

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091819\
 Data File : BF117137.D
 Acq On : 18 Sep 2019 10:15
 Operator : HP/JU
 Sample : PB123129BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB123129BS

Manual Integrations
 APPROVED

Jagrut
 9/19/2019 10:42:02 AM

Quant Time: Sep 18 13:32:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	72167	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	299823	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	155097	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	245575	20.00	ng	0.00
76) Chrysene-d12	13.97	240	155042	20.00	ng	0.00
87) Perylene-d12	15.42	264	115302	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	470980	101.89	ng	0.02
7) Phenol-d6	6.47	99	613746	102.74	ng	0.01
23) Nitrobenzene-d5	7.39	82	369719	64.79	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	141250	108.97	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	655760	66.11	ng	0.00
79) Terphenyl-d14	12.91	244	594345	71.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	75170	38.863	ng	98
3) Pyridine	3.40	79	173958	32.858	ng	98
4) n-Nitrosodimethylamine	3.35	42	66154	36.412	ng	100
6) Aniline	6.49	93	233727	30.389	ng	# 88
8) 2-Chlorophenol	6.62	128	206474	41.405	ng	96
9) Benzaldehyde	6.37	77	126707	43.206	ng	98
10) Phenol	6.48	94	260582	39.568	ng	91
11) bis(2-Chloroethyl)ether	6.56	93	191907	39.649	ng	99
12) 1,3-Dichlorobenzene	6.76	146	215949	38.592	ng	99
13) 1,4-Dichlorobenzene	6.84	146	224344	39.286	ng	99
14) 1,2-Dichlorobenzene	6.99	146	206475	38.835	ng	97
15) Benzyl Alcohol	6.97	79	186270	40.661	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.10	45	183470	38.367	ng	90
17) 2-Methylphenol	7.08	107	169206	40.487	ng	95
18) Hexachloroethane	7.33	117	84838	38.618	ng	97
19) n-Nitroso-di-n-propylamine	7.23	70	150012	41.239	ng	94
20) 3+4-Methylphenols	7.23	107	212182	40.759	ng	96
22) Acetophenone	7.23	105	301619	39.595	ng	98
24) Nitrobenzene	7.40	77	225578	40.030	ng	98
25) Isophorone	7.65	82	411081	40.687	ng	99
26) 2-Nitrophenol	7.72	139	105809	43.713	ng	96
27) 2,4-Dimethylphenol	7.76	122	180568	47.042	ng	96
28) bis(2-Chloroethoxy)methane	7.85	93	243396	39.101	ng	99
29) 2,4-Dichlorophenol	7.96	162	167991	41.768	ng	98
30) 1,2,4-Trichlorobenzene	8.05	180	167810	38.619	ng	98
31) Naphthalene	8.13	128	584984	40.569	ng	99
32) Benzoic acid	7.89	122	86461	32.327	ng	98
33) 4-Chloroaniline	8.17	127	156488	24.468	ng	99
34) Hexachlorobutadiene	8.25	225	95242	38.634	ng	98
35) Caprolactam	8.55	113	51977m	38.181	ng	
36) 4-Chloro-3-methylphenol	8.66	107	193284	41.215	ng	99
37) 2-Methylnaphthalene	8.82	142	400184	41.213	ng	98
38) 1-Methylnaphthalene	8.92	142	380236	41.811	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	157761	39.391	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.97	237	135929	77.098	nq	96
43) 2,4,6-Trichlorophenol	9.09	196	109171	40.167	nq	99
44) 2,4,5-Trichlorophenol	9.14	196	118698	40.876	nq	97
46) 1,1'-Biphenyl	9.27	154	473559	38.925	nq	99
47) 2-Chloronaphthalene	9.30	162	365131	38.029	nq	99
48) 2-Nitroaniline	9.40	65	118274	38.405	nq	96
49) Acenaphthylene	9.72	152	583848	40.592	nq	99
50) Dimethylphthalate	9.57	163	428446	39.015	nq	99
51) 2,6-Dinitrotoluene	9.64	165	94028	42.040	nq	94
52) Acenaphthene	9.89	154	332940	39.049	nq	99
53) 3-Nitroaniline	9.81	138	77262	27.379	nq	98
54) 2,4-Dinitrophenol	9.92	184	71040	74.297	nq	97
55) Dibenzofuran	10.06	168	507160	40.315	nq	97
56) 4-Nitrophenol	9.98	139	126628	69.236	nq	96
57) 2,4-Dinitrotoluene	10.05	165	126124	43.442	nq	97
58) Fluorene	10.40	166	409468	40.407	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.18	232	88564m	39.855	nq	
60) Diethylphthalate	10.27	149	430415	39.838	nq	98
61) 4-Chlorophenyl-phenylether	10.39	204	179030	40.833	nq	97
62) 4-Nitroaniline	10.42	138	110393	37.617	nq	97
63) Azobenzene	10.55	77	442994	40.146	nq	97
65) 4,6-Dinitro-2-methylphenol	10.46	198	50356	45.434	nq	89
66) n-Nitrosodiphenylamine	10.51	169	358651	42.288	nq	98
67) 4-Bromophenyl-phenylether	10.88	248	103421	41.238	nq	93
68) Hexachlorobenzene	10.95	284	109581	39.949	nq	94
69) Atrazine	11.03	200	98269	47.176	nq	99
70) Pentachlorophenol	11.15	266	102162	79.454	nq	99
71) Phenanthrene	11.36	178	566477	40.612	nq	100
72) Anthracene	11.42	178	587504	41.256	nq	97
73) Carbazole	11.57	167	550000	40.689	nq	99
74) Di-n-butylphthalate	11.89	149	665445	41.377	nq	99
75) Fluoranthene	12.54	202	558478	38.967	nq	99
77) Benzidine	12.66	184	207492	41.546	nq	99
78) Pyrene	12.77	202	561916	43.194	nq	100
80) Butylbenzylphthalate	13.38	149	257687	41.920	nq	98
81) Benzo(a)anthracene	13.96	228	406558	39.249	nq	99
82) 3,3'-Dichlorobenzidine	13.92	252	100231	30.027	nq	96
83) Chrysene	14.00	228	405572	40.144	nq	98
84) Bis(2-ethylhexyl)phthalate	13.94	149	343968	43.399	nq	98
85) Di-n-octyl phthalate	14.54	149	525200	42.102	nq	99
86) Indeno(1,2,3-cd)pyrene	16.86	276	294314	39.930	nq	98
88) Benzo(b)fluoranthene	15.00	252	313485	44.885	nq	99
89) Benzo(k)fluoranthene	15.02	252	303195	45.827	nq	100
90) Benzo(a)pyrene	15.36	252	287894	46.974	nq	98
91) Dibenzo(a,h)anthracene	16.86	278	248582	50.914	nq	99
92) Benzo(g,h,i)perylene	17.29	276	234316	48.161	nq	97

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

