

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052121\
 Data File : BF124027.D
 Acq On : 21 May 2021 19:59
 Operator : JU/CG
 Sample : M2482-01MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-052021MS

Manual Integrations
 APPROVED

mohammad
 5/24/2021 2:36:21 PM

Quant Time: May 22 01:39:21 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 14 19:05:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	58258	20.00	ng	-0.01
21) Naphthalene-d8	8.175	136	230842	20.00	ng	#-0.01
39) Acenaphthene-d10	9.933	164	141712	20.00	ng	-0.01
64) Phenanthrene-d10	11.422	188	238365	20.00	ng	0.00
76) Chrysene-d12	14.063	240	143002	20.00	ng	0.00
86) Perylene-d12	15.551	264	123631	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	236305	56.95	ng	0.00
7) Phenol-d6	6.522	99	577420	107.27	ng	0.00
23) Nitrobenzene-d5	7.457	82	431899	80.61	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	31162	18.31	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	825798	71.71	ng	-0.01
79) Terphenyl-d14	13.004	244	770436	81.60	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.710	88	55068	30.22	ng	99
3) Pyridine	3.469	79	143696	30.55	ng	88
4) n-Nitrosodimethylamine	3.428	42	114595	37.51	ng	# 86
6) Aniline	6.551	93	183714	30.06	ng	94
8) 2-Chlorophenol	6.675	128	176228	40.81	ng	95
9) Benzaldehyde	6.439	77	138200	59.34	ng	94
10) Phenol	6.539	94	232591	42.75	ng	95
11) bis(2-Chloroethyl)ether	6.628	93	173705	41.08	ng	91
12) 1,3-Dichlorobenzene	6.834	146	195430	38.85	ng	96
13) 1,4-Dichlorobenzene	6.910	146	200418	38.94	ng	96
14) 1,2-Dichlorobenzene	7.063	146	193861	40.31	ng	95
15) Benzyl Alcohol	7.034	79	187026	40.23	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.169	45	339869	39.93	ng	98
17) 2-Methylphenol	7.145	107	143940	41.11	ng	96
18) Hexachloroethane	7.404	117	76573	40.14	ng	96
19) n-Nitroso-di-n-propyla...	7.310	70	150423	41.36	ng	98
20) 3+4-Methylphenols	7.298	107	192359	41.70	ng	88
22) Acetophenone	7.304	105	266714	40.69	ng	# 97
24) Nitrobenzene	7.475	77	212337	38.76	ng	93
25) Isophorone	7.716	82	371989	36.09	ng	96
26) 2-Nitrophenol	7.792	139	95453	46.32	ng	89
27) 2,4-Dimethylphenol	7.828	122	160147	40.58	ng	96
28) bis(2-Chloroethoxy)met...	7.922	93	215720	41.62	ng	98
29) 2,4-Dichlorophenol	8.034	162	164721	41.24	ng	93
30) 1,2,4-Trichlorobenzene	8.116	180	169221	37.68	ng	99
31) Naphthalene	8.198	128	521283	38.16	ng	98
32) Benzoic acid	7.957	122	86235	39.82	ng	90
33) 4-Chloroaniline	8.239	127	31856	6.13	ng	94
34) Hexachlorobutadiene	8.316	225	114946	39.10	ng	94
35) Caprolactam	8.622	113	52548m	48.32	ng	
36) 4-Chloro-3-methylphenol	8.728	107	184282	42.19	ng	88
37) 2-Methylnaphthalene	8.892	142	369272	39.88	ng	99
38) 1-Methylnaphthalene	8.992	142	338919	39.22	ng	96
40) 1,2,4,5-Tetrachloroben...	9.057	216	190283	37.23	ng	98
41) Hexachlorocyclopentadiene	9.045	237	199104	61.25	ng	96
43) 2,4,6-Trichlorophenol	9.169	196	122429	36.92	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	130113	37.77	ng	# 90
46) 1,1'-Biphenyl	9.357	154	475910	38.13	ng	99
47) 2-Chloronaphthalene	9.380	162	360595	36.37	ng	95
48) 2-Nitroaniline	9.475	65	141051	42.13	ng	98
49) Acenaphthylene	9.798	152	550839	37.61	ng	99
50) Dimethylphthalate	9.657	163	583950	51.60	ng	100
51) 2,6-Dinitrotoluene	9.716	165	96609	42.22	ng	95
52) Acenaphthene	9.969	154	402700	40.16	ng	94
53) 3-Nitroaniline	9.880	138	55543	24.42	ng	93
54) 2,4-Dinitrophenol	9.992	184	84242	69.53	ng	95
55) Dibenzofuran	10.139	168	524086	36.39	ng	99
56) 4-Nitrophenol	10.045	139	140862	74.88	ng	86
57) 2,4-Dinitrotoluene	10.122	165	128201	45.68	ng	# 94
58) Fluorene	10.480	166	418158	38.53	ng	92
59) 2,3,4,6-Tetrachlorophenol	10.257	232	107906	38.09	ng	94
60) Diethylphthalate	10.357	149	455190	40.62	ng	94
61) 4-Chlorophenyl-phenyle...	10.475	204	212566	39.09	ng	90
62) 4-Nitroaniline	10.498	138	91317	40.37	ng	88
63) Azobenzene	10.633	77	460111	36.92	ng	91
65) 4,6-Dinitro-2-methylph...	10.527	198	60999	40.66	ng	85
66) n-Nitrosodiphenylamine	10.592	169	367294	40.33	ng	97
67) 4-Bromophenyl-phenylether	10.963	248	128833	41.13	ng	90
68) Hexachlorobenzene	11.033	284	138986	39.74	ng	94
69) Atrazine	11.122	200	128124	49.81	ng	97
70) Pentachlorophenol	11.222	266	143128	69.03	ng	98
71) Phenanthrene	11.445	178	613645	40.62	ng	99
72) Anthracene	11.498	178	589428	38.50	ng	98
73) Carbazole	11.651	167	468242	34.96	ng	97
74) Di-n-butylphthalate	11.980	149	681743	40.67	ng	99
75) Fluoranthene	12.633	202	536097	34.27	ng	99
77) Benzidine	12.751	184	179426	52.00	ng	# 95
78) Pyrene	12.863	202	545228	42.95	ng	98
80) Butylbenzylphthalate	13.480	149	224233	45.04	ng	97
81) Benzo(a)anthracene	14.051	228	404061	39.07	ng	98
82) 3,3'-Dichlorobenzidine	14.010	252	81499	25.37	ng	99
83) Chrysene	14.092	228	392228	39.43	ng	97
84) Bis(2-ethylhexyl)phtha...	14.039	149	321641	45.78	ng	98
85) Di-n-octyl phthalate	14.657	149	499399	47.03	ng	100
87) Indeno(1,2,3-cd)pyrene	17.051	276	310899	32.56	ng	96
88) Benzo(b)fluoranthene	15.115	252	397504	41.56	ng	98
89) Benzo(k)fluoranthene	15.145	252	373243	42.82	ng	98
90) Benzo(a)pyrene	15.492	252	353126	42.71	ng	99
91) Dibenzo(a,h)anthracene	17.068	278	257400	32.01	ng	98
92) Benzo(g,h,i)perylene	17.509	276	239395	29.69	ng	# 94

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

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