

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121919\
 Data File : BF118276.D
 Acq On : 19 Dec 2019 13:42
 Operator : JU
 Sample : PB125579BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB125579BS

Manual Integrations
 APPROVED

Sohil
 12/20/2019 4:26:15 PM

Quant Time: Dec 20 02:50:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 06 17:34:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	114734	20.00	ng	-0.01
21) Naphthalene-d8	8.15	136	468690	20.00	ng	-0.01
39) Acenaphthene-d10	9.91	164	252749	20.00	ng	-0.01
64) Phenanthrene-d10	11.40	188	467932	20.00	ng	0.00
76) Chrysene-d12	14.06	240	311745	20.00	ng	-0.01
87) Perylene-d12	15.54	264	243167	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	808680	123.45	ng	0.00
7) Phenol-d6	6.50	99	990491	125.01	ng	0.00
23) Nitrobenzene-d5	7.43	82	638905	76.69	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	350513	114.16	ng	0.00
45) 2-Fluorobiphenyl	9.23	172	1351359	81.36	ng	-0.01
79) Terphenyl-d14	12.99	244	1552288	85.63	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	113274	41.010	ng	97
3) Pyridine	3.46	79	267580	37.548	ng	99
4) n-Nitrosodimethylamine	3.42	42	135223	44.109	ng	99
6) Aniline	6.52	93	350898	34.621	ng	96
8) 2-Chlorophenol	6.65	128	342607	46.383	ng	98
9) Benzaldehyde	6.41	77	210275	58.233	ng	97
10) Phenol	6.51	94	404459	44.761	ng	97
11) bis(2-Chloroethyl)ether	6.60	93	293552	44.918	ng	94
12) 1,3-Dichlorobenzene	6.80	146	374439	43.394	ng	99
13) 1,4-Dichlorobenzene	6.87	146	381222	43.875	ng	100
14) 1,2-Dichlorobenzene	7.03	146	358932	43.989	ng	98
15) Benzyl Alcohol	7.00	79	281666	45.540	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.13	45	332730	42.344	ng	94
17) 2-Methylphenol	7.12	107	271399	46.196	ng	98
18) Hexachloroethane	7.37	117	138365	43.217	ng	93
19) n-Nitroso-di-n-propylamine	7.27	70	220905	45.882	ng	98
20) 3+4-Methylphenols	7.27	107	344201	46.644	ng	94
22) Acetophenone	7.27	105	459006	40.764	ng	100
24) Nitrobenzene	7.45	77	335863	41.033	ng	99
25) Isophorone	7.69	82	607991	43.050	ng	100
26) 2-Nitrophenol	7.76	139	191049	43.669	ng	96
27) 2,4-Dimethylphenol	7.80	122	295801	50.689	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	377544	40.911	ng	99
29) 2,4-Dichlorophenol	8.01	162	291473	42.844	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	325139	40.896	ng	99
31) Naphthalene	8.17	128	1016526	43.185	ng	99
32) Benzoic acid	7.94	122	199773	37.915	ng	99
33) 4-Chloroaniline	8.22	127	223299	21.970	ng	99
34) Hexachlorobutadiene	8.28	225	186953	39.215	ng	99
35) Caprolactam	8.59	113	89530m	42.675	ng	
36) 4-Chloro-3-methylphenol	8.71	107	307439	42.943	ng	97
37) 2-Methylnaphthalene	8.86	142	697539	43.772	ng	99
38) 1-Methylnaphthalene	8.96	142	647721	43.289	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	306603	40.044	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.02	237	297626	86.757	ng	98
43) 2,4,6-Trichlorophenol	9.15	196	208188	40.724	ng	98
44) 2,4,5-Trichlorophenol	9.19	196	224167	42.324	ng	98
46) 1,1'-Biphenyl	9.33	154	823938	41.333	ng	99
47) 2-Chloronaphthalene	9.36	162	643293	40.118	ng	98
48) 2-Nitroaniline	9.45	65	175713	38.426	ng	95
49) Acenaphthylene	9.77	152	1022915	43.251	ng	100
50) Dimethylphthalate	9.63	163	762946	40.867	ng	100
51) 2,6-Dinitrotoluene	9.70	165	171493	42.245	ng	94
52) Acenaphthene	9.95	154	588447	40.764	ng	99
53) 3-Nitroaniline	9.87	138	117765	24.650	ng	95
54) 2,4-Dinitrophenol	9.98	184	156967	77.720	ng	96
55) Dibenzofuran	10.12	168	899275	42.515	ng	100
56) 4-Nitrophenol	10.04	139	251374	75.819	ng	97
57) 2,4-Dinitrotoluene	10.10	165	227024	41.899	ng	97
58) Fluorene	10.46	166	705885	41.993	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.24	232	186682	42.722	ng	99
60) Diethylphthalate	10.33	149	765832	41.635	ng	98
61) 4-Chlorophenyl-phenylether	10.45	204	351620	41.574	ng	98
62) 4-Nitroaniline	10.49	138	175652	35.253	ng	99
63) Azobenzene	10.62	77	661890	42.153	ng	94
65) 4,6-Dinitro-2-methylphenol	10.52	198	114393	44.267	ng	95
66) n-Nitrosodiphenylamine	10.57	169	625367	42.796	ng	99
67) 4-Bromophenyl-phenylether	10.95	248	218188	40.847	ng	98
68) Hexachlorobenzene	11.02	284	253230	40.278	ng	# 93
69) Atrazine	11.10	200	214019	65.182	ng	97
70) Pentachlorophenol	11.21	266	244331	82.638	ng	97
71) Phenanthrene	11.43	178	1041771	41.881	ng	100
72) Anthracene	11.48	178	1093891	44.083	ng	99
73) Carbazole	11.64	167	915984	41.142	ng	99
74) Di-n-butylphthalate	11.96	149	1202011	43.139	ng	99
75) Fluoranthene	12.62	202	1055013	41.484	ng	97
77) Benzidine	12.75	184	389342	47.442	ng	98
78) Pyrene	12.85	202	1070219	43.563	ng	99
80) Butylbenzylphthalate	13.46	149	475807	42.766	ng	99
81) Benzo(a)anthracene	14.04	228	879216	40.787	ng	99
82) 3,3'-Dichlorobenzidine	14.00	252	243951	34.459	ng	99
83) Chrysene	14.09	228	852503	41.402	ng	99
84) Bis(2-ethylhexyl)phthalate	14.02	149	647116	44.109	ng	100
85) Di-n-octyl phthalate	14.64	149	976543	43.942	ng	100
86) Indeno(1,2,3-cd)pyrene	17.06	276	678761	39.177	ng	99
88) Benzo(b)fluoranthene	15.11	252	706646	43.939	ng	99
89) Benzo(k)fluoranthene	15.14	252	720588m	46.641	ng	
90) Benzo(a)pyrene	15.49	252	625516	44.039	ng	98
91) Dibenzo(a,h)anthracene	17.07	278	575587	47.246	ng	99
92) Benzo(g,h,i)perylene	17.51	276	572346	48.890	ng	99

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

