 1,3-Dichlorobenzene	6.839	146	828482	43.17	ng	96
13) 1,4-Dichlorobenzene	6.916	146	860509	44.53	ng	97
14) 1,2-Dichlorobenzene	7.069	146	845746	46.91	ng	100
15) Benzyl Alcohol	7.039	79	753599	50.66	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.169	45	1582006	52.89	ng	89
17) 2-Methylphenol	7.145	107	685028	48.70	ng	97
18) Hexachloroethane	7.410	117	323872	46.37	ng	90
19) n-Nitroso-di-n-propyla...	7.310	70	611507	48.91	ng	# 83
20) 3+4-Methylphenols	7.298	107	814290	46.28	ng	94
22) Acetophenone	7.304	105	1083321	44.52	ng	# 91
24) Nitrobenzene	7.475	77	886597	46.01	ng	93
25) Isophorone	7.722	82	1528291	45.55	ng	99
26) 2-Nitrophenol	7.792	139	437572	50.26	ng	97
27) 2,4-Dimethylphenol	7.828	122	743868	53.66	ng	97
28) bis(2-Chloroethoxy)met...	7.928	93	1047842	46.20	ng	99
29) 2,4-Dichlorophenol	8.033	162	671680	46.41	ng	96
30) 1,2,4-Trichlorobenzene	8.122	180	676695	43.89	ng	92
31) Naphthalene	8.204	128	2332112	46.88	ng	99
32) Benzoic acid	7.957	122	567221m	48.70	ng	
33) 4-Chloroaniline	8.245	127	612920	28.24	ng	# 93
34) Hexachlorobutadiene	8.322	225	355984	42.24	ng	100
35) Caprolactam	8.622	113	232397m	47.24	ng	
36) 4-Chloro-3-methylphenol	8.727	107	788094	51.19	ng	94
37) 2-Methylnaphthalene	8.898	142	1667668	48.60	ng	99
38) 1-Methylnaphthalene	8.998	142	1470087	45.36	ng	95
40) 1,2,4,5-Tetrachloroben...	9.063	216	515833	37.24	ng	# 95
41) Hexachlorocyclopentadiene	9.051	237	681756	101.31	ng	97
43) 2,4,6-Trichlorophenol	9.169	196	437736	42.39	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)	
44) 2,4,5-Trichlorophenol	9.204	196	451274	43.07	ng	#	93
46) 1,1'-Biphenyl	9.363	154	1779972	43.37	ng		97
47) 2-Chloronaphthalene	9.386	162	1392849	40.68	ng		98
48) 2-Nitroaniline	9.474	65	524010	44.90	ng	#	78
49) Acenaphthylene	9.804	152	2044101	40.92	ng		100
50) Dimethylphthalate	9.663	163	1604456	40.96	ng		99
51) 2,6-Dinitrotoluene	9.722	165	360283	42.06	ng	#	82
52) Acenaphthene	9.974	154	1509147	46.74	ng		98
53) 3-Nitroaniline	9.886	138	286652	27.07	ng		85
54) 2,4-Dinitrophenol	9.992	184	336036	70.47	ng	#	76
55) Dibenzofuran	10.151	168	1853371	41.18	ng		99
56) 4-Nitrophenol	10.045	139	695856	89.15	ng	#	75
57) 2,4-Dinitrotoluene	10.127	165	477110	43.30	ng		92
58) Fluorene	10.492	166	1360161	38.64	ng		99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	364039	43.53	ng	#	80
60) Diethylphthalate	10.363	149	1516574	39.84	ng		98
61) 4-Chlorophenyl-phenyle...	10.480	204	613740	38.77	ng		89
62) 4-Nitroaniline	10.504	138	385715	36.93	ng		98
63) Azobenzene	10.639	77	1710823	44.92	ng		90
65) 4,6-Dinitro-2-methylph...	10.533	198	205942	43.27	ng	#	77
66) n-Nitrosodiphenylamine	10.598	169	1314100	41.66	ng		98
67) 4-Bromophenyl-phenylether	10.974	248	374536	36.25	ng		97
68) Hexachlorobenzene	11.039	284	419396	35.50	ng	#	94
69) Atrazine	11.127	200	276260	32.89	ng		97
70) Pentachlorophenol	11.233	266	454197	73.10	ng		97
71) Phenanthrene	11.457	178	2114431	41.76	ng		99
72) Anthracene	11.510	178	1975037	39.46	ng		100
73) Carbazole	11.657	167	1885505	38.87	ng		99
74) Di-n-butylphthalate	11.992	149	2170217	37.80	ng		99
75) Fluoranthene	12.645	202	1877091	36.64	ng		95
77) Benzidine	12.762	184	921866	65.74	ng		100
78) Pyrene	12.874	202	1954054	46.08	ng		100
80) Butylbenzylphthalate	13.492	149	967772	50.92	ng		89
81) Benzo(a)anthracene	14.062	228	1575760	44.46	ng		98
82) 3,3'-Dichlorobenzidine	14.027	252	477369	41.14	ng		99
83) Chrysene	14.104	228	1480150	41.41	ng		100
84) Bis(2-ethylhexyl)phtha...	14.057	149	1206012	47.60	ng		99
85) Di-n-octyl phthalate	14.668	149	1964501	47.93	ng	#	86
87) Indeno(1,2,3-cd)pyrene	17.050	276	1759949	57.87	ng	#	93
88) Benzo(b)fluoranthene	15.121	252	1722082m	52.85	ng		
89) Benzo(k)fluoranthene	15.156	252	1504222	45.03	ng		99
90) Benzo(a)pyrene	15.498	252	1273989	43.28	ng		98
91) Dibenzo(a,h)anthracene	17.074	278	1438050	57.71	ng		97
92) Benzo(g,h,i)perylene	17.509	276	1265501	52.46	ng	#	95

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

