

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG121718\
 Data File : BG038681.D
 Acq On : 18 Dec 2018 3:36
 Operator : JU/SJ
 Sample : J6403-08MSD
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-1-BMSD

Manual Integrations
 APPROVED

Sohil
 12/18/2018 5:37:14 PM

Quant Time: Dec 18 05:08:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	20674	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	83619	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	53821	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	137905	20.00	ng	0.00
76) Chrysene-d12	21.73	240	137726	20.00	ng	0.00
87) Perylene-d12	25.03	264	150488	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	170158	145.98	ng	0.01
7) Phenol-d6	7.22	99	203178	126.97	ng	0.01
23) Nitrobenzene-d5	9.24	82	166056	101.45	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	121019	163.15	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	383728	97.81	ng	0.00
79) Terphenyl-d14	20.04	244	718989	113.61	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.50	88	21069	38.838	ng	# 1
3) Pyridine	3.92	79	52428	35.102	ng	88
4) n-Nitrosodimethylamine	3.83	42	39007	53.300	ng	# 84
6) Aniline	7.39	93	28555	13.979	ng	99
8) 2-Chlorophenol	7.62	128	66488	49.291	ng	98
9) Benzaldehyde	7.20	77	19015	18.318	ng	93
10) Phenol	7.24	94	70008	41.181	ng	92
11) bis(2-Chloroethyl)ether	7.48	93	60937	45.022	ng	95
12) 1,3-Dichlorobenzene	7.95	146	68580	43.000	ng	98
13) 1,4-Dichlorobenzene	8.09	146	69481	42.537	ng	97
14) 1,2-Dichlorobenzene	8.42	146	66633	43.270	ng	96
15) Benzyl Alcohol	8.30	79	62701	50.201	ng	94
16) 2,2'-oxybis(1-Chloropropan	8.58	45	138900	44.799	ng	99
17) 2-Methylphenol	8.50	107	57748	50.701	ng	98
18) Hexachloroethane	9.14	117	25095	42.044	ng	91
19) n-Nitroso-di-n-propylamine	8.86	70	50255	44.716	ng	94
20) 3+4-Methylphenols	8.83	107	77023	48.966	ng	99
22) Acetophenone	8.89	105	101465	48.580	ng	# 99
24) Nitrobenzene	9.28	77	81115	48.988	ng	97
25) Isophorone	9.79	82	145801	49.578	ng	99
26) 2-Nitrophenol	9.99	139	38743	45.715	ng	99
27) 2,4-Dimethylphenol	10.04	122	66765	54.969	ng	99
28) bis(2-Chloroethoxy)methane	10.27	93	80491	48.500	ng	97
29) 2,4-Dichlorophenol	10.52	162	73270	54.226	ng	97
30) 1,2,4-Trichlorobenzene	10.73	180	74189	45.455	ng	97
31) Naphthalene	10.93	128	204262	49.578	ng	98
32) Benzoic acid	10.19	122	26015m	36.745	ng	
33) 4-Chloroaniline	11.06	127	4897	2.738	ng	# 86
34) Hexachlorobutadiene	11.19	225	47656	43.152	ng	94
35) Caprolactam	11.83	113	17354m	31.034	ng	
36) 4-Chloro-3-methylphenol	12.15	107	75689	49.086	ng	95
37) 2-Methylnaphthalene	12.53	142	154075	52.420	ng	97
38) 1-Methylnaphthalene	12.75	142	147261	51.631	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	92684	46.070	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.86	237	93867	84.926	ng	97
43) 2,4,6-Trichlorophenol	13.13	196	65381	52.855	ng	97
44) 2,4,5-Trichlorophenol	13.21	196	69415	53.001	ng	99
46) 1,1'-Biphenyl	13.53	154	203775	45.164	ng	99
47) 2-Chloronaphthalene	13.58	162	164546	47.212	ng	96
48) 2-Nitroaniline	13.79	65	56844	51.205	ng	95
49) Acenaphthylene	14.42	152	271616	52.131	ng	100
50) Dimethylphthalate	14.15	163	225019	51.196	ng	99
51) 2,6-Dinitrotoluene	14.28	165	50290	50.491	ng	96
52) Acenaphthene	14.76	154	154181	49.487	ng	94
53) 3-Nitroaniline	14.62	138	13109	12.911	ng	# 90
54) 2,4-Dinitrophenol	14.82	184	34958	60.348	ng	89
55) Dibenzofuran	15.10	168	261923	49.044	ng	99
56) 4-Nitrophenol	14.91	139	63112	81.936	ng	93
57) 2,4-Dinitrotoluene	15.07	165	72440	51.637	ng	94
58) Fluorene	15.74	166	212666	53.748	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.32	232	64682	54.894	ng	99
60) Diethylphthalate	15.50	149	230607	47.794	ng	99
61) 4-Chlorophenyl-phenylether	15.73	204	118010	47.802	ng	97
62) 4-Nitroaniline	15.78	138	49815	47.656	ng	97
63) Azobenzene	16.03	77	200131	49.651	ng	97
65) 4,6-Dinitro-2-methylphenol	15.83	198	34064	34.627	ng	95
66) n-Nitrosodiphenylamine	15.95	169	197233	48.676	ng	97
67) 4-Bromophenyl-phenylether	16.62	248	78872	46.830	ng	94
68) Hexachlorobenzene	16.74	284	81648	46.766	ng	98
69) Atrazine	16.89	200	81588	54.257	ng	96
70) Pentachlorophenol	17.09	266	74086	71.743	ng	93
71) Phenanthrene	17.49	178	371290	51.919	ng	99
72) Anthracene	17.58	178	377431	53.462	ng	99
73) Carbazole	17.85	167	340274	48.736	ng	99
74) Di-n-butylphthalate	18.38	149	403541	50.370	ng	100
75) Fluoranthene	19.50	202	466714	52.747	ng	97
77) Benzidine	19.68	184	75379	18.506	ng	96
78) Pyrene	19.86	202	466608	51.284	ng	99
80) Butylbenzylphthalate	20.73	149	180616	50.810	ng	99
81) Benzo(a)anthracene	21.71	228	447084	53.273	ng	100
82) 3,3'-Dichlorobenzidine	21.62	252	61418	18.453	ng	98
83) Chrysene	21.77	228	414686	51.301	ng	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	250611	49.709	ng	99
85) Di-n-octyl phthalate	22.80	149	426788	50.547	ng	98
86) Indeno(1,2,3-cd)pyrene	28.81	276	514308	54.118	ng	# 71
88) Benzo(b)fluoranthene	23.97	252	442631	53.572	ng	99
89) Benzo(k)fluoranthene	24.04	252	434777	52.844	ng	96
90) Benzo(a)pyrene	24.87	252	429180	53.842	ng	98
91) Dibenzo(a,h)anthracene	28.87	278	427327	53.843	ng	99
92) Benzo(g,h,i)perylene	30.00	276	422249	53.986	ng	97

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

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