

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
 Data File : BF125930.D
 Acq On : 26 Oct 2021 12:41
 Operator : JU/CG
 Sample : M4325-01MS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RED-BOX-1-225057MS

Manual Integrations
 APPROVED

mohammad
 10/27/2021 11:48:48 AM

Quant Time: Oct 26 14:42:27 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 25 18:49:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	136848	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	566804	20.000	ng	# 0.00
39) Acenaphthene-d10	9.904	164	326610	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	608151	20.000	ng	0.00
76) Chrysene-d12	14.033	240	379881	20.000	ng	# 0.00
86) Perylene-d12	15.498	264	315720	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	1069000	116.356	ng	0.00
7) Phenol-d6	6.492	99	1370232	108.738	ng	0.00
23) Nitrobenzene-d5	7.434	82	927267	84.503	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	350295	123.031	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1447986	78.401	ng	0.00
79) Terphenyl-d14	12.980	244	1631583	70.254	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.693	88	198042	44.115	ng	# 100
3) Pyridine	3.451	79	450325	37.885	ng	# 86
4) n-Nitrosodimethylamine	3.387	42	358897	47.075	ng	# 73
6) Aniline	6.528	93	533784	35.963	ng	# 90
8) 2-Chlorophenol	6.651	128	470310	50.716	ng	# 80
9) Benzaldehyde	6.416	77	298129	39.538	ng	93
10) Phenol	6.510	94	597661m	49.190	ng	
11) bis(2-Chloroethyl)ether	6.604	93	478757	48.930	ng	# 78
12) 1,3-Dichlorobenzene	6.810	146	478561m	46.717	ng	
13) 1,4-Dichlorobenzene	6.887	146	484133	47.245	ng	94
14) 1,2-Dichlorobenzene	7.039	146	472344	47.553	ng	96
15) Benzyl Alcohol	7.010	79	470438	49.314	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.145	45	1146043	47.826	ng	96
17) 2-Methylphenol	7.116	107	398346	49.732	ng	# 91
18) Hexachloroethane	7.387	117	191738	49.351	ng	# 88
19) n-Nitroso-di-n-propyla...	7.287	70	355906	49.435	ng	# 93
20) 3+4-Methylphenols	7.269	107	458282	44.759	ng	88
22) Acetophenone	7.281	105	656304	44.848	ng	# 88
24) Nitrobenzene	7.451	77	560938	49.510	ng	# 87
25) Isophorone	7.692	82	999281	47.819	ng	# 86
26) 2-Nitrophenol	7.763	139	226166	61.567	ng	# 62
27) 2,4-Dimethylphenol	7.804	122	424659	52.060	ng	88
28) bis(2-Chloroethoxy)met...	7.904	93	601167	45.854	ng	# 97
29) 2,4-Dichlorophenol	8.004	162	383131	47.915	ng	89
30) 1,2,4-Trichlorobenzene	8.092	180	421758	44.972	ng	97
31) Naphthalene	8.175	128	1314490	45.610	ng	98
32) Benzoic acid	7.916	122	304340	50.848	ng	# 86
33) 4-Chloroaniline	8.216	127	196305	16.529	ng	# 89
34) Hexachlorobutadiene	8.298	225	263350	44.681	ng	98
35) Caprolactam	8.592	113	139212m	53.184	ng	
36) 4-Chloro-3-methylphenol	8.698	107	458187	48.963	ng	89
37) 2-Methylnaphthalene	8.863	142	886415	47.782	ng	91
38) 1-Methylnaphthalene	8.963	142	828929	46.297	ng	89
40) 1,2,4,5-Tetrachloroben...	9.033	216	374823	41.660	ng	98
41) Hexachlorocyclopentadiene	9.022	237	499975	102.888	ng	99
43) 2,4,6-Trichlorophenol	9.139	196	287496	46.819	ng	92

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.175	196	298156	46.023	ng	95
46) 1,1'-Biphenyl	9.328	154	1030257	43.067	ng	98
47) 2-Chloronaphthalene	9.351	162	794232	44.035	ng	95
48) 2-Nitroaniline	9.445	65	353385	54.855	ng	# 70
49) Acenaphthylene	9.769	152	1240767	45.541	ng	98
50) Dimethylphthalate	9.633	163	931770	46.055	ng	98
51) 2,6-Dinitrotoluene	9.686	165	222215	55.686	ng	# 60
52) Acenaphthene	9.939	154	874378	49.383	ng	98
53) 3-Nitroaniline	9.851	138	134344	29.367	ng	# 62
54) 2,4-Dinitrophenol	9.957	184	181369	105.830	ng	# 74
55) Dibenzofuran	10.116	168	1149915	43.482	ng	97
56) 4-Nitrophenol	10.010	139	368467	103.651	ng	# 53
57) 2,4-Dinitrotoluene	10.092	165	297369	60.669	ng	# 96
58) Fluorene	10.457	166	826607	50.130	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.227	232	249510	47.184	ng	# 91
60) Diethylphthalate	10.333	149	963637	46.695	ng	98
61) 4-Chlorophenyl-phenyle...	10.445	204	406199	41.739	ng	92
62) 4-Nitroaniline	10.469	138	216897	51.674	ng	92
63) Azobenzene	10.610	77	1091978	45.878	ng	95
65) 4,6-Dinitro-2-methylph...	10.498	198	135011	50.232	ng	94
66) n-Nitrosodiphenylamine	10.569	169	796088	43.704	ng	96
67) 4-Bromophenyl-phenylether	10.939	248	287989	44.329	ng	95
68) Hexachlorobenzene	11.004	284	307301	44.868	ng	# 88
69) Atrazine	11.092	200	288945	50.086	ng	94
70) Pentachlorophenol	11.192	266	352715	91.221	ng	95
71) Phenanthrene	11.416	178	1375932	43.603	ng	98
72) Anthracene	11.469	178	1426217	45.267	ng	98
73) Carbazole	11.622	167	1288011	43.004	ng	99
74) Di-n-butylphthalate	11.957	149	1530649	47.969	ng	99
75) Fluoranthene	12.604	202	1490734	46.861	ng	98
77) Benzidine	12.721	184	475467	40.125	ng	98
78) Pyrene	12.833	202	1521863	43.049	ng	97
80) Butylbenzylphthalate	13.457	149	637328	55.421	ng	# 83
81) Benzo(a)anthracene	14.021	228	1123302	43.934	ng	99
82) 3,3'-Dichlorobenzidine	13.980	252	309654	38.196	ng	99
83) Chrysene	14.057	228	1196085	46.237	ng	97
84) Bis(2-ethylhexyl)phtha...	14.021	149	747497	59.999	ng	98
85) Di-n-octyl phthalate	14.633	149	1357880	70.353	ng	# 93
87) Indeno(1,2,3-cd)pyrene	16.968	276	1076086	47.311	ng	# 90
88) Benzo(b)fluoranthene	15.068	252	1049952	50.383	ng	# 95
89) Benzo(k)fluoranthene	15.104	252	955233	47.721	ng	99
90) Benzo(a)pyrene	15.439	252	954219	56.719	ng	# 95
91) Dibenzo(a,h)anthracene	16.992	278	898297	48.751	ng	# 94
92) Benzo(g,h,i)perylene	17.415	276	921245	47.853	ng	# 88

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

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