

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG042920\
 Data File : BG045202.D
 Acq On : 29 Apr 2020 13:00
 Operator : CG/JU
 Sample : PB128519BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB128519BS

Manual Integrations
 APPROVED

mohammad
 4/30/2020 1:37:38 PM

Quant Time: Apr 29 14:14:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 29 12:24:27 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	58938	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	235071	20.00	ng	0.00
39) Acenaphthene-d10	14.63	164	167617	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	436173	20.00	ng	0.00
76) Chrysene-d12	21.65	240	439595	20.00	ng	0.00
87) Perylene-d12	24.82	264	491461	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.56	112	312458	87.52	ng	0.00
7) Phenol-d6	7.16	99	434566	81.93	ng	0.00
23) Nitrobenzene-d5	9.15	82	390552	75.81	ng	0.00
42) 2,4,6-Tribromophenol	16.12	330	218068	84.21	ng	0.00
45) 2-Fluorobiphenyl	13.25	172	927185	81.11	ng	0.00
79) Terphenyl-d14	19.99	244	1800886	89.29	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.45	88	57718	36.380	ng	98
3) Pyridine	3.84	79	159070	32.564	ng	96
4) n-Nitrosodimethylamine	3.76	42	62121	41.513	ng	89
6) Aniline	7.31	93	205889	29.855	ng	99
8) 2-Chlorophenol	7.56	128	173587	45.243	ng	99
9) Benzaldehyde	7.12	77	58050	14.492	ng	99
10) Phenol	7.18	94	234348	44.153	ng	96
11) bis(2-Chloroethyl)ether	7.41	93	160152	37.898	ng	98
12) 1,3-Dichlorobenzene	7.88	146	185898	41.118	ng	97
13) 1,4-Dichlorobenzene	8.02	146	188192	41.774	ng	96
14) 1,2-Dichlorobenzene	8.35	146	172357	39.991	ng	96
15) Benzyl Alcohol	8.23	79	167220	37.603	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.52	45	143985	34.710	ng	96
17) 2-Methylphenol	8.44	107	156684	38.261	ng	96
18) Hexachloroethane	9.07	117	70806	41.809	ng	94
19) n-Nitroso-di-n-propylamine	8.79	70	137762	36.726	ng	98
20) 3+4-Methylphenols	8.76	107	210129	36.659	ng	97
22) Acetophenone	8.81	105	254608	40.052	ng	97
24) Nitrobenzene	9.19	77	211818	40.111	ng	97
25) Isophorone	9.71	82	390019	36.968	ng	99
26) 2-Nitrophenol	9.90	139	95006	41.767	ng	97
27) 2,4-Dimethylphenol	9.97	122	157986	43.772	ng	95
28) bis(2-Chloroethoxy)methane	10.20	93	218464	39.411	ng	98
29) 2,4-Dichlorophenol	10.45	162	180901	43.407	ng	94
30) 1,2,4-Trichlorobenzene	10.66	180	190631	40.400	ng	99
31) Naphthalene	10.85	128	522486	40.630	ng	99
32) Benzoic acid	10.11	122	109502	38.943	ng	96
33) 4-Chloroaniline	10.95	127	127484	21.257	ng	98
34) Hexachlorobutadiene	11.15	225	135420	41.859	ng	96
35) Caprolactam	11.71	113	62301m	37.561	ng	
36) 4-Chloro-3-methylphenol	12.08	107	203352	41.536	ng	97
37) 2-Methylnaphthalene	12.46	142	411079	41.185	ng	99
38) 1-Methylnaphthalene	12.68	142	383693m	40.720	ng	
40) 1,2,4,5-Tetrachlorobenzene	12.83	216	223810	41.471	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.82	237	320278	94.951	nq	99
43) 2,4,6-Trichlorophenol	13.06	196	165760	43.519	nq	98
44) 2,4,5-Trichlorophenol	13.14	196	173563	40.033	nq	97
46) 1,1'-Biphenyl	13.46	154	514266	40.366	nq	98
47) 2-Chloronaphthalene	13.50	162	399852	41.075	nq	99
48) 2-Nitroaniline	13.70	65	131965	41.202	nq	99
49) Acenaphthylene	14.36	152	658086	41.503	nq	98
50) Dimethylphthalate	14.09	163	551089	40.378	nq	98
51) 2,6-Dinitrotoluene	14.20	165	126106	41.310	nq	96
52) Acenaphthene	14.70	154	512123m	47.735	nq	
53) 3-Nitroaniline	14.53	138	91720	27.395	nq	95
54) 2,4-Dinitrophenol	14.74	184	163286	89.954	nq	95
55) Dibenzofuran	15.03	168	655461	41.906	nq	99
56) 4-Nitrophenol	14.85	139	218912	84.327	nq	98
57) 2,4-Dinitrotoluene	14.99	165	187540	42.038	nq	# 93
58) Fluorene	15.68	166	550006	43.050	nq	100
59) 2,3,4,6-Tetrachlorophenol	15.26	232	175194	45.415	nq	98
60) Diethylphthalate	15.45	149	584539	41.079	nq	99
61) 4-Chlorophenyl-phenylether	15.67	204	304197	41.367	nq	99
62) 4-Nitroaniline	15.69	138	137350	39.552	nq	93
63) Azobenzene	15.97	77	554148	42.240	nq	98
65) 4,6-Dinitro-2-methylphenol	15.75	198	124946	43.581	nq	96
66) n-Nitrosodiphenylamine	15.88	169	492060	39.264	nq	99
67) 4-Bromophenyl-phenylether	16.57	248	210234	37.362	nq	95
68) Hexachlorobenzene	16.69	284	231869	39.732	nq	98
69) Atrazine	16.83	200	212001	46.545	nq	99
70) Pentachlorophenol	17.03	266	301281	84.155	nq	99
71) Phenanthrene	17.42	178	937326	41.819	nq	98
72) Anthracene	17.52	178	968385	43.071	nq	99
73) Carbazole	17.78	167	918974	40.277	nq	98
74) Di-n-butylphthalate	18.34	149	1116044	44.925	nq	100
75) Fluoranthene	19.43	202	1254665	48.042	nq	98
77) Benzidine	19.61	184	681922	54.863	nq	99
78) Pyrene	19.80	202	1233728	40.709	nq	99
80) Butylbenzylphthalate	20.68	149	518081	41.241	nq	99
81) Benzo(a)anthracene	21.62	228	1189752	41.757	nq	99
82) 3,3'-Dichlorobenzidine	21.54	252	360283	31.770	nq	# 98
83) Chrysene	21.69	228	1159876	42.369	nq	99
84) Bis(2-ethylhexyl)phthalate	21.54	149	714483	41.354	nq	99
85) Di-n-octyl phthalate	22.74	149	1248844	41.800	nq	100
86) Indeno(1,2,3-cd)pyrene	28.42	276	1549529	42.633	nq	# 96
88) Benzo(b)fluoranthene	23.81	252	1308166	42.494	nq	99
89) Benzo(k)fluoranthene	23.88	252	1261489	43.089	nq	99
90) Benzo(a)pyrene	24.67	252	1209843	41.624	nq	98
91) Dibenzo(a,h)anthracene	28.49	278	1259007	42.987	nq	98
92) Benzo(g,h,i)perylene	29.55	276	1264811	43.075	nq	99

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

