

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG091219\
 Data File : BG042781.D
 Acq On : 12 Sep 2019 14:22
 Operator : HP/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG091219

Quant Time: Sep 12 18:45:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG091219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 12 14:06:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	75781	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	303518	20.00	ng	0.00
39) Acenaphthene-d10	14.75	164	212252	20.00	ng	0.00
64) Phenanthrene-d10	17.48	188	504670	20.00	ng	0.00
76) Chrysene-d12	21.77	240	495443	20.00	ng	0.00
87) Perylene-d12	25.05	264	544658	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.68	112	346610	79.52	ng	0.00
7) Phenol-d6	7.27	99	494409	79.02	ng	0.00
23) Nitrobenzene-d5	9.28	82	477683	81.94	ng	0.00
42) 2,4,6-Tribromophenol	16.23	330	225679	79.92	ng	0.00
45) 2-Fluorobiphenyl	13.37	172	1075639	80.22	ng	0.00
79) Terphenyl-d14	20.09	244	1881379	84.16	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.59	88	74127	37.263	ng	98
3) Pyridine	3.98	79	233171	38.930	ng	90
4) n-Nitrosodimethylamine	3.89	42	111856	38.380	ng	100
6) Aniline	7.44	93	332168	39.351	ng	99
8) 2-Chlorophenol	7.68	128	186415	39.564	ng	100
9) Benzaldehyde	7.26	77	149322	36.676	ng	94
10) Phenol	7.30	94	257131	40.075	ng	98
11) bis(2-Chloroethyl)ether	7.53	93	192828	39.159	ng	100
12) 1,3-Dichlorobenzene	8.01	146	223342	38.863	ng	99
13) 1,4-Dichlorobenzene	8.16	146	225042	39.114	ng	97
14) 1,2-Dichlorobenzene	8.48	146	218420	39.881	ng	99
15) Benzyl Alcohol	8.35	79	212000	39.485	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.65	45	418933	38.741	ng	99
17) 2-Methylphenol	8.55	107	167528	39.028	ng	100
18) Hexachloroethane	9.22	117	81764	38.492	ng	95
19) n-Nitroso-di-n-propylamine	8.92	70	181552	39.381	ng	96
20) 3+4-Methylphenols	8.88	107	233874	39.525	ng	96
22) Acetophenone	8.94	105	306257	39.953	ng	# 97
24) Nitrobenzene	9.32	77	245577	40.132	ng	98
25) Isophorone	9.85	82	482824	39.732	ng	99
26) 2-Nitrophenol	10.03	139	99089	41.179	ng	94
27) 2,4-Dimethylphenol	10.09	122	168643	39.620	ng	96
28) bis(2-Chloroethoxy)methane	10.33	93	270132	39.861	ng	99
29) 2,4-Dichlorophenol	10.57	162	197786	39.986	ng	98
30) 1,2,4-Trichlorobenzene	10.79	180	241092	39.922	ng	96
31) Naphthalene	10.98	128	597495	39.871	ng	100
32) Benzoic acid	10.20	122	133900	40.025	ng	95
33) 4-Chloroaniline	11.07	127	280001	39.975	ng	99
34) Hexachlorobutadiene	11.28	225	168113	39.732	ng	94
35) Caprolactam	11.84	113	76648	41.671	ng	88
36) 4-Chloro-3-methylphenol	12.18	107	222914	41.506	ng	95
37) 2-Methylnaphthalene	12.58	142	444444	39.592	ng	100
38) 1-Methylnaphthalene	12.80	142	426077	40.469	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.94	216	278541	39.713	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.93	237	152868	38.046	ng	100
43) 2,4,6-Trichlorophenol	13.17	196	190072	40.706	ng	91
44) 2,4,5-Trichlorophenol	13.24	196	192233	40.192	ng	97
46) 1,1'-Biphenyl	13.58	154	589336	39.662	ng	99
47) 2-Chloronaphthalene	13.62	162	478866	38.958	ng	98
48) 2-Nitroaniline	13.81	65	179253	42.039	ng	97
49) Acenaphthylene	14.46	152	751930	40.031	ng	99
50) Dimethylphthalate	14.19	163	642634	40.416	ng	99
51) 2,6-Dinitrotoluene	14.31	165	136009	41.687	ng	93
52) Acenaphthene	14.81	154	489903	39.922	ng	100
53) 3-Nitroaniline	14.63	138	152733	41.711	ng	96
54) 2,4-Dinitrophenol	14.83	184	67152	41.810	ng	# 81
55) Dibenzofuran	15.14	168	724349	39.712	ng	97
56) 4-Nitrophenol	14.92	139	118593	41.888	ng	96
57) 2,4-Dinitrotoluene	15.09	165	185902	42.748	ng	# 97
58) Fluorene	15.79	166	601566	39.795	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.36	232	187419	40.423	ng	99
60) Diethylphthalate	15.56	149	652285	40.414	ng	99
61) 4-Chlorophenyl-phenylether	15.78	204	354296	39.639	ng	97
62) 4-Nitroaniline	15.79	138	163898	41.525	ng	98
63) Azobenzene	16.07	77	619062	39.831	ng	98
65) 4,6-Dinitro-2-methylphenol	15.85	198	89183	41.300	ng	96
66) n-Nitrosodiphenylamine	15.99	169	542527	39.950	ng	98
67) 4-Bromophenyl-phenylether	16.67	248	238721	39.081	ng	99
68) Hexachlorobenzene	16.80	284	256432	39.288	ng	99
69) Atrazine	16.93	200	169878	35.922	ng	98
70) Pentachlorophenol	17.13	266	160320	41.405	ng	96
71) Phenanthrene	17.53	178	975421	40.415	ng	99
72) Anthracene	17.62	178	977413	40.536	ng	99
73) Carbazole	17.88	167	923127	40.280	ng	98
74) Di-n-butylphthalate	18.45	149	1121536	40.852	ng	99
75) Fluoranthene	19.54	202	1277136	40.829	ng	100
77) Benzidine	19.71	184	466818	33.725	ng	97
78) Pyrene	19.89	202	1282108	41.334	ng	98
80) Butylbenzylphthalate	20.79	149	520612	41.472	ng	98
81) Benzo(a)anthracene	21.75	228	1278493	40.586	ng	98
82) 3,3'-Dichlorobenzidine	21.66	252	482946	39.450	ng	100
83) Chrysene	21.81	228	1222435	40.968	ng	99
84) Bis(2-ethylhexyl)phthalate	21.67	149	691778	40.331	ng	100
85) Di-n-octyl phthalate	22.92	149	1177897	40.283	ng	100
86) Indeno(1,2,3-cd)pyrene	28.79	276	1459329	39.493	ng	# 93
88) Benzo(b)fluoranthene	24.01	252	1293439	40.444	ng	99
89) Benzo(k)fluoranthene	24.08	252	1227875	40.645	ng	98
90) Benzo(a)pyrene	24.89	252	1207649	40.328	ng	100
91) Dibenzo(a,h)anthracene	28.87	278	1187681	40.324	ng	98
92) Benzo(g,h,i)perylene	29.96	276	1152953	40.132	ng	97

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

