

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF030121\
 Data File : BF123327.D
 Acq On : 1 Mar 2021 15:29
 Operator : JU/CG
 Sample : M1509-01MSD
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-022621MSD

Manual Integrations
 APPROVED

mohammad
 3/2/2021 9:43:02 AM

Quant Time: Mar 01 15:58:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 19 16:08:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	278401	20.00	ng	0.00
21) Naphthalene-d8	8.139	136	1115028	20.00	ng	# 0.00
39) Acenaphthene-d10	9.904	164	570979	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	1022782	20.00	ng	0.00
76) Chrysene-d12	14.051	240	657350	20.00	ng	0.00
86) Perylene-d12	15.521	264	618083	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	2085088	107.80	ng	0.05
7) Phenol-d6	6.516	99	2723948	103.67	ng	0.02
23) Nitrobenzene-d5	7.428	82	1908883	78.37	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	667225	114.76	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	2831613	71.24	ng	0.00
79) Terphenyl-d14	12.986	244	2486734	72.74	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.952	88	344132m	30.83	ng	Qvalue
3) Pyridine	3.651	79	944898	34.74	ng	93
4) n-Nitrosodimethylamine	3.628	42	516028	41.03	ng	82
6) Aniline	6.528	93	788002	25.39	ng	# 1
8) 2-Chlorophenol	6.651	128	948780	45.66	ng	92
9) Benzaldehyde	6.410	77	340969	21.82	ng	91
10) Phenol	6.528	94	1251021	43.69	ng	86
11) bis(2-Chloroethyl)ether	6.598	93	981255	47.77	ng	95
12) 1,3-Dichlorobenzene	6.798	146	946233	43.78	ng	# 93
13) 1,4-Dichlorobenzene	6.875	146	959217	43.54	ng	94
14) 1,2-Dichlorobenzene	7.028	146	907952m	44.31	ng	
15) Benzyl Alcohol	7.010	79	935710	45.67	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.134	45	1183576	42.01	ng	72
17) 2-Methylphenol	7.128	107	788539	45.12	ng	# 85
18) Hexachloroethane	7.369	117	384112	48.27	ng	97
19) n-Nitroso-di-n-propyla...	7.287	70	730606	46.74	ng	95
20) 3+4-Methylphenols	7.281	107	952982	41.52	ng	91
22) Acetophenone	7.275	105	1336833	46.22	ng	# 88
24) Nitrobenzene	7.445	77	1115846	43.45	ng	91
25) Isophorone	7.687	82	2040841	44.81	ng	96
26) 2-Nitrophenol	7.757	139	512501	47.64	ng	# 85
27) 2,4-Dimethylphenol	7.804	122	842291	50.58	ng	95
28) bis(2-Chloroethoxy)met...	7.892	93	1191983	50.00	ng	99
29) 2,4-Dichlorophenol	8.004	162	763912	44.43	ng	94
30) 1,2,4-Trichlorobenzene	8.081	180	839229	44.89	ng	99
31) Naphthalene	8.163	128	2640535	43.21	ng	98
32) Benzoic acid	7.981	122	621236	50.57	ng	86
33) 4-Chloroaniline	8.210	127	347182	13.96	ng	# 93
34) Hexachlorobutadiene	8.281	225	471718	44.73	ng	99
35) Caprolactam	8.628	113	284374m	49.29	ng	
36) 4-Chloro-3-methylphenol	8.710	107	932899	45.64	ng	89
37) 2-Methylnaphthalene	8.857	142	1797285	44.15	ng	98
38) 1-Methylnaphthalene	8.957	142	1672747	43.94	ng	99
40) 1,2,4,5-Tetrachloroben...	9.022	216	791721	46.40	ng	# 97
41) Hexachlorocyclopentadiene	9.010	237	704781	88.17	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	577363	47.71	ng	93

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44) 2,4,5-Trichlorophenol	9.192	196	503692	38.90	ng	93
46) 1,1'-Biphenyl	9.322	154	2125779	44.88	ng	98
47) 2-Chloronaphthalene	9.351	162	1613870	43.18	ng	99
48) 2-Nitroaniline	9.451	65	621809	46.44	ng	# 79
49) Acenaphthylene	9.769	152	2392344	42.20	ng	98
50) Dimethylphthalate	9.633	163	2039538	47.77	ng	99
51) 2,6-Dinitrotoluene	9.692	165	470322	47.79	ng	# 68
52) Acenaphthene	9.939	154	1862833m	47.66	ng	
53) 3-Nitroaniline	9.863	138	275841	25.26	ng	# 79
54) 2,4-Dinitrophenol	9.975	184	438870	84.90	ng	# 82
55) Dibenzofuran	10.116	168	2205641	41.68	ng	97
56) 4-Nitrophenol	10.045	139	719503	82.67	ng	# 64
57) 2,4-Dinitrotoluene	10.104	165	575131	43.46	ng	# 75
58) Fluorene	10.457	166	1789758	42.98	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.233	232	435317	42.14	ng	# 84
60) Diethylphthalate	10.333	149	1945955	46.02	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	857667	43.26	ng	96
62) 4-Nitroaniline	10.492	138	431890	39.28	ng	85
63) Azobenzene	10.610	77	2237911	45.94	ng	97
65) 4,6-Dinitro-2-methylph...	10.516	198	306152	44.47	ng	84
66) n-Nitrosodiphenylamine	10.569	169	1573932	44.50	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	521722	46.79	ng	98
68) Hexachlorobenzene	11.010	284	548179	46.67	ng	99
69) Atrazine	11.104	200	419790	37.56	ng	98
70) Pentachlorophenol	11.210	266	659228	88.88	ng	99
71) Phenanthrene	11.427	178	2992930	49.48	ng	100
72) Anthracene	11.475	178	2599772	43.08	ng	100
73) Carbazole	11.627	167	2311970	40.68	ng	98
74) Di-n-butylphthalate	11.957	149	2855585	44.94	ng	98
75) Fluoranthene	12.621	202	3192083	49.79	ng	99
77) Benzidine	12.739	184	812108	38.25	ng	99
78) Pyrene	12.851	202	3005960	58.57	ng	98
80) Butylbenzylphthalate	13.463	149	1191372	54.92	ng	93
81) Benzo(a)anthracene	14.039	228	2222711	49.75	ng	98
82) 3,3'-Dichlorobenzidine	14.004	252	452651	30.48	ng	100
83) Chrysene	14.080	228	2083163	47.48	ng	98
84) Bis(2-ethylhexyl)phtha...	14.021	149	1598031	53.51	ng	99
85) Di-n-octyl phthalate	14.639	149	2452634	48.07	ng	97
87) Indeno(1,2,3-cd)pyrene	17.033	276	2258286	52.87	ng	98
88) Benzo(b)fluoranthene	15.098	252	2470205	56.90	ng	99
89) Benzo(k)fluoranthene	15.127	252	1594123	40.50	ng	99
90) Benzo(a)pyrene	15.468	252	1873281	48.39	ng	99
91) Dibenzo(a,h)anthracene	17.051	278	1775794	48.46	ng	98
92) Benzo(g,h,i)perylene	17.486	276	1879482	55.68	ng	99

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

