

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122819\
 Data File : BF118358.D
 Acq On : 28 Dec 2019 11:50
 Operator : JU
 Sample : PB125767BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB125767BS

Manual Integrations
 APPROVED

mohammad
 12/30/2019 2:08:29 PM

Quant Time: Dec 29 07:19:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 15:58:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	179424	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	782608	20.00	ng	0.00
39) Acenaphthene-d10	9.91	164	431713	20.00	ng	0.01
64) Phenanthrene-d10	11.40	188	810146	20.00	ng	0.01
76) Chrysene-d12	14.06	240	543903	20.00	ng	0.00
87) Perylene-d12	15.54	264	387881	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.49	112	877607	83.49	ng	0.02
7) Phenol-d6	6.50	99	1066410	81.78	ng	0.01
23) Nitrobenzene-d5	7.43	82	1124834	82.21	ng	0.00
42) 2,4,6-Tribromophenol	10.70	330	387845	72.75	ng	0.01
45) 2-Fluorobiphenyl	9.22	172	2327792	82.31	ng	0.00
79) Terphenyl-d14	12.99	244	2730111	83.27	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.70	88	154052	36.868	ng	98
3) Pyridine	3.46	79	332826	29.896	ng	98
4) n-Nitrosodimethylamine	3.42	42	187095	40.648	ng	95
6) Aniline	6.52	93	545874	32.968	ng	# 57
8) 2-Chlorophenol	6.65	128	574580	48.265	ng	98
9) Benzaldehyde	6.40	77	248447	32.499	ng	96
10) Phenol	6.51	94	661098	45.109	ng	86
11) bis(2-Chloroethyl)ether	6.59	93	484710	46.245	ng	99
12) 1,3-Dichlorobenzene	6.80	146	557597	40.670	ng	97
13) 1,4-Dichlorobenzene	6.87	146	577761	41.488	ng	99
14) 1,2-Dichlorobenzene	7.03	146	546659	41.578	ng	98
15) Benzyl Alcohol	7.00	79	469968	47.337	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.13	45	494202	44.187	ng	74
17) 2-Methylphenol	7.12	107	454721	47.624	ng	96
18) Hexachloroethane	7.36	117	211219	41.326	ng	95
19) n-Nitroso-di-n-propylamine	7.27	70	366017	45.566	ng	98
20) 3+4-Methylphenols	7.27	107	578559	47.344	ng	# 84
22) Acetophenone	7.27	105	751716	39.191	ng	# 94
24) Nitrobenzene	7.45	77	548210	39.995	ng	99
25) Isophorone	7.69	82	1015685	42.592	ng	100
26) 2-Nitrophenol	7.76	139	324093	44.640	ng	97
27) 2,4-Dimethylphenol	7.80	122	501310	50.609	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	635854	40.771	ng	99
29) 2,4-Dichlorophenol	8.00	162	482895	42.354	ng	99
30) 1,2,4-Trichlorobenzene	8.09	180	522645	39.352	ng	98
31) Naphthalene	8.16	128	1643657	41.322	ng	99
32) Benzoic acid	7.96	122	327133	40.160	ng	96
33) 4-Chloroaniline	8.22	127	424588	24.899	ng	98
34) Hexachlorobutadiene	8.28	225	298106	37.707	ng	99
35) Caprolactam	8.61	113	152449m	42.829	ng	
36) 4-Chloro-3-methylphenol	8.71	107	520322	42.507	ng	98
37) 2-Methylnaphthalene	8.86	142	1156253	42.456	ng	99
38) 1-Methylnaphthalene	8.96	142	1078111	42.018	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.03	216	503038	38.781	ng	99

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41) Hexachlorocyclopentadiene	9.01	237	358523	67.260	nq	98
43) 2,4,6-Trichlorophenol	9.15	196	350690	40.469	nq	99
44) 2,4,5-Trichlorophenol	9.19	196	362329	40.646	nq	98
46) 1,1'-Biphenyl	9.33	154	1356956	39.740	nq	100
47) 2-Chloronaphthalene	9.35	162	1063082	38.683	nq	99
48) 2-Nitroaniline	9.45	65	297483	38.323	nq	98
49) Acenaphthylene	9.77	152	1643451	40.322	nq	99
50) Dimethylphthalate	9.63	163	1291157	39.843	nq	99
51) 2,6-Dinitrotoluene	9.70	165	283530	40.583	nq	95
52) Acenaphthene	9.95	154	982825	39.588	nq	99
53) 3-Nitroaniline	9.87	138	212527	25.918	nq	90
54) 2,4-Dinitrophenol	9.99	184	273978	77.921	nq	99
55) Dibenzofuran	10.12	168	1456698	39.710	nq	97
56) 4-Nitrophenol	10.05	139	408123	80.621	nq	93
57) 2,4-Dinitrotoluene	10.11	165	375875	40.263	nq	# 87
58) Fluorene	10.46	166	1171034	39.545	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.24	232	315051	41.595	nq	98
60) Diethylphthalate	10.33	149	1283633	40.039	nq	99
61) 4-Chlorophenyl-phenylether	10.45	204	592243	40.063	nq	93
62) 4-Nitroaniline	10.49	138	294652	34.589	nq	98
63) Azobenzene	10.61	77	1104698	40.229	nq	98
65) 4,6-Dinitro-2-methylphenol	10.52	198	202329	46.295	nq	85
66) n-Nitrosodiphenylamine	10.57	169	1057345	40.722	nq	98
67) 4-Bromophenyl-phenylether	10.94	248	373533	39.398	nq	98
68) Hexachlorobenzene	11.02	284	426587	38.225	nq	91
69) Atrazine	11.10	200	356404	44.475	nq	96
70) Pentachlorophenol	11.22	266	395092	81.745	nq	99
71) Phenanthrene	11.43	178	1730732	39.188	nq	100
72) Anthracene	11.48	178	1765599	39.940	nq	100
73) Carbazole	11.64	167	1535368	39.142	nq	99
74) Di-n-butylphthalate	11.96	149	2024456	40.703	nq	99
75) Fluoranthene	12.62	202	1757951	39.441	nq	96
77) Benzdine	12.74	184	528463	35.459	nq	99
78) Pyrene	12.85	202	1757914	39.579	nq	98
80) Butylbenzylphthalate	13.46	149	836532	41.830	nq	99
81) Benzo(a)anthracene	14.04	228	1482766	38.947	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	443668	34.371	nq	98
83) Chrysene	14.09	228	1385920	38.018	nq	99
84) Bis(2-ethylhexyl)phthalate	14.02	149	1208412	45.953	nq	99
85) Di-n-octyl phthalate	14.63	149	1834538	47.427	nq	100
86) Indeno(1,2,3-cd)pyrene	17.07	276	1123155	36.159	nq	99
88) Benzo(b)fluoranthene	15.11	252	1295151	50.302	nq	99
89) Benzo(k)fluoranthene	15.14	252	1081365	43.741	nq	99
90) Benzo(a)pyrene	15.49	252	1018661	44.912	nq	98
91) Dibenzo(a,h)anthracene	17.08	278	959523	48.750	nq	98
92) Benzo(g,h,i)perylene	17.53	276	927932	48.804	nq	99

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

