

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081519\
 Data File : BF116275.D
 Acq On : 15 Aug 2019 13:23
 Operator : HP/JU
 Sample : PB122172BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122172BS

Manual Integrations
 APPROVED

Jagrut
 8/16/2019 5:32:44 PM

Quant Time: Aug 15 15:09:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	147245	20.00	ng	-0.03
21) Naphthalene-d8	8.10	136	622467	20.00	ng	-0.03
39) Acenaphthene-d10	9.85	164	339158	20.00	ng	-0.04
64) Phenanthrene-d10	11.34	188	576808	20.00	ng	-0.03
76) Chrysene-d12	13.98	240	394130	20.00	ng	-0.03
87) Perylene-d12	15.44	264	345276	20.00	ng	-0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	1018936	115.85	ng	-0.01
7) Phenol-d6	6.46	99	1264931	113.99	ng	-0.02
23) Nitrobenzene-d5	7.39	82	849867	78.81	ng	-0.03
42) 2,4,6-Tribromophenol	10.65	330	365631	111.19	ng	-0.03
45) 2-Fluorobiphenyl	9.17	172	1478721	81.67	ng	-0.03
79) Terphenyl-d14	12.92	244	1570904	83.99	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	177371	39.917	ng	99
3) Pyridine	3.43	79	416909	33.588	ng	97
4) n-Nitrosodimethylamine	3.39	42	189483	38.682	ng	100
6) Aniline	6.49	93	462283	29.088	ng	# 91
8) 2-Chlorophenol	6.61	128	443997	45.649	ng	100
9) Benzaldehyde	6.37	77	161071	21.036	ng	99
10) Phenol	6.47	94	550598	42.133	ng	89
11) bis(2-Chloroethyl)ether	6.56	93	426920	44.434	ng	99
12) 1,3-Dichlorobenzene	6.76	146	464557	41.329	ng	100
13) 1,4-Dichlorobenzene	6.83	146	475854	42.280	ng	99
14) 1,2-Dichlorobenzene	6.99	146	437341	42.201	ng	99
15) Benzyl Alcohol	6.97	79	349936	41.552	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.09	45	553623	42.245	ng	90
17) 2-Methylphenol	7.08	107	369319	44.844	ng	98
18) Hexachloroethane	7.32	117	175107	42.258	ng	99
19) n-Nitroso-di-n-propylamine	7.23	70	300852	43.749	ng	99
20) 3+4-Methylphenols	7.23	107	445033	45.664	ng	99
22) Acetophenone	7.23	105	592029	43.367	ng	97
24) Nitrobenzene	7.40	77	496486	42.639	ng	98
25) Isophorone	7.64	82	918092	44.524	ng	99
26) 2-Nitrophenol	7.72	139	247885	45.163	ng	94
27) 2,4-Dimethylphenol	7.76	122	401812	48.659	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	562667	44.025	ng	100
29) 2,4-Dichlorophenol	7.96	162	397063	45.540	ng	98
30) 1,2,4-Trichlorobenzene	8.04	180	423296	42.590	ng	99
31) Naphthalene	8.12	128	1296935	44.276	ng	99
32) Benzoic acid	7.92	122	233422	40.094	ng	99
33) 4-Chloroaniline	8.17	127	356047	27.204	ng	98
34) Hexachlorobutadiene	8.23	225	235351	41.111	ng	99
35) Caprolactam	8.55	113	117112m	41.530	ng	
36) 4-Chloro-3-methylphenol	8.66	107	413299	45.565	ng	98
37) 2-Methylnaphthalene	8.81	142	859335	44.545	ng	100
38) 1-Methylnaphthalene	8.91	142	804676	44.091	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	398576	43.959	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	174600	45.641	ng	100
43) 2,4,6-Trichlorophenol	9.10	196	256433	41.089	ng	99
44) 2,4,5-Trichlorophenol	9.15	196	269823	41.825	ng	99
46) 1,1'-Biphenyl	9.27	154	1056607	44.127	ng	99
47) 2-Chloronaphthalene	9.30	162	832839	41.791	ng	97
48) 2-Nitroaniline	9.40	65	258561	39.252	ng	96
49) Acenaphthylene	9.72	152	1221438	42.011	ng	99
50) Dimethylphthalate	9.57	163	991317	42.927	ng	100
51) 2,6-Dinitrotoluene	9.65	165	217212	43.282	ng	94
52) Acenaphthene	9.89	154	751168	42.114	ng	99
53) 3-Nitroaniline	9.82	138	169560	27.344	ng	98
54) 2,4-Dinitrophenol	9.94	184	182738	72.414	ng	90
55) Dibenzofuran	10.06	168	1051221	40.526	ng	100
56) 4-Nitrophenol	10.00	139	270542	68.106	ng	90
57) 2,4-Dinitrotoluene	10.06	165	257089	38.477	ng	94
58) Fluorene	10.40	166	851610	42.672	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.19	232	230295	42.431	ng	99
60) Diethylphthalate	10.27	149	936515	43.437	ng	100
61) 4-Chlorophenyl-phenylether	10.39	204	438558	43.391	ng	97
62) 4-Nitroaniline	10.44	138	212676	33.187	ng	99
63) Azobenzene	10.55	77	962470	43.402	ng	99
65) 4,6-Dinitro-2-methylphenol	10.47	198	142166	46.510	ng	98
66) n-Nitrosodiphenylamine	10.51	169	777041	44.757	ng	99
67) 4-Bromophenyl-phenylether	10.88	248	270943	43.879	ng	98
68) Hexachlorobenzene	10.96	284	302659	42.389	ng	99
69) Atrazine	11.04	200	242449	42.449	ng	99
70) Pentachlorophenol	11.16	266	258550	74.586	ng	97
71) Phenanthrene	11.37	178	1240220	42.059	ng	100
72) Anthracene	11.42	178	1286574	43.112	ng	99
73) Carbazole	11.57	167	1163917	42.989	ng	99
74) Di-n-butylphthalate	11.89	149	1427566	45.199	ng	99
75) Fluoranthene	12.56	202	1293609	41.904	ng	100
77) Benzidine	12.67	184	420392	31.572	ng	98
78) Pyrene	12.79	202	1295135	43.592	ng	99
80) Butylbenzylphthalate	13.39	149	608117	48.388	ng	97
81) Benzo(a)anthracene	13.97	228	1082491	42.971	ng	99
82) 3,3'-Dichlorobenzidine	13.93	252	342051	39.083	ng	99
83) Chrysene	14.01	228	1084242	44.333	ng	99
84) Bis(2-ethylhexyl)phthalate	13.94	149	860925	52.716	ng	# 97
85) Di-n-octyl phthalate	14.55	149	1376812	53.077	ng	99
86) Indeno(1,2,3-cd)pyrene	16.90	276	911616	44.380	ng	100
88) Benzo(b)fluoranthene	15.02	252	950981	47.634	ng	99
89) Benzo(k)fluoranthene	15.04	252	977440	50.266	ng	99
90) Benzo(a)pyrene	15.38	252	907177	50.832	ng	99
91) Dibenzo(a,h)anthracene	16.90	278	765285	49.539	ng	99
92) Benzo(g,h,i)perylene	17.34	276	762743	50.275	ng	98

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

