

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG102518\
 Data File : BG037562.D
 Acq On : 26 Oct 2018 2:49
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 26 08:08:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG101818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 18 14:06:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	61770	20.00	ng	0.00
21) Naphthalene-d8	10.93	136	287442	20.00	ng	0.00
39) Acenaphthene-d10	14.74	164	190025	20.00	ng	0.00
64) Phenanthrene-d10	17.49	188	433867	20.00	ng	0.00
76) Chrysene-d12	21.77	240	379041	20.00	ng	0.00
87) Perylene-d12	25.10	264	402634	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.64	112	302744	81.94	ng	0.00
7) Phenol-d6	7.24	99	452499	80.99	ng	0.00
23) Nitrobenzene-d5	9.28	82	422075	82.26	ng	0.00
42) 2,4,6-Tribromophenol	16.22	330	169265	81.67	ng	0.00
45) 2-Fluorobiphenyl	13.36	172	975253	77.39	ng	0.00
79) Terphenyl-d14	20.08	244	1420196	80.06	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.53	88	74429	39.723	ng	97
3) Pyridine	3.94	79	189120	40.046	ng	95
4) n-Nitrosodimethylamine	3.85	42	69115	41.089	ng	# 92
6) Aniline	7.43	93	275322	40.396	ng	98
8) 2-Chlorophenol	7.66	128	172754	39.969	ng	98
9) Benzaldehyde	7.24	77	130771	37.419	ng	97
10) Phenol	7.27	94	238160	40.402	ng	98
11) bis(2-Chloroethyl)ether	7.52	93	180615	40.343	ng	96
12) 1,3-Dichlorobenzene	7.99	146	192239	39.951	ng	98
13) 1,4-Dichlorobenzene	8.14	146	192123	39.033	ng	98
14) 1,2-Dichlorobenzene	8.46	146	187054	39.507	ng	99
15) Benzyl Alcohol	8.34	79	169059	40.953	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.63	45	327959	40.940	ng	97
17) 2-Methylphenol	8.53	107	158950	40.787	ng	100
18) Hexachloroethane	9.19	117	68858	41.035	ng	93
19) n-Nitroso-di-n-propylamine	8.91	70	157497	40.938	ng	93
20) 3+4-Methylphenols	8.87	107	227083	40.756	ng	99
22) Acetophenone	8.94	105	283284	39.296	ng	# 97
24) Nitrobenzene	9.32	77	218275	40.948	ng	95
25) Isophorone	9.84	82	384390	41.000	ng	96
26) 2-Nitrophenol	10.03	139	110886	46.177	ng	98
27) 2,4-Dimethylphenol	10.08	122	171458	39.990	ng	98
28) bis(2-Chloroethoxy)methane	10.32	93	247101	40.227	ng	96
29) 2,4-Dichlorophenol	10.56	162	197029	41.273	ng	99
30) 1,2,4-Trichlorobenzene	10.78	180	206890	39.853	ng	98
31) Naphthalene	10.98	128	561177	39.748	ng	99
32) Benzoic acid	10.21	122	141825	50.928	ng	99
33) 4-Chloroaniline	11.09	127	268849	40.528	ng	97
34) Hexachlorobutadiene	11.24	225	122404	39.783	ng	95
35) Caprolactam	11.88	113	80859	42.233	ng	94
36) 4-Chloro-3-methylphenol	12.19	107	218030	40.498	ng	97
37) 2-Methylnaphthalene	12.57	142	412743	40.104	ng	99
38) 1-Methylnaphthalene	12.79	142	401282	40.149	ng	99
40) 1,2,4,5-Tetrachlorobenzene	12.93	216	240550	39.246	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.90	237	114024	38.945	nq	98
43) 2,4,6-Trichlorophenol	13.17	196	170826	41.717	nq	99
44) 2,4,5-Trichlorophenol	13.24	196	181472	40.703	nq	97
46) 1,1'-Biphenyl	13.57	154	618710	39.315	nq	99
47) 2-Chloronaphthalene	13.62	162	470310	39.502	nq	97
48) 2-Nitroaniline	13.83	65	154129	42.689	nq	99
49) Acenaphthylene	14.46	152	715647	39.958	nq	99
50) Dimethylphthalate	14.19	163	582197	39.603	nq	99
51) 2,6-Dinitrotoluene	14.32	165	133518	41.056	nq	96
52) Acenaphthene	14.80	154	435108	39.447	nq	98
53) 3-Nitroaniline	14.65	138	149165	41.317	nq	# 97
54) 2,4-Dinitrophenol	14.85	184	44424	44.380	nq	94
55) Dibenzofuran	15.14	168	710711	38.890	nq	100
56) 4-Nitrophenol	14.94	139	106578	39.882	nq	100
57) 2,4-Dinitrotoluene	15.10	165	182829	42.828	nq	96
58) Fluorene	15.78	166	535015	39.094	nq	99
59) 2,3,4,6-Tetrachlorophenol	15.35	232	154654	40.514	nq	99
60) Diethylphthalate	15.54	149	599089	39.694	nq	99
61) 4-Chlorophenyl-phenylether	15.77	204	306552	38.431	nq	99
62) 4-Nitroaniline	15.81	138	148319	41.344	nq	92
63) Azobenzene	16.06	77	568306	39.466	nq	99
65) 4,6-Dinitro-2-methylphenol	15.86	198	92423	41.907	nq	99
66) n-Nitrosodiphenylamine	15.99	169	524337	40.177	nq	98
67) 4-Bromophenyl-phenylether	16.66	248	191463	40.185	nq	97
68) Hexachlorobenzene	16.77	284	193734	39.233	nq	97
69) Atrazine	16.93	200	187540	39.745	nq	98
70) Pentachlorophenol	17.12	266	135113	40.848	nq	98
71) Phenanthrene	17.53	178	868788	39.400	nq	100
72) Anthracene	17.62	178	861959	39.833	nq	99
73) Carbazole	17.89	167	858109	39.643	nq	99
74) Di-n-butylphthalate	18.43	149	989852	41.108	nq	99
75) Fluoranthene	19.53	202	979633	39.171	nq	98
77) Benzidine	19.71	184	459203	35.629	nq	99
78) Pyrene	19.89	202	1000866	40.393	nq	99
80) Butylbenzylphthalate	20.76	149	439399	43.330	nq	98
81) Benzo(a)anthracene	21.75	228	902489	40.254	nq	98
82) 3,3'-Dichlorobenzidine	21.66	252	359005	41.312	nq	98
83) Chrysene	21.82	228	871759	40.437	nq	98
84) Bis(2-ethylhexyl)phthalate	21.62	149	611580	43.025	nq	99
85) Di-n-octyl phthalate	22.87	149	1017436	43.894	nq	99
86) Indeno(1,2,3-cd)pyrene	28.93	276	905292	39.152	nq	100
88) Benzo(b)fluoranthene	24.04	252	861725	39.038	nq	99
89) Benzo(k)fluoranthene	24.11	252	884037	40.878	nq	97
90) Benzo(a)pyrene	24.94	252	851816	40.610	nq	99
91) Dibenzo(a,h)anthracene	28.99	278	740423	38.268	nq	99
92) Benzo(g,h,i)perylene	30.13	276	728523	37.218	nq	99

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

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