

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051419\
 Data File : BF114257.D
 Acq On : 14 May 2019 17:06
 Operator : JU/SJ
 Sample : K2766-12MS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P026-SS008-0206-01MS

Manual Integrations
 APPROVED

mohammad
 5/16/2019 9:10:02 AM

Quant Time: May 15 02:20:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 10 15:56:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	231916	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	888249	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	421913	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	785791	20.00	ng	0.00
76) Chrysene-d12	14.03	240	524006	20.00	ng	0.01
87) Perylene-d12	15.53	264	562330	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	1580169	109.65	ng	0.01
7) Phenol-d6	6.51	99	2080471	122.06	ng	0.02
23) Nitrobenzene-d5	7.42	82	1347055	85.11	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	452627	103.77	ng	0.01
45) 2-Fluorobiphenyl	9.21	172	2072018	74.29	ng	0.00
79) Terphenyl-d14	12.96	244	1852854	72.89	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.62	88	246508	31.120	ng	99
3) Pyridine	3.43	79	136743	6.265	ng	96
4) n-Nitrosodimethylamine	3.33	42	376515	35.775	ng	# 92
8) 2-Chlorophenol	6.65	128	640111	41.606	ng	97
9) Benzaldehyde	6.40	77	220490	18.478	ng	99
10) Phenol	6.52	94	796180	39.708	ng	86
11) bis(2-Chloroethyl)ether	6.59	93	650595	42.739	ng	98
12) 1,3-Dichlorobenzene	6.80	146	643441	37.258	ng	98
13) 1,4-Dichlorobenzene	6.87	146	656144	37.705	ng	96
14) 1,2-Dichlorobenzene	7.03	146	612481	38.524	ng	99
15) Benzyl Alcohol	7.00	79	532999	40.094	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.13	45	1142714	38.716	ng	53
17) 2-Methylphenol	7.12	107	502500	39.761	ng	# 92
18) Hexachloroethane	7.37	117	240291	36.786	ng	96
19) n-Nitroso-di-n-propylamine	7.27	70	448816	41.543	ng	98
20) 3+4-Methylphenols	7.27	107	624223	42.750	ng	# 60
22) Acetophenone	7.26	105	874314	44.432	ng	# 89
24) Nitrobenzene	7.44	77	696229	43.189	ng	97
25) Isophorone	7.67	82	1257769	42.848	ng	98
26) 2-Nitrophenol	7.76	139	361422	46.211	ng	98
27) 2,4-Dimethylphenol	7.80	122	529848	43.134	ng	99
28) bis(2-Chloroethoxy)methane	7.88	93	806350	42.337	ng	100
29) 2,4-Dichlorophenol	8.00	162	523134	42.769	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	541731	38.948	ng	97
31) Naphthalene	8.16	128	1756180	42.326	ng	100
32) Benzoic acid	7.93	122	295698	35.843	ng	98
34) Hexachlorobutadiene	8.27	225	300483	37.968	ng	100
35) Caprolactam	8.59	113	138239m	34.660	ng	
36) 4-Chloro-3-methylphenol	8.70	107	552202	44.850	ng	97
37) 2-Methylnaphthalene	8.85	142	1171903	43.777	ng	97
38) 1-Methylnaphthalene	8.95	142	1105558	43.854	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	530037	41.472	ng	99
41) Hexachlorocyclopentadiene	9.00	237	258028	45.317	ng	99
43) 2,4,6-Trichlorophenol	9.13	196	361282	41.097	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.18	196	375873	41.692	ng	95
46) 1,1'-Biphenyl	9.31	154	1429459	41.938	ng	97
47) 2-Chloronaphthalene	9.34	162	1083470	39.498	ng	99
48) 2-Nitroaniline	9.44	65	403985	41.526	ng	96
49) Acenaphthylene	9.76	152	1708997	40.962	ng	99
50) Dimethylphthalate	9.61	163	1660473	53.611	ng	97
51) 2,6-Dinitrotoluene	9.68	165	292777	42.619	ng	88
52) Acenaphthene	9.93	154	997366	38.957	ng	99
53) 3-Nitroaniline	9.85	138	69040	8.096	ng	98
54) 2,4-Dinitrophenol	9.96	184	190257	57.180	ng	96
55) Dibenzofuran	10.10	168	1475772	39.311	ng	98
56) 4-Nitrophenol	10.03	139	406163	67.350	ng	88
57) 2,4-Dinitrotoluene	10.09	165	378785	40.624	ng	94
58) Fluorene	10.44	166	1174269	40.496	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.22	232	300473	38.627	ng	99
60) Diethylphthalate	10.31	149	1253284	39.825	ng	99
61) 4-Chlorophenyl-phenylether	10.43	204	573528	40.270	ng	97
62) 4-Nitroaniline	10.46	138	110918	12.378	ng	98
63) Azobenzene	10.59	77	1228103	40.317	ng	98
65) 4,6-Dinitro-2-methylphenol	10.50	198	135622	35.918	ng	89
66) n-Nitrosodiphenylamine	10.55	169	1021717	42.841	ng	98
67) 4-Bromophenyl-phenylether	10.92	248	359128	41.509	ng	97
68) Hexachlorobenzene	11.00	284	402042	42.005	ng	93
69) Atrazine	11.08	200	278298	37.498	ng	99
70) Pentachlorophenol	11.19	266	334643	73.755	ng	98
71) Phenanthrene	11.40	178	1653522	43.252	ng	99
72) Anthracene	11.46	178	1623212	42.273	ng	99
73) Carbazole	11.61	167	1473531	40.242	ng	98
74) Di-n-butylphthalate	11.93	149	1320573	31.304	ng	100
75) Fluoranthene	12.60	202	1701601	41.333	ng	99
78) Pyrene	12.83	202	1776518	42.144	ng	99
80) Butylbenzylphthalate	13.43	149	713772	40.229	ng	95
81) Benzo(a)anthracene	14.02	228	1494061	44.082	ng	100
83) Chrysene	14.06	228	1304608	39.267	ng	99
84) Bis(2-ethylhexyl)phthalate	14.00	149	1090785	47.006	ng	100
85) Di-n-octyl phthalate	14.63	149	1853962	49.285	ng	98
86) Indeno(1,2,3-cd)pyrene	17.03	276	1393846	47.470	ng	99
88) Benzo(b)fluoranthene	15.10	252	1597368	44.339	ng	98
89) Benzo(k)fluoranthene	15.13	252	1476190	44.607	ng	99
90) Benzo(a)pyrene	15.47	252	1464381	46.186	ng	98
91) Dibenzo(a,h)anthracene	17.04	278	1146661	43.523	ng	99
92) Benzo(g,h,i)perylene	17.49	276	1163929	44.084	ng	99

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

