

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF120920\
 Data File : BF122619.D
 Acq On : 9 Dec 2020 19:49
 Operator : JU/CG
 Sample : L4982-01MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 R-LA-36-5-WCMS

Manual Integrations
 APPROVED

mohammad
 12/10/2020 2:43:12 PM

Quant Time: Dec 10 06:56:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF120520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 07 12:08:49 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.839	152	192286	20.00	ng	0.00
21) Naphthalene-d8	8.122	136	764628	20.00	ng	0.00
39) Acenaphthene-d10	9.880	164	414272	20.00	ng	0.00
64) Phenanthrene-d10	11.363	188	764925	20.00	ng	0.00
76) Chrysene-d12	14.010	240	609929	20.00	ng	0.00
86) Perylene-d12	15.462	264	527125	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	1086534	89.46	ng	0.00
7) Phenol-d6	6.475	99	1380408	85.70	ng	0.00
23) Nitrobenzene-d5	7.404	82	867880	56.87	ng	0.00
42) 2,4,6-Tribromophenol	10.669	330	414137	76.37	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	1408905	46.00	ng	0.00
79) Terphenyl-d14	12.951	244	1444106	41.44	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.651	88	158019	27.68	ng	100
3) Pyridine	3.398	79	362127	24.00	ng	98
4) n-Nitrosodimethylamine	3.340	42	193916	32.05	ng	100
6) Aniline	6.498	93	577433	30.45	ng	98
8) 2-Chlorophenol	6.628	128	448922	33.53	ng	95
9) Benzaldehyde	6.387	77	206270	21.16	ng	99
10) Phenol	6.487	94	620367	37.17	ng	94
11) bis(2-Chloroethyl)ether	6.575	93	421681	33.07	ng	99
12) 1,3-Dichlorobenzene	6.781	146	480999	30.37	ng	98
13) 1,4-Dichlorobenzene	6.857	146	485589	30.40	ng	99
14) 1,2-Dichlorobenzene	7.010	146	466808	30.28	ng	99
15) Benzyl Alcohol	6.981	79	372410	33.58	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.116	45	548139	30.14	ng	98
17) 2-Methylphenol	7.092	107	372776	35.03	ng	98
18) Hexachloroethane	7.351	117	169779	29.83	ng	96
19) n-Nitroso-di-n-propyla...	7.257	70	297815	32.28	ng	98
20) 3+4-Methylphenols	7.245	107	432784	32.19	ng	96
22) Acetophenone	7.251	105	574953	30.24	ng	98
24) Nitrobenzene	7.422	77	458972	30.23	ng	98
25) Isophorone	7.663	82	826618	31.45	ng	99
26) 2-Nitrophenol	7.739	139	236412	31.93	ng	95
27) 2,4-Dimethylphenol	7.775	122	389112	35.85	ng	99
28) bis(2-Chloroethoxy)met...	7.875	93	530695	29.48	ng	99
29) 2,4-Dichlorophenol	7.981	162	387519	30.58	ng	98
30) 1,2,4-Trichlorobenzene	8.069	180	404130	28.14	ng	98
31) Naphthalene	8.145	128	1274115	30.20	ng	100
33) 4-Chloroaniline	8.192	127	506150	28.70	ng	98
34) Hexachlorobutadiene	8.263	225	235669	26.66	ng	99
35) Caprolactam	8.557	113	106762m	27.97	ng	
36) 4-Chloro-3-methylphenol	8.675	107	397310	31.82	ng	97
37) 2-Methylnaphthalene	8.839	142	850319	30.47	ng	98
38) 1-Methylnaphthalene	8.939	142	790528	29.94	ng	98
40) 1,2,4,5-Tetrachloroben...	9.004	216	396193	28.87	ng	99
41) Hexachlorocyclopentadiene	8.992	237	406746	67.04	ng	98
43) 2,4,6-Trichlorophenol	9.116	196	269111	28.40	ng	99
44) 2,4,5-Trichlorophenol	9.157	196	292205	29.83	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	9.304	154	1012038	29.44	ng	98
47) 2-Chloronaphthalene	9.328	162	810669	28.46	ng	98
48) 2-Nitroaniline	9.422	65	245431	29.55	ng	94
49) Acenaphthylene	9.739	152	1269044	30.16	ng	99
50) Dimethylphthalate	9.604	163	1074459	32.48	ng	100
51) 2,6-Dinitrotoluene	9.663	165	223276	31.95	ng	95
52) Acenaphthene	9.916	154	761548m	27.98	ng	
53) 3-Nitroaniline	9.833	138	229253	28.39	ng	98
54) 2,4-Dinitrophenol	9.939	184	41307	12.09	ng	95
55) Dibenzofuran	10.086	168	1136647	29.68	ng	99
56) 4-Nitrophenol	9.998	139	310698	53.97	ng	93
57) 2,4-Dinitrotoluene	10.069	165	290808	31.24	ng	95
58) Fluorene	10.427	166	879268	28.87	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.204	232	234047	27.19	ng	99
60) Diethylphthalate	10.298	149	923167	28.69	ng	99
61) 4-Chlorophenyl-phenyle...	10.422	204	422728	28.19	ng	95
62) 4-Nitroaniline	10.445	138	238314	29.08	ng	98
63) Azobenzene	10.580	77	876537	29.63	ng	97
65) 4,6-Dinitro-2-methylph...	10.475	198	72535	15.55	ng	96
66) n-Nitrosodiphenylamine	10.539	169	801568	29.66	ng	99
67) 4-Bromophenyl-phenylether	10.910	248	291773	28.15	ng	95
68) Hexachlorobenzene	10.980	284	317863	27.74	ng	94
69) Atrazine	11.063	200	224123	25.53	ng	99
70) Pentachlorophenol	11.174	266	226792	37.36	ng	98
71) Phenanthrene	11.392	178	1318483	29.00	ng	99
72) Anthracene	11.445	178	1374097	30.48	ng	99
73) Carbazole	11.598	167	1252455	30.64	ng	98
74) Di-n-butylphthalate	11.927	149	1412488	27.50	ng	98
75) Fluoranthene	12.580	202	1395623	29.28	ng	99
77) Benzidine	12.698	184	370258	28.07	ng	99
78) Pyrene	12.810	202	1401935	29.73	ng	100
80) Butylbenzylphthalate	13.421	149	582922	27.44	ng	97
81) Benzo(a)anthracene	13.998	228	1192005	29.24	ng	98
82) 3,3'-Dichlorobenzidine	13.957	252	362772	24.85	ng	99
83) Chrysene	14.033	228	1192045	29.39	ng	99
84) Bis(2-ethylhexyl)phtha...	13.986	149	769969	26.47	ng	98
85) Di-n-octyl phthalate	14.598	149	1244513	25.79	ng	99
87) Indeno(1,2,3-cd)pyrene	16.921	276	1026003	29.65	ng	99
88) Benzo(b)fluoranthene	15.045	252	1115572	30.16	ng	98
89) Benzo(k)fluoranthene	15.074	252	1035372	30.62	ng	99
90) Benzo(a)pyrene	15.404	252	928532	28.70	ng	98
91) Dibenzo(a,h)anthracene	16.939	278	850401	30.03	ng	99
92) Benzo(g,h,i)perylene	17.362	276	876011	31.96	ng	99

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

