

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121918\
 Data File : BG038732.D
 Acq On : 19 Dec 2018 18:25
 Operator : JU/SJ
 Sample : PB115792BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB115792BS

Manual Integrations
 APPROVED

Sohil
 12/20/2018 1:06:38 PM

Quant Time: Dec 20 00:22:46 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	23833	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	93979	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	63182	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	160138	20.00	ng	0.00
76) Chrysene-d12	21.72	240	161639	20.00	ng	0.00
87) Perylene-d12	25.02	264	172747	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	188531	140.30	ng	0.00
7) Phenol-d6	7.21	99	256644	139.13	ng	0.00
23) Nitrobenzene-d5	9.23	82	124562	67.71	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	128136	147.15	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	317849	69.01	ng	0.00
79) Terphenyl-d14	20.04	244	633538	85.30	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.49	88	21937	35.078	ng	95
3) Pyridine	3.89	79	59644	34.640	ng	99
4) n-Nitrosodimethylamine	3.82	42	36186	42.892	ng	# 90
6) Aniline	7.39	93	71178	30.226	ng	96
8) 2-Chlorophenol	7.62	128	67178	43.201	ng	96
9) Benzaldehyde	7.20	77	21026	17.570	ng	95
10) Phenol	7.24	94	84373	43.052	ng	97
11) bis(2-Chloroethyl)ether	7.48	93	55049	35.281	ng	95
12) 1,3-Dichlorobenzene	7.94	146	70590	38.393	ng	99
13) 1,4-Dichlorobenzene	8.09	146	71396	37.915	ng	98
14) 1,2-Dichlorobenzene	8.41	146	69252	39.010	ng	98
15) Benzyl Alcohol	8.30	79	61276	42.558	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.57	45	119762	33.507	ng	97
17) 2-Methylphenol	8.50	107	58117	44.262	ng	98
18) Hexachloroethane	9.13	117	25629	37.247	ng	90
19) n-Nitroso-di-n-propylamine	8.86	70	46875	36.180	ng	98
20) 3+4-Methylphenols	8.82	107	79488	43.835	ng	98
22) Acetophenone	8.89	105	95678	40.759	ng	# 96
24) Nitrobenzene	9.27	77	74188	39.865	ng	98
25) Isophorone	9.79	82	132827	40.187	ng	99
26) 2-Nitrophenol	9.98	139	39015	40.961	ng	94
27) 2,4-Dimethylphenol	10.03	122	66936	49.035	ng	94
28) bis(2-Chloroethoxy)methane	10.27	93	75925	40.705	ng	95
29) 2,4-Dichlorophenol	10.52	162	72911	48.011	ng	98
30) 1,2,4-Trichlorobenzene	10.73	180	76609	41.764	ng	97
31) Naphthalene	10.93	128	203095	43.861	ng	98
32) Benzoic acid	10.18	122	28060m	35.673	ng	
33) 4-Chloroaniline	11.04	127	49038	24.393	ng	96
34) Hexachlorobutadiene	11.18	225	51145	41.206	ng	97
35) Caprolactam	11.82	113	24188	38.487	ng	94
36) 4-Chloro-3-methylphenol	12.15	107	75101	43.336	ng	94
37) 2-Methylnaphthalene	12.52	142	154329	46.718	ng	100
38) 1-Methylnaphthalene	12.74	142	144169	44.975	ng	93
40) 1,2,4,5-Tetrachlorobenzene	12.89	216	95621	40.488	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.85	237	110106	84.859	ng	94
43) 2,4,6-Trichlorophenol	13.13	196	64547	44.450	ng	94
44) 2,4,5-Trichlorophenol	13.20	196	71753	46.669	ng	97
46) 1,1'-Biphenyl	13.53	154	204822	38.670	ng	98
47) 2-Chloronaphthalene	13.57	162	165575	40.469	ng	98
48) 2-Nitroaniline	13.79	65	52781	40.501	ng	92
49) Acenaphthylene	14.42	152	271602	44.405	ng	98
50) Dimethylphthalate	14.14	163	214352	41.544	ng	98
51) 2,6-Dinitrotoluene	14.28	165	48979	41.889	ng	96
52) Acenaphthene	14.76	154	152965	41.823	ng	99
53) 3-Nitroaniline	14.61	138	36197	30.368	ng	97
54) 2,4-Dinitrophenol	14.82	184	43061	63.067	ng	98
55) Dibenzofuran	15.09	168	261977	41.786	ng	97
56) 4-Nitrophenol	14.91	139	73024	80.759	ng	96
57) 2,4-Dinitrotoluene	15.06	165	70244	42.653	ng	95
58) Fluorene	15.74	166	210281	45.271	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.32	232	62763	45.373	ng	100
60) Diethylphthalate	15.50	149	221125	39.039	ng	99
61) 4-Chlorophenyl-phenylether	15.73	204	118564	40.911	ng	100
62) 4-Nitroaniline	15.78	138	51284	41.793	ng	90
63) Azobenzene	16.02	77	187655	39.658	ng	96
65) 4,6-Dinitro-2-methylphenol	15.82	198	42546	37.244	ng	92
66) n-Nitrosodiphenylamine	15.95	169	194499	41.337	ng	98
67) 4-Bromophenyl-phenylether	16.62	248	80700	41.263	ng	99
68) Hexachlorobenzene	16.74	284	84546	41.703	ng	96
69) Atrazine	16.89	200	79899	45.757	ng	96
70) Pentachlorophenol	17.09	266	84023	70.069	ng	99
71) Phenanthrene	17.49	178	366478	44.132	ng	99
72) Anthracene	17.58	178	375703	45.829	ng	99
73) Carbazole	17.85	167	330625	40.779	ng	98
74) Di-n-butylphthalate	18.38	149	382058	41.068	ng	99
75) Fluoranthene	19.49	202	457566	44.534	ng	97
77) Benzidine	19.67	184	188784	39.492	ng	99
78) Pyrene	19.86	202	460544	43.129	ng	100
80) Butylbenzylphthalate	20.72	149	174013	41.711	ng	99
81) Benzo(a)anthracene	21.70	228	445056	45.186	ng	100
82) 3,3'-Dichlorobenzidine	21.62	252	120409	30.824	ng	99
83) Chrysene	21.77	228	413488	43.585	ng	99
84) Bis(2-ethylhexyl)phthalate	21.57	149	242437	40.974	ng	98
85) Di-n-octyl phthalate	22.79	149	413894	41.768	ng	98
86) Indeno(1,2,3-cd)pyrene	28.80	276	506402	45.403	ng	# 76
88) Benzo(b)fluoranthene	23.96	252	448571	47.295	ng	99
89) Benzo(k)fluoranthene	24.03	252	430084	45.538	ng	97
90) Benzo(a)pyrene	24.86	252	434823	47.521	ng	98
91) Dibenzo(a,h)anthracene	28.86	278	420216	46.124	ng	97
92) Benzo(g,h,i)perylene	29.99	276	417734	46.526	ng	97

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

