

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062519\
 Data File : BF115189.D
 Acq On : 25 Jun 2019 16:01
 Operator : HP/JU
 Sample : PB120652BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120652BS

Manual Integrations
 APPROVED

mohammad
 6/26/2019 2:03:25 PM

Quant Time: Jun 26 06:57:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	275605	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1109988	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	558417	20.00	ng	0.00
64) Phenanthrene-d10	11.11	188	1079843	20.00	ng	0.00
76) Chrysene-d12	13.74	240	739785	20.00	ng	-0.01
87) Perylene-d12	15.09	264	777546	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	1962138	109.86	ng	0.02
7) Phenol-d6	6.26	99	2423286	111.79	ng	0.00
23) Nitrobenzene-d5	7.18	82	1496843	82.95	ng	0.00
42) 2,4,6-Tribromophenol	10.42	330	719040	122.10	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	2794556	85.01	ng	-0.01
79) Terphenyl-d14	12.69	244	2935388	84.36	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.30	88	300328	35.631	ng	99
3) Pyridine	2.94	79	819082	34.478	ng	98
4) n-Nitrosodimethylamine	2.89	42	342423	36.767	ng	96
6) Aniline	6.27	93	739665	24.827	ng	# 1
8) 2-Chlorophenol	6.40	128	827899	41.225	ng	99
9) Benzaldehyde	6.15	77	114649	8.107	ng	98
10) Phenol	6.27	94	948289	37.346	ng	94
11) bis(2-Chloroethyl)ether	6.35	93	751135	40.130	ng	98
12) 1,3-Dichlorobenzene	6.55	146	872245	39.136	ng	99
13) 1,4-Dichlorobenzene	6.63	146	882007	39.464	ng	99
14) 1,2-Dichlorobenzene	6.78	146	821776	39.705	ng	99
15) Benzyl Alcohol	6.76	79	472638	29.943	ng	98
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1060879	39.078	ng	96
17) 2-Methylphenol	6.88	107	720051	43.438	ng	98
18) Hexachloroethane	7.12	117	316966	39.339	ng	96
19) n-Nitroso-di-n-propylamine	7.04	70	535836	38.610	ng	99
20) 3+4-Methylphenols	7.03	107	815893	42.627	ng	92
22) Acetophenone	7.02	105	1096742	43.160	ng	# 97
24) Nitrobenzene	7.20	77	839657	42.239	ng	99
25) Isophorone	7.44	82	1547940	41.877	ng	99
26) 2-Nitrophenol	7.51	139	471171	47.995	ng	99
27) 2,4-Dimethylphenol	7.56	122	804329	50.041	ng	98
28) bis(2-Chloroethoxy)methane	7.65	93	968025	41.434	ng	99
29) 2,4-Dichlorophenol	7.75	162	724481	43.676	ng	97
30) 1,2,4-Trichlorobenzene	7.84	180	785510	42.872	ng	99
31) Naphthalene	7.91	128	2335239	42.859	ng	99
32) Benzoic acid	7.70	122	592769	51.665	ng	97
33) 4-Chloroaniline	7.97	127	564503	22.669	ng	98
34) Hexachlorobutadiene	8.04	225	422494	42.078	ng	99
35) Caprolactam	8.34	113	243329m	43.632	ng	
36) 4-Chloro-3-methylphenol	8.46	107	752633	44.803	ng	95
37) 2-Methylnaphthalene	8.61	142	1595481	43.429	ng	99
38) 1-Methylnaphthalene	8.71	142	1523420	43.418	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	676426	42.866	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.77	237	699599	74.768	ng	99
43) 2,4,6-Trichlorophenol	8.88	196	512234	42.046	ng	100
44) 2,4,5-Trichlorophenol	8.93	196	584627	46.599	ng	99
46) 1,1'-Biphenyl	9.07	154	1854155	41.497	ng	98
47) 2-Chloronaphthalene	9.09	162	1512179	41.723	ng	97
48) 2-Nitroaniline	9.18	65	456297	43.183	ng	98
49) Acenaphthylene	9.50	152	2329625	41.255	ng	98
50) Dimethylphthalate	9.38	163	1722398	41.924	ng	99
51) 2,6-Dinitrotoluene	9.43	165	416507	43.912	ng	97
52) Acenaphthene	9.68	154	1384068	39.170	ng	99
53) 3-Nitroaniline	9.59	138	314897	27.205	ng	97
54) 2,4-Dinitrophenol	9.70	184	447268	90.572	ng	96
55) Dibenzofuran	9.85	168	2034775	40.121	ng	99
56) 4-Nitrophenol	9.77	139	728070	78.460	ng	94
57) 2,4-Dinitrotoluene	9.83	165	552313	45.097	ng	95
58) Fluorene	10.19	166	1405146	41.034	ng	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	484822	46.086	ng	99
60) Diethylphthalate	10.08	149	1701390	41.198	ng	100
61) 4-Chlorophenyl-phenylether	10.18	204	683418	42.162	ng	97
62) 4-Nitroaniline	10.21	138	467680	40.276	ng	98
63) Azobenzene	10.34	77	1568880	40.672	ng	99
65) 4,6-Dinitro-2-methylphenol	10.24	198	270366	47.153	ng	93
66) n-Nitrosodiphenylamine	10.30	169	1446314	42.991	ng	99
67) 4-Bromophenyl-phenylether	10.67	248	491866	42.012	ng	97
68) Hexachlorobenzene	10.73	284	546489	43.004	ng	95
69) Atrazine	10.83	200	464952	42.963	ng	99
70) Pentachlorophenol	10.92	266	683274	84.397	ng	97
71) Phenanthrene	11.14	178	2308447	39.705	ng	99
72) Anthracene	11.19	178	2367926	40.347	ng	98
73) Carbazole	11.34	167	2171906	41.443	ng	98
74) Di-n-butylphthalate	11.69	149	2488336	41.089	ng	98
75) Fluoranthene	12.32	202	2389729	41.196	ng	97
77) Benzidine	12.44	184	809388	24.639	ng	99
78) Pyrene	12.54	202	2428234	42.711	ng	98
80) Butylbenzylphthalate	13.18	149	1143710	45.585	ng	92
81) Benzo(a)anthracene	13.72	228	2243143	42.258	ng	99
82) 3,3'-Dichlorobenzidine	13.69	252	702325	36.639	ng	99
83) Chrysene	13.76	228	2116462	43.527	ng	99
84) Bis(2-ethylhexyl)phthalate	13.74	149	1447122	44.018	ng	100
85) Di-n-octyl phthalate	14.35	149	2500119	45.262	ng	99
86) Indeno(1,2,3-cd)pyrene	16.37	276	2403183	41.768	ng	99
88) Benzo(b)fluoranthene	14.72	252	2403507	49.540	ng	98
89) Benzo(k)fluoranthene	14.75	252	2112171	48.546	ng	98
90) Benzo(a)pyrene	15.05	252	2200773	50.328	ng	98
91) Dibenzo(a,h)anthracene	16.39	278	1958320	47.970	ng	99
92) Benzo(g,h,i)perylene	16.75	276	1987796	48.109	ng	98

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

