

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038580.D
 Acq On : 14 Dec 2018 23:04
 Operator : JU/SJ
 Sample : PB115428BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115428BS

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:11:52 PM

Quant Time: Dec 15 00:03:27 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	23415	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	93508	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	60008	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	153680	20.00	ng	0.00
76) Chrysene-d12	21.72	240	162167	20.00	ng	0.00
87) Perylene-d12	25.02	264	175577	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	175688	133.08	ng	0.00
7) Phenol-d6	7.21	99	228459	126.06	ng	0.00
23) Nitrobenzene-d5	9.24	82	115667	63.19	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	103129	124.70	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	266195	60.86	ng	0.00
79) Terphenyl-d14	20.04	244	552430	74.14	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	20617	33.555	ng	# 89
3) Pyridine	3.89	79	57939	34.250	ng	90
4) n-Nitrosodimethylamine	3.82	42	37436	45.165	ng	# 88
6) Aniline	7.39	93	79158	34.215	ng	98
8) 2-Chlorophenol	7.62	128	66280	43.385	ng	98
9) Benzaldehyde	7.20	77	26706	22.715	ng	99
10) Phenol	7.24	94	82254	42.720	ng	99
11) bis(2-Chloroethyl)ether	7.48	93	57327	37.397	ng	100
12) 1,3-Dichlorobenzene	7.94	146	67147	37.173	ng	94
13) 1,4-Dichlorobenzene	8.09	146	69756	37.706	ng	96
14) 1,2-Dichlorobenzene	8.42	146	67028	38.432	ng	97
15) Benzyl Alcohol	8.30	79	57973	40.982	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.58	45	131818	37.538	ng	99
17) 2-Methylphenol	8.50	107	56685	43.942	ng	97
18) Hexachloroethane	9.14	117	25030	37.026	ng	97
19) n-Nitroso-di-n-propylamine	8.86	70	46947	36.883	ng	97
20) 3+4-Methylphenols	8.83	107	75209	42.216	ng	95
22) Acetophenone	8.89	105	94266	40.360	ng	# 97
24) Nitrobenzene	9.28	77	75875	40.977	ng	98
25) Isophorone	9.79	82	134584	40.924	ng	99
26) 2-Nitrophenol	9.99	139	36337	38.342	ng	96
27) 2,4-Dimethylphenol	10.03	122	63547	46.787	ng	99
28) bis(2-Chloroethoxy)methane	10.28	93	77782	41.911	ng	95
29) 2,4-Dichlorophenol	10.52	162	66461	43.985	ng	98
30) 1,2,4-Trichlorobenzene	10.74	180	71592	39.225	ng	96
31) Naphthalene	10.93	128	197981	42.972	ng	98
32) Benzoic acid	10.18	122	28811m	36.488	ng	
33) 4-Chloroaniline	11.05	127	41934	20.964	ng	98
34) Hexachlorobutadiene	11.19	225	46521	37.669	ng	95
35) Caprolactam	11.82	113	22915m	36.645	ng	
36) 4-Chloro-3-methylphenol	12.15	107	70163	40.691	ng	94
37) 2-Methylnaphthalene	12.53	142	145971	44.411	ng	99
38) 1-Methylnaphthalene	12.75	142	137478	43.103	ng	98
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	84387	37.621	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121418\
 Data File : BG038580.D
 Acq On : 14 Dec 2018 23:04
 Operator : JU/SJ
 Sample : PB115428BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115428BS

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:11:52 PM

Quant Time: Dec 15 00:03:27 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)
41) Hexachlorocyclopentadiene	12.86	237	113978	92.489	ng	93
43) 2,4,6-Trichlorophenol	13.13	196	58516	42.428	ng	97
44) 2,4,5-Trichlorophenol	13.20	196	62639	42.896	ng	94
46) 1,1'-Biphenyl	13.53	154	186164	37.006	ng	99
47) 2-Chloronaphthalene	13.58	162	151521	38.993	ng	95
48) 2-Nitroaniline	13.79	65	50696	40.958	ng	95
49) Acenaphthylene	14.42	152	245557	42.270	ng	99
50) Dimethylphthalate	14.14	163	198138	40.433	ng	99
51) 2,6-Dinitrotoluene	14.28	165	44418	39.997	ng	97
52) Acenaphthene	14.76	154	140116	40.336	ng	98
53) 3-Nitroaniline	14.61	138	31815	28.104	ng	89
54) 2,4-Dinitrophenol	14.82	184	49483	75.218	ng	98
55) Dibenzofuran	15.10	168	235021	39.470	ng	96
56) 4-Nitrophenol	14.91	139	73032	85.039	ng	94
57) 2,4-Dinitrotoluene	15.07	165	64592	41.296	ng	96
58) Fluorene	15.74	166	188208	42.662	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.31	232	57189	43.531	ng	98
60) Diethylphthalate	15.50	149	206145	38.320	ng	98
61) 4-Chlorophenyl-phenylether	15.73	204	107810	39.168	ng	98
62) 4-Nitroaniline	15.78	138	48930	41.983	ng	97
63) Azobenzene	16.02	77	180845	40.241	ng	97
65) 4,6-Dinitro-2-methylphenol	15.82	198	40019	36.504	ng	96
66) n-Nitrosodiphenylamine	15.95	169	173762	38.482	ng	98
67) 4-Bromophenyl-phenylether	16.62	248	69648	37.109	ng	96
68) Hexachlorobenzene	16.74	284	73153	37.599	ng	99
69) Atrazine	16.89	200	74101	44.220	ng	96
70) Pentachlorophenol	17.09	266	82224	71.450	ng	98
71) Phenanthrene	17.49	178	333917	41.900	ng	99
72) Anthracene	17.58	178	342643	43.552	ng	99
73) Carbazole	17.85	167	315963	40.609	ng	98
74) Di-n-butylphthalate	18.38	149	384596	43.078	ng	99
75) Fluoranthene	19.49	202	438175	44.439	ng	98
77) Benzidine	19.68	184	217310	45.311	ng	99
78) Pyrene	19.86	202	432078	40.331	ng	99
80) Butylbenzylphthalate	20.72	149	172935	41.317	ng	99
81) Benzo(a)anthracene	21.70	228	426749	43.186	ng	99
82) 3,3'-Dichlorobenzidine	21.62	252	110189	28.116	ng	98
83) Chrysene	21.77	228	390523	41.030	ng	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	244747	41.229	ng	99
85) Di-n-octyl phthalate	22.79	149	417908	42.036	ng	99
86) Indeno(1,2,3-cd)pyrene	28.80	276	483169	43.179	ng	# 86
88) Benzo(b)fluoranthene	23.96	252	420307	43.601	ng	99
89) Benzo(k)fluoranthene	24.03	252	409721	42.682	ng	98
90) Benzo(a)pyrene	24.86	252	408301	43.903	ng	99
91) Dibenzo(a,h)anthracene	28.86	278	401786	43.390	ng	98
92) Benzo(g,h,i)perylene	30.00	276	395522	43.342	ng	98

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

